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NAS PENSACOLA
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DRAFT FINAL REMEDIAL INVESTIGATION REPORT

APPENDICES
VOLUME II
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NAVAL AIR STATION PENSACOLA
OPERABLE UNIT 10 AND SITE 13

CONTRACT N62467-89-D-0318
CTO - 048

Submitted to:

SOUTHNAVFACENGCOM

Submitted by:

EnSafe/Allen & Hoshall
5720 Summer Trees Drive, Suite 8
Memphis, Tennessee 38134
(901) 383-9115

October 14, 1994

Release of this document requires prior notification of the Commanding Officer of the Naval Air Station, Pensacola, FL

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S02002	13S02003	13S02004	13S02004DE	13S02006	13S03002
	LAB ID:						
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92
NAME	UNITS						
I Aluminum	MG/KG	9410.0000	.	8360.0000	.	.	5570.0000
I Antimony	MG/KG	11.9000 W	.	12.1000 U	.	.	11.8000 U
I Arsenic	MG/KG	2.5000	.	1.5000 J			1.8000 J
I Barium	MG/KG	14.4000 J	.	8.3000 J		.	4.9000 J
I Beryllium	MG/KG	0.4200 U	.	0.4200 W			0.4100 U
I Cadmium	MG/KG	1.5000	.	1.1000 U	.	.	1.0000 U
I Calcium	MG/KG	1060.0000	.	343.0000 J	.	.	3480.0000
I Chromium	MG/KG	10.1000	.	8.8000	.	.	5.6000
I Cobalt	MG/KG	2.3000 W	.	2.3000 W	.	.	2.3000 W
I Copper	MG/KG	12.7000	.	5.9000	.	.	5.2000 W
I Cyanide	MG/KG	0.5200 W	.	0.5300 W		.	0.5000 W
I Iron	MG/KG	5700.0000	.	3830.0000	.	.	2520.0000
I Lead	MG/KG	32.7000	.	16.1000	.	.	10.6000
I Magnesium	MG/KG	110.0000 J	.	90.0000 W	.	.	88.2000 W
I Manganese	MG/KG	54.8000	.	60.2000	.	.	52.7000
I Mercury	MG/KG	0.1000 U	.	0.1100 W	.	.	0.1000 W
I Nickel	MG/KG	8.4000 U	.	8.5000 W	.	.	8.3000 W
I Potassium	MG/KG	1010.0000 U	.	1030.0000 W	.	.	1010.0000 W
I Selenium	MG/KG	0.8400 U	.	0.8500 U	.	.	0.8300 U
I Silver	MG/KG	2.1000 U	.	2.1000 U	.	.	2.1000 U
I Sodium	MG/KG	125.0000 J	.	109.0000 J	.	.	122.0000 J
I Thallium	MG/KG	0.4200 U	.	0.4200 U	.	.	0.4100 U
I Vanadium	MG/KG	15.5000	.	14.2000	.	.	9.4000 J
I Zinc	MG/KG	29.4000	.	18.8000		.	12.0000
P 4,4'-DDD	UG/KG	.	0.1400 J	.	.	3.4000 U	.
P 4,4'-DDE	UG/KG	.	3.5000 U	.	.	0.0690 J	.
P 4,4'-DDT	UG/KG	.	0.8400 J	.	.	3.4000 U	.
P Aldrin	UG/KG	.	1.8000 W	.	.	1.7000 U	.
P Aroclor 1016	UG/KG	.	35.0000 W	.	.	34.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	13802002	13S02003	13802004	13S02004RE	13802006	13803002
		12/09/92	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92
P Aroclor 1221	UG/KG		71.0000 U			68.0000 U	
P Aroclor 1232	UG/KG		35.0000 U			34.0000 U	
P Aroclor 1242	UG/KG		35.0000 U			34.0000 U	
P Aroclor 1248	UG/KG		35.0000 U	.	.	34.0000 U	
P Aroclor 1254	UG/KG	.	35.0000 U	.	.	34.0000 U	.
P Aroclor 1260	UG/KG		35.0000 U	.	.	34.0000 U	
P Dieldrin	UG/KG		0.7800 J	.	.	0.1400 J	.
P Endosulfan I	UG/KG		1.8000 U	.		1.7000 U	.
P Endosulfan II	UG/KG	.	3.5000 U	.		3.4000 U	.
P Endosulfan sulfate	UG/KG		3.5000 U	.	.	3.4000 U	.
P Endrin	UG/KG		1.1000 J	.		3.4000 U	.
P Endrin aldehyde	UG/KG		3.5000 U	.	.	3.4000 U	.
P Endrin ketone	UG/KG	.	1.1000 J	.	.	3.4000 U	.
P Heptachlor	UG/KG		0.3600 U	.	.	0.1600 U	.
P Heptachlor epoxide	UG/KG		0.1400 J	.	.	0.0870 J	.
P Methoxychlor	UG/KG	.	0.5900 J	.		17.0000 U	.
P Toxaphene	UG/KG	.	180.0000 U	.	.	170.0000 U	.
P alpha-BHC	UG/KG	.	1.8000 U	.	.	1.7000 U	.
P alpha-Chlordane	UG/KG	.	1.8000 U	.	.	0.0860 J	.
P beta-BHC	UG/KG	.	1.8000 U	.	.	1.7000 U	.
P delta-BHC	UG/KG	.	1.8000 U	.		0.0960 J	.
P gamma-BHC (Lindane)	UG/KG	.	1.8000 U	.	.	1.7000 U	.
P gamma-Chlordane	UG/KG	.	1.1000 J	.	.	1.7000 U	.
S 1,2,4-Trichlorobenzene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 1.2-Dichlorobenzene	UG/KG	.	340.0000 U		.	330.0000 U	.
S 1,3-Dichlorobenzene	UG/KG	.	340.0000 U		.	330.0000 U	.
S 1,4-Dichlorobenzene	UG/KG	.	340.0000 U	.	.	330.0000 U	
S 2,4,5-Trichlorophenol	UG/KG	.	820.0000 U		.	800.0000 U	.
S 2,4,6-Trichlorophenol	UG/KG	.	340.0000 U	.	.	330.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S02002	13802003	13S02004	13S02004RE	13802006	13503002
	LAB ID:						
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92
NAME	UNITS						
S 2,4-Dichlorophenol	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2,4-Dimethylphenol	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2,4-Dinitrophenol	UG/KG	.	820.0000 U	.	.	800.0000 U	
S 2,4-Dinitrotoluene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2,6-Dinitrotoluene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2-Chloronaphthalene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2-Chlorophenol	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2-Methylnaphthalene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2-Methylphenol	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 2-Nitroaniline	UG/KG	.	820.0000 U	.	.	800.0000 U	.
S 2-Nitrophenol	UG/KG		340.0000 U	.	.	330.0000 U	.
S 3,3'-Dichlorobenzidine	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 3-Nitroaniline	UG/KG	.	820.0000 U	.	.	800.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/KG	.	820.0000 U	.	.	800.0000 U	.
S 4-Bromophenyl-phenylethet	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 4-Chloro-3-methylphenol	UG/KG	.	240.0000 J	.	.	330.0000 U	.
S 4-Chloroaniline	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 4-Chlorophenyl-phenylether	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S 4-Methylphenol	UG/KG		340.0000 U		.	330.0000 U	.
S 4-Nitroaniline	UG/KG	.	820.0000 U	.	.	800.0000 U	.
S 4-Nitrophenol	UG/KG	.	820.0000 U	.	.	800.0000 U	.
S Acenaphthene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Acenaphthylene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Anthracene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Benzo(a)anthracene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Benzo(a)pyrene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Benzo(b,k)fluoranthene	UG/KG	.	340.0000 U	.	.	65.0000 J	.
S Benzo(g,h,i)perylene	UG/KG	.	340.0000 U	.	.	330.0000 U	
S Butylbenzylphthalate	UG/KG	.	340.0000 U	.	.	330.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S02002	13S02003	13S02004	13S02004RE	13S02006	13S03002
	LAB ID:						
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92
NAME	UNITS						
S Carbazole	UG/KG		340.0000 U	.	.	330.0000 U	
S Chrysene	UG/KG	.	340.0000 U	.	.	35.0000 J	
S Di-n-butylphthalate	UG/KG		340.0000 U	.		330.0000 U	
S Di-n-octylphthalate	UG/KG		340.0000 U	.		330.0000 U	.
S Dibenz (a,h) anthracene	UG/KG		340.0000 U	.	.	330.0000 U	.
S Dibenzofuran	UG/KG		340.0000 U	.	.	330.0000 U	.
S Diethylphthalate	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Dimethylphthalate	UG/KG		340.0000 U	.	.	330.0000 U	
S Fluoranthene	UG/KG	.	35.0000 J	.	.	330.0000 U	.
S Fluorene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Hexachlorobenzene	UG/KG	.	340.0000 U	.	.	330.0000 U	
S Hexachlorobutadiene	UG/KG		340.0000 U	.		330.0000 U	
S Hexachlorocyclopentadiene	UG/KG		340.0000 U	.	.	330.0000 U	.
S Hexachloroethane	UG/KG		340.0000 U	.	.	330.0000 U	
S Indeno(1,2,3-cd)pyrene	UG/KG		340.0000 U	.		330.0000 U	.
S Isophorone	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S N-Nitroso-di-n-propylamine	UG/KG		340.0000 U		.	330.0000 U	.
S N-Nitrosodiphenylamine	UG/KG	.	340.0000 U	.		330.0000 U	.
S Naphthalene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Nitrobenzene	UG/KG	.	340.0000 U	.	.	330.0000 U	.
S Pentachlorophenol	UG/KG		820.0000 U	.	.	800.0000 U	.
S Phenanthrene	UG/KG		340.0000 U			330.0000 U	.
S Phenol	UG/KG	.	99.0000 J			330.0000 U	
S Pyrene	UG/KG	.	37.0000 J	.		330.0000 U	
S bis(2-Chloroethoxy)methane	UG/KG	.	340.0000 U	.		330.0000 U	
S bis(2-Ethlyhexy)phthalate	UG/KG		340.0000 U	.		44.0000 U	
S bis(2-chloroethyl)ether	UG/KG		340.0000 U			330.0000 U	
S bis(2-chloroisopropyl)ether	UG/KG		340.0000 U			330.0000 U	
V 1,1,1-Trichloroethane	UG/KG	10.0000 U		11.0000 U	15.0000 U		10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13802002	13802003	13802004	13S02004RE	13S02006	13803002
	LAB ID:						
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92
NAME	UNITS						
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 2-Butanone	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Acetone	UG/KG	130.0000 U	.	240.0000 U	170.0000 U	.	52.0000 U
V Benzene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Bromoform	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Bromomethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Chloroethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Chloroform	UG/KG	2.0000 U	.	2.0000 U	15.0000 U	.	3.0000 U
V Chloromethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Methylene Chloride	UG/KG	61.0000 U	.	59.0000 U	41.0000 U	.	51.0000 U
V Styrene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Toluene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Trichloroethene	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	.	11.0000 U	15.0000 U	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S02002	13S02003	13S02004	13S02004RE	13S02006	13S03002
	LAB ID:						
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92	12/09/92
NAME	UNITS						
V Xylene (total)	UG/KG	10.0000 U		11.0000 U	15.0000 U		10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U		11.0000 U	15.0000 U		10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U		11.0000 U	15.0000 U		10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13503003	13503004	13503005	13S04002*	13S04004	13504006
	LAB ID:					532980	532983
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/12/92	01/20/93	01/20/93
NAME	UNITS						
I Aluminum	MG/KG	.	3120.0000	.	3480.0000	12700.0000	64.5000
I Antimony	MG/KG		12.2000 U	.	12.2000 U	12.2000 U	11.9000 U
I Arsenic	MG/KG	.	0.9200 J	.	0.8500 U	2.3000	0.6300 U
I Barium	MG/KG		4.4000 J	.	5.7000 J	8.2000 J	1.3000 U
I Beryllium	MG/KG	.	0.4300 U	.	0.4300 U	0.6400 U	0.6300 U
I Cadmium	MG/KG	.	1.1000 U	.	1.1000 U	1.1000 U	1.0000 U
I Calcium	MG/KG	.	685.0000 J	.	504.0000 J	323.0000 J	27.2000 J
I Chromium	MG/KG	.	5.4000	.	15.1000	7.7000	2.1000 U
I Cobalt	MG/KG	.	2.3000 U	.	2.4000 U	2.6000 U	2.5000 U
I Copper	MG/KG	.	5.3000 U		5.3000 U	10.2000	5.2000 U
I Cyanide	MG/KG		0.5300 U	.	0.5300 U	0.5300 U	0.5200 U
I Iron	MG/KG	.	1460.0000	.	1830.0000	5380.0000	52.0000
I Lead	MG/KG		13.4000	.	5.2000	9.4000	0.7500
I Magnesium	MG/KG	.	90.6000 U	.	90.8000 U	118.0000 U	115.0000 U
I Manganese	MG/KG	.	23.9000	.	31.4000	101.0000	1.3000 U
I Mercury	MG/KG	.	0.1100 U	.	0.1100 U	0.1100 U	0.1000 U
I Nickel	MG/KG	.	8.5000 U	.	8.5000 U	8.1000 U	7.9000 U
I Potassium	MG/KG	.	1040.0000 U	.	1040.0000 U	844.0000 U	824.0000 U
I Selenium	MG/KG		0.8500 U		0.8500 U	0.6400 U	0.6300 U
I Silver	MG/KG	.	2.1000 U	.	2.1000 U	2.1000 U	2.1000 U
I sodium	MG/KG	.	117.0000 J	.	116.0000 J	116.0000 J	107.0000 J
I Thallium	MG/KG	.	0.4300 U		0.4300 U	0.6400 U	0.6300 U
I Vanadium	MG/KG	.	5.0000 J	.	6.4000 J	20.1000	3.1000 U
I Zinc	MG/KG		14.8000		4.9000,	7.6000	2.7000 J
P 4,4'-DDD	UG/KG	3.6000 U		3.5000 U	3.5000 U	3.5000 U	3.4000 U
P 4,4'-DDE	UG/KG	3.6000 U		6.1000	3.5000 U	3.5000 U	3.4000 U
P 4,4'-DDT	UG/KG	0.8700 J		4.1000	0.2400 J	3.5000 U	3.4000 U
P Aldrin	UG/KG	1.9000 U		1.8000 U	0.4600 J	1.8000 U	1.7000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S03003	13S03004	13S03005	13S04002*	13S04004	13S04006
	LAB ID:					532980	532983
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/12/92	01/20/93	01/20/93
NAME	UNITS						
P Aroclor 1016	UG/KG	36.0000 U		35.0000 U	35.0000 U	35.0000 U	34.0000 U
P Aroclor 1221	UG/KG	73.0000 U		71.0000 U	71.0000 U	71.0000 U	69.0000 U
P Aroclor 1232	UG/KG	36.0000 U		35.0000 U	35.0000 U	35.0000 U	34.0000 U
P Aroclor 1242	UG/KG	36.0000 U		35.0000 U	35.0000 U	35.0000 U	34.0000 U
P Aroclor 1248	UG/KG	36.0000 U		35.0000 U	35.0000 U	35.0000 U	34.0000 U
P Aroclor 1254	UG/KG	36.0000 U		35.0000 U	35.0000 U	35.0000 U	34.0000 U
P Aroclor 1260	UG/KG	36.0000 U		35.0000 U	35.0000 U	35.0000 U	34.0000 U
P Dieldrin	UG/KG	0.9900 J		0.3400 J	3.5000 U	3.5000 U	3.4000 U
P Endosulfan I	UG/KG	0.6900 J		0.1900 J	0.2500 J	0.2900 J	1.7000 U
P Endosulfan 11	UG/KG	0.3300 J		3.5000 U	3.5000 U	2.4000 J	3.4000 U
P Endosulfan sulfate	UG/KG	0.2400 J		3.5000 U	1.5000 J	3.5000 U	3.4000 U
P Endrin	UG/KG	3.7000 J		0.5800 J	0.5000 J	3.2000 J	3.4000 U
P Endrin aldehyde	UG/KG	3.6000 U		3.5000 U	0.6500 J	3.5000 U	3.4000 U
P Endrin ketone	UG/KG	5.4000 J		0.5800 J	3.5000 U	0.8100 J	3.4000 U
P Heptachlor	UG/KG	0.3100 U		0.3900 U	1.8000 U	0.4200 U	1.7000 U
P Heptachlor epoxide	UG/KG	1.9000 U		0.1400 J	1.8000 U	0.7000 J	1.7000 U
P Methoxychlor	UG/KG	22.0000 J		18.0000 U	0.6400 J	18.0000 U	17.0000 U
P Toxaphene	UG/KG	190.0000 U		180.0000 U	180.0000 U	180.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.9000 U		1.8000 U	1.8000 U	1.8000 U	1.7000 U
P alpha-Chlordane	UG/KG	0.1700 J		1.8000 U	0.1200 J	1.1000 J	1.7000 U
P beta-BHC	UG/KG	1.9000 U		1.8000 U	1.8000 U	1.8000 U	1.7000 U
P delta-BHC	UG/KG	1.9000 U		0.1000 J	0.1000 J	1.8000 U	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.9000 U		1.8000 U	1.8000 U	1.8000 U	1.7000 U
P gamma-Chlordane	UG/KG	0.3800 J		0.1900 J	0.2500 J	0.3800 J	1.7000 U
8 1,2,4-Trichlorobenzene	UG/KG	350.0000 U		340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 1,2-Dichlorobenzene	UG/KG	350.0000 U		340.0000 U	350.0000 'U	350.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG	350.0000 U		340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG	350.0000 U		340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	860.0000 U		830.0000 U	850.0000 U	850.0000 U	820.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S03003	13S03004	13S03005	13S04002*	13S04004	13S04006
	LAB ID:					532980	532983
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/12/92	01/20/93	01/20/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4-Dichlorophenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4-Dimethylphenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4-Dinitrophenol	UG/KG	860.0000 U	.	830.0000 U	850.0000 U	850.0000 U	820.0000 U
S 2,4-Dinitrotoluene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,6-Dinitrotoluene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Chloronaphthalene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Chlorophenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Methylnaphthalene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Methylphenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Nitroaniline	UG/KG	860.0000 U	.	830.0000 U	850.0000 U	850.0000 U	820.0000 U
S 2-Nitrophenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 3-Nitroaniline	UG/KG	860.0000 U	.	830.0000 U	850.0000 U	850.0000 U	820.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	860.0000 U	.	830.0000 U	850.0000 U	850.0000 U	820.0000 U
S 4-Bromophenyl-phenylether	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Chloro-3-methylphenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Chloroaniline	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Methylphenol	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Nitroaniline	UG/KG	860.0000 U	.	830.0000 U	850.0000 U	850.0000 U	820.0000 U
S 4-Nitrophenol	UG/KG	860.0000 U	.	830.0000 U	850.0000 U	850.0000 U	820.0000 U
S Acenaphthene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Acenaphthylene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Anthracene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Benzo(a)anthracene	UG/KG	350.0000 U	.	160.0000 J	350.0000 U	350.0000 U	340.0000 U
S Benzo(a)pyrene	UG/KG	350.0000 U	.	94.0000 J	350.0000 U	350.0000 U	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG	93.0000 J	.	310.0000 J	350.0000 U	350.0000 U	340.0000 U
S Benzo(g,h,i)perylene	UG/KG	350.0000 U	.	83.0000 J	350.0000 U	350.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S03003	13S03004	13S03005	13S04002*	13S04004	13S04006
	LAB ID:					532980	532983
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/12/92	01/20/93	01/20/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	350.0000 U	.	340.0000 U	86.0000 U	460.0000 U	280.0000 U
S Carbazole	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Chrysene	UG/KG	52.0000 J	.	180.0000 J	350.0000 U	350.0000 U	340.0000 U
S Di-n-butylphthalate	UG/KG	350.0000 U	.	340.0000 U	61.0000 U	350.0000 U	340.0000 U
S Di-n-octylphthalate	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Dibenz (a,h) anthracene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Dibenzofuran	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Diethylphthalate	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Dimethylphthalate	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Fluoranthene	UG/KG	75.0000 J	.	520.0000	350.0000 U	350.0000 U	340.0000 U
S Fluorene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachloroethane	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	350.0000 U	.	57.0000 J	350.0000 U	350.0000 U	340.0000 U
S Isophorone	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Naphthalene	UG/KG	350.0006 U	.	340.0'000 U	350.0000 U	350.0000 U	340.0000 U
S Nitrobenzene	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	860.0000 U	.	830.0000 U	850.0000 U	850.0000 U	820.0000 U
S Phenanthrene	UG/KG	350.0000 U	.	94.0000 J	350.0000 U	350.0000 U	340.0000 U
S Phenol	UG/KG	350.0000 U	.	340.0000 U	50.0000 J	350.0000 U	340.0000 U
S Pyrene	UG/KG	49.0000 J	.	260.0000 J	350.0000 U	350.0000 U	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	350.0000 U	.	340.0000 U	80.0000 U	55.0000 U	340.0000 U
S bis(2-chloroethyl)ether	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	350.0000 U	.	340.0000 U	350.0000 U	350.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S03003	13S03004	13S03005	13S04002*	13S04004	13S04006
	LAB ID:					532980	532983
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/12/92	01/20/93	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 2-Butanone	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 2-Hexanone	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Acetone	UG/KG	.	28.0000 U	.	120.0000 U	120.0000 U	34.0000 U
V Benzene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Brmodichlormethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Bromoform	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Bromomethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Chlorobenzene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Chloroethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Chloroform	UG/KG	.	2.0000 U	.	1.0000 U	2.0000 U	10.0000 U
V Chloromethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Ethylbenzene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Methylene Chloride	UG/KG	.	75.0000 U	.	29.0000 U	24.0000 U	18.0000 U
V Styrene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Toluene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U
V Trichloroethene	UG/KG	.	11.0000 U	.	11.0000 U	11.0000 U	10.0000 U

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13803003	13S03004	13S03005	13S04002*	13S04004	13804006
	LAB ID:					532980	532983
	DATE RCVD:	12/09/92	12/09/92	12/09/92	12/12/92	01/20/93	01/20/93
NAME	UNITS						
V Vinyl Chloride	UG/KG		11.0000 U		11.0000 U	11.0000 U	10.0000 U
V Xylene (total)	UG/KG		11.0000 U		11.0000 U	11.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG		11.0000 U		11.0000 U	11.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG		11.0000 U		11.0000 U	11.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 ANI) SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S04008	13505002	13S05002RE	13S05004	33806002	13506004
	LAB ID:	532982				529923	529883
	DATE RCVD:	01/20/93	12/12/92	12/12/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
I Aluminum	MG/KG	83.0000	63.2000	.	831.0000	1530.0000	254.0000
I Antimony	MG/KG	11.8000 U	11.8000 U		12.9000 U	12.9000 U	11.6000 U
I Arsenic	MG/KG	0.6200 U	0.8300 U	.	0.9000 U	0.6800 U	0.6100 U
I Barium	MG/KG	1.2000 U	0.7100 J	.	4.6000 J	3.9000 J	1.4000 J
I Beryllium	MG/KG	0.6200 U	0.4100 U	.	0.4500 U	0.4500 U	0.4100 U
I Cadmium	MG/KG	1.0000 U	1.0000 U	.	1.3000	1.1000 U	1.0000 U
I Calcium	MG/KG	83.7000 J	192.0000 U	.	657.0000 J	5020.0000	13600.0000
I Chromium	MG/KG	2.1000 U	2.1000 U	.	2.3000 U	2.3000 U	2.0000 U
I Cobalt	MG/KG	2.5000 U	2.3000 U	.	2.5000 U	2.5000 U	2.2000 U
I Copper	MG/KG	5.2000 U	5.2000 U		5.7000 U	5.7000 U	5.4000
I Cyanide	MG/KG	0.5200 U	0.5200 U	.	1.7000	0.5700 U	0.5100 U
I Iron	MG/KG	36.5000	23.3000	.	500.0000	793.0000	242.0000
I Lead	MG/KG	0.9800	0.7500	.	16.9000	1.5000	2.3000
I Magnesium	MG/KG	115.0000 U	88.2000 U	.	96.2000 U	96.2000 U	86.8000 U
I Manganese	MG/KG	1.2000 U	1.8000 J	.	3.7000	9.3000	4.9000
I Mercury	MG/KG	0.1000 U	0.1000 U	.	0.1100 U	0.1100 U	0.1000 U
I Nickel	MG/KG	7.9000 U	8.3000 U	.	9.0000 U	9.0000 U	8.2000 U
I Potassium	MG/KG	821.0000 U	1010.0000 U	.	1100.0000 U	1100.0000 U	992.0000 U
I Selenium	MG/KG	0.6200 U	0.8300 U	.	0.9000 U	0.6800 U	0.6100 U
I Silver	MG/KG	2.1000 U	2.1000 U	.	2.3000 U	2.3000 U	2.0000 U
I Sodium	MG/KG	109.0000 J	109.0000 J	.	117.0000 J	118.0000 J	221.0000 J
I Thallium	MG/KG	0.6200 U	0.4100 U	.	0.4500 U	0.6800 U	0.4100 U
I Vanadium	MG/KG	3.1000 U	3.9000 U	.	4.3000 U	4.3000 U	3.9000 U
I Zinc	MG/KG	2.0000 J	1.0000 U	.	1.5000 J	3.1000 J	7.0000
P 4,4'-DDD	UG/KG	3.3000 U	3.4000 U	.	3.7000 'U	2.3000 J	3.3000 U
P 4,4'-DDE	UG/KG	3.3000 U	3.4000 U	.	0.0690 J	6.6000 J	3.3000 U
P 4,4'-DDT	UG/KG	3.3000 U	3.4000 U	.	3.7000 U	2.0000 J	0.0870 J
P Aldrin	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	6.1000	1.7000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

NAME	ENSAFE ID:	13804008	13805002	13S05002RE	13805004	13806002	13806004
	LAB ID:	532982				529923	529883
	DATE RCVD:	01/20/93	12/12/92	12/12/92	12/12/92	12/18/92	12/18/92
	UNITS						
P Aroclor 1016	UG/KG	33.0000 U	34.0000 U	.	37.0000 U	36.0000 U	33.0000 U
P Aroclor 1221	UG/KG	68.0000 U	69.0000 U	.	76.0000 U	74.0000 U	68.0000 U
P Aroclor 1232	UG/KG	33.0000 U	34.0000 U	.	37.0000 U	36.0000 U	33.0000 U
P Aroclor 1242	UG/KG	33.0000 U	34.0000 U	.	37.0000 U	36.0000 U	33.0000 U
P Aroclor 1248	UG/KG	33.0000 U	34.0000 U	.	37.0000 U	36.0000 U	33.0000 U
P Aroclor 1254	UG/KG	33.0000 U	34.0000 U	.	37.0000 U	36.0000 U	33.0000 U
P Aroclor 1260	UG/KG	33.0000 U	34.0000 U	.	37.0000 U	36.0000 U	33.0000 U
P Dieldrin	UG/KG	3.3000 U	3.4000 U	.	3.7000 U	1.6000 J	0.1400 J
P Endosulfan I	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	1.9000 U	1.7000 U
P Endosulfan II	UG/KG	3.3000 U	3.4000 U	.	3.7000 U	3.6000 U	3.3000 U
P Endosulfan sulfate	UG/KG	3.3000 U	3.4000 U	.	3.7000 U	9.2000	3.3000 U
P Endrin	UG/KG	3.3000 U	3.4000 U	.	3.7000 U	3.6000 U	3.3.000 U
P Endrin aldehyde	UG/KG	3.3000 U	0.1100 J	.	3.7000 U	3.6000 U	3.3000 U
P Endrin ketone.	UG/KG	3.3000 U	3.4000 U	.	3.7000 U	0.8000 J	3.3000 U
P Heptachlor	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	1.9000 U	0.1200 J
P Heptachlor epoxide	UG/KG	1.7000 U	0.1500 J	.	1.9000 U	0.5400 J	1.7000 U
P Methoxychlor	UG/KG	17.0000 U	0.4100 J	.	19.0000 U	19.0000 U	17.0000 U
P Toxaphene	UG/KG	170.0000 U	180.0000 U	.	190.0000 U	190.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	0.2100 J	1.7000 U
P alpha-Chlordane	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	1.9000 U	1.7000 U
P beta-BHC	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	1.9000 U	1.7000 U
P delta-BHC	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	2.6000 J	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	1.9000 U	1.7000 U
P gamma-Chlordane	UG/KG	1.7000 U	1.8000 U	.	1.9000 U	1.1000 J	1.7000 U
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	.	370.0000 'U	11000.0000 U	330.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2,4,5-Trichlorophenol	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	13804008	13505002	13S05002RE	13S05004	13806002	13506004
	LAB ID:	532982				529923	529883
	DATE RCVD:	01/20/93	12/12/92	12/12/92	12/12/92	12/18/92	12/18/92
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2,4-Dinitrophenol	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	2100.0000 J	330.0000 U
S 2-Methylphenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 2-Nitroaniline	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 3-Nitroaniline	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S 4-Nitroaniline	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U
S 4-Nitrophenol	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U
S Acenaphthene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Acenaphthylene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Anthracene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Benzo(a)anthracene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Benzo(a)pyrene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S04008	13S05002	13S05002RE	33805004	13S06002	13S06004
	LAB ID:	532982				529923	529883
	DATE RCVD:	01/20/93	12/12/92	12/12/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	250.0000 J	340.0000 U	.	50.0000 U	11000.0000 U	330.0000 U
S Carbazole	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Chrysene	UG/KG	340.0000 U	340.0000 U		370.0000 U	11000.0000 U	330.0000 U
S Di-n-butylphthalate	UG/KG	340.0000 U	48.0000 U	.	85.0000 U	11000.0000 U	330.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Dibenz(ah)anthracene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Dibenzofuran	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Diethylphthalate	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Fluoranthene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Fluorene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Hexachloroethane	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Isophorone	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Naphthalene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Nitrobenzene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Pentachlorophenol	UG/KG	820.0000 U	820.0000 U	.	900.0000 U	27000.0000 U	810.0000 U
S Phenanthrene	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	3600.0000 J	330.0000 U
S Phenol	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S Pyrene	UG/KG	340.0000 U	340.0000 U		370.0000 U	11000.0000 U	330.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	340.0000 U	340.0000 U	.	62.0000 U	11000.0000 U	330.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	340.0000 U	.	370.0000 U	11000.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S04008	13S05002	13S05002RE	33805004	13S06002	13806004
	LAB ID:	532982				529923	529883
	DATE RCVD:	01/20/93	12/12/92	12/12/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10,0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10,0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Acetone	UG/KG	27.0000 U	490.0000	800.0000 J	190.0000 U	78.0000 U	48.0000 U
V Benzene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Chloroform	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	1.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Methylene Chloride	UG/KG	11.0000 U	14.0000 U	75.0000 U	26.0000 U	33.0000 U	24.0000 U
V Styrene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	1.0000 J	10.0000 U
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	13S04008	13S05002	13S05002RE	13S05004	13S06002	13S06004
	LAB ID:	532982				529923	529883
	DATE RCVD:	01/20/93	12/12/92	12/12/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	180.0000	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	35.0000 U	11.0000 U	11.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S61002	13S61004	13S61006	13S61007	13S62002	13S62004
	LAB ID:	532082	532083	532084	532085	532086	532087
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
I Aluminum	MG/KG	201.0000	220.0000	207.0000	122.0000	259.0000	185.0000
I Antimony	MG/KG	12.0000 U	12.0000 U	11.9000 W	11.5000 U	11.9000 W	12.0000 W
I Arsenic	MG/KG	0.6200 U	0.6300 U	0.6300 W	0.6100 U	0.6300 W	0.6300 W
I Barium	MG/KG	1.3000 U	1.3000 U	1.3000 W	1.2000 U	1.3000 W	1.3000 W
I Beryllium	MG/KG	0.6300 U	0.6300 U	0.6300 W	0.6100 U	0.6300 W	0.6300 W
I Cadmium	MG/KG	1.1000 U	1.1000 U	1.0000 W	1.0000 U	1.0000 W	1.1000 W
I Calcium	MG/KG	36.4000 J	92.3000 J	32.4000 W	26.3000 J	27.1000 W	27.4000 W
I Chromium	MG/KG	2.1000 U	2.1000 U	2.1000 W	2.0000 U	2.1000 W	2.1000 W
I Cobalt	MG/KG	2.5000 U	2.5000 U	2.5000 W	2.4000 U	2.5000 W	2.5000 W
I Copper	MG/KG	5.3000 U	5.3000 U	5.2000 W	5.0000 U	5.2000 W	5.3000 W
I Cyanide	MG/KG	0.5300 U	0.5300 U	0.5200 W	0.5100 U	0.5300 W	0.5300 W
I Iron	MG/KG	123.0000	134.0000	119.0000	84.6000	156.0000	106.0000
I Lead	MG/KG	1.4000	1.3000	0.7600	0.7600	0.9600	0.7600
I Magnesium	MG/KG	116.0000 W	116.0000 W	116.0000 W	111.0000 W	115.0000 W	116.0000 W
I Manganese	MG/KG	1.3000 W	1.3000 W	1.3000 W	1.6000 W	1.3000 W	1.3000 W
I Mercury	MG/KG	0.1100 W	0.1000 W	0.1000 W	0.1000 W	0.1100 W	0.1100 W
I Nickel	MG/KG	8.0000 W	8.0000 W	8.0000 W	7.7000 W	7.9000 W	8.0000 W
I Potassium	MG/KG	830.0000 W	833.0000 W	829.0000 W	798.0000 W	824.0000 W	833.0000 W
I Selenium	MG/KG	0.6200 W	0.6000 W	0.6300 W	0.6100 W	0.6300 W	0.6300 W
I Silver	MG/KG	2.1000 W	2.1000 W	2.1000 W	2.0000 W	2.1000 W	2.1000 W
I Sodium	MG/KG	108.0000 W	122.0000 W	121.0000 W	108.0000 W	110.0000 W	109.0000 W
I Thallium	MG/KG	0.6200 W	0.6000 W	0.6300 W	0.6100 W	0.6300 W	0.6300 W
I Vanadium	MG/KG	3.2000 W	3.2000 W	3.1000 W	3.0000 W	3.1000 W	3.2000 W
I Zinc	MG/KG	1.3000 W	3.2000 W	2.6000 W	2.3000 W	1.9000 W	2.6000 W
P 4,4'-DDD	UG/KG	3.5000 W	3.5000 W	3.5000 W	3.4000 W	3.5000 W	3.5000 W
P 4,4'-DDE	UG/KG	0.2800 W	0.2000 W	0.1700 W	3.4000 W	0.0410 W	3.5000 W
P 4,4'-DDT	UG/KG	0.2000 W	0.1700 W	3.5000 W	0.0770 W	3.5000 W	3.5000 W
P Aldrin	UG/KG	1.8000 W	1.8000 W	1.8000 W	1.7000 W	1.8000 W	1.8000 W

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13861002	13861004	13861006	13S61007	13862002	13862004
	LAB ID:	532082 R1	532083 R1	532084 R1	532085 R1	532086 R1	532087 R1
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
P Aroclor 1016	UG/KG	35.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1221	UG/KG	71.0000 U	71.0000 U	71.0000 U	68.0000 U	71.0000 U	71.0000 U
P Aroclor 1232	UG/KG	35.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1242	UG/KG	35.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1248	UG/KG	35.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1254	UG/KG	35.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1260	UG/KG	35.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Dieldrin	UG/KG	3.5000 U	3.5000 U	3.5000 U	3.4000 U	3.5000 U	3.5000 U
P Endosulfan I	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P Endosulfan II	UG/KG	0.0820 J	3.5000 U	3.5000 U	3.4000 U	3.5000 U	3.5000 U
P Endosulfan sulfate	UG/KG	3.5000 U	0.5200 J	0.1400 J	3.4000 U	3.5000 U	3.5000 U
P Endrin	UG/KG	3.5000 U	3.5000 U	0.1900 J	3.4000 U	3.5000 U	3.5000 U
P Endrin aldehyde	UG/KG	3.5000 U	3.5000 U	3.5000 U	3.4000 U	3.5000 U	3.5000 U
P Endrin ketone	UG/KG	3.5000 U	3.5000 U	3.5000 U	3.4000 U	3.5000 U	3.5000 U
P Heptachlor	UG/KG	1.8000 U	1.8000 U	0.0690 J	1.7000 U	1.8000 U	1.8000 U
P Heptachlor epoxide	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P Methoxychlor	UG/KG	18.0000 U	18.0000 U	18.0000 U	17.0000 U	18.0000 U	18.0000 U
P Toxaphene	UG/KG	180.0000 U	180.0000 U	180.0000 U	170.0000 U	180.0000 U	180.0000 U
P alpha-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P alpha-Chlordane	UG/KG	1.8000 U	0.1400 J	0.0740 J	1.7000 U	1.8000 U	1.8000 U
P beta-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P delta-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P gamma-Chlordane	UG/KG	0.0680 J	1.8000 U	0.1900 J	0.1500 J	0.2000 J	0.1000 J
S 1,2,4-Trichlorobenzene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 1,2-Dichlorobenzene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 1,3-Dichlorobenzene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 1,4-Dichlorobenzene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4,5-Trichlorophenol	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13961002	13861004	13861006	13861007	13862002	13862004
ENSAFE ID:		13961002	13861004	13861006	13861007	13862002	13862004
LAB ID:		532082	532083	532084	532085	532086	532087
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
S 2,4,6-Trichlorophenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4-Dichlorophenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4-Dimethylphenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4-Dinitrophenol	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U
S 2,4-Dinitrotoluene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,6-Dinitrotoluene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Chloronaphthalene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Chlorophenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Methylnaphthalene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Methylphenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Nitroaniline	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U
S 2-Nitrophenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 3-Nitroaniline	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U
S 4-Bromophenyl-phenylether	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Chloro-3-methylphenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Chloroaniline	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Methylphenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Nitroaniline	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U
S 4-Nitrophenol	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U
S Acenaphthene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Acenaphthylene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Anthracene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(a)anthracene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(a)pyrene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(b,k)fluoranthene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(g,h,i)perylene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data **not** shown in **significant** digits.

	ENSAFE ID:	13S61002	13861004	13861006	13861007	13862002	13862004
	LAB ID:	532082	532083	532084	532085	532086	532087
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Carbazole	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Chrysene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Di-n-butylphthalate	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Di-n-octylphthalate	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Dibenz(a,h)anthracene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Dibenzofuran	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Diethylphthalate	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	48.0000 J
S Dimethylphthalate	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Fluoranthene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Fluorene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachlorobenzene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachlorobutadiene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachlorocyclopentadiene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachloroethane	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Isophorone	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S N-Nitrosodiphenylamine	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Naphthalene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Nitrobenzene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Pentachlorophenol	UG/KG	840.0000 U	840.0000 U	840.0000 U	820.0000 U	840.0000 U	850.0000 U
S Phenanthrene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Phenol	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S Pyrene	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-chloroethyl)ether	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	350.0000 U	350.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digiti.

NAME	UNITS	13861002	13861004	13861006	13861007	13862002	13862004
ENSAFE ID:		13861002	13861004	13861006	13861007	13862002	13862004
LAB ID:		532082	532083	532084	532085	532086	532087
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	4.0000 U	11.0000 U
V 2-Hexanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	34.0000 U	23.0000 U	22.0000 U	32.0000 U	21.0000 U	35.0000 U
V Benzene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	11.0000 U	11.0000 U	11.0000 U,	10.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chloroform	UG/KG	11.0000 U	11.0000 U	2.0000 U	10.0000 U	2.0000 U	2.0000 U
V Chloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	16.0000 U	12.0000 U	9.0000 U	8.0000 U	9.0000 U	12.0000 U
V Styrene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Trichloroethene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S61002	13S61004	13S61006	13S61007	13S62002	13S62004
	LAB ID:	532082	532083	532084	532085	532086	532087
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U	11.0000 U'	10.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13562006	13862007	13863002	13863004	13863006	13864002
ENSAFE ID:		13562006	13862007	13863002	13863004	13863006	13864002
LAB ID:		532088	532089	532090	532091	532092	532093
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
I Aluminum	MG/KG	321.0000	191.0000	118.0000	200.0000	1640.0000	122.0000
I Antimony	MG/KG	11.8000 U	12.1000 U	11.6000 U	11.8000 U	12.0000 U	12.0000 U
I Arsenic	MG/KG	0.6200 U	0.6400 U	0.6100 U	0.6200 U	0.6300 U	0.6300 U
I Barium	MG/KG	1.2000 U	1.3000 U	1.2000 U	1.9000 J	2.6000 J	1.3000 U
I Beryllium	MG/KG	0.6200 U	0.6400 U	0.6100 U	0.6200 U	0.6300 U	0.6300 U
I Cadmium	MG/KG	1.0000 U	1.1000 U	1.0000 U	1.0000 U	1.1000 U	1.1000 U
I Calcium	MG/KG	26.9000 U	27.6000 U	26.4000 U	1780.0000	24100.0000	119.0000 J
I Chromium	MG/KG	2.6000	2.1000 U	2.0000 U	2.1000 U	3.2000	2.9000
I Cobalt	MG/KG	2.5000 U	2.6000 U	2.4000 U	2.5000 U	2.5000 U	2.5000 U
I Copper	MG/KG	5.2000 U	5.3000 U	5.1000 U	5.2000 U	5.3000 U	5.3000 U
I Cyanide	MG/KG	0.5200 U	0.5300 U	0.5100 U	0.5300 U	3.3000	0.5300 U
I Iron	MG/KG	177.0000	101.0000	59.9000	134.0000	1050.0000	57.3000
I Lead	MG/KG	1.1000	1.0000	0.4100 U	1.0000	6.3000	0.5500 J
I Magnesium	MG/KG	114.0000 U	117.0000 U	112.0000 U	114.0000 U	179.0000 J	116.0000 U
I Manganese	MG/KG	1.4000 J	1.3000 U	1.2000 U	1.6000 J	13.6000	1.3000 U
I Mercury	MG/KG	0.1000 U	0.1100 U	0.1000 U	0.1100 U	0.1100 U	0.1100 U
I Nickel	MG/KG	7.9000 U	8.1000 U	7.7000 U	7.9000 U	8.0000 U	8.0000 U
I Potassium	MG/KG	819.0000 U	840.0000 U	803.0000 U	818.0000 U	831.0000 U	830.0000 U
I Selenium	MG/KG	0.6200 U	0.6400 U	0.6100 U	0.6200 U	0.6300 U	0.6300 U
I Silver	MG/KG	2.1000 U	2.1000 U	2.0000 U	2.1000 U	2.1000 U	2.1000 U
I Sodium	MG/KG	107.0000 J	124.0000 J	122.0000 J	130.0000 J	335.0000 J	125.0000 J
I Thallium	MG/KG	0.6200 U	0.6400 U	0.6100 U	0.6200 U	0.6300 U	0.6300 U
I Vanadium	MG/KG	3.1000 U	3.2000 U	3.0000 U	3.1000 U	3.2000 U	3.2000 U
I Zinc	MG/KG	2.5000 J	3.3000 J	1.7000 J	4.5000	5.4000	1.7000 J
P 4,4'-DDD	UG/KG	3.4000 U	3.5000 U	3.3000 U	3.5000 U	3.5000 U	3.5000 U
P 4,4'-DDE	UG/KG	3.4000 U	0.0640 U	0.0810 U	3.5000 U	0.7000 U	0.1900 U
P 4,4'-DDT	UG/KG	3.4000 U	3.5000 U	3.3000 U	3.5000 U	3.5000 U	0.1500 J
P Aldrin	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S62006	13S62007	13S63002	13S63004	13S63006	13S64002
	LAB ID:	532088 R1	532089 R1	532090 R1	532091 R1	532092 R1	532093 R1
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
P Aroclor 1016	UG/KG	34.0000 U	35.0000 U	33.0000 U	35.0000 U	35.0000 U	35.0000 U
P Aroclor 1221	UG/KG	70.0000 U	71.0000 U	68.0000 U	71.0000 U	71.0000 U	71.0000 U
P Aroclor 1232	UG/KG	34.0000 U	35.0000 U	33.0000 U	35.0000 U	35.0000 U	35.0000 U
P Aroclor 1242	UG/KG	34.0000 U	35.0000 U	33.0000 U	35.0000 U	35.0000 U	35.0000 U
P Aroclor 1248	UG/KG	34.0000 U	35.0000 U	33.0000 U	35.0000 U	35.0000 U	35.0000 U
P Aroclor 1254	UG/KG	34.0000 U	35.0000 U	33.0000 U	35.0000 U	35.0000 U	35.0000 U
P Aroclor 1260	UG/KG	34.0000 U	35.0000 U	33.0000 U	35.0000 U	10.0000 J	35.0000 U
P Dieldrin	UG/KG	3.4000 U	3.5000 U	3.3000 U	3.5000 U	0.0520 J	3.5000 U
P Endosulfan I	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U
P Endosulfan II	UG/KG	0.1300 J	3.5000 U	3.3000 U	3.5000 U	3.5000 U	3.5000 U
P Endosulfan sulfate	UG/KG	3.4000 U	3.5000 U	3.3000 U	3.5000 U	0.2400 J	3.5000 U
P Endrin	UG/KG	3.4000 U	3.5000 U	3.3000 U	3.5000 U	3.5000 U	3.5000 U
P Endrin aldehyde	UG/KG	3.4000 U	3.5000 U	3.3000 U	3.5000 U	3.5000 U	3.5000 U
P Endrin ketone	UG/KG	3.4000 U	3.5000 U	3.3000 U	3.5000 U	3.5000 U	3.5000 U
P Heptachlor	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U
P Heptachlor epoxide	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U
P Methoxychlor	UG/KG	18.0000 U	18.0000 U	17.0000 U	18.0000 U	18.0000 U	18.0000 U
P Toxaphene	UG/KG	180.0000 U	180.0000 U	170.0000 U	180.0000 U	180.0000 U	180.0000 U
P alpha-BHC	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U
P alpha-Chlordane	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	0.0550 J	1.8000 U
P beta-BHC	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U
P delta-BHC	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U
P gamma-Chlordane	UG/KG	0.2400 J	0.1400 J	0.0710 U	0.1900 J	0.2000 J	0.2000 J
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S62006	13S62007	13S63002	13S63004	13S63006	13S64002
	LAB ID:	532088	532089	532090	532091	532092	532093
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,4-Dinitrophenol	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Methylphenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 2-Nitroaniline	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 3-Nitroaniline	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S 4-Nitroaniline	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U
S 4-Nitrophenol	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U
S Acenaphthene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Acenaphthylene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Anthracene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Benzo(a)anthracene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Benzo(a)pyrene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key present d at end of Appendix. Data not shown in significan digits.

NAME	UNITS						
	ENSAFE ID:	13862006	13862007	13863002	13863004	13S63006	13864002
	LAB ID:	532088	532089	532090	532091	532092	532093
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
S Butylbenzylphthalate	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	240.0000 U	300.0000 U
S Carbazole	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Chrysene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Di-n-butylphthalate	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Dibenz (a,h) anthracene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Dibenzofuran	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Diethylphthalate	UG/KG	41.0000 U	42.0000 U	42.0000 U	43.0000 U	350.0000 U	340.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Fluoranthene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Fluorene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Hexachloroethane	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Isophorone	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Naphthalene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Nitrobenzene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	830.0000 U	850.0000 U	820.0000 U	840.0000 U	850.0000 U	830.0000 U
S Phenanthrene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Phenol	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S Pyrene	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000,U	350.0000 U	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S bis (2-Ethlyhexy) phthalate	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	320.0000 U	340.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	13662006	13862007	13863002	13663004	13S63006	13864002
	LAB ID:	532088	532089	532090	532091	532092	532093
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 2-Eutanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	40.0000 U	32.0000 U	47.0000 U	26.0000 U	63.0000 U	69.0000 U
V Benzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Chloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Chloroform	UG/KG	1.0000 U	2.0000 U	10.0000 U	1.0000 U	11.0000 U	11.0000 U
V Chloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	17.0000 U	21.0000 U	12.0000 U	16.0000 U	30.0000 U	27.0000 U
V Styrene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Trichloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	13S62006	13S62007	13S63002	13S63004	13S63006	13S64002
	LAB ID:	532088	532089	532090	532091	532092	532093
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S64004	13S64006	13S64008	13S65002	13S65004	13S65006
	LAB ID:	532094	532095	532096	532098	532099	532100
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
I Aluminum	MG/KG	119.0000	2080.0000	9250.0000	80.8000	149.0000	275.0000
I Antimony	MG/KG	12.0000 U	11.8000 U	12.0000 U	12.0000 U	11.7000 U	11.6000 U
I Arsenic	MG/KG	0.6300 U	0.6200 U	2.4000	0.6300 U	0.6100 U	0.6100 U
I Barium	MG/KG	1.3000 U	5.2000 J	6.8000 J	1.3000 U	1.2000 U	1.2000 U
I Beryllium	MG/KG	0.6300 U	0.6200 U	0.6300 U	0.6300 U	0.6100 U	0.6100 U
I Cadmium	MG/KG	1.0000 U	1.0000 U	1.1000 U	1.1000 U	1.0000 U	1.0000 U
I Calcium	MG/KG	47.3000 J	34700.0000	2760.0000	27.4000 U	70.6000 J	71.9000 J
I Chromium	MG/KG	2.1000 U	2.3000	7.4000	2.1000 U	2.0000 U	2.0000 U
I Cobalt	MG/KG	2.5000 U	2.5000 U	2.5000 U	2.5000 U	2.5000 U	2.4000 U
I Copper	MG/KG	5.2000 U	5.2000 U	5.3000 U	5.3000 U	5.1000 U	5.1000 U
I Cyanide	MG/KG	0.5300 U	1.9000	2.7000	0.5300 U	0.5200 U	0.5100 U
I Iron	MG/KG	67.0000	1400.0000	5560.0000	27.7000	69.6000	177.0000
I Lead	MG/KG	0.4200 U	2.6000	11.9000	0.5300 J	0.5700 J	1.8000
I Magnesium	MG/KG	116.0000 U	522.0000 J	135.0000 J	116.0000 U	113.0000 U	112.0000 U
I Manganese	MG/KG	1.3000 U	20.3000	23.2000	1.3000 U	1.2000 U	1.6000 J
I Mercury	MG/KG	0.1100 U	0.1000 U	0.1100 U	0.1100 U	0.1000 U	0.1000 U
I Nickel	MG/KG	8.0000 U	7.9000 U	8.0000 U	8.0000 U	7.8000 U	7.7000 U
I Potassium	MG/KG	829.0000 U	819.0000 U	833.0000 U	833.0000 U	810.0000 U	803.0000 U
I Selenium	MG/KG	0.6300 U	0.6200 U	0.6300 U	0.6300 U	0.6100 U	0.6100 U
I Silver	MG/KG	2.1000 U	2.1000 U	2.1000 U	2.1000 U	2.0000 U	2.0000 U
I Sodium	MG/KG	128.0000 J	425.0000 J	180.0000 J	118.0000 J	95.5000 J	106.0000 J
I Thallium	MG/KG	0.6300 U	0.6200 U	0.6300 U	0.6300 U	0.6100 U	0.6100 U
I Vanadium	MG/KG	3.1000 U	4.1000 J	15.3000	3.2000 U	3.1000 U	3.0000 U
I Zinc	MG/KG	2.2000 J	4.7000	11.3000	0.8500 J	1.8000 J	6.4000
P 4,4'-DDD	UG/KG	3.5000 U	0.3700 J	2.6000 J	3.5000 'U	3.4000 U	3.4000 U
P 4,4'-DDE	UG/KG	3.5000 U	0.5600 U	0.4400 U	3.5000 U	0.0920 U	0.0690 U
P 4,4'-DDT	UG/KG	3.5000 U	2.7000 J	0.1900 J	0.0880 J	0.3600 J	0.1700 J
P Aldrin	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in eignificant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	13964004 / 532094 / 01/15/93	13864006 / 532095 / 01/15/93	13864008 / 532096 / 01/15/93	13865002 / 532098 / 01/15/93	13865004 / 532099 / 01/15/93	13865006 / 532100 / 01/15/93
P Aroclor 1016	UG/KG	35.0000 U	34.0000 U	35.0000 U	35.0000 U	34.0000 U	34.0000 U
P Aroclor 1221	UG/KG	70.0000 U	69.0000 U	71.0000 U	71.0000 U	69.0000 U	68.0000 U
P Aroclor 1232	UG/KG	35.0000 U	34.0000 U	35.0000 U	35.0000 U	34.0000 U	34.0000 U
P Aroclor 1242	UG/KG	35.0000 U	34.0000 U	35.0000 U	35.0000 U	34.0000 U	34.0000 U
P Aroclor 1248	UG/KG	35.0000 U	34.0000 U	35.0000 U	35.0000 U	34.0000 U	34.0000 U
P Aroclor 1254	UG/KG	35.0000 U	34.0000 U	35.0000 U	35.0000 U	34.0000 U	34.0000 U
P Aroclor 1260	UG/KG	35.0000 U	3.7000 J	35.0000 U	35.0000 U	34.0000 U	34.0000 U
P Dieldrin	UG/KG	0.1100 J	1.0000 J	0.2400 J	3.5000 U	0.1200 J	0.1300 J
P Endosulfan I	UG/KG	1.8000 U	0.0580 J	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P Endosulfan II	UG/KG	3.5000 U	3.4000 U	3.5000 U	3.5000 U	3.4000 U	3.4000 U
P Endosulfan sulfate	UG/KG	3.5000 U	3.4000 U	3.5000 U	3.5000 U	3.4000 U	3.4000 U
P Endrin	UG/KG	3.5000 U	3.4000 U	3.5000 U	3.5000 U	3.4000 U	3.4000 U
P Endrin aldehyde	UG/KG	3.5000 U	3.4000 U	0.0960 J	3.5000 U	3.4000 U	3.4000 U
P Endrin ketone	UG/KG	3.5000 U	0.1300 J	3.5000 U	3.5000 U	3.4000 U	3.4000 U
P Heptachlor	UG/KG	0.0740 J	0.0650 J	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P Heptachlor epoxide	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P Methoxychlor	UG/KG	18.0000 U	17.0000 U	18.0000 U	18.0000 U	0.3600 J	17.0000 U
P Toxaphene	UG/KG	180.0000 U	170.0000 U	180.0000 U	180.0000 U	180.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P alpha-Chlordane	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P beta-BHC	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P delta-BHC	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P gamma-Chlordane	UG/KG	1.8000 U	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
S 1,2,4-Trichlorobenzene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 1,2-Dichlorobenzene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13564004	13564006	13S64008	13S65002	13865004	13565006
ENSAFE ID:		13564004	13564006	13S64008	13S65002	13865004	13565006
LAB ID:		532094	532095	532096	532098	532099	532100
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
S 2,4,6-Trichlorophenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2,4-Dichlorophenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2,4-Dimethylphenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2,4-Dinitrophenol	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U
S 2,4-Dinitrotoluene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2,6-Dinitrotoluene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2-Chloronaphthalene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2-Chlorophenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2-Methylnaphthalene	UG/KG	350.0000 U	340.0000 U	86,0000 J	350.0000 U	340.0000 U	340.0000 U
S 2-Methylphenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 2-Nitroaniline	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U
S 2-Nitrophenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 3,3*-Dichlorobenzidine	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 3-Nitroaniline	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U
S 4-Bromophenyl-phenylether	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 4-Chloro-3-methylphenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 4-Chloroaniline	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 4-Methylphenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S 4-Nitroaniline	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U
S 4-Nitrophenol	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U
S Acenaphthene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Acenaphthylene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Anthracene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Benzo(a)anthracene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Benzo(a)pyrene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Benzo(g,h,i)perylene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Kox presented at end of Appendix. Data not shown in significant digits

	ENSAFE ID:	13S54004	13S64006	13S54008	13S65002	13S55004	13S65006
	LAB ID:	532094	532095	532096	532098	532099	532100
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	190.0000 U	280.0000 U	160.0000 U	240.0000 U	280.0000 U	240.0000 U
S Carbazole	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Chrysene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Di-n-butylphthalate	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Di-n-octylphthalate	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Dibenz(a,h)anthracene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Dibenzofuran	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Diethylphthalate	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Dimethylphthalate	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Fluoranthene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Fluorene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Hexachloroethane	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Isophorone	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Naphthalene	UG/KG	350.0000 U	340.0000 U	39.0000 J	350.0000 U	340.0000 U	340.0000 U
S Nitrobenzene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	850.0000 U	830.0000 U	840.0000 U	840.0000 U	820.0000 U	820.0000 U
S Phenanthrene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Phenol	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S Pyrene	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	47.0000 U	36.0000 U	50.0000 U	33.0000 U	89.0000 U	60.0000 U
S bis(2-chloroethyl)ether	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	350.0000 U	340.0000 U	350.0000 U	350.0000 U	340.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13864004	13864006	13864008	13S65002	13865004	13865006
	LAB ID:	532094	532095	532096	532098	532099	532100
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Acetone	UG/KG	58.0000 U	57.0000 U	38.0000 U	20.0000 U	21.0000 U	17.0000 U
V Benzene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/KG	11.0000 U	10.0000 U	1.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	19.0000 U	20.0000 U	9.0000 U	7.0000 U	8.0000 U	7.0000 U
V Styrene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Toluene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Trichloroethene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13864004	13864006	13864008	13865002	13865004	13565006
	LAB ID:	532094	532095	532096	532098	532099	532100
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	WITS						
V Vinyl Chloride	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	10.0000 U	11.0000 U	11.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits

NAME	UNITS	13S65007	13S66002	13S66004	13S66006	13S66008	13S67002
ENSAFE ID:		13S65007	13S66002	13S66004	13S66006	13S66008	13S67002
LAB ID:		532112	532113	532114	532115	532116	532118
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
I Aluminum	MG/KG	650.0000	111.0000	135.0000	382.0000	216.0000	107.0000
I Antimony	MG/KG	11.9000 U	11.9000 U	11.9000 U	12.5000 U	11.9000 U	11.9000 U
I Arsenic	MG/KG	0.6300 U	0.6300 U	0.6800 J	0.6600 U	0.6200 U	0.6200 U
I Barium	MG/KG	1.3000 U	1.3000 U	1.3000 U	2.1000 J	1.2000 U	1.2000 U
I Beryllium	MG/KG	0.6300 U	0.6300 U	0.6300 U	0.6600 U	0.6200 U	0.6200 U
I Cadmium	MG/KG	1.0000 U	1.0000 U	1.0000 U	1.1000 U	1.0000 U	1.0000 U
I Calcium	MG/KG	40.8000 J	175.0000 J	≤36.0000 J	55.0000 J	≤96.0000 J	536.0000 J
I Chromium	MG/KG	2.6000	2.1000 U	2.1000 U	2.4000	2.7000	2.1000 U
I Cobalt	MG/KG	2.5000 U	2.5000 U	2.5000 U	2.6000 U	2.5000 U	2.5000 U
I Copper	MG/KG	5.2000 U	5.2000 U	5.2000 U	5.5000 U	5.2000 U	5.2000 U
I Cyanide	MG/KG	1.3000	0.5200 U	0.5200 U	0.5500 U	0.5200 U	0.5200 U
I Iron	MG/KG	351.0000	104.0000	95.5000	318.0000	173.0000	191.0000
I Lead	MG/KG	2.2000	0.4200 U	0.5600 J	3.1000	1.1000	0.5900 J
I Magnesium	MG/KG	115.0000 U	115.0000 U	115.0000 U	175.0000 J	115.0000 U	115.0000 U
I Manganese	MG/KG	2.6000 J	1.4000 J	1.5000 J	7.0000	1.6000 J	1.6000 J
I Mercury	MG/KG	0.1000 U	0.1000 U	0.1000 U	0.1100 ω	0.1000 U	0.1000 U
I Nickel	MG/KG	7.9000 U	7.9000 U	7.9000 U	8.3000 ω	7.9000 U	7.9000 U
I Potassium	MG/KG	825.0000 U	≥6.0000 U	826.0000 U	≤7.0000 ω	24.0000 U	22.0000 U
I Selenium	MG/KG	0.6300 U	0.6300 U	0.6300 U	0.6600 ω	0.6200 U	0.6200 U
I Silver	MG/KG	2.1000 U	2.1000 U	2.1000 U	2.2000 ω	2.1000 U	2.1000 U
I Sodium	MG/KG	123.0000 J	115.0000 J	121.0000 J	132.0000 J	17.0000 J	113.0000 J
I Thallium	MG/KG	0.6300 U	0.6300 U	0.6300 U	0.6600 ω	0.6200 U	0.6200 U
I Vanadium	MG/KG	3.1000 U	3.1000 U	3.1000 U	3.3000 ω	3.1000 U	3.1000 U
I Zinc	MG/KG	6.4000	3.6000 J	2.7000 J	3.9000 J	2.7000 J	1.3000 J
P 4,4'-DDD	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.6000 U	3.4000 U	0.3000 J
P 4,4'-DDE	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.6000 ω	0.1500 J	3.4000 U
P 4,4'-DDT	UG/KG	3.4000 U	3.4000 U	0.3800 J	3.6000 ω	3.4000 U	0.3900 J
P Aldrin	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.9000 ω	1.8000 U	1.8000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13S65007	13866002	13S66004	13866006	13866008	13867002
ENSAFE ID:		13S65007	13866002	13S66004	13866006	13866008	13867002
LAB ID:		532112	532113	532114	532115	532116	532118
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
P Aroclor 1016	UG/KG	34.0000 U	34.0000 U	34.0000 U	36.0000 U	34.0000 U	34.0000 U
P Aroclor 1221	UG/KG	70.0000 U	70.0000 U	70.0000 U	74.0000 U	70.0000 U	70.0000 U
P Aroclor 1232	UG/KG	34.0000 U	34.0000 U	34.0000 U	36.0000 U	34.0000 U	34.0000 U
P Aroclor 1242	UG/KG	34.0000 U	34.0000 U	34.0000 U	36.0000 U	34.0000 U	34.0000 U
P Aroclor 1248	UG/KG	34.0000 U	34.0000 U	34.0000 U	36.0000 U	34.0000 U	34.0000 U
P Aroclor 1254	UG/KG	34.0000 U	34.0000 U	34.0000 U	36.0000 U	34.0000 U	34.0000 U
P Aroclor 1260	UG/KG	34.0000 U	34.0000 U	34.0000 U	6.4000 J	15.0000 J	34.0000 U
P Dieldrin	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.6000 U	3.4000 U	0.1900 J
P Endosulfan I	UG/KG	0.1300 J	1.8000 U	1.8000 U	1.9000 U	1.8000 U	1.8000 U
P Endosulfan II	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.6000 U	3.4000 U	3.4000 U
P Endosulfan sulfate	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.6000 U	3.4000 U	3.4000 U
P Endrin	UG/KG	3.4000 U	3.4000 U	0.0990 J	3.6000 U	3.4000 U	3.4000 U
P Endrin aldehyde	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.6000 U	3.4000 U	3.4000 U
P Endrin ketone	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.6000 U	3.4000 U	3.4000 U
P Heptachlor	UG/KG	0.0800 U	1.8000 U	0.1600 U	0.1000 U	0.0790 U	0.1000 U
P Heptachlor epoxide	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.9000 U	1.8000 U	1.8000 U
P Methoxychlor	UG/KG	18.0000 U	18.0000 U	18.0000 U	19.0000 U	18.0000 U	18.0000 U
P Toxaphene	UG/KG	180.0000 U	180.0000 U	180.0000 U	190.0000 U	180.0000 U	180.0000 U
P alpha-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.9000 U	1.8000 U	1.8000 U
P alpha-Chlordane	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.9000 U	1.8000 U	1.8000 U
P beta-BHC	UG/KG	1.8000 U	0.1800 J	0.1700 J	1.9000 U	1.8000 U	1.8000 U
P delta-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.9000 U	1.8000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.9000 U	1.8000 U	1.8000 U
P gamma-Chlordane	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.9000 U	1.8000 U	0.0770 U
S 1,2,4-Trichlorobenzene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 1,2-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S65007	13866002	13866004	13866006	13866008	13867002
	LAB ID:	532112	532113	532114	532115	532116	532118
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2,4-Dichlorophenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2,4-Dimethylphenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2,4-Dinitrophenol	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U
S 2,4-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2,6-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2-Chloronaphthalene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2-Chlorophenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2-Methylnaphthalene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2-Methylphenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 2-Nitroaniline	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U
S 2-Nitrophenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 3-Nitroaniline	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U
S 4-Bromophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 4-Chloro-3-methylphenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 4-Chloroaniline	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 4-Methylphenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S 4-Nitroaniline	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U
S 4-Nitrophenol	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U
S Acenaphthene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Acenaphthylene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Anthracene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Benzo(a)anthracene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Benzo(a)pyrene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Benzo(g,h,i)perylene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S65007	13S66002	13S66004	13S66006	13S66008	13S67002
	LAB ID:	532112	532113	532114	532115	532116	532118
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	330.0000 U	330.0000 U	140.0000 U	360.0000 U	340.0000 U	78.0000 U
S Carbazole	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0600 U	340.0000 U
S Chrysene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Di-n-butylphthalate	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Di-n-octylphthalate	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Dibenz(a,h)anthracene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Dibenzofuran	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Diethylphthalate	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Dimethylphthalate	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Fluoranthene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Fluorene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Hexachloroethane	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Isophorone	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Naphthalene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Nitrobenzene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	810.0000 U	810.0000 U	820.0000 U	860.0000 U	830.0000 U	820.0000 U
S Phenanthrene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Phenol	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S Pyrene	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S bis(2-Ethylhexyl)phthalate	UG/KG	82.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S bis(2-chloroethyl)ether	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	330.0000 U	330.0000 U	340.0000 U	360.0000 U	340.0000 U	340.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13865007	13866002	13866004	13S66006	13S66008	13867002
ENSAFE ID:		13865007	13866002	13866004	13S66006	13S66008	13867002
LAB ID:		532112	532113	532114	532115	532116	532118
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	lo.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 u	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Acetone	UG/KG	50.0000 U	28.0000 U	34.0000 U	50.0000 U	120.0000 U	53.0000 U
V Benzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/KG	1.0000 U	1.0000 U	1.0000 U	11.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	13.0000 U	13.0000 U	10.0000 U	11.0000 U	21.0000 U	27.0000 U
V Styrene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 'u	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13365007	13366002	13866004	13866006	13866008	13867002
	LAB ID:	532112	532113	532114	532115	532116	532118
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	11.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:		13867004	13867006	13S67008	13868002	13868004	13S68005
	LAB ID:		532119	532120	532121	532967	532969	532970
	DATE RCVD:		01/15/93	01/15/93	01/15/93	01/20/93	01/20/93	01/20/93
NAME	WITS							
I Aluminum		MG/KG	114.0000	130.0000	171.0000	95.8000	105.0000	104.0000
I Antimony		MG/KG	11.9000 U	11.8000 U	11.9000 U	11.9000 U	12.0000 U	12.1000 U
I Arsenic		MG/KG	0.6300 U	0.6200 U	0.6300 U	0.6200 U	0.6300 U	0.6300 U
I Barium		MG/KG	1.3000 U	1.2000 U	1.3000 U	1.2000 U	1.3000 U	1.3000 U
I Beryllium		MG/KG	0.6300 U	0.6200 U	0.6300 U	0.6200 U	0.6300 U	0.6300 U
I Cadmium		MG/KG	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.1000 U	1.1000 U
I Calcium		MG/KG	135.0000 J	282.0000 J	195.0000 J	103.0000 J	97.1000 J	81.7000 J
I Chromium		MG/KG	2.1000 U	2.3000	3.4000	2.1000 U	2.1000 U	2.1000 U
I Cobalt		MG/KG	2.5000 u	2.5000 U	2.5000 U	2.5000 U	2.5000 U	2.5000 U
I Copper		MG/KG	5.2000 U	5.2000 U	5.2000 U	5.2000 U	5.3000 U	5.3000 U
I Cyanide		MG/KG	0.5200 U	0,5200 U	0.5200 U	0.5200 U	0.5300 U	0.5300 U
I Iron		MG/KG	83.5000	84.5000	174.0000	67.6000	202.0000	68.0000
I Lead		MG/KG	0.6200 J	0.8300	1.3000	0.4200 U	0.4200 U	0.5600 J
I Magnesium		MG/KG	115.0000 U	114.0000 U	115.0000 U	115.0000 U	116.0000 U	117.0000 U
I Manganese		MG/KG	1.5000 J	1.2000 U	1.6000 J	1.2000 U	1.3000 U	1.3000 U
I Mercury		MG/KG	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1100 U	0.1100 U
I Nickel		MG/KG	7.9000 U	7.8000 U	7.9000 U	7.9000 U	8.0000 U	8.0000 U
I Potassium		MG/KG	827.0000 U	815.0000 U	824.0000 U	824.0000 U	830.0000 U	837.0000 U
I Selenium		MG/KG	0.6300 U	0.6200 U	0.6300 U	0.6200 U	0.6300 U	0.6300 U
I Silver		MG/KG	2.1000 U	2.1000 U	2.1000 U	2.1000 U	2.1000 U	2.1000 U
I Sodium		MG/KG	118.0000 J	104.0000 J	113.0000 J	102.0000 J	113.0000 J	108.0000 J
I Thallium		MG/KG	0.6300 U	0.6200 U	0.6300 U	0.6200 U	0.6300 U	0.6300 U
I Vanadium		MG/KG	3.1000 U	3.1000 U	3.1000 U	3.1000 U	3.2000 U	3.2000 U
I Zinc		MG/KG	1.5000 J	2.7000 J	4.5000	1.6000 J	1.4000 J	1.7000 J
P 4,4'-DDD		UG/KG	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U
P 4,4'-DDE		UG/KG	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U
P 4,4'-DDT		UG/KG	0.1900 J	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U
P Aldrin		UG/KG	1.8000 U	1.8000 U	1.8000 U	0.0410 J	1.8000 u	1.7000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13S67004	13S67006	13S67008	13S68002	13S68004	13S68005
ENSAFE ID:		13S67004	13S67006	13S67008	13S68002	13S68004	13S68005
LAB ID:		532119	532120	532121	532967	532969	532970
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/20/93	01/20/93	01/20/93
P Aroclor 1016	UG/KG	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U
P Aroclor 1221	UG/KG	70.0000 U	69.0000 U	70.0000 U	70.0000 U	70.0000 U	69.0000 U
P Aroclor 1232	UG/KG	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U
P Aroclor 1242	UG/KG	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U
P Aroclor 1248	UG/KG	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U
P Aroclor 1254	UG/KG	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U
P Aroclor 1260	UG/KG	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U	34.0000 U
P Dieldrin	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.4000 U	0.0990 J	0.0620 J
P Endosulfan I	UG/KG	0.0890 J	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P Endosulfan II	UG/KG	3.4000 U	3.4000 U	0.1500 J	3.4000 U	3.4000 U	3.4000 U
P Endosulfan sulfate	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U
P Endrin	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U
P Endrin aldehyde	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U
P Endrin ketone	UG/KG	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U	3.4000 U
P Heptachlor	UG/KG	1.8000 U	1.8000 U	1.8000 U	0.0700 U	1.8000 U	0.0550 U
P Heptachlor epoxide	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P Methoxychlor	UG/KG	18.0000 U	18.0000 U	18.0000 U	18.0000 U	18.0000 U	17.0000 U
P Toxaphene	UG/KG	180.0000 U	180.0000 U	180.0000 U	180.0000 U	180.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P alpha-Chlordane	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P beta-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P delta-BHC	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
P gamma-Chlordane	UG/KG	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.8000 U	1.7000 U
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4,5-Trichlorophenol	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S67004	13S67006	13S67008	13S68002	13S68004	13568005
	LAB ID:	532119	532120	532121	532967	532969	532970
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,4-Dinitrophenol	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Methylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 2-Nitroaniline	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 3-Nitroaniline	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S 4-Nitroaniline	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U
S 4-Nitrophenol	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U
S Acenaphthene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Acenaphthylene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Anthracene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(a)anthracene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(a)pyrene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S67004	13S67006	13S67008	13S68002	13S68004	13868005
	LAB ID:	532119	532120	532121	532967	532969	532970
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	36.0000 U	43.0000 U	340.0000 U	80.0000 U	350.0000 U	94.0000 U
S Carbazole	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Chrysene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Di-n-butylphthalate	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Dibenz (a,h) anthracene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Dibenzofuran	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Diethylphthalate	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Fluoranthene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Fluorene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Hexachloroethane	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Isophorone	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Naphthalene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0600 U	350.0000 U
S Nitrobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Pentachlorophenol	UG/KG	820.0000 U	810.0000 U	830.0000 U	820.0000 U	840.0000 U	850.0000 U
S Phenanthrene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Phenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S Pyrene	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13567004	13S67006	13S67008	13868002	13568004	13S68005
	LAB ID:	532119	532120	532121	532967	532969	532970
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	62.0000 U	99.0000 U	78.0000 U	44.0000 U	35.0000 U	40.0000 U
V Benzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromofo m	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chlorofo m	UG/KG	10.0000 U	10.0000 U	1.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	23.0000 U	21.0000 U	18.0000 U	12.0000 U	13.0000 U	11.0000 U
V Styrene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S67004	13S67006	13S67008	13S68002	13S68004	13S68005
	LAB ID:	532119	532120	532121	532967	532969	532970
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S69002	13S69004	13S70002	13S70004	33S01001	33S01002
	LAB ID:	532985	532987	532962	532979		
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	12/10/92	12/10/92
NAME	UNITS						
I Aluminum	MG/KG	109.0000	108.0000	635.0000	116.0000	3000.0000	1280.0000
I Antimony	MG/KG	11.8000 U	11.9000 U	11.8000 U	12.2000 U	13.0000 U	21.8000 U
I Arsenic	MG/KG	0.6200 U	0.6300 U	0.6200 U	0.6400 U	0.9200 J	3.2000 J
I Barium	MG/KG	1.2000 U	1.3000 U	1.2000 U	1.3000 U	4.5000 J	2.6000 J
I Beryllium	MG/KG	0.6200 U	0.6300 U	0.6200 U	0.6400 U	0.4600 U	0.7600 U
I Cadmium	MG/KG	1.0000 U	1.0000 U	1.0000 U	1.1000 U	1.1000 U	6.8000
I Calcium	MG/KG	637.0000 J	53.4000 J	43.6000 J	36.0000 J	1650.0000	1170.0000 J
I Chromium	MG/KG	2.1000 U	2.1000 U	2.1000 U	2.1000 U	7.3000	289.0000
I Cobalt	MG/KG	2.5000 U	2.5000 U	2.5000 U	2.6000 U	2.5000 U	4.2000 U
I Copper	MG/KG	16.7000	5.2000 U	5.2000 U	5.4000 U	5.7000 U	9.5000 U
I Cyanide	MG/KG	0.5200 U	0.5200 U	0.5200 U	0.5400 U	7.5000	169.0000
I Iron	MG/KG	140.0000	63.0000	410.0000	41.9000	1670.0000	250.0000
I Lead	MG/KG	1.3000	1.8000	2.8000	0.4500 J	8.1000	12.5000
I Magnesium	MG/KG	361.0000 J	116.0000 U	115.0000 U	119.0000 U	186.0000 J	162.0000 U
I Manganese	MG/KG	2.5000 J	1.5000 J	1.6000 J	1.3000 U	9.1000	15.3000
I Mercury	MG/KG	0.1000 U	0.1000 U	0.1000 U	0.1100 U	0.1100 U	0.1900 U
I Nickel	MG/KG	7.8000 U	8.0000 U	7.9000 U	8.2000 U	9.1000 U	15.3000 U
I Potassium	MG/KG	816.0000 U	829.0000 U	821.0000 U	849.0000 U	1100.0000 U	1850.0000 U
I Selenium	MG/KG	0.6200 U	0.6300 U	0.6200 U	0.6400 U	0.9100 U	1.5000 U
I Silver	MG/KG	2.1000 U	2.1000 U	2.1000 U	2.1000 U	2.3000 U	3.8000 U
I Sodium	MG/KG	112.0000 J	73.1000 J	128.0000 J	129.0000 J	119.0000 J	197.0000 J
I Thallium	MG/KG	0.6200 U	0.6300 U	0.6200 U	0.6400 U	0.4600 U	0.7600 U
I Vanadium	MG/KG	3.1000 U	3.1000 U	3.1000 U	3.2000 U	4.6000 J	7.3000 U
I Zinc	MG/KG	2.2000 J	1.3000 J	1.2000 J	1.6000 J	7.8000	12.7000
P 4,4'-DDD	UG/KG	3.4000 U	3.5000 U	3.4000 U	3.5000 U	3.7000 U	1.2000 J
P 4,4'-DDE	UG/KG	3.4000 U	1.4000 J	0.2600 J	3.5000 U	2.5000 J	19.0000 J
P 4,4'-DDT	UG/KG	3.4000 U	1.4000 J	3.4000 U	3.5000 U	1.0000 J	46.0000 J
P Aldrin	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.9000 U	6.5000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13869002	13869004	13870002	13870004	33601001	33S01002
	LAB ID:	532985	532987	532962	532979		
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	12/10/92	12/10/92
NAME	UNITS						
P Aroclor 1016	UG/KG	34.0000 U	35.0000 U	34.0000 U	35.0000 U	37.0000 U	130,0000 U
P Aroclor 1221	UG/KG	69.0000 U	72.0000 U	70.0000 U	71.0000 U	76.0000 U	260.0000 U
P Aroclor 1232	UG/KG	34.0000 U	35.0000 U	34.0000 U	35.0000 U	37.0000 U	130.0000 U
P Aroclor 1242	UG/KG	34.0000 U	35.0000 U	34.0000 U	35.0000 U	37.0000 U	130.0000 U
P Aroclor 1248	UG/KG	34.0000 U	35.0000 U	34.0000 U	35.0000 U	37.0000 U	130.0000 U
P Aroclor 1254	UG/KG	34.0000 U	35.0000 U	34.0000 U	35.0000 U	37.0000 U	130.0000 U
P Aroclor 1260	UG/KG	34.0000 U	3.5000 J	34.0000 U	35.0000 U	20.0000 J	130.0000 U
P Dieldrin	UG/KG	0.0730 J	3.5000 U	3.4000 U	3.5000 U	4.8000	18.0000 J
P Endosulfan I	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.9000 U	6.5000 U
P Endosulfan II	UG/KG	3.4000 U	3.5000 U	3.4000 U	3.5000 U	3.7000 U	2.2000 J
P Endosulfan sulfate	UG/KG	3.4000 U	3.5000 U	3.4000 U	3.5000 U	3.7000 U	110.0000 J
P Endrin	UG/KG	3.4000 U	3.5000 U	3.4000 U	3.5000 U	3.7000 U	27.0000 J
P Endrin aldehyde	UG/KG	3.4000 U	3.5000 U	3.4000 U	3.5000 U	0.4600 J	21.0000
P Endrin ketone	UG/KG	3.4000 U	3.5000 U	3.4000 U	3.5000 U	3.7000 U	13.0000 U
P Heptachlor	UG/KG	1.7000 U	0.2000 U	1.8000 U	1.8000 U	1.9000 U	6.8000 U
P Heptachlor epoxide	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.9000 U	26.0000 J
P Methoxychlor	UG/KG	17.0000 U	18.0000 U	18.0000 U	18.0000 U	0.2400 J	25.0000 J
P Toxaphene	UG/KG	170.0000 U	180.0000 U	180.0000 U	180.0000 U	190.0000 U	650.0000 U
P alpha-BHC	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.9000 U	2.6000 J
P alpha-Chlordane	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	0.1800 J	7.2000 J
P beta-BHC	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.9000 U	6.5000 U
P delta-BHC	UG/KG	1.7000 U	1.8000 U	0.0760 U	1.8000 U	0.1500 J	1.2000 J
P gamma-BHC (Lindane)	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.9000 U	6.5000 U
P gamma-Chlordane	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U	1.1000 J	6.5000 U
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	17000.0000 J
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	4200.0000 J
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	10000.0000 J
S 2,4,5-Trichlorophenol	UG/KG	810.0000 U	830.0000 U	820.0000 U	850.0000 U	900.0000 U	46000.0000 U

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13869002	13869004	13870002	13870004	33801061	33801002
ENSAFE ID:		13869002	13869004	13870002	13870004	33801061	33801002
LAB ID:		532985	532987	532962	532979		
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/20/93	12/10/92	12/10/92
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2,4-Dinitrophenol	UG/KG	810.0000 U	830.0000 U	820.0000 U	850.0000 U	900.0000 U	46000.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	36000.0000
S 2-Methylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 2-Nitroaniline	UG/KG	810.0000 U	830.0000 U	820.0000 U	850.0000 U	900.0000 U	46000.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 3-Nitroaniline	UG/KG	810.0000 U	830.0000 U	820.0000 U	850.0000 U	900.0000 U	46000.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	810.0000 U	830.0000 U	820.0000 U	850.0000 U	900.0000 U	46000.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S 4-Nitroaniline	UG/KG	810.0000 U	830.0000 U	820.0000 U	850.0000 U	900.0000 U	46000.0000 U
S 4-Nitrophenol	UG/KG	810.0000 U	830.0000 U	820.0000 U	850.0000 U	900.0000 U	46000.0000 U
S Acenaphthene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	3900.0000 J
S Acenaphthylene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S Anthracene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	4100.0000 J
S Benzo(a)anthracene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S Benzo(a)pyrene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	370.0000 U	19000.0000 U
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	55.0000 J	19000.0000 U
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	340.0000 U	340.0000 U	350.0000 U	51.0000 J	19000.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S69002	13S69004	13S70002	13S70004	33S01001	33S01002
	LAB ID:	532983	532987	532962	532979		
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	12/10/92	12/10/92
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	150.0000 J	110.0000 J	140.0000 J	120.0000 U	370.0000 U	19000.0000 U
S Carbazole	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Chrysene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Di-n-butylphthalate	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	72.0000 U	19000.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Dibenz(a,h)anthracene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Dibenzofuran	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Diethylphthalate	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Dimethylphthalate	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Fluoranthene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	33.0000 J	14000.0000 J
S Fluorene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	5200.0000 J
S Hexachlorobenzene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Hexachloroethane	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Isophorone	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Naphthalene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	26000.0000
S Nitrobenzene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Pentachlorophenol	UG/KG	810.0000 J	830.0000 J	820.0000 J	850.0000 U	900.0000 U	46000.0000 U
S Phenanthrene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	8200.0000 J
S Phenol	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S Pyrene	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	2400.0000 J
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	43.0000 U	43000.0000
S bis(2-chloroethyl)ether	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 J	340.0000 J	340.0000 J	350.0000 U	370.0000 U	19000.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13869002	13869004	13870002	13870004	33501001	33801002
ENSAFE ID:		13869002	13869004	13870002	13870004	33501001	33801002
LAB ID:		532985	532987	532962	532979		
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/20/93	12/10/92	12/10/92
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 2-Butanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 2-Hexanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Acetone	UG/KG	26.0000 U	130.0000 U	25.0000 U	40.0000 U	12.0000 U	54.0000 U
V Benzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Bromoform	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Bromomethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Chlorobenzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	7.0000 J	110.0000
V Chlotoethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Chloroform	UG/KG	10.0000 U	3.0000 U	10.0000 U	11.0000 U	2.0000 U	2.0000 U
V Chloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Ethylbenzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Methylene Chloride	UG/KG	8.0000 U	7.0000 U	20.0000 U	13.0000 U	19.0000 U	79.0000 U
V Styrene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Toluene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Trichloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13869002	13S69004	13S70002	13870004	33S01001	33501002
	LAB ID:	532985	532987	532962	532979		
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	12/10/92	12/10/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V Xylene (total)	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	10.0000 U	11.0000 U	11.0000 U	19.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33502002	33S02002DL	33S02002MS	33S02002MSD	33S02002RE	33902003
	LAB ID:						
	DATE RCM:	12/10/92	12/10/92	12/18/92	12/18/92	12/10/92	12/10/92
NAME	UNITS						
I Aluminum	MG/KG	1630.0000			.	.	434.0000
I Antimony	MG/KG	12.0000 U			.	.	13.4000 U
I Arsenic	MG/KG	0.8400 U			.	.	0.9400 U
I Barium	MG/KG	9.0000 J			.	.	1.7000 J
I Beryllium	MG/KG	0.4200 U			.	.	0.4700 U
I Cadmium	MG/KG	2.4000			.	.	1.2000 U
I Calcium	MG/KG	724.0000 J			.	.	308.0000 J
I Chromium	MG/KG	71.4000			.	.	8.3000
I Cobalt	MG/KG	2.3000 U			.	.	2.6000 U
I Copper	MG/KG	7.6000			.	.	5.9000 U
I Cyanide	MG/KG	0.5200 U			.	.	0.5900 U
I Iron	MG/KG	772.0000			.	.	161.0000
I Lead	MG/KG	12.3000			.	.	6.1000
I Magnesium	MG/KG	89.6000 U			.	.	100.0000 U
I Manganese	MG/KG	6.3000			.	.	2.1000 J
I Mercury	MG/KG	0.1300			.	.	0.1200 U
I Nickel	MG/KG	8.4000 U			.	.	9.4000 U
I Potassium	MG/KG	1020.0000 U			.	.	1140.0000 U
I Selenium	MG/KG	0.8400 U			.	.	0.9400 U
I Silver	MG/KG	2.4000 U			.	.	2.4000 U
I Sodium	MG/KG	119.0000 J			.	.	112.0000 J
I Thallium	MG/KG	0.4200 U			.	.	0.4700 U
I Vanadium	MG/KG	4.0000 U			.	.	4.5000 U
I Zinc	MG/KG	19.2000			.	.	4.4000 J
P 4,4'-DDD	UG/KG	6.9000 U	69.0000 U		.	.	3.9000 U
P 4,4'-DDE	UG/KG	160.0000 J	200.0000 J		.	.	7.6000
P 4,4'-DDT	UG/KG	56.0000	54.0000 J		.	.	2.6000 J
P Aldrin	UG/KG	3.5000 U	35.0000 U		.	.	2.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33802002	33S02002DL	33S02002MS	33S02002MSD	33S02002RE	33502003
		12/10/92	12/10/92	12/18/92	12/18/92	12/10/92	12/10/92
P Aroclor 1016	UG/KG	69.0000 U	690.0000 U	.	.	.	39.0000 U
P Aroclor 1221	UG/KG	140.0000 U	1400.0000 U	.	.	.	78.0000 U
P Aroclor 1232	UG/KG	69.0000 U	690.0000 U	.	.	.	39.0000 U
P Aroclor 1242	UG/KG	69.0000 U	690.0000 U	.	.	.	39.0000 U
P Aroclor 1248	UG/KG	69.0000 U	690.0000 U	.	.	.	39.0000 U
P Aroclor 1254	UG/KG	69.0000 U	690.0000 U	.	.	.	39.0000 U
P Aroclor 1260	UG/KG	69.0000 U	690.0000 U	.	.	.	39.0000 U
P Dieldrin	UG/KG	5.0000 J	5.2000 J	.	.	.	1.5000 J
P Endosulfan I	UG/KG	3.5000 U	35.0000 U	.	.	.	2.0000 U
P Endosulfan II	UG/KG	6.9000 U	69.0000 U	.	.	.	3.9000 U
P Endosulfan sulfate	UG/KG	1.3000 J	69.0000 U	.	.	.	3.9000 U
P Endrin	UG/KG	6.9000 U	3.9000 J	.	.	.	3.9000 U
P Endrin aldehyde	UG/KG	6.9000 U	69.0000 U	.	.	.	3.9000 U
P Endrin ketone	UG/KG	6.9000 U	69.0000 U	.	.	.	3.9000 U
P Heptachlor	UG/KG	3.5000 U	35.0000 U	.	.	.	0.1200 J
P Heptachlor epoxide	UG/KG	7.8000 J	4.0000 J	.	.	.	0.0760 J
P Methoxychlor	UG/KG	35.0000 U	350.0000 U	.	.	.	20.0000 U
P Toxaphene	UG/KG	350.0000 U	3500.0000 U	.	.	.	200.0000 U
P alpha-BHC	UG/KG	3.5000 U	35.0000 U	.	.	.	2.0000 U
P alpha-Chlordane	UG/KG	130.0000 J	170.0000 J	.	.	.	2.8000
P beta-BHC	UG/KG	3.5000 U	35.0000 U	.	.	.	2.0000 U
P delta-BHC	UG/KG	3.5000 U	35.0000 U	.	.	.	2.0000 U
P gamma-BHC (Lindane)	UG/KG	3.5000 U	35.0000 U	.	.	.	2.0000 U
P gamma-Chlordane	UG/KG	100.0000 J	130.0000 J	.	.	.	1.9000 J
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U		.	.	.	380.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U		.	.	.	380.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U		.	.	.	380.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U		.	.		380.0000 U
S 2,4,5-Trichlorophenol	UG/KG	830.0000 U		.	.	.	930.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33S02002	33S02002DL	33S02002MS	33S02002MSD	33S02002RE	33802003
		12/10/92	12/10/92	12/18/92	12/18/92	12/10/92	12/10/92
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2,4-Dinitrophenol	UG/KG	830.0000 U	.	.	.	.	930.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2-Methylphenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 2-Nitroaniline	UG/KG	830.0000 U		.	.	.	930.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 3-Nitroaniline	UG/KG	830.0000 U	.	.	.	.	930.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	830.0000 U	.	.	.	.	930.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S 4-Nitroaniline	UG/KG	830.0000 U	.	.	.	.	930.0000 U
S 4-Nitrophenol	UG/KG	830.0000 U	.	.	.	.	930.0000 U
S Acenaphthene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S Acenaphthylene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S Anthracene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S Benzo(a)anthracene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S Benzo(a)pyrene	UG/KG	340.0000 U	.	.	.	.	380.0000 U
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	.		.	.	380.0000 U
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	.	.	.	.	380.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / WITS	33902002	33S02002DL	33S02002MS	33S02002MSD	33S02002RE	33S02003
		12/10/92	12/10/92	12/18/92	12/18/92	12/10/92	12/10/92
S Butylbenzylphthalate	UG/KG	340.0000 U		.	.		380.0000 U
S Carbazole	UG/KG	340.0000 U		.	.		380.0000 U
S Chrysene	UG/KG	340.0000 U	.	.			380.0000 U
S Di-n-butylphthalate	UG/KG	84.0000 U	.	.	.		63.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	.	.	.		380.0000 U
S Dibenz(a,h)anthracene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Dibenzofuran	UG/KG	340.0000 U	.	.	.		380.0000 U
S Diethylphthalate	UG/KG	340.0000 U	.	.	.		380.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	.	.	.		380.0000 U
S Fluoranthene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Fluorene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Hexachloroethane	UG/KG	340.0000 U	.	.	.		380.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Ieophorone	UG/KG	340.0000 U	.	.	.		380.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	.	.	.		380.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	.	.	.		380.0000 U
S Naphthalene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Nitrobenzene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Pentachlorophenol	UG/KG	830.0000 U	.	.	.		930.0000 U
S Phenanthrene	UG/KG	340.0000 U	.	.	.		380.0000 U
S Phenol	UG/KG	340.0000 U	.	.	.		380.0000 U
S Pyrene	UG/KG	340.0000 U	.	.	.		380.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	.	.	.		380.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	44.0000 U	.	.	.	. .	41.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	.	.	.		380.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	.	.	.		380.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S02002	33S02002DL	33S02002MS	33S02002MSD	33S02002RE	33S02003
	LAB ID:						
	DATE RCVD:	12/10/92	12/10/92	12/18/92	12/18/92	12/10/92	12/10/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	.	11000.0000	11000.0000	1300.0000 U	12.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 2-Butanone	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1100.0000 U	12.0000 U
V 2-Hexanone	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Acetone	UG/KG	2000.0000 J	.	9100.0000 U	9500.0000 U	11000.0000 U	530.0000 U
V Benzene	UG/KG	11.0000 U	.	7600.0000	7400.0000	1300.0000 U	12.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Bromoform	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Bromomethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Chlorobenzene	UG/KG	11.0000 U	.	6400.0000	6500.0000	1300.0000 U	12.0000 U
V Chloroethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Chloroform	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	2.0000 U
V Chloromethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Ethylbenzene	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Methylene Chloride	UG/KG	14.0000 U	.	780.0000 U	640.0000 U	780.0000 U	19.0000 U
V Styrene	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	.	1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V Toluene	UG/KG	11.0000 U	.	6700.0000	6900.0000	1300.0000 U	12.0000 U
V Trichloroethene	UG/KG	11.0000 U	.	7000.0000	6800.0000	1300.0000 U	12.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S02002	33S02002DL	33S02002MS	33S02002MSD	33S02002RE	33S02003
	LAB ID:						
	DATE RCVD:	12/10/92	12/10/92	12/18/92	12/18/92	12/10/92	12/10/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U		1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V xylene (total)	UG/KG	11.0000 U		1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U		1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U		1300.0000 U	1300.0000 U	1300.0000 U	12.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID:	33S02003RE	33S03002	33503004	33S04002	33804004	33505002
	LAB ID:						528712
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92
	UNITS						
I Aluminum	MG/KG		1650.0000	20100.0000	5120.0000	169.0000	2350.0000
I Antimony	MG/KG		13.5000 U	19.2000 U	12.4000 U	13.1000 U	12.0000 U
I Arsenic	MG/KG		2.2000 J	7.7000	1.3000 J	0.9200 U	0.8400 U
I Barium	MG/KG		12.6000 J	127.0000	6.4000 J	0.8100 J	2.4000 J
I Beryllium	MG/KG		0.4800 U	0.6700 U	0.4300 U	0.4600 U	0.4200 U
I Cadmium	MG/KG		6.5000	93.4000	1.6000	1.1000 U	1.1000 U
I Calcium	MG/KG		635.0000 J	14900.0000	1080.0000 J	807.0000 J	382.0000 J
I Chromium	MG/KG		174.0000	6180.0000	31.4000	2.3000 U	4.5000
I Cobalt	MG/KG		2.6000 U	3.7000 U	2.4000 U	2.5000 U	2.3000 U
I Copper	MG/KG		16.4000	196.0000	5.4000 U	5.7000 U	5.3000 U
I Cyanide	MG/KG		0.5900 U	11.9000	0.5400 U	0.5700 U	0.5200 U
I Iron	MG/KG		969.0000	3370.0000	2340.0000	70.5000	1210.0000
I Lead	MG/KG		44.0000	461.0000	9.6000	0.6700 J	6.8000
I Magnesium	MG/KG		101.0000 U	575.0000 J	92.1000 U	97.7000 U	89.3000 U
I Manganese	MG/KG		20.0000	66.7000	136.0000	4.8000	38.3000
I Mercury	MG/KG		0.3100	1.6000	0.1100 U	0.1100 U	0.1100 U
I Nickel	MG/KG		10.1000	83.3000	8.7000 U	9.2000 U	8.4000 U
I Potassium	MG/KG		1150.0000 U	1640.0000 U	1050.0000 U	1120.0000 U	1020.0000 U
I Selenium	MG/KG		0.9500 U	1.3000 U	0.8700 U	0.9200 U	0.8400 U
I Silver	MG/KG		2.6000	42.9000	2.2000 U	2.3000 U	2.1000 U
I Sodium	MG/KG		123.0000 J	282.0000 J	125.0000 J	117.0000 J	110.0000 J
I Thallium	MG/KG		0.4800 U	0.6700 U	0.4300 U	0.4600 U	0.4200 U
I Vanadium	MG/KG		4.5000 U	12.7000 J	7.3000 J	4.4000 U	4.0000 U
I Zinc	MG/KG		32.7000	352.0000	7.9000	1.4000 J	6.4000
P 4,4'-DDD	UG/KG		3.9000 U	5.5000 U	3.6000 U	3.8000 U	3.5000 U
P 4,4'-DDE	UG/KG		9.9000	25.0000	2.8000 J	0.0910 J	1.0000 J
P 4,4'-DDT	UG/KG		1.9000 J	8.6000 J	1.8000 J	3.8000 U	0.3700 J
P Aldrin	UG/KG		2.0000 U	0.3700 J	1.8000 U	1.9000 U	0.0840 J

NAS PENSACOLA OU-10 ANI) SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID:	33S02003RE	33803002	33903004	33804002	33804004	33805002
	LAB ID:						528712
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92
	WITS						
P Aroclor 1016	UG/KG		39.0000 U	55.0000 U	36.0000 U	38.0000 U	35.0000 U
P Atoclor 1221	UG/KG	.	79.0000 U	110.0000 U	73.0000 U	77.0000 U	70.0000 U
P Aroclor 1232	UG/KG		39.0000 U	55.0000 U	36.0000 U	38.0000 U	35.0000 U
P Aroclor 1242	UG/KG		39.0000 U	55.0000 U	36.0000 U	38.0000 U	35.0000 U
P Aroclor 1248	UG/KG	.	39.0000 U	55.0000 U	36.0000 U	38.0000 U	35.0000 U
P Aroclor 1254	UG/KG	.	39.0000 U	55.0000 U	36.0000 U	38.0000 U	35.0000 U
P Aroclor 1260	UG/KG	.	39.0000 U	55.0000 U	36.0000 U	38.0000 U	35.0000 U
P Dieldrin	UG/KG	.	5.2000 J	38.0000 J	0.4300 J	3.8000 U	3.5000 U
P Endosulfan I	UG/KG	.	2.0000 U	2.9000 U	1.8000 U	1.9000 U	1.8000 U
P Endosulfan II	UG/KG	.	0.4700 J	8.3000 J	3.6000 U	3.8000 U	3.5000 U
P Endosulfan sulfate	UG/KG	.	3.9000 U	0.8800 J	3.6000 U	3.8000 U	0.1600 J
P Endrin	UG/KG	.	0.9800 J	15.0000	0.1300 J	3.8000 U	3.5000 U
P Endrin aldehyde	UG/KG	.	3.9000 U	23.0000 J	3.6000 U	3.8000 U	3.5000 U
P Endrin ketone	UG/KG	.	3.9000 U	0.2300 J	3.6000 U	3.8000 U	3.5000 U
P Heptachlor	UG/KG	.	2.0000 U	2.9000 U	0.1000 U	0.1500 J	0.2500 U
P Heptachlor epoxide	UG/KG	.	2.0000 U	5.4000 J	0.2200 J	1.9000 U	1.8000 U
P Methoxychlor	UG/KG	.	20.0000 U	5.8000 J	18.0000 U	19.0000 U	18.0000 U
P Toxaphene	UG/KG	.	200.0000 U	290.0000 U	180.0000 U	190.0000 U	180.0000 U
P alpha-BHC	UG/KG	.	2.0000 U	2.9000 U	1.8000 U	1.9000 U	1.8000 U
P alpha-Chlordane	UG/KG	.	3.5000	20.0000	0.3100 J	0.1200 J	1.8000 U
P beta-BHC	UG/KG	.	2.0000 U	2.9000 U	1.8000 U	1.9000 U	1.8000 U
P delta-BHC	UG/KG	.	0.2500 J	1.3000 J	1.8000 U	1.9000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	.	2.0000 U	2.9000 U	1.8000 U	1.9000 U	1.8000 U
P gamma-Chlordane	UG/KG	.	4.3000	26.0000 J	1.8000 U	1.9000 U	0.0640 J
S 1,2,4-Trichlorobenzene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 1,2-Dichlorobenzene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 'U	'370.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	.	950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S02003RE	33S03002	33S03004	33S04002	33S04004	33S05002
	LAB ID:						528712
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2,4-Dichlorophenol	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2,4-Dimethylphenol	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2,4-Dinitrophenol	UG/KG	.	950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U
S 2,4-Dinitrotoluene	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2,6-Dinitrotoluene	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2-Chloronaphthalene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2-Chlorophenol	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2-Methylnaphthalene	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2-Methylphenol	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 2-Nitroaniline	UG/KG		950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U
S 2-Nitrophenol	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 3,3'-Dichlorobenzidine	UG/KG		390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 3-Nitroaniline	UG/KG	.	950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG		950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U
S 4-Bromophenyl-phenylether	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 4-Chloro-3-methylphenol	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 4-Chloroaniline	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 4-Methylphenol	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S 4-Nitroaniline	UG/KG	.	950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U
S 4-Nitrophenol	UG/KG	.	950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U
S Acenaphthene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Acenaphthylene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Anthracene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Benzo(a)anthracene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Benzo(a)pyrene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Benzo(g,h,i)perylene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S02003RE	33803002	33S03004	33S04002	33804004	33805002
	LAB ID:						528712
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	.	390.0000 U	1800.0000 U	360.0000 U	370.0000 U	340.0000 U
S Carbazole	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Chrysene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Di-n-butylphthalate	UG/KG	.	89.0000 U	1200.0000 U	53.0000 U	66.0000 U	340.0000 U
S Di-n-octylphthalate	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Dibenz(a,h)anthracene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Dibenzofuran	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Diethylphthalate	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Dimethylphthalate	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Fluoranthene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Fluorene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Hexachloroethane	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Isophorone	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Naphthalene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Nitrobenzene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	.	950.0000 U	27000.0000 U	860.0000 U	900.0000 U	830.0000 U
S Phenanthrene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Phenol	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S Pyrene	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S bis(2-Ethylhexy)phthalate	UG/KG	.	47.0000 U	3800.0000 U	40.0000 U	370.0000 U	40.0000 U
S bis(2-chloroethyl)ether	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	.	390.0000 U	11000.0000 U	360.0000 U	370.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S02003RE	33S03002	33S03004	33S04002	33S04004	33S05002
	LAB ID:						528712
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	810.0000	15.0000 U	120.0000 U	27.0000 U	11.0000 U	22.0000 U
V Benzene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Chloroethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Chloroform	UG/KG	59.0000 U	12.0000 U	4.0000 U	1.0000 U	2.0000 U	11.0000 U
V Chloromethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Ethylbeneene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	43.0000 U	19.0000 U	52.0000 U	37.0000 U	51.0000 U	12.0000 U
V Styrene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Trichloroethene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S02003RE	33S03002	33S03004	33S04002	33S04004	33S05002
	LAB ID:						528712
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	59.0000 U	12.0000 U	17.0000 U	11.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S05004	33S06002	33S07002	33S08002	33S08004	33S09002
	LAB ID:	528715	528730	528721		528718	528693
	DATE RCVD:	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92
NAME	UNITS						
I Aluminum	MG/KG	2440.0000	1450.0000	1570.0000	989.0000	3950.0000	992.0000
I Aluminum	MG/KG	2440.0000	133.0000	1570.0000	989.0000	3950.0000	992.0000
I Antimony	MG/KG	13.9000 U	12.8000 U	12.6000 U	11.7000 U	12.0000 U	12.1000 U
I Arsenic	MG/KG	1.3000 J	0.6400 U	0.6300 U	0.8200 U	0.8400 U	0.8500 U
I Barium	MG/KG	2.2000 J	13.6000 J	26.6000 J	9.9000 J	15.7000 J	2.2000 J
I Beryllium	MG/KG	0.4900 U	0.2100 U	0.2100 U	0.4100 U	0.4200 U	0.4200 U
I Cadmium	MG/KG	1.2000 U	1.1000 U	1.1000 U	4.5000	6.7000	1.1000 U
I Calcium	MG/KG	225.0000 U	239.0000 J	1520.0000	534.0000 J	944.0000 J	4620.0000
I Chromium	MG/KG	20.8000	1.8000 J	69.7000	85.0000	81.8000	2.1000 U
I Cobalt	MG/KG	2.7000 U	1.9000 U	1.9000 U	2.3000 U	2.3000 U	2.3000 U
I Copper	MG/KG	6.1000 U	2.7000 J	7.7000	11.4000	18.0000	5.3000 U
I Cyanide	MG/KG	0.6100 U	0.5300 U	0.5300 U	0.5100 U	0.5200 U	0.5300 U
I Iron	MG/KG	1170.0000	940.0000	684.0000	1420.0000	2440.0000	563.0000
I Lead	MG/KG	7.5000	15.6000	9.5000	14.6000	45.1000	5.8000
I Magnesium	MG/KG	103.0000 U	58.4000 J	61.3000 J	87.4000 U	98.6000 J	90.0000 U
I *Manganese*	MG/KG	4.4000	6.6000	7.5000	14.7000	14.4000	4.8000
I Mercury	MG/KG	0.1200 U	0.1100 U	0.1100 U	0.1000	0.1500	0.1100 U
I Nickel	MG/KG	9.7000 U	2.3000 U	2.3000 U	8.2000 U	9.8000	8.5000 U
I Potassium	MG/KG	1180.0000 U	290.0000 U	286.0000 U	999.0000 U	1020.0000 U	1030.0000 U
I Selenium	MG/KG	0.9700 U	0.6400 U	0.6300 U	0.8200 U	0.8400 U	0.8500 U
I Silver	MG/KG	2.4000 U	1,3000 U	1.2000 U	2.1000 U	2.1000 U	2.1000 U
I Sodium	MG/KG	137.0000 J	91.9000 J	89.3000 J	115.0000 J	120.0000 J	146.0000 J
I Thallium	MG/KG	0.4900 U	0.4300 U	0.4200 U	0.4100 U	0.4200 U	0.4200 U
I Vanadium	MG/KG	4.6000 U	1.9000 J	1.4000 J	3.9000 U	6.2000 J	4.0000 U
I Zinc	MG/KG	2.4000 J	12.4000	12.3000	25.2000	49.9000	3.1000 J
P 4,4'-DDD	UG/KG	4.0000 U	0.0990 J	3.5000 U	3.4000 U	3.5000 U	3.5000 U
P 4,4'-DDE	UG/KG	1.2000 J	0.3200 J	3.5000 U	0.1600 J	0.5900 J	0.0900 J
P 4,4'-DDT	UG/KG	0.3400 U	3.5000 J	0.2000 U	0.9400 U	3.5000 J	0.0790 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S05004	33S06002	33807002	33808002	33S08004	33909002
ENSAFE ID:		33S05004	33S06002	33807002	33808002	33S08004	33909002
LAB ID:		528715	528730	528721		528718	528693
DATE RCVD:		12/12/92	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92
P Aldrin	UG/KG	2.0000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P Aroclor 1016	UG/KG	40.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1221	UG/KG	81.0000 U	71.0000 U	70.0000 U	69.0000 U	70.0000 U	71.0000 U
P Aroclor 1232	UG/KG	40.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1242	UG/KG	40.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1248	UG/KG	40.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclor 1254	UG/KG	40.0000 U	35.0000 U	35.0000 U	34.0000 U	35.0000 U	35.0000 U
P Aroclot 1260	UG/KG	40.0000 U	35.0000 U	35.0000 U	34.0000 U	13.0000 J	35.0000 U
P Dieldrin	UG/KG	4.0000 U	3.5000 U	0.1600 J	5.1000	1.8000 J	0.0620 J
P Endosulfan I	UG/KG	2.0000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P Endosulfan II	UG/KG	4.0000 U	3.5000 U	3.5000 U	3.4000 U	3.5000 U	0.1000 J
P Endosulfan sulfate	UG/KG	4.0000 U	3.5000 U	3.5000 U	0.1300 J	3.5000 U	3.5000 U
P Endrin	UG/KG	4.0000 U	3.5000 U	3.5000 U	0.2700 J	3.5000 U	3.5000 U
P Endrin aldehyde	UG/KG	4.0000 U	3.5000 U	3.5000 U	3.3000 J	3.5000 U	3.5000 U
P Endrin ketone	UG/KG	4.0000 U	0.2200 J	3.5000 U	3.4000 U	3.5000 U	3.5000 U
P Heptachlor	UG/KG	0.1600 U	0.2300 U	0.2000 U	1.7000 U	0.1700 U	0.1600 U
P Heptachlor epoxide	UG/KG	2.0000 U	0.1000 U	1.8000 U	0.2000 J	1.8000 U	1.8000 U
P Methoxychlor	UG/KG	20.0000 U	18.0000 U	18.0000 U	0.3300 J	18.0000 U	18.0000 U
P Toxaphene	UG/KG	200.0000 U	180.0000 U	180.0000 U	170.0000 U	180.0000 U	180.0000 U
P alpha-BHC	UG/KG	2.0000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P alpha-Chlordane	UG/KG	2.0000 U	1.8000 U	1.8000 U	1.8000 J	1.1000 J	1.8000 U
P beta-BHC	UG/KG	2.0000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P delta-BHC	UG/KG	2.0000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	2.0000 U	1.8000 U	1.8000 U	1.7000 U	1.8000 U	1.8000 U
P gamma-Chlordane	UG/KG	2.0000 U	1.8000 U	0.1000 J	1.7000 U	1.4000 J	1.8000 U
S 1,2,4-Trichlorobenzene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 1,2-Dichlorobenzene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 1,3-Dichlorobenzene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 1,4-Dichlorobenzene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33805004 / 528715 / 12/12/92	33806002 / 528730 / 12/12/92	33807002 / 528721 / 12/12/92	33S08002 / 12/12/92	33808004 / 528718 / 12/12/92	33809002 / 528693 / 12/12/92
S 2,4,5-Trichlorophenol	UG/KG	960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S 2,4,6-Trichlorophenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2,4-Dichlorophenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2,4-Dimethylphenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2,4-Dinitrophenol	UG/KG	960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S 2,4-Dinitrotoluene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2,6-Dinitrotoluene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2-Chloronaphthalene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2-Chlorophenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2-Methylnaphthalene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2-Methylphenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 2-Nitroaniline	UG/KG	960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S 2-Nitrophenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 3-Nitroaniline	UG/KG	960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S 4-Bromophenyl-phenylether	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 4-Chloro-3-methylphenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 4-Chloroaniline	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 4-Methylphenol	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S 4-Nitroaniline	UG/KG	960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S 4-Nitrophenol	UG/KG	960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S Acenaphthene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Acenaphthylene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Anthracene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Benzo(a)anthracene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Benzo(a)pyrene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Benzo(b,k)fluoranthene	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME		ENSAFE ID:	33S05004	33806002	33807002	33S08002	33SO0004	33809002
		LAB ID:	528715	528730	520721		520710	520693
		DATE RCVD:	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92
		UNITS						
S Benzo(g,h,i)perylene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Butylbenzylphthalate	UG/KG		400.0000 U	350.0000 U	340.0000 U	38.0000 U	38.0000 U	350.0000 U
S Carbazole	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Chrysene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Di-n-butylphthalate	UG/KG		400.0000 U	350.0000 U	340.0000 U	77.0000 U	340.0000 U	350.0000 U
S Di-n-octylphthalate	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Dibenz (a,h)anthracene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Dibenzofuran	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Diethylphthalate	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Dimethylphthalate	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Fluoranthene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Fluorene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Hexachlorobenzene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Hexachlorobutadiene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Hexachlorocyclopentadiene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Hexachloroethane	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Isophorone	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S N-Nitroso-di-n-propylamine	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S N-Nitrosodiphenylamine	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Naphthalene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Nitrobenzene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Pentachlorophenol	UG/KG		960.0000 U	840.0000 U	830.0000 U	820.0000 U	830.0000 U	840.0000 U
S Phenanthrene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Phenol	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S Pyrene	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S bis(2-Chloroethoxy)methane	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG		41.0000 U	81.0000 U	39.0000 U	74.0000 U	66.0000 U	36.0000 U
S bis(2-chloroethyl)ether	UG/KG		400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S05004	33S06002	33S07002	33S08002	33S08004	33809002
	LAB ID:	528715	528730	528721		528718	528693
	DATE RCVD:	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92
NAME	UNITS						
S bis(2-chloroisopropyl)ether	UG/KG	400.0000 U	350.0000 U	340.0000 U	340.0000 U	340.0000 U	350.0000 U
V 1,1,1-Trichloroethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	44.0000 U	38.0000 U	71.0000 U	1600.0000 U	180.0000 U	89.0000 U
V Benzene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Chloroethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Chlorofonn	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Chloromethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	20.0000 U	34.0000 U	26.0000 U	740.0000 U	16.0000 U	13.0000 U
V Styrene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S05004	33S06002	33S07002	33S08002	33S08004	33S09002
	LAB ID:	528715	528730	528721		528718	528693
	DATE RCVD:	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92	12/12/92
NAME	UNITS						
V Trichloroethene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Vinyl Chloride	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	12.0000 U	11.0000 U	11.0000 U	1200.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33810002	33S10004	33S11002	33511004	33512002	33S12002RE
	LAB ID:	528736	528733	529539	529537	530135	530135
	DATE RCVD:	12/12/92	12/12/92	12/17/92	12/17/92	12/19/92	12/19/92
NAME	UNITS						
I Aluminum	MG/KG	2760.0000	2770.0000	511.0000	983.0000	2270.0000	.
I Antimony	MG/KG	13.2000 U	13.5000 U	12.5000 U	12.7000 U	11.8000 U	.
I Arsenic	MG/KG	0.8300 J	0.6700 U	0.6300 U	0.6300 U	0.6200 U	.
I Barium	MG/KG	5.9000 J	6.5000 J	1.4000 J	2.8000 J	16.2000 J	.
I Beryllium	MG/KG	0.2200 U	0.2200 U	0.2100 U	0.2100 U	0.4100 U	.
I Cadmium	MG/KG	2.2000	2.3000	1.0000 U	1.1000 U	1.0000 U	.
I Calcium	MG/KG	456.0000 J	750.0000 J	385.0000 J	393.0000 J	191.0000 U	.
I Chromium	MG/KG	83.9000	90.6000	3.6000	1.7000 J	8.5000	.
I Cobalt	MG/KG	2.0000 U	2.0000 U	1.9000 U	1.9000 U	2.3000 U	.
I Copper	MG/KG	23.6000	7.9000	1.0000 U	1.9000 J	5.2000 U	.
I Cyanide	MG/KG	0.5500 U	0.5600 U	0.5200 U	0.5300 U	0.5200 U	.
I Iron	MG/KG	1480.0000	1390.0000	366.0000	424.0000	1140.0000	.
I Lead	MG/KG	18.9000	20.6000	2.8000	5.4000	48.4000	.
I Magnesium	MG/KG	71.4000 J	81.1000 J	50.6000 J	33.8000 J	88.6000 J	.
I Manganese	MG/KG	5.3000	5.4000	6.4000	2.7000 J	4.3000	.
I Mercury	MG/KG	0.1100 U	0.1100 U	0.1000 U	0.1100 U	0.1000 U	.
I Nickel	MG/KG	6.6000 J	3.6000 J	2.3000 U	2.3000 U	8.2000 U	.
I Potassium	MG/KG	299.0000 U	306.0000 U	285.0000 U	288.0000 U	1000.0000 U	.
I Selenium	MG/KG	0.6600 U	0.6700 U	0.6300 U	0.6300 U	0.6200 U	.
I Silver	MG/KG	1.3000 U	1.4000 U	1.3000 U	1.3000 U	2.1000 U	.
I Sodium	MG/KG	84.7000 J	92.8000 J	81.2000 J	79.9000 J	95.7000 J	.
I Thallium	MG/KG	0.4400 U	0.4500 U	0.4200 U	0.4200 U	0.6200 U	.
I Vanadium	MG/KG	3.2000 J	3.8000 J	1.0000 U	1.2000 J	3.9000 U	.
I Zinc	MG/KG	28.1000	24.1000	5.4000	4.9000,	13.8000	.
P 4,4'-DDD	UG/KG	3.6000 U	0.3800 J	0.1400 J	5.3000		.
P 4,4'-DDE	UG/KG	0.6600 J	4.1000	0.7800 J	9.6000	.	.
P 4,4'-DDT	UG/KG	1.1000 U	3.7000 U	1.2000 U	0.4800 U	.	.
P Aldrin	UG/KG	1.8000 U	1.9000 U	0.0820 J	0.2300 J	.	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33810002	33810004	33811002	33811004	33612002	33812002RE
ENSAFE ID:		33810002	33810004	33811002	33811004	33612002	33812002RE
LAB ID:		528736	528733	529539	529537	530135	530135
DATE RCVD:		12/12/92	12/12/92	12/17/92	12/17/92	12/19/92	12/19/92
P Aroclor 1016	UG/KG	36.0000 U	37.0000 U	34.0000 U	34.0000 U	.	.
P Aroclor 1221	UG/KG	72.0000 U	75.0000 U	70.0000 U	70.0000 U	.	.
P Aroclor 1232	UG/KG	36.0000 U	37.0000 U	34.0000 U	34.0000 U	.	.
P Aroclor 1242	UG/KG	36.0000 U	37.0000 U	34.0000 U	34.0000 U	.	.
P Aroclor 1248	UG/KG	36.0000 U	37.0000 U	34.0000 U	34.0000 U	.	.
P Aroclor 1254	UG/KG	36.0000 U	37.0000 U	34.0000 U	34.0000 U	.	.
P Aroclor 1260	UG/KG	36.0000 U	17.0000 J	34.0000 U	34.0000 U	.	.
P Dieldrin	UG/KG	9.5000	9.8000	0.4400 J	0.3000 J	.	.
P Endosulfan I	UG/KG	1.8000 U	1.9000 U	1.8000 U	1.8000 U	.	.
P Endosulfan II	UG/KG	0.4500 J	3.7000 U	3.4000 U	3.4000 U	.	.
P Endosulfan sulfate	UG/KG	3.6000 U	3.7000 U	0.5200 J	1.4000 J	.	.
P Endrin	UG/KG	3.6000 U	3.7000 U	3.4000 U	3.4000 U	.	.
P Endrin aldehyde	UG/KG	0.6100 J	3.7000 U	3.4000 U	0.3100 J	.	.
P Endrin ketone	UG/KG	0.4900 J	0.2500 U	3.4000 U	3.4000 U	.	.
P Heptachlor	UG/KG	0.3200 U	0.3100 U	0.0810 U	1.8000 U	.	.
P Heptachlor epoxide	UG/KG	0.1600 J	1.9000 U	1.8000 U	1.8000 U	.	
P Methoxychlor	UG/KG	18.0000 U	19.0000 U	18.0000 U	18.0000 U	.	.
P Toxaphene	UG/KG	180.0000 U	190.0000 U	180.0000 U	180.0000 U	.	
P alpha-BHC	UG/KG	0.2200 J	0.1000 J	1.8000 U	1.8000 U	.	.
P alpha-Chlordane	UG/KG	1.8000 U	0.8500 J	1.8000 U	1.8000 U	.	.
P beta-BHC	UG/KG	1.8000 U	1.9000 U	1.8000 U	1.8000 U	.	.
P delta-BHC	UG/KG	1.8000 U	1.9000 U	1.8000 U	1.8000 U	.	.
P gamma-BHC (Lindane)	UG/KG	2.0000 J	1.9000 U	1.8000 U	1.8000 U	.	.
P gamma-Chlordane	UG/KG	0.4900 J	1.5000 J	1.8000 U	1.8000 U		.
S 1,2,4-Trichlorobenzene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 1,2-Dichlorobenzene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 1,3-Dichlorobenzene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 1,4-Dichlorobenzene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2,4,5-Trichlorophenol	UG/KG	870.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key present d at end of Appendix. Data not shown in significan digits.

	ENSAFE ID:	33810002	33810004	33811002	33811004	33812002	33S12002RE
	LAB ID:	528736	528733	529539	529537	530135	530135
	DATE RCVD:	12/12/92	12/12/92	12/17/92	12/17/92	12/19/92	12/19/92
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2,4-Dichlorophenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2,4-Dimethylphenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2,4-Dinitrophenol	UG/KG	870.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.
S 2,4-Dinitrotoluene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2,6-Dinitrotoluene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2-Chloronaphthalene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2-Chlorophenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2-Methylnaphthalene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2-Methylphenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 2-Nitroaniline	UG/KG	870.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.
S 2-Nitrophenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 3,3'-Dichlorobenzidine	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 3-Nitroaniline	UG/KG	870.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/KG	870.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.
S 4-Bromophenyl-phenylether	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 4-Chloro-3-methylphenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 4-Chloroaniline	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 4-Chlorophenyl-phenylether	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 4-Methylphenol	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S 4-Nitroaniline	UG/KG	870.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.
S 4-Nitrophenol	UG/KG	870.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.
S Acenaphthene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Acenaphthylene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Anthracene	UG/KG	360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Benzo(a)anthracene	UG/KG	360.0000 U	120.0000 J	340.0000 U	130.0000 J	340.0000 U	.
S Beneo(a)pyrene	UG/KG	360.0000 U	140.0000 J	340.0000 U	160.0000 J	340.0000 U	.
S Benzo(b,k)fluoranthene	UG/KG	57.0000 J	360.0000 J	62.0000 J	360.0000 J	59.0000 J	.
S Benzo(g,h,i)perylene	UG/KG	120.0000 J	280.0000 J	340.0000 U	350.0000 U	340.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME		ENSAFE ID:	33810002	33810004	33811002	33511004	33812002	33S12002RE
		LAB ID:	528736	528733	529539	529537	530135	530135
		DATE RCVD:	12/12/92	12/12/92	12/17/92	12/17/92	12/19/92	12/19/92
		UNITS						
S Butylbenzylphthalate	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Carbazole	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Chrysene	UG/KG		360.0000 U	100.0000 J	340.0000 U	140.0000 J	340.0000 U	.
S Di-n-butylphthalate	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Di-n-octylphthalate	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Dibenz (a,h) anthracene	UG/KG		360.0000 U	69.0000 J	340.0000 U	350.0000 U	340.0000 U	.
S Dibenzofuran	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Diethylphthalate	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Dimethylphthalate	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Fluoranthene	UG/KG		360.0000 U	160.0000 J	39.0000 J	140.0000 J	58.0000 J	.
S Fluorene	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Hexachlorobenzene	UG/KG		360,0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Hexachlorobutadiene	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Hexachlorocyclopentadiene	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Hexachloroethane	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Indeno(1,2,3-cd) pyrene	UG/KG		44.0000 J	190.0000 J	340.0000 U	260.0000 J	340.0000 U	.
S Isophorone	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S N-Nitroso-di-n-propylamine	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S N-Nitrosodiphenylamine	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Naphthalene	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Nitrobenzene	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Pentachlorophenol	UG/KG		55.0000 U	890.0000 U	830.0000 U	840.0000 U	820.0000 U	.
S Phenanthrene	UG/KG		360.0000 U	82.0000 J	36.0000 J	39.0000 J	340.0000 U	.
S Phenol	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S Pyrene	UG/KG		41.0000 J	190.0000 J	56.0000 J	150.0000 J	53.0000 J	.
S bis(2-Chloroethoxy)methane	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S bis(2-Ethylhexy)phthalate	UG/KG		240.0000 U	300.0000 U	40.0000 U	48.0000 U	41.0000 U	.
S bis(2-chloroethyl) ether	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.
S bis(2-chloroisopropyl)ether	UG/KG		360.0000 U	370.0000 U	340.0000 U	350.0000 U	340.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33510002	33810004	33S11002	33S11004	33512002	33S12002RE
ENSAFE ID:		33510002	33810004	33S11002	33S11004	33512002	33S12002RE
LAB ID:		528736	528733	529539	529537	530135	530135
DATE RCVD:		12/12/92	12/12/92	12/17/92	12/17/92	12/19/92	12/19/92
V 1,1,1-Trichloroethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 1,1,2-Trichloroethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 1,1-Dichloroethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 1,1-Dichloroethene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 1,2-Dichloroethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 1,2-Dichloroethene (total)	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 1,2-Dichloropropane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 2-Butanone	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 2-Hexanone	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V 4-Methyl-2-Pentanone	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Acetone	UG/KG	1000.0000 U	37.0000 U	20.0000 U	38.0000 U	560.0000 U	710.0000 U
V Benzene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Bromodichloromethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Bromoform	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Bromomethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Carbon Disulfide	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Carbon Tetrachloride	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Chlorobenzene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Chloroethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Chloroform	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Chloromethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Dibtomochloromethane	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Ethylbenzene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Methylene Chloride	UG/KG	190.0000 U	49.0000 U	30.0000 U	30.0000 U	5.0000 U	52.0000 U
V Styrene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Tetrachloroethene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Toluene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Trichloroethene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S10002	33510004	33511002	33511004	33912002	33S12002RE
	LAB ID:	528736	528733	529539	529537	530135	530135
	DATE RCVD:	12/12/92	12/12/92	12/17/92	12/17/92	12/19/92	12/19/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V Xylene (total)	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V cis-1,3-Dichloropropene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U
V trans-1,3-Dichloropropene	UG/KG	55.0000 U	11.0000 U	10.0000 U	11.0000 U	10.0000 U	34.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33812004	33S12004MS	33S12004MSD	33812006	33S12006RE	33812008
	LAB ID:	530144	530145	530146	530132	530132	530117
	DATE RCM:	12/19/92	02/29/92	02/29/92	12/19/92	12/19/92	12/19/92
NAME	UNITS						
I Aluminum	MG/KG	1990.0000			1660.0000	.	1780.0000
I Antimony	MG/KG	12.0000 u			11.8000 u	.	12.2000 u
I Arsenic	MG/KG	0.6300 U			0.6200 U	.	0.6400 U
I Barium	MG/KG	7.4000 J			3.2000 J	.	2.8000 J
I Beryllium	MG/KG	0.4200 U			0.4100 U	.	0.4300 U
I Cadmium	MG/KG	1.0000 U			1.0000 U	.	1.1000 U
I Calcium	MG/KG	194.0000 U			191.0000 U	.	198.0000 u
I Chromium	MG/KG	28.9000			73.8000	.	67.8000
I Cobalt	MG/KG	2.3000 U			2.3000 U	.	2.4000 U
I Copper	MG/KG	5.2000 u			5.2000 U	.	5.4000 U
I Cyanide	MG/KG	0.5200 U			0.5200 U	.	0.5400 U
I Iron	MG/KG	1110.0000			1040.0000	.	1040.0000
I Lead	MG/KG	9.3000			6.4000	.	5.0000
I Magnesium	MG/KG	89.2000 U			87.8000 U	.	91.1000 U
I Manganese	MG/KG	4.3000			2.2000 J	.	2.6000 J
I Mercury	MG/KG	0.1000 u			0.1000 u	.	0.1100 u
I Nickel	MG/KG	8.4000 U			8.3000 U	.	8.6000 U
I Potassium	MG/KG	1020.0000 u			1000.0000 U	.	1040.0000 U
I Selenium	MG/KG	0.6300 U			0.6200 U	.	0.6400 U
I Silver	MG/KG	2.1000 u			2.1000 u	.	2.1000 U
I *Sodium*	MG/KG	92.1000 J			106.0000 J	.	81.3000 J
I Thallium	MG/KG	0.6300 U			0.6200 U	.	0.4300 U
I Vanadium	MG/KG	4.0000 U			3.9000 U	.	4.1000 U
I Zinc	MG/KG	21.3000			8.8000,	.	4.5000
P 4,4'-DDD	UG/KG	3.0000 U	*3.0000* U	3.0000 U	3.4000 U	.	3.5000 U
P 4,4'-DDE	UG/KG	0.2000 J	0.2000 J	0.1000 J	3.4000 U	.	3.5000 U
P 4,4'-DDT	UG/KG	3.0000 U	3.0000	3.0000	3.4000 U	.	3.5000 U
P Aldrin	UG/KG	1.0000 u	1.0000	1.0000	1.7000 U	.	0.4100 J

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33512004	33S12004MS	33S12004MSD	33512006	33S12006RE	33512008
	LAB ID:	530144	530151	530152	530132	530132	530117
	DATE RCVD:	12/23/09	12/23/09	12/23/09	12/19/92	12/19/92	12/19/92
NAME	UNITS						
P Aroclor 1016	UG/KG	3.0000 U	3.0000 U	3.0000 U	34.0000 U	.	35.0000 U
P Aroclor 1221	UG/KG	7.0000 U	7.0000 U	7.0000 U	69.0000 U	.	72.0000 U
P Aroclor 1232	UG/KG	3.0000 U	3.0000 U	3.0000 U	34.0000 U	.	35.0000 U
P Aroclor 1242	UG/KG	3.0000 U	3.0000 U	3.0000 U	34.0000 U	.	35.0000 U
P Aroclor 1248	UG/KG	3.0000 U	3.0000 U	3.0000 U	34.0000 U	.	35.0000 U
P Aroclor 1254	UG/KG	3.0000 U	3.0000 U	3.0000 U	34.0000 U	.	35.0000 U
P Aroclor 1260	UG/KG	5.0000 J	6.0000 J	5.0000 J	16.0000 J	.	8.4000 J
P Dieldrin	UG/KG	2.0000 J	3.0000	2.0000	4.5000 J	.	1.9000 J
P Endosulfan I	UG/KG	1.0000 U	1.0000 U	1.0000 U	1.7000 U	.	1.8000 U
P Endosulfan II	UG/KG	3.0000 U	3.0000 U	3.0000 U	3.4000 U	.	3.5000 U
P Endosulfan sulfate	UG/KG	0.2000 U	0.0900 U	3.0000 U	0.1200 J	.	3.5000 U
P Endrin	UG/KG	3.0000 U	3.0000	3.0000	3.4000 U	.	3.5000 U
P Endrin aldehyde	UG/KG	3.0000 U	3.0000 U	3.0000 U	0.8100 J	.	0.4800 J
P Endrin ketone	UG/KG	3.0000 U	3.0000 U	3.0000 U	0.2900 J	.	3.5000 U
P Heptachlor	UG/KG	1.0000 U	1.0000	1.0000	1.7000 U	.	2.1000 J
P Heptachlor epoxide	UG/KG	1.0000 U	1.0000 U	1.0000 U	1.7000 U	.	1.8000 U
P Methoxychlor	UG/KG	1.0000 U	1.0000 U	1.0000 U	17.0000 U	.	18.0000 U
P Toxaphene	UG/KG	8.0000 U	18.0000 U	18.0000 U	170.0000 U	.	180.0000 U
P alpha-BHC	UG/KG	1.0000 U	1.0000 U	1.0000 U	1.7000 U	.	1.8000 U
P alpha-Chlordane	UG/KG	1.0000 U	1.0000 U	1.0000 U	1.7000 U	.	0.1200 J
P beta-BHC	UG/KG	0.8000 U	0.6000 U	1.0000 U	1.7000 U	.	1.8000 U
P delta-BHC	UG/KG	1.0000 U	1.0000 U	1.0000 U	1.7000 U	.	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.0000 U	1.0000	1.0000	1.7000 U	.	1.8000 U
P gamma-Chlordane	UG/KG	0.9000 J	0.9000 J	0.9000 J	0.7800 J	.	1.1000 J
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	900.0000	700.0000	340.0000 U	.	350.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	500.0000	400.0000	340.0000 U	.	350.0000 U
S 2,4,5-Trichlorophenol	UG/KG	820.0000 U	840.0000 U	850.0000 U	820.0000 U	.	860.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S12004	33S12004MS	33S12004MSD	33S12006	33S12006RE	33812008
	LAB ID:	530144	530148	530149	530132	530132	530117
	DATE RCVD:	12/19/92	12/29/92	12/29/92	12/19/92	12/19/92	12/19/92
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 2,4-Dinitrophenol	UG/KG	820.0000 U	840.0000 U	850.0000 U	820.0000 U	.	860.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	500.0000	400.0000	340.0000 U	.	350.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	700.0000	400.0000	340.0000 U	.	350.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 2-Methylphenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 2-Nitroaniline	UG/KG	820.0000 U	840.0000 U	850.0000 U	820.0000 U	.	860.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 3-Nitroaniline	UG/KG	820.0000 U	840.0000 U	850.0000 U	820.0000 U	.	860.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	820.0000 U	840.0000 U	850.0000 U	820.0000 U	.	860.0000 U
S 4-Brmophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	300.0000 J	100.0000 J	340.0000 U	.	350.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S 4-Nitroaniline	UG/KG	820.0000 U	840.0000 U	850.0000 U	820.0000 U	.	860.0000 U
S 4-Nitrophenol	UG/KG	820.0000 U	300.0000 J	200.0000 J	820.0000 U	.	860.0000 U
S Acenaphthene	UG/KG	340.0000 U	800.0000	600.0000	340.0000 U	.	350.0000 U
S Acenaphthylene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Anthracene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Benzo(a)anthracene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Bento(a)pyrene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33812004	33S12004MS	33S12004MSD	33512006	33S12006RE	33S12008
	LAB ID:	530144	530148	530149	530132	530132	530117
	DATE RCVD:	12/19/92	12/29/92	12/29/92	12/19/92	12/19/92	12/19/92
NAME	UNITS						
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Butylbenzylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	100.0000 U	.	350.0000 U
S Carbazole	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Chrysene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Di-n-butylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Dibenz (a,h) anthracene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Dibenzofuran	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Diethylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Fluoranthene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Fluorene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Hexachloroethane	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S fndeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Isophorone	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	600.0000	400.0000	340.0000 U	.	350.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Naphthalene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Nitrobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Pentachlorophenol	UG/KG	820.0000 U	700.0000 J	500.0000 J	820.0000 U	.	860.0000 U
S Phenanthrene	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S Phenol	UG/KG	340.0000 U	900.0000 J	600.0000 J	340.0000 U	.	350.0000 U
S Pyrene	UG/KG	340.0000 U	900.0000	800.0000	340.0000 'U	.	350.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	60.0000 U	46.0000 J	56.0000 U	43.0000 U	.	37.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U	.	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	33S12004	33S12004MS	33S12004MSD	33512006	33S12006RE	33S12008
	LAB ID:	530144	530148	530149	530132	530132	530117
	DATE RCVD:	12/19/92	12/29/92	12/29/92	12/19/92	12/19/92	12/19/92
NAME	UNITS						
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	350.0000 U	350.0000 U	340.0000 U		350.0000 U
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	62.0000	67.0000	10.0000 U	34.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 2-Butanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 2-Hexanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Acetone	UG/KG	110.0000 U	6.0000 U	11.0000 U	620.0000 U	850.0000 U	110.0000 U
V Benzene	UG/KG	11.0000 U	58.0000	57.0000	10.0000 U	34.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Bromoform	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Bromomethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Chlorobenzene	UG/KG	11.0000 U	55.0000	56.0000	10.0000 U	34.0000 U	11.0000 U
V Chloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Chloroform	UG/KG	11.0000 U	11.0000 U	2.0000 U	10.0000 U	34.0000 U	11.0000 U
V Chloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Ethylbeneene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Methylene Chloride	UG/KG	6.0000 U	25.0000 U	44.0000 U	6.0000 U	49.0000 U	5.0000 U
V Styrene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Toluene	UG/KG	11.0000 U	57.0000	59.0000	10.0000 U	34.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33912004	33S12004MS	33S12004MSD	33S12006	33S12006RE	33912008
	LAB ID:	530144	530145	530146	530132	530132	530117
	DATE RCVD:	12/19/92	02/29/92	02/29/92	12/19/92	12/19/92	12/19/92
NAME	UNITS						
V Trichloroethene	UG/KG	11.0000 U	51.0000	53.0000	10.0000 U	34.0000 U	11.0000 U
V Vinyl Chloride	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V Xylene (total)	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	34.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD:	UNITS	33812010 / 530126 / 12/19/92	33S12010RE / 530126 / 12/19/92	33813002 / 546852 / 04/21/93	33S13004 / 546855 / 04/21/93	33813006 / 546859 / 04/21/93	33814002 / 529896 / 12/18/92
I Aluminum		MG/KG	1790.0000	.	2280.0000 J	656.0000 J	497.0000	3600.0000
I Antimony		MG/KG	11.9000 U	.	9.6000 U	10.5000 U	11.5000 U	12.4000 U
I Arsenic		MG/KG	0.6300 U	.	0.6300 U	0.6800 U	0.7500 U	1.7000 J
I Barium		MG/KG	3.4000 J	.	5.8000 J	1.4000 J	1.2000 U	7.3000 J
I Beryllium		MG/KG	0.4200 U	.	0.4200 U	0.4500 U	0.5000 U	0.4400 U
I Cadmium		MG/KG	1.0000 U	.	1.0000 U	1.1000 U	1.2000 U	1.1000 U
I Calcium		MG/KG	193.0000 U	.	216.0000 J	71.6000 J	78.9000 J	304.0000 J
I Chromium		MG/KG	64.9000	.	5.4000	2.3000 U	2.5000 U	9.5000
I Cobalt		MG/KG	2.3000 U	.	1.9000 U	2.0000 U	2.2000 U	2.4000 U
I Copper		MG/KG	5.2000 U	.	5.2000 U	5.7000 U	6.2000 U	14.0000
I Cyanide		MG/KG	0.5200 U	.	0.5200 U	0.5700 U	0.6200 U	0.5500 U
I Iron		MG/KG	917.0000	.	1380.0000	407.0000	306.0000	1870.0000
I Lead		MG/KG	7.9000	.	6.9000	1.4000	0.5300 J	77.1000
I Magnesium		MG/KG	88.8000 U		99.6000 U	108.0000 U	119.0000 U	119.0000 J
I Manganese		MG/KG	2.2000 J	.	20.6000	3.1000 J	3.6000 J	72.5000
I Mercury		MG/KG	0.1000 U	.	0.1000 U	0.1100 U	0.1200 U	0.1100 U
I Nickel		MG/KG	8.4000 U	.	6.5000 U	7.0000 U	7.7000 U	8.7000 U
I Potassium		MG/KG	1010.0000 U		468.0000 U	508.0000 U	558.0000 U	1060.0000 U
I Selenium		MG/KG	0.6300 U	.	0.6300 U	0.6800 U	0.7500 U	0.6600 U
I Silver		MG/KG	2.1000 U		2.1000 U	2.3000 U	2.5000 U	2.2000 U
I sodium		MG/KG	97.0000 J	.	82.5000 J	59.9000 J	110.0000 J	115.0000 J
I Thallium		MG/KG	0.6300 U		0.8400 U	0.9100 U	1.0000 U	0.4400 U
I Vanadium		MG/KG	4.0000 U	.	3.9000 J	1.6000 U	1.7000 U	4.6000 J
I Zinc		MG/KG	6.5000		7.4000 J	2.4000 J	2.3000 J	30.3000
P 4,4'-DDD		UG/KG	3.4000 U	.	3.4000 U	3.7000 U	4.1000 U	3.6000 U
P 4,4'-DDE		UG/KG	3.4000 U		0.1800 J	3.7000 U	4.1000 U	1.4000 J
P 4,4'-DDT		UG/KG	3.4000 U		0.4300 U	3.7000 U	0.6900 J	0.8700 J
P Aldrin		UG/KG	1.7000 U	.	1.8000 U	1.9000 U	2.1000 U	1.8000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33812010	33S12010RE	33813002	33813004	33813006	33814002
ENSAFE ID:		33812010	33S12010RE	33813002	33813004	33813006	33814002
LAB ID:		530126	530126	546852 R1	546855	546859	529896
DATE RCVD:		12/19/92	12/19/92	04/21/93	04/21/93	04/21/93	12/18/92
P Aroclor 1016	UG/KG	34.0000 U	.	34.0000 U	37.0000 U	41.0000 U	36.0000 U
P Aroclor 1221	UG/KG	69.0000 U	.	70.0000 U	76.0000 U	83.0000 U	72.0000 U
P Aroclor 1232	UG/KG	34.0000 U	.	34.0000 U	37.0000 U	41.0000 U	36.0000 U
P Aroclor 1242	UG/KG	34.0000 U	.	34.0000 U	37.0000 U	41.0000 U	36.0000 U
P Aroclor 1248	UG/KG	34.0000 U	.	34.0000 U	37.0000 U	41.0000 U	36.0000 U
P Aroclor 1254	UG/KG	34.0000 U	.	34.0000 U	37.0000 U	41.0000 U	36.0000 U
P Aroclor 1260	UG/KG	7.0000 J	.	34.0000 U	37.0000 U	41.0000 U	36.0000 U
P Dieldrin	UG/KG	1.4000 J	.	0.8200 J	3.7000 U	4.1000 U	0.4900 J
P Endosulfan I	UG/KG	1.7000 U	.	1.8000 U	1.9000 U	2.1000 U	1.8000 U
P Endosulfan II	UG/KG	3.4000 U	.	3.4000 U	3.7000 U	4.1000 U	3.6000 U
P Endosulfan sulfate	UG/KG	0.1800 J	.	3.4000 U	3.7000 U	4.1000 U	3.6000 U
P Endrin	UG/KG	3.4000 U	.	3.4000 U	3.7000 U	4.1000 U	3.6000 U
P Endrin aldehyde	UG/KG	0.4700 J	.	0.1700 J	3.7000 U	4.1000 U	3.6000 U
P Endrin ketone	UG/KG	0.1500 J	.	3.4000 U	3.7000 U	4.1000 U	3.6000 U
P Heptachlor	UG/KG	1.7000 U	.	0.1100 J	1.9000 U	2.1000 U	0.0870 J
P Heptachlor epoxide	UG/KG	1.7000 U	.	1.8000 U	1.9000 U	2.1000 U	1.8000 U
P Methoxychlor	UG/KG	17.0000 U	.	18.0000 U	19.0000 U	21.0000 U	18.0000 U
P Toxaphene	UG/KG	170.0000 U	.	180.0000 U	190.0000 U	210.0000 U	180.0000 U
P alpha-BHC	UG/KG	1.7000 U	.	1.8000 U	1.9000 U	2.1000 U	1.8000 U
P alpha-Chlordane	UG/KG	0.4800 J	.	0.0560 J	1.9000 U	2.1000 U	1.8000 U
P beta-BHC	UG/KG	1.7000 U	.	1.8000 U	1.9000 U	2.1000 U	1.8000 U
P delta-BHC	UG/KG	1.7000 U	.	1.8000 U	1.9000 U	2.1000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.7000 U	.	1.8000 U	1.9000 U	2.1000 U	1.8000 U
P gamma-Chlordane	UG/KG	0.5300 J	.	1.8000 U	1.9000 U	2.1000 U	0.1600 J
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2,4,5-Trichlorophenol	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	33812010	33S12010RE	33813002	33813004	33813006	33814002
	LAB ID:	530126	530126	546852	546855	546859	529896
	DATE RCVD:	12/19/92	12/19/92	04/21/93	04/21/93	04/21/93	12/18/92
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2,4-Dinitrophenol	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2-Methylphenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 2-Nitroaniline	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 3-Nitroaniline	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S 4-Nitroaniline	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U
S 4-Nitrophenol	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U
S Acenaphthene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Acenaphthylene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Anthracene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Benzo(a)anthracene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Benzo(a)pyrene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	WITS						
	ENSAFE ID:	33512010	33S12010RE	33513002	33513004	33513006	33514002
	LAB ID:	530126	530126	546852	546855	546859	529896
	DATE RCVD:	12/19/92	12/19/92	04/21/93	04/21/93	04/21/93	12/18/92
S Butylbenzylphthalate	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Carbazole	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Chrysene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Di-n-butylphthalate	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Dibenz(a,h)anthracene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Dibenzofuran	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Diethylphthalate	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Fluoranthene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Fluorene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Hexachloroethane	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Ieophorone	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Naphthalene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Nitrobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Pentachlorophenol	UG/KG	830.0000 U	.	820.0000 U	900.0000 U	990.0000 U	860.0000 U
S Phenanthrene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Phenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S Pyrene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	340.0000 U	.	51.0000 U	430.0000 U	410.0000 U	54.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	410.0000 U	360.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33812010	33S12010RE	33813002	33813004	33813006	33814002
	LAB ID:	530126	530126	546852	546855	546059	529896
	DATE RCVD:	12/19/92	12/19/92	04/21/93	04/21/93	04/21/93	12/18/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 2-Butanone	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 2-Hexanone	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Acetone	UG/KG	1900.0000 U	4800.0000 U	49.0000 U	200.0000 U	150.0000 U	3100.0000
V Benzene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Bromoform	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Bromomethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Chlorobenzene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Chloroethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Chloroform	UG/KG	10.0000 U	1300.0000 U	1.0000 J	11.0000 U	2.0000 J	1300.0000 U
V Chloromethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Ethylbenzene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Methylene Chloride	UG/KG	6.0000 U	230.0000 U	9.0000 U	10.0000 U	18.0000 U	220.0000 U
V Styrene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Toluene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Trichloroethene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S12010	33S12010RE	33S13002	33S13004	33S13006	33S14002
	LAB ID:	530126	530126	546852	546855	546859	529896
	DATE RCVD:	12/19/92	12/19/92	04/21/93	04/21/93	04/21/93	12/18/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V Xylene (total)	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	1300.0000 U	10.0000 U	11.0000 U	12.0000 U	1300.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33814004	33S14004RE	33S14S02	33S15002	33816002	33817002
	LAB ID:	529900	529900	530665	529535	528731	529322
	DATE RCVD:	12/18/92	12/18/92	12/18/92	12/17/92	12/18/92	12/16/92
NAME	UNITS						
I Aluminum	MG/KG	1040.0000	.	.	193.0000	133.0000 U	447.0000
I Antimony	MG/KG	12.2000 U	.	.	13.9000 U	12.3000 U	13.1000 U
I Arsenic	MG/KG	0.6400 U	.	.	0.7000 U	0.6200 U	0.6500 U
I Barium	MG/KG	2.6000 J		.	1.1000 J	0.9000 J	0.5100 J
I Beryllium	MG/KG	0.4300 U	.	.	0.2300 U	0.2100 U	0.2200 U
I Cadmium	MG/KG	1.1000 U	.	.	1.2000 U	1.0000 U	1.1000 U
I Calcium	MG/KG	198.0000 U	.	.	98.6000 J	9610.0000	90.4000 J
I Chromium	MG/KG	2.4000	.	.	1.8000 J	1.2000 U	1.7000 J
I Cobalt	MG/KG	2.4000 U	.	.	2.1000 U	1.8000 U	2.0000 U
I Copper	MG/KG	5.4000 U	.	.	1.2000 U	3.1000 J	1.1000 U
I Cyanide	MG/KG	0.5400 U	.	.	0.5800 U	0.5100 U	0.5400 U
I Iron	MG/KG	593.0000	.	.	74.9000	363.0000	187.0000
I Lead	MG/KG	3.3000	.	.	2.1000	1.7000	3.4000
I Magnesium	MG/KG	91.1000 U	.	.	10.7000 U	86.9000 J	34.0000 J
I Manganese	MG/KG	8.8000	.	.	1.6000 J	4.2000	1.2000 J
I Mercury	MG/KG	0.1100 U	.	.	0.1200 U	0.1000 U	0.1100 U
I Nickel	MG/KG	8.6000 U	.	.	2.6000 U	2.3000 U	2.4000 U
I Potassium	MG/KG	1040.0000 U	.	.	317.0000 U	280.0000 U	297.0000 U
I Selenium	MG/KG	0.6400 U	.	.	0.7000 U	0.6200 U	0.6500 U
I Silver	MG/KG	2.1000 U	.	.	1.4000 U	1.3000 U	1.3000 U
I Sodium	MG/KG	91.5000 J	.	.	89.3000 J	183.0000 J	96.8000 J
I Thallium	MG/KG	0.6400 U	.	.	0.4600 U	0.4100 U	0.4400 U
I Vanadium	MG/KG	4.1000 U	.	.	1.2000 U	1.0000 U	1.1000 U
I Zinc	MG/KG	11.0000	.	.	2.4000,J	6.0000	2.4000 J
P 4,4'-DDD	UG/KG	3.5000 U	.	.	3.8000 U	0.0900 J	3.5000 U
P 4,4'-DDE	UG/KG	3.5000 J	.	.	0.0830 J	3.4000 U	4.4000
P 4,4'-DDT	UG/KG	2.6000 J	.	.	3.8000 U	0.1400 J	7.6000 J
P Aldrin	UG/KG	1.8000 U	.		0.0460 J	1.7000 U	1.8000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	33S14004	33S14004RE	33S14S02	33815002	33816002	33S17002
	LAB ID:	529900	529900	530665	529535	529865	529322
	DATE RCVD:	12/18/92	12/18/92	12/18/92	12/17/92	12/18/92	12/16/92
NAME	UNITS						
P Aroclor 1016	UG/KG	35.0000 U		.	38.0000 U	34.0000 U	35.0000 U
P Aroclor 1221	UG/KG	72.0000 U			78.0000 U	68.0000 U	72.0000 U
P Aroclor 1232	UG/KG	35.0000 U		.	38.0000 U	34.0000 U	35.0000 U
P Aroclor 1242	UG/KG	35.0000 U		.	38.0000 U	34.0000 U	35.0000 U
P Aroclor 1248	UG/KG	35.0000 U	.	.	38.0000 U	34.0000 U	35.0000 U
P Aroclor 1254	UG/KG	35.0000 U			38.0000 U	34.0000 U	35.0000 U
P Aroclor 1260	UG/KG	35.0000 U			38.0000 U	34.0000 U	9.6000 J
P Dieldrin	UG/KG	0.4100 J	.		3.8000 U	0.1600 J	3.5000 U
P Endosulfan I	UG/KG	0.1600 J	.	.	2.0000 U	0.1800 J	1.8000 U
P Endosulfan II	UG/KG	3.5000 U	.		3.8000 U	3.4000 U	3.5000 U
P Endosulfan sulfate	UG/KG	3.5000 U	.		3.8000 U	3.4000 U	0.7200 J
P Endrin	UG/KG	3.5000 U	.		2.2000 J	3.4000 U	3.5000 U
P Endrin aldehyde	UG/KG	3.5000 U	.		3.8000 U	3.4000 U	3.5000 U
P Endrin ketone	UG/KG	3.5000 U	.	.	3.8000 U	3.4000 U	3.5000 U
P Heptachlor	UG/KG	1.8000 U	.		2.0000 U	1.7000 U	0.1900 U
P Heptachlor epoxide	UG/KG	1.8000 U	.	.	2.0000 U	1.7000 U	0.0980 J
P Methoxychlor	UG/KG	18.0000 U	.	.	20.0000 U	17.0000 U	18.0000 U
P Toxaphene	UG/KG	180.0000 U	.	.	200.0000 U	170.0000 U	180.0000 U
P alpha-BHC	UG/KG	1.8000 U	.	.	2.0000 U	1.7000 U	1.8000 U
P alpha-Chlordane	UG/KG	1.8000 U	.	.	0.2700 J	1.7000 U	0.1700 J
P beta-BHC	UG/KG	1.8000 U	.	.	2.0000 U	1.7000 U	1.8000 U
P delta-BHC	UG/KG	1.8000 U	.	.	2.0000 U	1.7000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	.	.	2.0000 U	1.7000 U	1.8000 U
P gamma-Chlordane	UG/KG	1.8000 U	.	.	2.0000 U	1.7000 U	1.8000 U
S 1,2,4-Trichlorobenzene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 1,2-Dichlorobenzene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 1,3-Dichlorobenzene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 1,4-Dichlorobenzene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 2,4,5-Trichlorophenol	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S14004	33S14004RE	33S14S02	33S15002	33516002	33817002
	LAB ID:	529900	529900	530665	529535	529865	529322
	DATE RCVD:	12/18/92	12/18/92	12/18/92	12/17/92	12/18/92	12/16/92
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	350.0000 U			380.0000 U	340.0000 U	360.0000 U
S 2,4-Dichlorophenol	UG/KG	350.0000 U			380.0000 U	340.0000 U	360.0000 U
S 2,4-Dimethylphenol	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 2,4-Dinitrophenol	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U
S 2,4-Dinitrotoluene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 2,6-Dinitrotoluene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 2-Chloronaphthalene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 2-Chlorophenol	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 2-Methylnaphthalene	UG/KG	350.0000 U	.		380.0000 U	340.0000 U	360.0000 U
S 2-Methylphenol	UG/KG	350.0000 U	.		380.0000 U	340.0000 U	360.0000 U
S 2-Nitroaniline	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U
S 2-Nitrophenol	UG/KG	350.0000 U	.		380.0000 U	340.0000 U	360.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	350.0000 U	.		380.0000 U	340.0000 U	360.0000 U
S 3-Nitroaniline	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U
S 4-Bromophenyl-phenylether	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 4-Chloro-3-methylphenol	UG/KG	350.0000 U	.		380.0000 U	340.0000 U	360.0000 U
S 4-Chloroaniline	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	350.0000 U	.		380.0000 U	340.0000 U	360.0000 U
S 4-Methylphenol	UG/KG	350.0000 U			380.0000 U	340.0000 U	360.0000 U
S 4-Nitroaniline	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U
S 4-Nitrophenol	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U
S Acenaphthene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Acenaphthylene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Anthracene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Benzo(a)anthracene	UG/KG	89.0000 J			380.0000 U	340.0000 U	360.0000 U
S Benzo(a)pyrene	UG/KG	110.0000 J	.	.	380.0000 U	340.0000 U	360.0000 U
S Benzo(b,k)fluoranthene	UG/KG	210.0000 J			380.0000 U	340.0000 U	360.0000 U
S Benzo(g,h,i)perylene	UG/KG	350.0000 U			380.0000 U	340.0000 U	360.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

N—	ENSAFE ID:	33914004	33S14004RE	33S14S02	33815002	33816002	33S17002
	LAB ID:	529900	529900	530665	529535	529865	529322
	DATE RCVD:	12/18/92	12/18/92	12/18/92	12/17/92	12/18/92	12/16/92
	UNITS						
S Butylbenzylphthalate	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Carbazole	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Chrysene	UG/KG	56.0000 J	.	.	380.0000 U	340.0000 U	360.0000 U
S Di-n-butylphthalate	UG/KG	350.0000 U	.	.	63.0000 U	340.0000 U	360.0000 U
S Di-n-octylphthalate	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Dibenz(a,h)anthracene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Dibenzofuran	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Diethylphthalate	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Dimethylphthalate	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Fluoranthene	UG/KG	38.0000 J	.	.	380.0000 U	340.0000 U	360.0000 U
S Fluorene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Hexachlorobenzene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Hexachlorobutadiene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Hexachlorocyclopentadiene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Hexachloroethane	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Isophorone	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S N-Nitrosodiphenylamine	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Naphthalene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Nitrobenzene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Pentachlorophenol	UG/KG	850.0000 U	.	.	920.0000 U	810.0000 U	860.0000 U
S Phenanthrene	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Phenol	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S Pyrene	UG/KG	110.0000 J	.	.	380.0000 U	340.0000 U	360.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	62.0000 U	.	.	58.0000 U	57.0000 U	37.0000 U
S bis(2-chloroethyl)ether	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	350.0000 U	.	.	380.0000 U	340.0000 U	360.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33814004	33S14004RE	33S14S02	33515002	33816002	331517002
	LAB ID:	529900	529900	530665	529535	529865	529322
	DATE RCVD:	12/18/92	12/18/92	12/18/92	12/17/92	12/18/92	12/16/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	54.0000 U	11000.0000	12.0000 U	10.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 1,2-Dichloroethene (total	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 2-Butanone	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 2-Hexanone	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Acetone	UG/KG	1200.0000 U	1200.0000 U	1300.0000 U	36.0000 U	71.0000 U	28.0000 U
V Benzene	UG/KG	11.0000 U	54.0000 U	6900.0000	12.0000 U	10.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Bromoform	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Bromomethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Chlorobenzene	UG/KG	11.0000 U	54.0000 U	6700.0000	12.0000 U	10.0000 U	11.0000 U
V Chloroethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Chloroform	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	1.0000 U
V Chloromethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Ethylbenzene	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Methylene Chloride	UG/KG	21.0000 U	110.0000 U	360.0000 U	36.0000 U	31.0000 U	32.0000 U
V Styrene	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Toluene	UG/KG	11.0000 U	54.0000 U	6600.0000	12.0000 U	10.0000 U	11.0000 U
V Trichloroethene	UG/KG	11.0000 U	54.0000 U	6700.0000	12.0000 U	10.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S14004	33S14004RE	33S14S02	33S15002	33S16002	33S17002
	LAB ID:	529900	529900	530665	529535	529865	529322
	DATE RCVD:	12/18/92	12/18/92	12/18/92	12/17/92	12/18/92	12/16/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V Xylene (total)	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	54.0000 U	1300.0000 U	12.0000 U	10.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33918002	33919002	33S19002RE	33820002	33821002	33S22002
	LAB ID:	529334	529337	529337	529346	529341	529538
	DATE RCVD:	12/16/92	12/16/92	12/16/92	12/16/92	12/16/92	12/17/92
NAME	UNITS						
I Aluminum	MG/KG	33.4000 J	1060.0000	.	17500.0000	332.0000	314.0000
I Antimony	MG/KG	14.7000 U	15.1000 U	.	15.5000 U	13.4000 U	13.1000 U
I Arsenic	MG/KG	0.7300 U	0.7500 U	.	3.5000	0.6700 U	0.6600 U
I Barium	MG/KG	0.4900 U	2.8000 J	.	15.8000 J	0.8200 J	1.5000 J
I Beryllium	MG/KG	0.2400 U	0.2500 U	.	0.2600 U	0.2200 U	0.2200 U
I Cadmium	MG/KG	1.2000 U	1.3000 U	.	1.3000 U	1.1000 U	1.1000 U
I Calcium	MG/KG	21.2000 J	189.0000 J	.	374.0000 J	178.0000 J	423.0000 J
I Chromium	MG/KG	1.5000 U	37.0000	.	22.1000	74.5000	2.4000
I Cobalt	MG/KG	2.2000 U	2.3000 U	.	2.3000 U	2.0000 U	2.0000 U
I Copper	MG/KG	1.2000 U	3.1000 J	.	10.2000	3.2000 J	1.2000 J
I Cyanide	MG/KG	0.6100 U	0.6300 U	.	0.6500 U	0.5600 U	0.5500 U
I Iron	MG/KG	18.2000 J	498.0000	.	7500.0000	131.0000	226.0000
I Lead	MG/KG	0.5500 J	8.7000	.	28.9000	3.7000	7.4000
I Magnesium	MG/KG	11.2000 U	59.1000 J	.	265.0000 J	10.5000 J	67.3000 J
I Manganese	MG/KG	0.4900 U	2.1000 J	.	537.0000	1.8000 J	5.0000
I Mercury	MG/KG	0.1200 U	0.1300 U	.	0.1300 U	0.1100 U	0.1100 U
I Nickel	MG/KG	2.7000 U	2.8000 U	.	3.5000 J	2.5000 U	2.4000 U
I Potassium	MG/KG	333.0000 U	342.0000 U	.	352.0000 U	304.0000 U	299.0000 U
I Selenium	MG/KG	0.7300 U	0.7500 U	.	0.7800 U	0.6700 U	0.6600 U
I Silver	MG/KG	1.5000 U	1.5000 U	.	1.6000 U	1.3000 U	1.3000 U
I sodium	MG/KG	107.0000 J	118.0000 J	.	133.0000 J	81.5000 J	102.0000 J
I Thallium	MG/KG	0.4900 U	0.5000 U	.	0.5200 U	0.4700 J	0.4400 U
I Vanadium	MG/KG	1.2000 U	1.7000 J	.	21.8000	1.1000 U	1.1000 U
I Zinc	MG/KG	2.2000 J	6.7000	.	17.7000	7.7000	8.6000
P 4,4'-DDD	UG/KG	4.0000 U	4.1000 U	.	4.3000 U	3.7000 U	3.6000 U
P 4,4'-DDE	UG/KG	0.2100 J	1.2000 J	.	0.6500 J	0.2300 J	6.2000
P 4,4'-DDT	UG/KG	0.1900 J	4.1000 U	.	0.2600 U	3.7000 U	12.0000
P Aldrin	UG/KG	2.1000 U	2.1000 U	.	0.1500 J	1.9000 U	1.9000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33910002	33919002	33S19002RE	33S20002	33821002	33822002
	LAB ID:	529334	529337	529337	529346	529341	529538
	DATE RCVD:	12/16/92	12/16/92	12/16/92	12/16/92	12/16/92	12/17/92
NAME	UNITS						
P Aroclor 1016	UG/KG	40.0000 U	41.0000 U		43.0000 U	37.0000 U	36.0000 U
P Aroclor 1221	UG/KG	81.0000 U	84.0000 U		87.0000 U	74.0000 U	73.0000 U
P Aroclor 1232	UG/KG	40.0000 U	41.0000 U		43.0000 U	37.0000 U	36.0000 U
P Aroclor 1242	UG/KG	40.0000 U	41.0000 U		43.0000 U	37.0000 U	36.0000 U
P Aroclor 1248	UG/KG	40.0000 U	41.0000 U		43.0000 U	37.0000 U	36.0000 U
P Aroclor 1254	UG/KG	40.0000 U	41.0000 U		43.0000 U	37.0000 U	36.0000 U
P Aroclor 1260	UG/KG	40.0000 U	12.0000 J		43.0000 U	6.1000 J	4.1000 J
P Dieldrin	UG/KG	4.0000 U	4.1000 U		4.3000 U	0.4100 J	0.5900 J
P Endosulfan I	UG/KG	2.1000 U	2.1000 U		2.2000 U	1.9000 U	1.9000 U
P Endosulfan II	UG/KG	4.0000 U	4.1000 U		4.3000 U	3.7000 U	3.6000 U
P Endosulfan sulfate	UG/KG	0.3200 J	0.8400 J		2.3000 J	3.7000 U	0.7500 J
P Endrin	UG/KG	4.0000 U	4.1000 U		0.3600 J	3.7000 U	3.6000 U
P Endrin aldehyde	UG/KG	4.0000 U	4.1000 U		3.9000 J	3.7000 U	3.6000 U
P Endrin ketone	UG/KG	4.0000 U	4.1000 U		4.3000 U	3.7000 U	3.6000 U
P Heptachlor	UG/KG	2.1000 U	0.4900 U		2.2000 U	0.0820 U	1.9000 U
P Heptachlor epoxide	UG/KG	2.1000 U	0.1300 J		2.2000 U	1.9000 U	0.0950 J
P Methoxychlor	UG/KG	21.0000 U	21.0000 U		22.0000 U	19.0000 U	19.0000 U
P Toxaphene	UG/KG	210.0000 U	210.0000 U		220.0000 U	190.0000 U	190.0000 U
P alpha-BHC	UG/KG	2.1000 U	2.1000 U		2.2000 U	1.9000 U	1.9000 U
P alpha-Chlordane	UG/KG	2.1000 U	2.1000 U		29.0000	0.0900 J	1.9000 U
P beta-BHC	UG/KG	2.1000 U	2.1000 U		2.2000 U	1.9000 U	1.9000 U
P de1ta-BHC	UG/KG	2.1000 U	2.1000 U		2.2000 U	1.9000 U	0.0770 J
P gamma-BHC (Lindane)	UG/KG	2.1000 U	2.1000 U		2.2000 U	1.9000 U	1.9000 U
P gamma-Chlordane	UG/KG	0.0490 J	2.1000 U		2.2000 U	0.4800 J	1.9000 U
S 1,2,4-Trichlorobenzene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 1,2-Dichlorobenzene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	'360.0000 U	360.0000 U
S 1,3-Dichlorobenzene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 1,4-Dichlorobenzene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2,4,5-Trichlorophenol	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33818002	33519002	33S19002RE	33820002	33S21002	33522002
ENSAFE ID:		33818002	33519002	33S19002RE	33820002	33S21002	33522002
LAB ID:		529334	529337	529337	529346	529341	529538
DATE RCVD:		12/16/92	12/16/92	12/16/92	12/16/92	12/16/92	12/17/92
S 2,4,6-Trichlorophenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2,4-Dichlorophenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2,4-Dimethylphenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2,4-Dinitrophenol	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U
S 2,4-Dinitrotoluene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2,6-Dinitrotoluene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2-Chloronaphthalene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2-Chlorophenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2-Methylnaphthalene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2-Methylphenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 2-Nitroaniline	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U
S 2-Nitrophenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 3-Nitroaniline	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U
S 4-Bromophenyl-phenylether	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 4-Chloro-3-methylphenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 4-Chloroaniline	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 4-Methylphenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S 4-Nitroaniline	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U
S 4-Nitrophenol	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U
S Acenaphthene	UG/KG	400.0000 U	410.0000 U	410.0000 U	99.0000 J	360.0000 U	360.0000 U
S Acenaphthylene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Anthracene	UG/KG	400.0000 U	410.0000 U	410.0000 U	230.0000 J	360.0000 U	360.0000 U
S Benzo(a)anthracene	UG/KG	400.0000 U	410.0000 U	410.0000 U	7500.0000 J	360.0000 U	360,0000 U
S Benzo(a)pyrene	UG/KG	400.0000 U	410.0000 U	410.0000 U	6200.0000 J	360.0000 U	360.0000 U
S Benzo(b,k)fluoranthene	UG/KG	400.0000 U	410.0000 U	410.0000 U	14000.0000 J	360.0000 U	360.0000 U
S Benzo(g,h,i)perylene	UG/KG	400.0000 U	410.0000 U	410.0000 U	5300.0000 J	360.0000 U	360.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33518002	33S19002	33S19002RE	33520002	33821002	33522002
	LAB ID:	529334	529337	529337	529346	529341	529538
	DATE RCVD:	12/16/92	12/16/92	12/16/92	12/16/92	12/16/92	12/17/92
NAME	WITS						
S Butylbenzylphthalate	UG/KG	400.0000 U	60.0000 U	62.0000 U	420.0000 U	43.0000 U	360.0000 U
S Carbazole	UG/KG	400.0000 U	410.0000 U	410.0000 U	59.0000 J	360.0000 U	360.0000 U
S Chrysene	UG/KG	400.0000 U	410.0000 U	410.0000 U	5600.0000 J	360.0000 U	360.0000 U
S Di-n-butylphthalate	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Di-n-octylphthalate	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Dibenz (a,h) anthracene	UG/KG	400.0000 U	410.0000 U	410.0000 U	1400.0000	360.0000 U	360.0000 U
S Dibenzofuran	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Diethylphthalate	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Dimethylphthalate	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Fluoranthene	UG/KG	400.0000 U	410.0000 U	410.0000 U	4800.0000 J	360.0000 U	360.0000 U
S Fluorene	UG/KG	400.0000 U	410.0000 U	410.0000 U	70.0000 J	360.0000 U	360.0000 U
S Hexachlorobenzene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Hexachlorobutadiene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Hexachlorocyclopentadiene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Hexachloroethane	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Indeno(1,2,3-cd) pyrene	UG/KG	400.0000 U	410.0000 U	410.0000 U	4800.0000 J	360.0000 U	360.0000 U
S Isophorone	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S N-Nitrosodiphenylamine	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Naphthalene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Nitrobenzene	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Pentachlorophenol	UG/KG	960.0000 U	990.0000 U	990.0000 U	1000.0000 U	880.0000 U	880.0000 U
S Phenanthrene	UG/KG	400.0000 U	410.0000 U	410.0000 U	1500.0000	360.0000 U	360.0000 U
S Phenol	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U
S Pyrene	UG/KG	400.0000 U	47.0000 J	410.0000 U	8600.0000 J	360.0000 U	360.0000 U
S bis(2-Chloroethoxy) methane	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 'U	360.0000 U	360.0000 U
S bis (2-Ethlyhexy) phthalate	UG/KG	400.0000 U	130.0000 U	110.0000 U	120.0000 U	100.0000 U	360.0000 U
S bis(2-chloroethyl) ether	UG/KG	400.0000 U	850.0000	810.0000	420.0000 U	360.0000 U	360.0000 U
S bis(2-chloroisopropyl) ether	UG/KG	400.0000 U	410.0000 U	410.0000 U	420.0000 U	360.0000 U	360.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33818002	33819002	33S19002RE	33820002	33821002	33822002
	LAB ID:	529334	529337	529337	529346	529341	529538
	DATE RCVD:	12/16/92	12/16/92	12/16/92	12/16/92	12/16/92	12/17/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	190.0000 U	31.0000 U	.	47.0000 U	36.0000 U	49.0000 U
V Benzene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Chloroethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Chloroform	UG/KG	1.0000 U	12.0000 U	.	13.0000 U	1.0000 U	11.0000 U
V Chloromethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	36.0000 U	31.0000 U	.	39.0000 U	41.0000 U	62.0000 U
V Styrene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U
V Trichloroethene	UG/KG	12.0000 U	12.0000 U	.	13.0000 U	11.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S18002	33S19002	33S19002DE	33S20002	33S21002	33S22002
	LAB ID:	529334	529337	529337	529346	529341	529538
	DATE RCVD:	12/16/92	12/16/92	12/16/92	12/16/92	12/16/92	12/17/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	12.0000 ω	12.0000 ω	.	13.0000 ω	11.0000 ω	11.0000 ω
V Xylene (total)	UG/KG	12.0000 ω	12.0000 ω	.	13.0000 ω	11.0000 ω	11.0000 ω
V cis-1,3-Dichloropropene	UG/KG	12.0000 ω	12.0000 ω	.	13.0000 ω	11.0000 ω	11.0000 ω
V trans-1,3-Dichloropropene	UG/KG	12.0000 ω	12.0000 ω	.	13.0000 ω	11.0000 ω	11.0000 ω

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	331522004	33823002	33823004	33524002	33S24004	33524005
	LAB ID:	529540	529541	529552	530123	530129	547123
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/19/92	12/19/92	04/22/93
NAME	WITS						
I Aluminum	MG/KG	114.0000	5810.0000	84.0000	1230.0000	953.0000	533.0000
I Antimony	MG/KG	13.1000 U	13.3000 U	13.6000 U	11.6000 U	13.8000 U	11.5000 U
I Arsenic	MG/KG	0.6600 U	0.9400 J	0.6800 U	0.6100 U	0.7300 U	0.7500 U
I Barium	MG/KG	0.5800 J	8.1000 J	0.4500 U	3.5000 J	19.6000 J	4.3000 J
I Beryllium	MG/KG	0.2200 U	0.2200 U	0.2300 U	0.4100 U	0.4800 U	0.5000 U
I Cadmium	MG/KG	1.1000 U	1.1000 U	1.1000 U	1.0000 U	5.9000	1.6000
I Calcium	MG/KG	160.0000 J	340.0000 J	26.1000 J	7630.0000	563.0000 J	2490.0000
I Chromium	MG/KG	1.3000 J	13.1000	1.4000 U	2.0000 U	70.7000	19.2000
I Cobalt	MG/KG	2.0000 U	2.0000 U	2.0000 U	2.2000 U	2.7000 U	2.2000 U
I Copper	MG/KG	1.1000 U	3.2000 J	1.1000 U	5.1000 U	21.3000	8.7000
I Cyanide	MG/KG	0.5500 U	0.5500 U	0.5700 U	0.5100 U	0.6100 U	0.6200 U
I Iron	MG/KG	70.6000	2860.0000	39.4000	684.0000	408.0000	342.0000
I Lead	MG/KG	2.8000	8.8000	0.4500 U	183.0000	43.7000	7.4000
I Magnesium	MG/KG	30.0000 J	138.0000 J	10.4000 U	116.0000 J	103.0000 U	198.0000 J
I Manganese	MG/KG	1.5000 J	49.0000	0.8900 J	15.5000	2.5000 J	4.4000
I Mercury	MG/KG	0.1100 U	0.1100 U	0.1100 U	0.1000 U	1.1000	0.1900
I Nickel	MG/KG	2.4000 U	2.4000 U	2.5000 U	8.2000 U	9.7000 U	7.7000 U
I Potassium	MG/KG	298.0000 U	302.0000 U	309.0000 U	991.0000 U	1180.0000 U	558.0000 U
I Selenium	MG/KG	0.6606 U	0.6700 U	0.6800 U	0.6100 U	0.7300 U	0.7500 U
I Silver	MG/KG	1.3000 U	1.3000 U	1.4000 U	2.0000 U	9.0000 U	2.5000 U
I Sodium	MG/KG	88.7000 J	95.4000 J	91.4000 J	187.0000 J	135.0000 J	97.3000 J
I Thallium	MG/KG	0.4400 U	0.4400 U	0.4500 U	0.6100 U	0.7300 U	1.0000 U
I Vanadium	MG/KG	1.1000 U	7.3000 J	1.1000 U	3.9000 U	4.6000 U	1.7000 U
I Zinc	MG/KG	5.3000	106.0000	2.0000 J	7.5000,	16.4000	7.5000
P 4,4'-DDD	UG/KG	3.6000 U	2.0000 J	1.7000 J			24.0000
P 4,4'-DDE	UG/KG	2.8000 J	42.0000	31.0000			7.9000
P 4,4'-DDT	UG/KG	5.2000	21.0000 J	11.0000			1.3000 J
P Aldrin	UG/KG	1.9000 U	0.0810 J	1.9000 U			2.1000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33822004	33S23002	33823004	33524002	33824004	33S24005
	LAB ID:	529540	529541	529552	530123	530129	547123
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/19/92	12/19/92	04/22/93
NAME	UNITS						
P Aroclor 1016	UG/KG	36.0000 U	36.0000 U	37.0000 U	.	.	41.0000 U
P Aroclor 1221	UG/KG	73.0000 U	74.0000 U	76.0000 U	.	.	83.0000 U
P Aroclor 1232	UG/KG	36.0000 U	36.0000 U	37.0000 U	.	.	41.0000 U
P Aroclor 1242	UG/KG	36.0000 U	36.0000 U	37.0000 U	.	.	41.0000 U
P Aroclor 1248	UG/KG	36.0000 U	36.0000 U	37.0000 U	.	.	41.0000 U
P Aroclor 1254	UG/KG	36.0000 U	36.0000 U	37.0000 U	.	.	41.0000 U
P Aroclor 1260	UG/KG	36.0000 U	36.0000 U	37.0000 U	.	.	11.0000 J
P Dieldrin	UG/KG	0.2000 J	2.7000 J	0.9100 J	.	.	1.1000 J
P Endosulfan I	UG/KG	1.9000 U	1.9000 U	0.3400 J	.	.	2.1000 U
P Endosulfan II	UG/KG	0.1100 J	3.6000 U	3.7000 U	.	.	4.1000 U
P Endosulfan sulfate	UG/KG	0.3800 J	3.6000 U	3.7000 U	.	.	4.1000 U
P Endrin	UG/KG	3.6000 U	3.6000 U	3.7000 U	.	.	0.2300 J
P Endrin aldehyde	UG/KG	3.6000 U	0.4900 J	3.7000 U	.	.	4.1000 U
P Endrin ketone	UG/KG	3.6000 U	3.6000 U	3.7000 U	.	.	4.1000 U
P Heptachlor	UG/KG	1.9000 U	0.1300 J	0.0570 J	.	.	2.1000 U
P Heptachlor epoxide	UG/KG	1.9000 U	1.9000 U	1.9000 U	.	.	2.1000 U
P Methoxychlor	UG/KG	1.0000 J	19.0000 U	19.0000 U	.	.	21.0000 U
P Toxaphene	UG/KG	190.0000 U	190.0000 U	190.0000 U	.	.	210.0000 U
P alpha-BHC	UG/KG	1.9000 U	1.9000 U	1.9000 U	.	.	2.1000 U
P alpha-Chlordane	UG/KG	1.9000 U	1.9000 U	1.9000 U	.	.	1.6000 J
P beta-BHC	UG/KG	1.9000 U	1.9000 U	1.9000 U	.	.	2.1000 U
P delta-BHC	UG/KG	1.9000 U	1.9000 U	1.9000 U	.	.	2.1000 U
P gamma-BHC (Lindane)	UG/KG	1.9000 U	0.1000 J	1.9000 U	.	.	2.1000 U
P gamma-Chlordane	UG/KG	0.1300 J	1.9000 U	1.9000 U	.	.	0.4700 J
S 1,2,4-Trichlorobenzene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 1,2-Dichlorobenzene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	77.0000 J	410.0000 U
S 1,3-Dichlorobenzene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	85.0000 J	410.0000 U
S 1,4-Dichlorobenzene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	120.0000 J	410.0000 U
S 2,4,5-Trichlorophenol	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S22004	33S23002	33823004	33824002	33824004	33S24005
	LAB ID:	529540	529541	529552	530123	530129	547123
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/19/92	12/19/92	04/22/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2,4-Dichlorophenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2,4-Dimethylphenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2,4-Dinitrophenol	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U
S 2,4-Dinitrotoluene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2,6-Dinitrotoluene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2-Chloronaphthalene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2-Chlorophenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2-Methylnaphthalene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2-Methylphenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 2-Nitroaniline	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U
S 2-Nitrophenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 3-Nitroaniline	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U
S 4-Bromophenyl-phenylether	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 4-Chloro-3-methylphenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 4-Chloroaniline	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 4-Methylphenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S 4-Nitroaniline	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U
S 4-Nitrophenol	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U
S Acenaphthene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Acenaphthylene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Anthracene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Benzo(a)anthracene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Benzo(a)pyrene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Benzo(b,k)fluoranthene	UG/KG	360.0000 U	76.0000 J	370.0000 U	340.0000 U	44.0000 J	410.0000 U
S Benzo(g,h,i)perylene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33822004	33823002	33823004	33824002	33824004	33S24005
	LAB ID:	529540	529541	529552	530123	530129	547123
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/19/92	12/19/92	04/22/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Carbazole	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Chrysene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Di-n-butylphthalate	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Di-n-octylphthalate	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Dibenz (a,h) anthracene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Dibenzofuran	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Diethylphthalate	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Dimethylphthalate	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Fluoranthene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Fluorene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Hexachlorobenzene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Hexachlorobutadiene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Hexachlorocyclopentadiene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Hexachloroethane	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	360.0000 U	40.0000 J	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Isophorone	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S N-Nitrosodiphenylamine	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Naphthalene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Nitrobenzene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Pentachlorophenol	UG/KG	870.0000 U	880.0000 U	900.0000 U	810.0000 U	970.0000 U	990.0000 U
S Phenanthrene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Phenol	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S Pyrene	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S bis (2-Chloroethoxy) methane	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S bis(2-Ethlyhexy) phthalate	UG/KG	48.0000 U	42.0000 U	47.0000 U	340.0000 U	85.0000 U	480.0000 U
S bis(2-chloroethyl) ether	UG/KG	82.0000 J	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U
S bis(2-chloroisopropyl) ether	UG/KG	360.0000 U	360.0000 U	370.0000 U	340.0000 U	400.0000 U	410.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33822004	33S23002	33823004	33824002	33824004	33821005
ENSAFE ID:		33822004	33S23002	33823004	33824002	33824004	33821005
LAB ID:		529540	529541	529552	530123	530129	547123
DATE RCVD:		12/17/92	12/17/92	12/17/92	12/19/92	12/19/92	04/22/93
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 2-Butanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 2-Hexanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Acetone	UG/KG	110.0000 U	50.0000 U	30.0000 U	32.0000 U	43.0000 U	97.0000 U
V Benzene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Bromoform	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Bromomethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Chlorobenzene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	6.0000 J	12.0000 U
V Chloroethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Chloroform	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	2.0000 J
V Chloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Bthylbenzene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Methylene Chloride	UG/KG	50.0000 U	32.0000 U	27.0000 U	5.0000 U	20.0000 U	28.0000 U
V Styrene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Toluene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V Trichloroethene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33922004	33823002	33823004	33824002	33924004	33824005
	LAB ID:	529540	529541	529552	530123	530129	547123
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/19/92	12/19/92	04/22/93
NAME	WITS						
V Vinyl Chloride	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V xylene (total)	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U	11.0000 U	10.0000 U	12.0000 U	12.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not ehown in significant digits.

	ENSAFE ID:	33825002	33825004	33S32002	33S32003	33833002	33S33002DL
	LAB ID:	530141	530120	532972	532975	532978	532978
	DATE RCVD:	12/19/92	12/19/92	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
I Aluminum	MG/KG	208.0000	94.4000	5030.0000	1490.0000	8010.0000	
I Antimony	MG/KG	11.7000 U	13.6000 U	12.4000 U	12.7000 U	14.3000 U	
I Arsenic	MG/KG	0.6200 U	0.7100 U	1.1000 J	0.6700 U	0.7500 U	
I Barium	MG/KG	1.2000 J	0.9300 J	18.4000 J	3.7000 J	43.0000 J	
I Beryllium	MG/KG	0.4100 U	0.4800 U	0.6600 U	0.6700 U	0.7500 U	
I Cadmium	MG/KG	1.0000 U	1.2000 U	31.2000	3.6000	22.1000	
I Calcium	MG/KG	1980.0000	2120.0000	2230.0000	2210.0000	6210.0000	
I Chromium	MG/KG	2.1000 U	2.4000 U	935.0000	148.0000	607.0000	
I Cobalt	MG/KG	2.3000 U	2.6000 U	2.6000 U	2.7000 U	3.0000 U	
I Copper	MG/KG	5.1000 U	5.9000 U	19.7000	5.5000 U	19.7000	
I Cyanide	MG/KG	0.5100 U	0.5900 U	0.5500 U	0.5500 U	0.6300 U	
I Iron	MG/KG	93.3000	33.7000	1700.0000	630.0000	3460.0000	
I Lead	MG/KG	1.9000	0.5100 J	54.7000	3.0000	83.3000	
I Magnesium	MG/KG	87.4000 U	101.0000 U	191.0000 J	123.0000 U	403.0000 J	
I Manganese	MG/KG	4.2000	4.8000	40.0000	11.3000	78.8000	
I Mercury	MG/KG	0.1000 U	0.1200 U	0.7400	0.1100 U	1.2000	
I Nickel	MG/KG	8.2000 U	9.5000 U	53.1000	8.4000 U	27.8000	
I Potassium	MG/KG	998.0000 U	1150.0000 U	863.0000 U	877.0000 U	992.0000 U	
I Selenium	MG/KG	0.6200 U	0.7100 U	0.8700 U	0.6700 U	0.7500 U	
I Silver	MG/KG	2.1000 U	2.4000 U	6.4000	2.2000 U	6.6000	
I Sodium	MG/KG	112.0000 J	117.0000 J	138.0000 J	135.0000 J	176.0000 J	
I Thallium	MG/KG	0.6200 U	0.7100 U	0.6600 U	0.6700 U	0.7500 U	
I Vanadium	MG/KG	3.9000 U	4.5000 U	5.6000 J	3.3000 U	8.8000 J	
I Zinc	MG/KG	2.3000 J	3.6000 J	136.0000	13.6000	120.0000	
P 4,4'-DDD	UG/KG			3.5000 U	3.5000 'U	0.5800 J	
P 4,4'-DDE	UG/KG			0.6200 J	3.5000 U	6.4000 J	
P 4,4'-DDT	UG/KG			3.5000 U	3.5000 U	4.1000 U	
P Aldrin	UG/KG			0.1000 J	1.8000 U	2.8000 J	

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	33825002	33825004	33832002	33S32003	33933002	33S33002DL
	LAB ID:	530141	530120	532972	532975	532978	532978
	DATE RCVD:	12/19/92	12/19/92	01/20/93	01/20/93	01/20/93	01/20/93
P Aroclor 1016	UG/KG	.		35.0000 U	35.0000 U	41.0000 U	.
P Aroclor 1221	UG/KG	.		72.0000 U	71.0000 U	82.0000 U	.
P Aroclor 1232	UG/KG	.	.	35.0000 U	35.0000 U	41,0000 U	.
P Aroclor 1242	UG/KG	.	.	35.0000 U	35.0000 U	41.0000 U	.
P Aroclor 1248	UG/KG	.	.	35.0000 U	35.0000 U	41.0000 U	.
P Aroclor 1254	UG/KG	.	.	35.0000 U	35.0000 U	41.0000 U	
P Aroclor 1260	UG/KG	.		39.0000	35.0000 U	50.0000 J	.
P Dieldrin	UG/KG			5.9000 J	3.5000 U	7.7000 J	.
P Endosulfan I	UG/KG	.	.	1.8000 U	1.8000 U	2.1000 U	.
P Endosulfan II	UG/KG	.	.	3.5000 U	3.5000 U	4.1000 U	.
P Endosulfan sulfate	UG/KG	.	.	3.5000 U	3.5000 U	4.1000 U	.
P Endrin	UG/KG	.	.	3.5000 U	3.5000 U	4.1000 U	.
P Endrin aldehyde	UG/KG	.	.	3.5000 U	3.5000 U	4.1000 U	.
P Endrin ketone	UG/KG	.	.	3.5000 U	3.5000 U	4.1000 U	.
P Heptachlor	UG/KG	.	.	1.8000 U	0.0620 U	2.1000 U	.
P Heptachlor epoxide	UG/KG	.	.	1.8000 U	1.8000 U	2.1000 U	.
P Methoxychlor	UG/KG	.	.	18.0000 U	18.0000 U	21.0000 U	.
P Toxaphene	UG/KG	.	.	180.0000 U	180.0000 U	210.0000 U	.
P alpha-BHC	UG/KG	.	.	0.1300 J	1.8000 U	2.1000 U	.
P alpha-Chlordane	UG/KG	.	.	1.5000 J	1.8000 U	2.1000 U	.
P beta-BHC	UG/KG	.	.	1.8000 U	1.8000 U	2.1000 U	.
P delta-BHC	UG/KG	.	.	1.8000 U	1.8000 U	0.4000 J	.
P gamma-BHC (Lindane)	UG/KG	.	.	1.8000 U	1.8000 U	2.1000 U	.
P gama-Chlordane	UG/KG			1.7000 J	0.0930 J	5.7000	
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	'260.0000 J	240.0000 J
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	1700.0000	1600.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	3000.0000	2800.0000 J
S 2,4,5-Trichlorophenol	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33525002	33S25004	33532002	33532003	33833002	33S33002DL
	LAB ID:	530141	530120	532972	532975	532978	532978
	DATE RCVD:	12/19/92	12/19/92	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2,4-Dinitrophenol	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	4100.0000 J	4200.0000 J
S 2-Methylphenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 2-Nitroaniline	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 3-Nitroaniline	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S 4-Nitroaniline	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U
S 4-Nitrophenol	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U
S Acenaphthene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Acenaphthylene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Anthracene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Benzo(a)anthracene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	230.0000 J	210.0000 J
S Benzo(a)pyrene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	160.0000 J	190.0000 J
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	330.0000 J	400.0000 J
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	100.0000 J	820.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33825002	33825004	33832002	33832003	33533002	33S33002DL
	LAB ID:	530141	530120	532972	532975	532978	532978
	DATE RCVD:	12/19/92	12/19/92	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	340.0000 U	390.0000 U	58.0000 U	360.0000 U	1100.0000 U	1300.0000 U
S Carbazole	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Chrysene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	260.0000 J	260.0000 J
S Di-n-butylphthalate	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Dibenz(a,h)anthracene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	45.0000 J	820.0000 U
S Dibenzofuran	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Diethylphthalate	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Fluoranthene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	1000.0000	1000.0000 U
S Fluorene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	750.0000	570.0000 J
S Hexachlorobenzene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Hexachloroethane	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	92.0000 J	820.0000 U
S Isophorone	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Naphthalene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	1400.0000	1400.0000 J
S Nitrobenzene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S Pentachlorophenol	UG/KG	820.0000 U	950.0000 U	860.0000 U	880.0000 U	990.0000 U	2000.0000 U
S Phenanthrene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	1800.0000	1600.0000 J
S Phenol	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	340.0000 J	290.0000 J
S Pyrene	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	490.0000	480.0000 J
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	1200.0000	1300.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	390.0000 U	350.0000 U	360.0000 U	410.0000 U	820.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S25002	33S25004	33S32002	33S32003	33S33002	33S33002DL
	LAB ID:	530141	530120	532972	532975	532978	532978
	DATE RCVD:	12/19/92	12/19/92	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 1,1-Dichloroethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 1,1-Dichloroethene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 1,2-Dichloroethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 1,2-Dichloropropane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 2-Butanone	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 2-Hexanone	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Acetone	UG/KG	6.0000 U	63.0000 U	23.0000 U	65.0000 U	23.0000 U	
V Benzene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Bromodichloromethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Bromoform	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Bromomethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Carbon Disulfide	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Carbon Tetrachloride	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Chlorobenzene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	28.0000	
V Chloroethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Chloroform	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	2.0000 U	
V Chloromethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Dibrmochloromethane	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Ethylbenzene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Methylene Chloride	UG/KG	7.0000 U	11.0000 U	17.0000 U	17.0000 U	23.0000 U	
V Styrene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Tetrachloroethene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Toluene	UG/KG	10.0000 U	12.0000 U	2.0000 J	1.0000 J	12.0000 U	
V Trichloroethene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S25002	33S25004	33S32002	33S32003	33S33002	33S33002DL
	LAB ID:	530141	530120	532972	532975	532978	532978
	DATE RCVD:	12/19/92	12/19/92	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V vinyl Chloride	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V Xylene (total)	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	12.0000 U	11.0000 U	11.0000 U	12.0000 U	

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33834002 / 532976 / 01/20/93	33S35002 / 532977 / 01/20/93	33835004 / 532936 / 01/20/93	33536002 / 532942 / 01/20/93	33836003 / 532943 / 01/20/93	33837002 / 532957 / 01/20/93
I Aluminum	MG/KG	2270.0000	4490.0000	409.0000	5580.0000	7490.0000	7350.0000
I Antimony	MG/KG	12.2000 U	13.9000 U	13.0000 U	12.9000 U	13.5000 U	14.3000 U
I Arsenic	MG/KG	0.7900 J	0.8700 J	0.6800 U	1.7000 J	2.4000	2.0000 J
I Barium	MG/KG	10.1000 J	82.8000	5.1000 J	20.4000 J	40.5000 J	33.8000 J
I Beryllium	MG/KG	0.6400 U	0.7300 U	0.6800 U	0.6800 U	0.7100 U	0.7500 U
I Cadmium	MG/KG	8.2000	27.4000	1.6000	3.5000	22.3000	36.6000
I Calcium	MG/KG	11600.0000	3820.0000	170.0000 J	2970.0000	4450.0000	1690.0000
I Chromium	MG/KG	248.0000	615.0000	29.0000	80.1000	1340.0000	1300.0000
I Cobalt	MG/KG	2.6000 U	2.9000 U	2.7000 U	2.7000 U	2.8000 U	3.0000 U
I Copper	MG/KG	19.4000	68.8000	7.0000	16.6000	58.4000	74.1000
I Cyanide	MG/KG	0.5400 U	0.6100 U	0.5700 U	0.5600 U	1.6000	0.7500
I Iron	MG/KG	1070.0000	1920.0000	97.1000	2800.0000	1740.0000	2420.0000
I Lead	MG/KG	17.7000	86.0000	1.8000	25.0000	105.0000	77.5000
I Magnesium	MG/KG	129.0000 J	179.0000 J	126.0000 U	309.0000 J	130.0000 U	156.0000 J
I Manganese	MG/KG	15.4000	23.4000	1.8000 J	22.1000	26.9000	24.0000
I Mercury	MG/KG	0.7200	0.9100	0.1100 U	0.3800	0.5100	0.7000
I Nickel	MG/KG	12.1000	11.9000	8.7000 U	8.6000 U	29.3000	49.4000
I Potassium	MG/KG	847.0000 U	961.0000 U	900.0000 U	892.0000 U	933.0000 U	993.0000 U
I Selenium	MG/KG	0.6400 U	0.7300 U	0.6800 U	0.6800 U	0.7100 U	0.7500 U
I Silver	MG/KG	2.9000	23.4000	2.3000 U	3.3000 U	14.5000	12.8000
I Sodium	MG/KG	267.0000 J	182.0000 J	141.0000 J	145.0000 J	159.0000 J	134.0000 J
I Thallium	MG/KG	0.6400 U	0.7300 U	0.6800 U	0.6800 U	0.7100 U	0.7500 U
I Vanadium	MG/KG	3.2000 U	4.8000 J	3.4000 U	7.0000 J	5.2000 J	6.4000 J
I Zinc	MG/KG	50.0000	185.0000	12.9000	31.1000,	94.7000	166.0000
P 4,4'-DDD	UG/KG	3.5000 U	4.0000 U	20.0000	18.0000 U	5.1000 J	4.1000 U
P 4,4'-DDE	UG/KG	0.5000 J	4.8000	4.4000 J	4.6000 J	3.7000 J	1.4000 J
P 4,4'-DDT	UG/KG	3.5000 U	4.0000 U	0.4900 J	18.0000 U	3.9000 U	4.1000 U
P Aldrin	UG/KG	0.0660 J	0.2500 J	0.0500 J	0.7800 J	0.2800 J	2.1000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33S34002 / 532976 / 01/20/93	33535002 / 532977 / 01/20/93	33835004 / 532936 R1 / 01/20/93	33536002 / 532942 R1D50 / 01/20/93	33836003 / 532943 / 01/20/93	33837002 / 532957 / 01/20/93
P Aroclor 1016	UG/KG	35.0000 U	40.0000 U	37.0000 U	180.0000 U	39.0000 U	41.0000 U
P Aroclor 1221	UG/KG	70.0000 U	80.0000 U	75.0000 U	370.0000 U	79.0000 U	83.0000 U
P Aroclor 1232	UG/KG	35.0000 U	40.0000 U	37.0000 U	180.0000 U	39.0000 U	41.0000 U
P Aroclor 1242	UG/KG	35.0000 U	40.0000 U	37.0000 U	180.0000 U	39.0000 U	41.0000 U
P Aroclor 1248	UG/KG	35.0000 U	40.0000 U	37.0000 U	180.0000 U	39.0000 U	41.0000 U
P Aroclor 1254	UG/KG	35.0000 U	40.0000 U	37.0000 U	180.0000 U	39.0000 U	41.0000 U
P Aroclor 1260	UG/KG	14.0000 J	95.0000	5.6000 J	350.0000 J	89.0000	150.0000
P Dieldrin	UG/KG	3.4000 J	28.0000 J	3.2000 J	77.0000	20.0000	46.0000
P Endosulfan I	UG/KG	1.8000 U	2.0000 U	1.9000 U	9.4000 U	2.0000 U	2.1000 U
P Endosulfan II	UG/KG	3.5000 U	4.0000 U	3.7000 U	18.0000 U	3.9000 U	4.1000 U
P Endosulfan sulfate	UG/KG	3.5000 U	4.0000 U	0.0530 J	18.0000 U	3.9000 U	4.1000 U
P Endrin	UG/KG	3.5000 U	0.4700 J	3.7000 U	18.0000 U	3.9000 U	0.2800 J
P Endrin aldehyde	UG/KG	3.5000 U	4.0000 U	3.7000 U	18.0000 U	3.9000 U	4.1000 U
P Endrin ketone	UG/KG	3.5000 U	4.0000 U	3.7000 U	18.0000 U	3.9000 U	4.1000 U
P Heptachlor	UG/KG	0.0580 U	2.0000 U	0.0410 U	0.9200 J	0.1300 U	0.5000 U
P Heptachlor epoxide	UG/KG	1.8000 U	0.3700 J	1.9000 U	9.4000 U	0.3400 J	4.4000 J
P Methoxychlor	UG/KG	18.0000 U	20.0000 U	0.1600 J	94.0000 U	20.0000 U	21.0000 U
P Toxaphene	UG/KG	180.0000 U	200.0000 U	190.0000 U	940.0000 U	200.0000 U	210.0000 U
P alpha-BHC	UG/KG	0.1200 J	2.0000 U	1.9000 U	9.4000 U	0.0710 J	2.1000 U
P alpha-Chlordane	UG/KG	1.3000 J	16.0000	1.0000 J	19.0000	4.1000 J	14.0000 J
P beta-BHC	UG/KG	1.8000 U	2.0000 U	1.9000 U	9.4000 U	2.0000 U	2.1000 U
P delta-BHC	UG/KG	1.8000 U	2.0000 U	0.1800 J	0.4600 J	0.5000 J	0.3400 J
P gamma-BHC (Lindane)	UG/KG	1.8000 U	2.0000 U	1.9000 U	9.4000 U	2.0000 U	2.1000 U
P gamma-Chlordane	UG/KG	0.8200 J	11.0000 J	1.9000 J	31.0000 J	5.0000 J	51.0000 J
S 1,2,4-Trichlorobenzene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 1,2-Dichlorobenzene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 1,3-Dichlorobenzene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 1,4-Dichlorobenzene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2,4,5-Trichlorophenol	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S34002	33S35002	33S35004	33S36002	33S36003	33S37002
	LAB ID:	532976	532977	532936	532942	532943	532957
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2,4-Dichlorophenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2,4-Dimethylphenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2,4-Dinitrophenol	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U
S 2,4-Dinitrotoluene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2,6-Dinitrotoluene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2-Chloronaphthalene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2-Chlorophenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2-Methylnaphthalene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2-Methylphenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 2-Nitroaniline	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U
S 2-Nitrophenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 3-Nitroaniline	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U
S 4-Bromophenyl-phenylether	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 4-Chloro-3-methylphenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 4-Chloroaniline	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 4-Methylphenol	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S 4-Nitroaniline	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U
S 4-Nitrophenol	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U
S Acenaphthene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Acenaphthylene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Anthracene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	50.0000 J
S Benzo(a)anthracene	UG/KG	350.0000 U	400.0000 U	370.0000 U	57.0000 J	72.0000 J	130.0000 J
S Benzo(a)pyrene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	46.0000 J	94.0000 J
S Benzo(b,k)fluoranthene	UG/KG	350.0000 U	64.0000 J	370.0000 U	160.0000 J	110.0000 J	230.0000 J
S Benzo(g,h,i)perylene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U

NAS PENSACOLA OU-10 ANI) SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	33S34002	33S35002	33S35004	33S36002	33S36003	33S37002
	LAB ID:	532976	532977	532936	532942	532943	532957
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	50.0000 U	120.0000 U	320.0000 U	110.0000 U	140.0000 U	91.0000 U
S Carbazole	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Chxysene	UG/KG	350.0000 U	400.0000 U	370.0000 U	50.0000 J	66.0000 J	99.0000 J
S Di-n-butylphthalate	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	73.0000 U	69.0000 U
S Di-n-octylphthalate	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Dibenz (a,h) anthracene	UG/KG	350,0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Dibenzofuran	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Diethylphthalate	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Dimethylphthalate	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Fluoranthene	UG/KG	350.0000 U	400.0000 U	370.0000 U	120.0000 J	200.0000 J	210.0000 J
S Fluorene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Hexachlorobenzene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Hexachlorobutadiene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Hexachlorocyclopentadiene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Hexachloroethane	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Isophorone	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S N-Nitrosodiphenylamine	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Naphthalene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Nitrobenzene	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S Pentachlorophenol	UG/KG	850.0000 U	970.0000 U	900.0000 U	890.0000 U	940.0000 U	990.0000 U
S Phenanthrene	UG/KG	350.0000 U	400.0000 U	370.0000 U	53.0000 J	65.0000 J	75.0000 J
S Phenol	UG/KG	350.0000 U	61.0000 J	370.0000 U	370.0000 U	64.0000 J	160.0000 J
S Pyrene	UG/KG	350.0000 U	400.0000 U	370.0000 U	67.0000 J	88.0000 J	110.0000 J
S bis(2-Chloroethoxy)methane	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S bis (2-Ethlyhexy)phthalate	UG/KG	50.0000 U	200.0000 U	110.0000 U	110.0000 U	620.0000	320.0000 U
S bis(2-chloroethyl)ether	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	350.0000 U	400.0000 U	370.0000 U	370.0000 U	390.0000 U	410.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33834002	33S35002	33835004	33836002	33836003	33837002
ENSAFE ID:		33834002	33S35002	33835004	33836002	33836003	33837002
LAB ID:		532976	532977	532936	532942	532943	532957
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 2-Butanone	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 2-Hexanone	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Acetone	UG/KG	61.0000 U	25.0000 U	430.0000	17.0000 U	23000.0000	32.0000 U
V Benzene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Bromoform	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Bromomethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Chlorobenzene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Chloroethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Chloroform	UG/KG	11.0000 U	12.0000 U	23.0000 U	2.0000 U	1400.0000 U	12.0000 U
V Chloromethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Ethylbenzene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Methylene Chloride	UG/KG	18.0000 U	21.0000 U	16.0000 U	28.0000,U	1100.0000 U	54.0000 U
V Styrene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Toluene	UG/KG	1.0000 J	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Trichloroethene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S34002	33S35002	33S35004	33S36002	33S36003	33S37002
	LAB ID:	532976	532977	532936	532942	532943	532957
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V Xylene (total)	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	12.0000 U	23.0000 U	11.0000 U	1400.0000 U	12.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not ehown in significant digits.

NAME	UNITS	33837004	33S37004RE	33838002	33S38004	33S38004DL	33S39002
ENSAFE ID:		33837004	33S37004RE	33838002	33S38004	33S38004DL	33S39002
LAB ID:		532966	532966	532959	532960	532960	532961
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
I Aluminum	MG/KG	2420.0000	.	4100.0000	8950.0000	.	10200.0000
I Antimony	MG/KG	12.7000 U		16.5000 U	13.8000 U	.	15.4000 U
I Arsenic	MG/KG	4.0000		1.6000 J	1.2000 J	.	1.4000 J
I Barium	MG/KG	8.8000 J	.	30.2000 J	61.2000	.	77.9000
I Beryllium	MG/KG	0.6700 U	.	0.8700 U	0.7200 U	.	0.8100 U
I Cadmium	MG/KG	18.3000	.	15.3000	31.8000	.	39.0000
I Calcium	MG/KG	1160.0000	.	1670.0000	2450.0000	.	2630.0000
I Chromium	MG/KG	561.0000	.	631.0000	1160.0000	.	1350.0000
I Cobalt	MG/KG	2.7000 U	.	3.5000 U	2.9000 U	.	3.2000 U
I Copper	MG/KG	19.7000	.	24.7000	114.0000	.	107.0000
I Cyanide	MG/KG	0.9200	.	0.7200 U	1.0000	.	0.7800
I Iron	MG/KG	575.0000	.	1590.0000	2000.0000	.	2610.0000
I Lead	MG/KG	86.9000	.	47.8000	70.0000	.	101.0000
I Magnesium	MG/KG	123.0000 U	.	160.0000 U	147.0000 J	.	280.0000 J
I Manganese	MG/KG	17.5000	.	20.7000	21.7000	.	48.8000
I Mercury	MG/KG	0.1500	.	0.6300	2.5000	.	0.3800
I Nickel	MG/KG	30.2000	.	19.0000	59.0000	.	62.6000
I Potassium	MG/KG	883.0000 U	.	1140.0000 U	955.0000 U	.	1070.0000 U
I Selenium	MG/KG	0.6700 U	.	0.8700 U	0.7200 U	.	0.8100 U
I Silver	MG/KG	2.4000	.	9.2000	29.7000	.	24.3000
I Sodium	MG/KG	124.0000 J	.	161.0000 J	143.0000 J	.	143.0000 J
I Thallium	MG/KG	0.6700 U	.	0.8700 U	0.7200 U	.	0.8100 U
I Vanadium	MG/KG	3.4000 U		4.5000 J	5.3000 J	.	7.3000 J
I Zinc	MG/KG	70.0000	.	92.5000	223.0000	.	240.0000
P 4,4'-DDD	UG/KG	3.7000 U	.	0.7500 J	1.3000 J	.	4.5000 U
P 4,4'-DDE	UG/KG	0.2900 J		6.9000	3.8000 J	.	3.8000 J
P 4,4'-DDT	UG/KG	3.7000 U	.	4.8000 U	4.0000 U	.	4.5000 U
P Aldrin	UG/KG	1.9000 U	.	1.5000 J	2.0000 U	.	0.7200 J

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	33S37004	33S37004RE	33S38002	33838004	33S38004DL	33839002
	LAB ID:	532966	532966	532959	532960	532960	532961
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
P Aroclor 1016	UG/KG	37.0000 U	.	48.0000 U	40.0000 U	.	45.0000 U
P Aroclor 1221	UG/KG	74.0000 U	.	97.0000 U	81.0000 U		91.0000 U
P Aroclor 1232	UG/KG	37.0000 U	.	48.0000 U	40.0000 U		45.0000 U
P Aroclor 1242	UG/KG	37.0000 U	.	48.0000 U	40.0000 U		45.0000 U
P Aroclor 1248	UG/KG	37.0000 U	.	48.0000 U	40.0000 U	.	45.0000 U
P Aroclor 1254	UG/KG	37.0000 U	.	48.0000 U	40.0000 U		45.0000 U
P Aroclor 1260	UG/KG	25.0000 J	.	210.0000	82.0000		210.0000
P Dieldrin	UG/KG	6.1000	.	4.8000 U	20.0000 J		28.0000 J
P Endosulfan I	UG/KG	1.9000 U	.	2.5000 U	2.0000 U		2.3000 U
P Endosulfan II	UG/KG	3.7000 U	.	4.8000 U	4.0000 U		4.5000 U
P Endosulfan sulfate	UG/KG	3.7000 U	.	4.8000 U	4.0000 U		4.5000 U
P Endrin	UG/KG	3.7000 U	.	4.8000 U	4.0000 U		1.0000 J
P Endrin aldehyde	UG/KG	3.7000 U	.	4.8000 U	4.0000 U		4.5000 U
P Endrin ketone	UG/KG	3.7000 U	.	4.8000 U	4.0000 U		4.5000 U
P Heptachlor	UG/KG	0.1800 U	.	2.5000 U	0.2400 U		2.3000 U
P Heptachlor epoxide	UG/KG	0.8300 J	.	2.5000 U	0.4200 J		0.5200 U
P Methoxychlor	UG/KG	19.0000 U	.	25.0000 U	20.0000 U		23.0000 U
P Toxaphene	UG/KG	190.0000 U	.	250.0000 U	200.0000 U		230.0000 U
P alpha-BHC	UG/KG	1.9000 U	.	0.1200 J	2.0000 U		0.1200 J
P alpha-Chlordane	UG/KG	2.1000 J	.	23.0000 J	8.1000 J		19.0000 J
P beta-BHC	UG/KG	1.9000 U	.	2.5000 U	2.0000 U		2.3000 U
P delta-BHC	UG/KG	1.9000 U	.	1.4000 J	0.5200 J		2.3000 U
P gamma-BHC (Lindane)	UG/KG	1.9000 U	.	2.5000 U	2.0000 U		2.3000 U
P gamma-Chlordane	UG/KG	11.0000 J	.	21.0000 J	14.0000 J		32.0000
S 1,2,4-Trichlorobenzene	UG/KG	360.0000 U	.	480.0000 U	1700.0000	1800.0000 U	440.0000 U
S 1,2-Dichlorobenzene	UG/KG	360.0000 U	.	480.0000 U	15000.0000 J	23000.0000 J	440.0000 U
S 1,3-Dichlorobenzene	UG/KG	360.0000 U	.	480.0000 U	1100.0000	1200.0000 J	440.0000 U
S 1,4-Dichlorobenzene	UG/KG	360.0000 U	.	480.0000 U	4100.0000 J	4600.0000 J	440.0000 U
S 2,4,5-Trichlorophenol	UG/KG	880.0000 U	.	1200.0000 U	950.0000 U	14000.0000 U	1100.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33837004	33S37004RE	33838002	33838004	33S38004DL	33839002
ENSAFE ID:		33837004	33S37004RE	33838002	33838004	33S38004DL	33839002
LAB ID:		532966	532966	532959	532960	532960	532961
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
S 2,4,6-Trichlorophenol	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2,4-Dichlorophenol	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2,4-Dimethylphenol	UG/KG	360.0000 U	.	480.0000 U	64.0000 J	5900.0000 U	440.0000 U
S 2,4-Dinitrophenol	UG/KG	880.0000 U	.	1200.0000 U	950.0000 U	14000.0000 U	1100.0000 U
S 2,4-Dinitrotoluene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2,6-Dinitrotoluene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2-Chloronaphthalene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2-Chlorophenol	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2-Methylnaphthalene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2-Methylphenol	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 2-Nitroaniline	UG/KG	880.0000 U	.	1200.0000 U	950.0000 U	14000.0000 U	1100.0000 U
S 2-Nitrophenol	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 3-Nitroaniline	UG/KG	880.0000 U	.	1200.0000 U	950.0000 U	14000.0000 U	1100.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	880.0000 U	.	1200.0000 U	950.0000 U	14000.0000 U	1100.0000 U
S 4-Bromophenyl-phenylether	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 4-Chloro-3-methylphenol	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 4-Chloroaniline	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 4-Methylphenol	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S 4-Nitroaniline	UG/KG	880.0000 U	.	1200.0000 U	950.0000 U	14000.0000 U	1100.0000 U
S 4-Nitrophenol	UG/KG	880.0000 U	.	1200.0000 U	950.0000 U	14000.0000 U	1100.0000 U
S Acenaphthene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Acenaphthylene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Anthracene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Benzo(a)anthracene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Benzo(a)pyrene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Benzo(b,k)fluoranthene	UG/KG	360.0000 U	.	81.0000 J	390.0000 U	5900.0000 U	440.0000 U
S Benzo(g,h,i)perylene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S37004	33S37004RE	33838002	33838004	33S38004DL	33839002
ENSAFE ID:		33S37004	33S37004RE	33838002	33838004	33S38004DL	33839002
LAB ID:		532966	532966	532959	532960	532960	532961
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
S Butylbenzylphthalate	UG/KG	110.0000 U		480.0000 U	160.0000 U	5900.0000 U	190.0000 U
S Carbazole	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Chrysene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Di-n-butylphthalate	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Di-n-octylphthalate	UG/KG	360.0000 U		480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Dibenz(a,h)anthracene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Dibenzofuran	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Diethylphthalate	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Dimethylphthalate	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Fluoranthene	UG/KG	360.0000 U	.	58.0000 J	390.0000 U	5900.0000 U	48.0000 J
S Fluorene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Hexachlorobenzene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Hexachlorobutadiene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Hexachlorocyclopentadiene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Hexachloroethane	UG/KG	360.0000 U	.	480.0000 U	390,0000 U	5900.0000 U	440.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Isophorone	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S N-Nitrosodiphenylamine	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Naphthalene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Nitrobenzene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Pentachlorophenol	UG/KG	880.0000 U	.	110.0000 J	950.0000 U	14000.0000 U	1100.0000 U
S Phenanthrene	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S Phenol	UG/KG	360.0000 U	.	160.0000 J	110.0000 J	5900.0000 U	170.0000 J
S Pyrene	UG/KG	360.0000 U		480.0000 U	390.0000,U	5900.0000 U	440.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	360.0000 U		480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	42.0000 U	.	280.0000 U	360.0000 U	5900.0000 U	730.0000
S bis(2-chloroethyl)ether	UG/KG	360.0000 U		480.0000 U	390.0000 U	5900.0000 U	440.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	360.0000 U	.	480.0000 U	390.0000 U	5900.0000 U	440.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33837004	33S37004RE	33838002	33S38004	33S38004DL	33839002
	LAB ID:	532966	532966	532959	532960	532960	532961
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 2-Butanone	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 2-Hexanone	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Acetone	UG/KG	760.0000 U	1600.0000 U	55.0000 U	2100.0000 U	.	77.0000 U
V Benzene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Bromoform	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Bromomethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Chlorobenzene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2900.0000	.	14.0000 U
V Chloroethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Chloroform	UG/KG	11.0000 U	56.0000 U	14.0000 U	300.0000 U	.	14.0000 U
V Chloromethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Ethylbenzene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Methylene Chloride	UG/KG	23.0000 U	64.0000 U	73.0000 U	4600.0000 U	.	29.0000 U
V Styrene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000'U	.	14.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Toluene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U
V Trichloroethene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U	.	14.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S37004	33S37004RE	33S38002	33S38004	33S38004DL	33S39002
	LAB ID:	532966	532966	532959	532960	532960	532961
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U		14.0000 U
V Xylene (total)	UG/KG	11.0000 U	56.0000 U	14.0000 U	290.0000 J		14.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U		14.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	56.0000 U	14.0000 U	2100.0000 U		14.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key pteeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S45002	33S45002DL	33S45004	33845006	33S45008	33S45008RE
	LAB ID:	546837	546837 D50	546840	546844	546847	546847
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
I Aluminum	MG/KG	2570.0000 J	.	949.0000 J	326.0000 J	262.0000 J	.
I Antimony	MG/KG	9.7000 U	.	9.6000 U	10.6000 U	11.4000 U	.
I Arsenic	MG/KG	1.2000 J	.	0.6300 U	0.6900 U	0.7500 U	.
I Barium	MG/KG	4.4000 J	.	58.9000	1.1000 U	1.2000 U	.
I Beryllium	MG/KG	0.4200 U	.	0.4200 U	0.4600 U	0.5000 U	.
I Cadmium	MG/KG	1.1000 U		1.0000 U	1.1000 U	1.2000 U	.
I Calcium	MG/KG	960.0000 J		2200.0000	866.0000 J	306.0000 J	.
I Chromium	MG/KG	3.6000		2.1000 U	5.2000	3.8000	.
I Cobalt	MG/KG	1.9000 U		1.9000 U	2.1000 U	2.2000 U	.
I Copper	MG/KG	5.3000 U		5.2000 U	5.7000 U	6.2000 U	.
I Cyanide	MG/KG	0.5300 U		0.5200 U	0.5700 U	0.6200 U	.
I Iron	MG/KG	1570.0000		1820.0000	237.0000	178.0000	.
I Lead	MG/KG	14.5000		2.8000	7.6000	2.9000	.
I Magnesium	MG/KG	100.0000 U		196.0000 J	109.0000 U	118.0000 U	.
I Manganese	MG/KG	17.4000		9.4000	1.9000 J	1.8000 J	.
I Mercury	MG/KG	0.1100 U		0.1000 U	0.1100 U	0.1200 U	.
I Nickel	MG/KG	6.5000 U		6.5000 U	7.1000 U	7.7000 U	.
I Potassium	MG/KG	471.0000 U		468.0000 U	514.0000 U	555.0000 U	.
I Selenium	MG/KG	0.6300 U		0.6300 U	0.6900 U	0.7500 U	.
I Silver	MG/KG	2.1000 U		2.1000 U	2.3000 U	2.5000 U	.
I Sodium	MG/KG	71.0000 J		84.6000 J	87.6000 J	70.4000 J	.
I Thallium	MG/KG	0.8400 U		0.8400 U	0.9200 U	0.9900 U	.
I Vanadium	MG/KG	3.8000 J		2.0000 J	1.6000 U	1.7000 U	.
I Zinc	MG/KG	18.1000 J		4.8000 J	3.8000 J	3.4000 J	.
P 4,4'-DDD	UG/KG	3.9000 J	18.0000 J	3.4000 U	0.3000 J	2.5000 J	.
P 4,4'-DDE	UG/KG	24.0000	26.0000 J	0.9800 J	2.6000 J	3.6000 J	.
P 4,4'-DDT	UG/KG	100.0000 J	120.0000 J	5.0000	11.0000 J	15.0000 J	.
P Aldrin	UG/KG	1.8000 U	8.9000 U	1.8000 U	2.0000 U	2.1000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S45002	33S45002DL	33S45004	33845006	33845008	33S45008RE
	LAB ID:	546837	546837 D50	546840	546844 R1	546847 R1	546847
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
P Aroclor 1016	UG/KG	35.0000 U	170.0000 U	34.0000 U	38.0000 U	41.0000 U	.
P Aroclor 1221	UG/KG	71.0000 U	350.0000 U	70.0000 U	77.0000 U	83.0000 U	.
P Aroclor 1232	UG/KG	35.0000 U	170.0000 U	34.0000 U	38.0000 U	41.0000 U	.
P Aroclor 1242	UG/KG	35.0000 U	170.0000 U	34.0000 U	38.0000 U	41.0000 U	.
P Aroclor 1248	UG/KG	35.0000 U	170.0000 U	34.0000 U	38.0000 U	41.0000 U	.
P Aroclor 1254	UG/KG	35.0000 U	170.0000 U	34.0000 U	38.0000 U	41.0000 U	.
P Aroclor 1260	UG/KG	35.0000 U	170.0000 U	34.0000 U	38.0000 U	41.0000 U	.
P Dieldrin	UG/KG	7.3000	7.3000 J	0.5700 J	0.7000 J	0.6100 J	.
P Endosulfan I	UG/KG	1.8000 U	8.9000 U	1.8000 U	2.0000 U	2.1000 U	.
P Endosulfan 11	UG/KG	3.5000 U	17.0000 U	3.4000 U	3.8000 U	4.1000 U	.
P Endosulfan sulfate	UG/KG	3.5000 U	17.0000 U	3.4000 U	3.8000 U	4.1000 U	.
P Endrin	UG/KG	3.5000 U	17.0000 U	3.4000 U	0.0780 J	4.1000 U	.
P Endrin aldehyde	UG/KG	3.5000 U	17.0000 U	3.4000 U	3.8000 U	4.1000 U	.
P Endrin ketone	UG/KG	3.5000 U	17.0000 U	3.4000 U	3.8000 U	4.1000 U	.
P Heptachlor	UG/KG	0.0890 U	8.9000 U	1.8000 U	0.1000 J	0.1400 U	.
P Heptachlor epoxide	UG/KG	1.8000 U	8.9000 U	1.8000 U	2.0000 U	2.1000 U	.
P Methoxychlor	UG/KG	18.0000 U	89.0000 U	18.0000 U	20.0000 U	21.0000 U	.
P Toxaphene	UG/KG	180.0000 U	890.0000 U	180.0000 U	200.0000 U	210.0000 U	.
P alpha-BHC	UG/KG	1.8000 U	8.9000 U	1.8000 U	2.0000 U	2.1000 U	.
P alpha-Chlordane	UG/KG	0.7700 J	0.7200 J	1.8000 U	2.0000 U	2.1000 U	.
P beta-BHC	UG/KG	1.8000 U	8.9000 U	1.8000 U	2.0000 U	2.1000 U	.
P delta-BHC	UG/KG	1.8000 U	8.9000 U	1.8000 U	2.0000 U	2.1000 U	.
P gamma-BHC (Lindane)	UG/KG	1.8000 U	8.9000 U	1.8000 U	2.0000 U	2.1000 U	
P gama-Chlordane	UG/KG	0.5900 J	8.9000 U	1.8000 U	0.0690 J	2.1000 U	.
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	
S 1,2-Dichlorobenzene	UG/KG	340.0000 U		340.0000 U	370.0000 U	A00.0000 U	
S 1,3-Dichlorobenzene	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S 1,4-Dichlorobenzene	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S 2,4,5-Trichlorophenol	UG/KG	830.0000 U		820.0000 U	910.0000 U	980.0000 U	

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33845002	33S45002DL	33S45004	33845006	33S45000	33S45008RE
ENSAFE ID:		33845002	33S45002DL	33S45004	33845006	33S45000	33S45008RE
LAB ID:		546837	546837 D50	546040	546844	546847	546047
DATE RCVD:		04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S 2,4-Dichlorophenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S 2,4-Dimethylphenol	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S 2,4-Dinitrophenol	UG/KG	830.0000 U	.	820.0000 U	910.0000 U	980.0000 U	.
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 2,6-Dinitrotoluene	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S 2-Chloronaphthalene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 2-Chlorophenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S 2-Methylnaphthalene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 2-Methylphenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 2-Nitroaniline	UG/KG	830.0000 U	.	820.0000 U	910.0000 U	980.0000 U	.
S 2-Nitrophenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 3-Nitroaniline	UG/KG	830.0000 U	.	820.0000 U	910.0000 U	980.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/KG	830.0000 U	.	820.0000 U	910.0000 U	980.0000 U	.
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S 4-Chloroaniline	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S 4-Methylphenol	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S 4-Nitroaniline	UG/KG	830.0000 U	.	820.0000 U	910.0000 U	980.0000 U	
S 4-Nitrophenol	UG/KG	87.0000 J	.	820.0000 U	910.0000 U	980.0000 U	.
S Acenaphthene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S Acenaphthylene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S Anthracene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S Benzo(a)anthracene	UG/KG	250.0000 J		340.0000 U	370.0000 U	400.0000 U	.
S Benzo(a)pyrene	UG/KG	99.0000 J	.	340.0000 U	370.0000 U	400.0000 U	.
S Benzo(b,k)fluoranthene	UG/KG	350.0000 J	.	43.0000 J	370.0000 U	400.0000 U	.
S Benzo(g,h,i)perylene	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33845002	33S45002DL	33845004	33845006	33S45008	33S45008RE
	LAB ID:	546837	546837 D50	546840	546844	546847	546847
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S Carbazole	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S Chrysene	UG/KG	130.0000 J		340.0000 U	370.0000 U	400.0000 U	.
S Di-n-butylphthalate	UG/KG	42.0000 J	.	340.0000 U	58.0000 J	400.0000 U	.
S Di-n-octylphthalate	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S Dibenz(a,h)anthracene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S Dibenzofuran	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S Diethylphthalate	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	
S Dimethylphthalate	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S Fluoranthene	UG/KG	170.0000 J		340.0000 U	370.0000 U	400.0000 U	.
S Fluorene	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	.
S Hexachlorobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S Hexachlorobutadiene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S Hexachloroethane	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/KG	110.0000 J	.	340.0000 U	370.0000 U	400.0000 U	.
S Isophorone	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S Naphthalene	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	
S Nitrobenzene	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	
S Pentachlorophenol	UG/KG	830.0000 U	.	820.0000 U	910.0000 U	980.0000 U	.
S Phenanthrene	UG/KG	81.0000 J	.	340.0000 U	370.0000 U	400.0000 U	.
S Phenol	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S Pyrene	UG/KG	200.0000 J		340.0000 U	370.0000 U	400.0000 U	
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/KG	81.0000 U		40.0000 U	75.0000 U	54.0000 U	
S bis(2-chloroethyl)ether	UG/KG	340.0000 U		340.0000 U	370.0000 U	400.0000 U	
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	.	340.0000 U	370.0000 U	400.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	33S45002	33S45002DL	33S45004	33S45006	33S45008	33S45008RE
	LAB ID:	546837	546837 D50	546840	546844	546847	546847
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U		10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 2-Butanone	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 2-Hexanone	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Acetone	UG/KG	11.0000 U	.	38.0000 U	46.0000 U	430.0000 U	2000.0000 U
V Benzene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Bromoform	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Bromomethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Chlorobenzene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Chloroethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Chloroform	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	7.0000 U
V Chloromethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Ethylbenzene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Methylene Chloride	UG/KG	31.0000 U	.	10.0000 U	34.0000 U	15.0000 U	150.0000 U
V Styrene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Toluene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Trichloroethene	UG/KG	11.0000 U	.	10.0000 U	11.0000 U	12.0000 U	62.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S45002	33S45002DL	33S45004	33S45006	33S45008	33S45008RE
	LAB ID:	546837	546837 D50	546840	546844	546847	546847
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U		10.0000 U	11.0000 U	12.0000 U	62.0000 U
V Xylene (total)	UG/KG	11.0000 U		10.0000 U	11.0000 U	12.0000 U	62.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U		10.0000 U	11.0000 U	12.0000 U	62.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U		10.0000 U	11.0000 U	12.0000 U	62.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S46002	33S46002DL	33S46004	33846006	33846008	33847002
	LAB ID:	546825	546825 D20	546828	546831	546834	546809
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
I Aluminum	MG/KG	1360.0000 J		475.0000 J	771.0000 J	646.0000 J	1190.0000 J
I Antimony	MG/KG	9.6000 U		10.0000 U	10.5000 U	12.3000 U	9.6000 U
I Arsenic	MG/KG	0.6300 J		0.6500 U	0.6800 U	0.8000 U	0.6200 U
I Barium	MG/KG	3.4000 J		2.8000 J	7.1000 J	2.7000 J	6.6000 J
I Beryllium	MG/KG	0.4200 U		0.4300 U	0.4600 U	0.5300 U	0.4200 U
I Cadmium	MG/KG	1.0000 U		1.1000 U	1.1000 U	1.3000 U	1.0000 U
I Calcium	MG/KG	265.0000 J		264.0000 J	194.0000 J	246.0000 J	958.0000 J
I Chromium	MG/KG	6.5000		2.2000 U	3.0000	2.7000	7.3000
I Cobalt	MG/KG	1.9000 U		2.0000 U	2.1000 U	2.4000 U	1.9000 U
I Copper	MG/KG	5.2000 U		5.4000 U	6.5000	6.7000 U	9.5000
I Cyanide	MG/KG	0.5200 U		0.5400 U	0.5700 U	0.6600 U	0.5200 U
I Iron	MG/KG	926.0000		372.0000	432.0000	495.0000	808.0000
I Lead	MG/KG	13.7000		16.0000	22.7000	9.3000	19.7000
I Magnesium	MG/KG	99.6000 U		103.0000 U	109.0000 U	127.0000 U	99.2000 U
I Manganese	MG/KG	7.6000		2.2000 J	2.2000 J	3.9000 J	9.3000
I Mercury	MG/KG	0.1000 U		0.1100 U	0.1100 U	0.1300 U	0.1000 U
I Nickel	MG/KG	6.5000 U		6.7000 U	7.1000 U	8.3000 U	6.5000 U
I Potassium	MG/KG	468.0000 U		485.0000 U	510.0000 U	595.0000 U	466.0000 U
I Selenium	MG/KG	0.6300 U		0.6500 U	0.6800 U	0.8000 U	0.6200 U
I Silver	MG/KG	2.1000 U		2.2000 U	2.3000 U	2.7000 U	2.1000 U
I Sodium	MG/KG	68.7000 J		69.8000 J	61.1000 J	77.8000 J	70.6000 J
I Thallium	MG/KG	0.8400 U		0.8700 U	0.9100 U	1.1000 U	0.8300 U
I Vanadium	MG/KG	2.1000 J		1.5000 U	1.6000 U	1.9000 U	2.6000 J
I Zinc	MG/KG	10.0000 J		8.7000 J	19.4000 J	8.1000 J	16.5000 J
P 4,4'-DDD	UG/KG	0.8600 J	2.0000 J	3.6000 U	3.1000 J	52.0000	3.4000 U
P 4,4'-DDE	UG/KG	18.0000	19.0000 J	2.1000 J	3.3000 J	6.5000 J	3.4000 U
P 4,4'-DDT	UG/KG	60.0000 J	70.0000	3.8000	2.5000 J	3.2000 J	3.4000 U
P Aldrin	UG/KG	0.0880 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U

NAS PENSACOLA OU-10 ANI) SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S46002	33S46002DL	33846004	33S46006	33846008	33847002
	LAB ID:	546825	546825 D20	546828	546831	546834	546809
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
P Aroclor 1016	UG/KG	34.0000 U	68.0000 U	36.0000 U	38.0000 U	44.0000 U	34.0000 U
P Aroclor 1221	UG/KG	69.0000 U	140.0000 U	73.0000 U	76.0000 U	89.0000 U	70.0000 U
P Aroclor 1232	UG/KG	34.0000 U	68.0000 U	36.0000 U	38.0000 U	44.0000 U	34.0000 U
P Aroclor 1242	UG/KG	34.0000 U	68.0000 U	36.0000 U	38.0000 U	44.0000 U	34.0000 U
P Aroclor 1248	UG/KG	34.0000 U	68.0000 U	36.0000 U	38.0000 U	44.0000 U	34.0000 U
P Aroclor 1254	UG/KG	34.0000 U	68.0000 U	36.0000 U	38.0000 U	44.0000 U	34.0000 U
P Aroclor 1260	UG/KG	34.0000 U	68.0000 U	36.0000 U	38.0000 U	44.0000 U	82.0000
P Dieldrin	UG/KG	33.0000	36.0000 J	0.8200 J	0.8800 J	0.3000 J	3.4000 U
P Endosulfan I	UG/KG	1.8000 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U
P Endosulfan II	UG/KG	3.4000 U	6.8000 U	'3.6000 U	3.8000 U	4.4000 U	3.4000 U
P Endosulfan sulfate	UG/KG	0.3700 U	0.5500 J	3.6000 U	3.8000 U	0.4500 J	3.4000 U
P Endrin	UG/KG	0.2700 J	6.8000 U	3.6000 U	3.8000 U	0.3800 J	3.4000 U
P Endrin aldehyde	UG/KG	3.4000 U	6.8000 U	3.6000 U	3.8000 U	4.4000 U	3.4000 U
P Endrin ketone	UG/KG	1.4000 J	1.4000 J	3.6000 U	3.8000 U	4.4000 U	3.4000 U
P Heptachlor	UG/KG	1.8000 U	3.5000 U	1.8000 U	0.1300 U	0.1900 U	1.8000 U
P Heptachlor epoxide	UG/KG	1.8000 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U
P Methoxychlor	UG/KG	18.0000 U	35.0000 U	18.0000 U	19.0000 U	23.0000 U	18.0000 U
P Toxaphene	UG/KG	180.0000 U	350.0000 U	180.0000 U	190.0000 U	230.0000 U	180.0000 U
P alpha-BHC	UG/KG	1.8000 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U
P alpha-Chlordane	UG/KG	0.2300 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U
P beta-BHC	UG/KG	1.8000 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U
P delta-BHC	UG/KG	1.8000 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	3.5000 U	1.8000 U	1.9000 U	2.3000 U	1.8000 U
P gamma-Chlordane	UG/KG	0.2300 J	0.2300 J	1.8000 U	0.0820 J	2.3000 U	1.8000 U
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U		350.0000 U	370.0000 U	430.0000 U	340.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U		350.0000 U	370.0000 U	t430.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U		350.0000 U	370.0000 U	430.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U		350.0000 U	370.0000 U	430.0000 U	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	820.0000 U		860.0000 U	900.0000 U	1100.0000 U	820.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	33846002	33S46002DL	33546004	33846006	33546008	33S47002
	LAB ID:	546825	546825 D20	546828	546831	546834	546809
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
2,4,6-Trichlorophenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2,4-Dichlorophenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2,4-Dimethylphenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2,4-Dinitrophenol	UG/KG	820.0000 U	.	860.0000 U	900.0000 U	1100.0000 U	820.0000 U
2,4-Dinitrotoluene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2,6-Dinitrotoluene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2-Chloronaphthalene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2-Chlorophenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2-Methylnaphthalene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2-Methylphenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
2-Nitroaniline	UG/KG	820.0000 U	.	860.0000 U	900.0000 U	1100.0000 U	820.0000 U
2-Nitrophenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
3,3'-Dichlorobenzidine	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
3-Nitroaniline	UG/KG	820.0000 U	.	860.0000 U	900.0000 U	1100.0000 U	820.0000 U
4,6-Dinitro-2-methylphenol	UG/KG	820.0000 U	.	860.0000 U	900.0000 U	1100.0000 U	820.0000 U
4-Bromophenyl-phenylether	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
4-Chloro-3-methylphenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
4-Chloroaniline	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
4-Chlorophenyl-phenylether	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
4-Methylphenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
4-Nitroaniline	UG/KG	820.0000 U	.	860.0000 U	900.0000 U	1100.0000 U	820.0000 U
4-Nitrophenol	UG/KG	820.0000 U	.	860.0000 U	900.0000 U	1100.0000 U	820.0000 U
Acenaphrhene	UG/KG	35.0000 J	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
Acenaphthylene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
Anthracene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
Benzo(a)anthracene	UG/KG	220.0000 J	.	350.0000 U	370.0000 U	190.0000 J	340.0000 U
Benzo(a)pyrene	UG/KG	110.0000 J	.	350.0000 U	370.0000 U	87.0000 J	340.0000 U
Benzo(b,k)fluoranthene	UG/KG	330.0000 J	.	350.0000 U	370.0000 U	190.0000 J	340.0000 U
Benzo(g,h,i)perylene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S46002	33S46002DL	33S46004	33S46006	33S46008	33S47002
	LAB ID:	546825	546825 D20	546828	546831	546834	546809
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	51.0000 J	45.0000 J
S Carbazole	UG/KG	50.0000 J	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Chrysene	UG/KG	250.0000 J	.	350.0000 U	370.0000 U	110.0000 J	340.0000 U
S Di-n-butylphthalate	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	52.0000 J	340.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Dibenz(a,h)anthracene	UG/KG	46.0000 J	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Dibenzofuran	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Diethylphthalate	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Fluoranthene	UG/KG	450.0000	.	350.0000 U	370.0000 U	150.0000 J	340.0000 U
S Fluorene	UG/KG	40.0000 J	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Hexachloroethane	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	110.0000 J	.	350.0000 U	370.0000 U	69.0000 J	340.0000 U
S Isophorone	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Naphthalene	UG/KG	55.0000 J	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Nitrobenzene	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	820.0000 U	.	860.0000 U	900.0000 U	1100.0000 U	820.0000 U
S Phenanthrene	UG/KG	410.0000	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Phenol	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S Pyrene	UG/KG	320.0000 J	.	350.0000 U	370.0000 U	140.0000 J	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	56.0000 U	.	350.0000 U	54.0000 U	53.0000 U	60.0000 U
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	.	350.0000 U	370.0000 U	430.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33846002	33846002DL	33S46004	33846006	33846008	33847002
	LAB ID:	546825	546825 D20	546028	546831	546834	546809
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 2-Butanone	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Acetone	UG/KG	17.0000 U		46.0000 U	37.0000 U	48.0000 U	18.0000 U
V Benzene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Chloroform	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Methylene Chloride	UG/KG	7.0000 U		15.0000 U	14.0000 U	35.0000 'U	9.0000 U
V Styrene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	. .	11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Trichloroethene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S46002	33S46002DL	33S46004	33S46006	33S46008	33S47002
	LAB ID:	546825	546825 D20	546828	546831	546834	546809
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U		11.0000 U	11.0000 U	13.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33847004	33347006	33850002	33350004	33851002	33851004
	LAB ID:	546812	546819	546892	546896	546886	546889
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
I Aluminum	MG/KG	593.0000 J	670.0000 J	1390.0000	662.0000	4250.0000	6000.0000
I Antimony	MG/KG	10.6000 U	11.0000 U	9.9000 U	10.5000 U	9.7000 U	10.2000 U
I Arsenic	MG/KG	0.6900 U	0.7200 U	0.6500 U	0.6800 U	0.9800 J	1.8000 J
I Barium	MG/KG	4.8000 J	4.6000 J	4.2000 J	1.6000 J	12.3000 J	40.9000 J
I Beryllium	MG/KG	0.4600 U	0.4800 U	0.4300 U	0.4500 U	0.4200 U	0.4400 U
I Cadmium	MG/KG	1.2000 U	1.2000 U	2.1000	1.1000 U	2.2000	7.5000
I Calcium	MG/KG	522.0000 J	765.0000 J	948.0000 J	327.0000 J	624.0000 J	1540.0000
I Chromium	MG/KG	2.3000 U	4.3000	45.1000	20.9000	48.3000	174.0000
I Cobalt	MG/KG	2.1000 U	2.2000 U	1.9000 U	2.0000 U	1.9000 U	2.0000 U
I Copper	MG/KG	5.8000 U	6.0000 U	7.0000	5.7000 U	13.8000	41.3000
I Cyanide	MG/KG	0.5800 U	0.6000 U	0.5400 U	0.5700 U	0.5300 U	0.6900
I Iron	MG/KG	641.0000	384.0000	755.0000	442.0000	2630.0000	3560.0000
I Lead	MG/KG	14.3000	15.8000	28.4000	4.1000	13.3000	82.8000
I Magnesium	MG/KG	110.0000 U	114.0000 U	103.0000 U	108.0000 U	101.0000 U	109.0000 J
I Manganese	MG/KG	4.3000	4.6000	12.9000	2.0000 J	33.0000	25.8000
I Mercury	MG/KG	0.1200 U	0.1200 U	0.1100 U	0.1100 U	0.1100 U	0.3600
I Nickel	MG/KG	7.1000 U	7.4000 U	8.8000	7.0000 U	6.6000 U	7.3000 J
I Potassium	MG/KG	514.0000 U	537.0000 U	482.0000 U	508.0000 U	474.0000 U	496.0000 U
I Selenium	MG/KG	0.6900 U	0.7200 U	0.6500 U	0.6800 U	0.6400 U	0.6700 U
I Silver	MG/KG	2.3000 U	2.4000 U	2.2000 U	2.3000 U	2.1000 U	6.7000
I Sodium	MG/KG	70.1000 J	74.5000 J	100.0000 J	94.8000 J	98.7000 J	106.0000 J
I Thallium	MG/KG	0.9200 U	0.9600 U	0.8600 U	0.9100 U	0.8500 U	0.8900 U
I Vanadium	MG/KG	1.6000 U	1.7000 J	2.0000 J	1.6000 U	5.9000 J	7.0000 J
I Zinc	MG/KG	10.4000 J	9.5000 J	20.4000	5.3000	27.3000	77.8000
P 4,4'-DDD	UG/KG	5.8000	3.9000 U	3.5000 U	3.7000 U	3.5000 U	67.0000 J
P 4,4'-DDE	UG/KG	3.2000 J	0.7300 J	3.5000 U	3.7000 U	2.7000 R	78.0000 J
P 4,4'-DDT	UG/KG	3.7000 U	1.1000 J	3.5000 U	0.8600 J	2.3000 R	4.5000 J
P Aldrin	UG/KG	1.9000 U	2.0000 U	1,8000 U	1.9000 U	0.1200 R	1.9000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33847004	33847006	33850002	33850004	33851002	33851004
ENSAFE ID:		33847004	33847006	33850002	33850004	33851002	33851004
LAB ID:		546812	546819	546892	546896	546886 R1	546889
DATE RCVD:		04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
P Aroclor 1016	UG/KG	37.0000 U	39.0000 U	35.0000 U	37.0000 U	35.0000 U	36.0000 U
P Aroclor 1221	UG/KG	76.0000 U	79.0000 U	72.0000 U	76.0000 U	71.0000 U	74.0000 U
P Aroclor 1232	UG/KG	37.0000 U	39.0000 U	35.0000 U	37.0000 U	35.0000 U	36.0000 U
P Aroclor 1242	UG/KG	37.0000 U	39.0000 U	35.0000 U	37.0000 U	35.0000 U	36.0000 U
P Aroclor 1248	UG/KG	37.0000 U	39.0000 U	35.0000 U	37.0000 U	35.0000 U	36.0000 U
P Aroclor 1254	UG/KG	37.0000 U	39.0000 U	35.0000 U	37.0000 U	35.0000 U	36.0000 U
P Aroclor 1260	UG/KG	10.0000 J	10.0000 J	180.0000	79.0000	35.0000 U	36.0000 U
P Dieldrin	UG/KG	3.7000 U	0.7500 J	3.5000 U	3.9000 J	0.6000 R	11.0000
P Endosulfan I	UG/KG	1.9000 U	2.0000 U	1.8000 U	1.9000 U	1.8000 U	1.9000 U
P Endosulfan II	UG/KG	3.7000 U	3.9000 U	3.5000 U	3.7000 U	3.5000 U	3.6000 U
P Endosulfan sulfate	UG/KG	3.7000 U	3.9000 U	3.5000 U	3.7000 U	3.5000 U	3.6000 U
P Endrin	UG/KG	3.7000 U	3.9000 U	3.5000 U	1.6000 J	3.5000 U	3.6000 U
P Endrin aldehyde	UG/KG	3.7000 U	3.9000 U	3.5000 U	3.7000 U	3.5000 U	3.6000 U
P Endrin ketone	UG/KG	3.7000 U	3.9000 U	3.5000 U	3.7000 U	3.5000 U	3.6000 U
P Heptachlor	UG/KG	1.9000 U	0.1900 U	1.8000 U	1.9000 U	0.1200 R	1.9000 U
P Heptachlor epoxide	UG/KG	1.9000 U	2.0000 U	0.3500 J	0.3200 J	1.8000 U	1.9000 U
P Methoxychlor	UG/KG	19.0000 U	20.0000 U	18.0000 U	19.0000 U	18.0000 U	19.0000 U
P Toxaphene	UG/KG	190.0000 U	200.0000 U	180.0000 U	190.0000 U	180.0000 U	190.0000 U
P alpha-BHC	UG/KG	1.9000 U	2.0000 U	1.8000 U	0.1600 J	1.8000 U	1.9000 U
P alpha-Chlordane	UG/KG	1.9000 U	2.0000 U	1.6000 J	1.9000 U	0.1300 R	3.0000 J
P beta-BHC	UG/KG	1.9000 U	2.0000 U	1.8000 U	2.9000 J	1.8000 U	1.9000 U
P delta-BHC	UG/KG	1.9000 U	2.0000 U	0.3400 J	0.4900 J	1.8000 U	1.9000 U
P gamma-BHC (Lindane	UG/KG	1.9000 U	2.0000 U	0.1000 J	1.9000 U	1.8000 U	1.9000 U
P gamma-Chlordane	UG/KG	1.9000 U	2.0000 U	3.1000 J	1.9000 U	1.2000 R	3.0000 J
S 1,2,4-Trichlorobenzene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 1,2-Dichlorobenzene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	12000.0000	350.0000 U	370.0000 U
S 1,3-Dichlorobenzene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	8600.0000	350.0000 U	370.0000 U
S 1,4-Dichlorobenzene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	12000.0000	350.0000 U	370.0000 U
S 2,4,5-Trichlorophenol	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
ENSAFE ID:		33S47004	33S47006	33850002	33550004	33651002	33S51004
LAB ID:		546812	546819	546892	546896	546886	546889
DATE RCVD:		04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
S 2,4,6-Trichlorophenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2,4-Dichlorophenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2,4-Dimethylphenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2,4-Dinitrophenol	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U
S 2,4-Dinitrotoluene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2,6-Dinitrotoluene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2-Chloronaphthalene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2-Chlorophenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2-Methylnaphthalene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	24000.0000	350.0000 U	370.0000 U
S 2-Methylphenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 2-Nitroaniline	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U
S 2-Nitrophenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 3-Nitroaniline	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U
S 4-Bromophenyl-phenylether	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 4-Chloro-3-methylphenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 4-Chloroaniline	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 4-Methylphenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S 4-Nitroaniline	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U
S 4-Nitrophenol	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U
S Acenaphthene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	1900.0000 J	350.0000 U	370.0000 U
S Acenaphthylene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Anthracene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	38.0000 J	370.0000 U
S Benzo(a)anthracene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	870.0000	81.0000 J
S Benzo(a)pyrene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	380.0000	85.0000 J
S Benzo(b,k)fluoranthene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	1000.0000 J	210.0000 J
S Benzo(g,h,i)perylene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	340.0000 J	370.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	33947004	33547006	33850002	33850004	33851002	33851004
	LAB ID:	546812	546819	546892	546896	546886	546889
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
S Butylbenzylphthalate	UG/KG	44.0000 J	390.0000 U	3500.0000 U	1800.0000 J	350.0000 U	87.0000 J
S Carbazole	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Chrysene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	690.0000	74.0000 J
S Di-n-butylphthalate	UG/KG	370.0000 U	390.0000 U	370.0000 J	7500.0000 U	350.0000 U	370.0000 U
S Di-n-octylphthalate	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Dibenz (a,h) anthracene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	84.0000 J	370.0000 U
S Dibenzofuran	UG/KG	370.0000 U	390.0000 U	3500.0000 U	1400.0000 J	350.0000 U	370.0000 U
S Diethylphthalate	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Dimethylphthalate	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Fluoranthene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	1900.0000 J	1800.0000	95.0000 J
S Fluorene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	3600.0000 J	350.0000 U	370.0000 U
S Hexachlorobenzene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Hexachlorobutadiene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Hexachlorocyclopentadiene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Hexachloroethane	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Indeno(1,2,3-cd) pyrene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	390.0000	370.0000 U
S Ieophorone	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S N-Nitrosodiphenylamine	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Naphthalene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	10000.0000	350.0000 U	370.0000 U
S Nitrobenzene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Pentachlorophenol	UG/KG	910.0000 U	950.0000 U	8600.0000 U	18000.0000 U	850.0000 U	890.0000 U
S Phenanthrene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	6600.0000 J	410.0000	370.0000 U
S Phenol	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S Pyrene	UG/KG	370.0000 U	390.0000 U	3500.0000 U	2000.0000.J	1100.0000	86.0000 J
S bis(2-Chloroethoxy) methane	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S bis(2-Ethlyhexy) phthalate	UG/KG	72.0000 U	54.0000 U	3000.0000 J	9600.0000 J	46.0000 U	1000.0000 J
S bis (2-chloroethyl) ether	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U
S bis(2-chloroisopropyl) ether	UG/KG	370.0000 U	390.0000 U	3500.0000 U	7500.0000 U	350.0000 U	370.0000 U

NAS YENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in signific nt digits.

	ENSAFE ID:	33547004	33S47006	33850002	33850004	33S51002	33851004
	LAB ID:	546812	546819	546892	546896	546886	546889
	DATE RCM:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400,0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	11.0000 U	12.0000 U	11.0000 U	720.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	51.0000 U	43.0000 U	65.0000 U	1400.0000 U	44.0000 U	41.0000 U
V Benzene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Chloroethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Chloroform	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	2.0000 J
V Chloromethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	15.0000 U	11.0000 U	36.0000 U	700.0000. U	37.0000 U	53.0000 U
V Styrene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Trichloroethene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33847004	33847006	33850002	33850004	33851002	33851004
	LAB ID:	546812	546819	546892	546896	546886	546889
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	11.0000 U	12.0000 U	11.0000 U	460.0000 J	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	12.0000 U	11.0000 U	1400.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preaented at end of Appendix. Data not shown in significant digita.

	ENSAFE ID:	33S51004DL	33852002	33852004	33853002	33853004	33854002
	LAB ID:	546889 D50	546880	546883	546755	546776	546779
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
I Aluminum	MG/KG	.	1100.0000	1090.0000	942.0000 J	58.9000 J	570.0000 J
I Antimony	MG/KG	.	9.6000 U	10.5000 U	10.5000 U	10.4000 U	9.5000 U
I Arsenic	MG/KG		1.4000 J	0.7200 J	0.8800 J	0.6800 U	0.6200 U
I Barium	MG/KG		12.7000 J	11.9000 J	35.1000 J	2.1000 J	7.9000 J
I Beryllium	MG/KG		0.4200 U	0.4600 U	0.4600 U	0.4500 U	0.4100 U
I Cadmium	MG/KG		1.0000 U	1.1000 U	16.9000	1.3000	1.3000
I Calcium	MG/KG	.	331.0000 J	202.0000 J	2570.0000	74.5000 J	812.0000 J
I Chromium	MG/KG	.	2.1000 U	2.3000 U	105.0000	20.5000	19.0000
I Cobalt	MG/KG	.	1.9000 U	2.1000 U	2.1000 U	2.0000 U	1.9000 U
I Copper	MG/KG		5.9000	6.4000	18.9000	5.7000 U	5.2000 U
I Cyanide	MG/KG	.	0.5200 U	0.5700 U	0.5700 U	0.5700 U	0.5200 U
I Iron	MG/KG	.	785.0000	812.0000	1630.0000	59.9000	244.0000
I Lead	MG/KG	.	38.7000	37.4000	21.0000	2.3000	3.1000
I Magnesium	MG/KG	.	99.4000 U	109.0000 U	406.0000 J	108.0000 U	174.0000 J
I Manganese	MG/KG	.	16.3000	7.3000	76.2000	2.2000 J	12.6000
I Mercury	MG/KG	.	0.1000 U	0.1100 U	0.2700	0.1100 U	0.1000 U
I Nickel	MG/KG	.	6.5000 U	7.1000 U	7.1000 U	7.0000 U	6.4000 U
I Potassium	MG/KG	.	467.0000 U	511.0000 U	510.0000 U	507.0000 U	461.0000 U
I Selenium	MG/KG	.	0.6300 U	0.6900 U	0.6800 U	0.6800 U	0.6200 U
I Silver	MG/KG	.	2.1000 U	2.3000 U	4.8000	2.3000 U	2.1000 U
I Sodium	MG/KG		131.0000 J	103.0000 J	67.1000 J	76.3000 J	88.0000 J
I Thallium	MG/KG		0.8400 U	0.9100 U	0.9100 U	0.9100 U	0.8300 U
I Vanadium	MG/KG		1.5000 U	1.6000 U	1.6000 U	1.6000 U	1.4000 U
I Zinc	MG/KG		31.1000	23.2000	63.6000 J	7.2000'J	11.7000 J
P 4,4'-DDD	UG/KG	96.0000	0.4900 J	1.7000 J	3.7000 U	3.7000 U	3.4000 U
P 4,4'-DDE	UG/KG	130.0000	1.1000 J	2.2000 J	6.6000	1.1000 J	0.3300 J
P 4,4'-DDT	UG/KG	5.4000 J	1.7000 J	1.3000 J	3.7000 U	3.7000 U	3.4000 U
P Aldrin	UG/KG	9.4000 U	1.8000 U	2.0000 U	1.9000 U	1.9000 U	1.7000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S51004DL	33S52002	33S52004	33S53002	33S53004	33854002
	LAB ID:	546889 D50	546880	546883	546755	546776	546779
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
P Aroclor 1016	UG/KG	180.0000 U	34.0000 U	38.0000 U	37.0000 U	37.0000 U	34.0000 U
P Aroclor 1221	UG/KG	370.0000 U	70.0000 U	77.0000 U	76.0000 U	75.0000 U	69.0000 U
P Aroclor 1232	UG/KG	180.0000 U	34.0000 U	38.0000 U	37.0000 U	37.0000 U	34.0000 U
P Aroclor 1242	UG/KG	180.0000 U	34.0000 U	38.0000 U	37.0000 U	37.0000 U	34.0000 U
P Aroclor 1248	UG/KG	180.0000 U	34.0000 U	38.0000 U	37.0000 U	37.0000 U	34.0000 U
P Aroclor 1254	UG/KG	180.0000 U	34.0000 U	38.0000 U	37.0000 U	37.0000 U	34.0000 U
P Aroclor 1260	UG/KG	180.0000 U	34.0000 U	38.0000 U	310.0000	50.0000	34.0000
P Dieldrin	UG/KG	12.0000 J	0.4200 J	0.3100 J	4.2000	0.2600 J	1.0000 J
P Endosulfan I	UG/KG	9.4000 U	1.8000 U	2.0000 U	1.9000 U	1.9000 U	1.7000 U
P Endosulfan II	UG/KG	18.0000 U	3.4000 U	3.8000 U	3.7000 U	3.7000 U	3.4000 U
P Endosulfan sulfate	UG/KG	18.0000 U	3.4000 U	3.8000 U	3.7000 U	3.7000 U	3.4000 U
P Endrin	UG/KG	18.0000 U	3.4000 U	3.8000 U	3.7000 U	3.7000 U	3.4000 U
P Endrin aldehyde	UG/KG	18.0000 U	3.4000 U	3.8000 U	3.7000 U	3.7000 U	3.4000 U
P Endrin ketone	UG/KG	18.0000 U	3.4000 U	3.8000 U	3.7000 U	3.7000 U	3.4000 U
P Heptachlor	UG/KG	9.4000 U	1.8000 U	2.0000 U	1.9000 U	1.9000 U	1.9000 U
P Heptachlor epoxide	UG/KG	9.4000 U	0.1500 J	2.0000 U	0.1700 J	1.9000 U	1.7000 U
P Methoxychlor	UG/KG	94.0000 U	18.0000 U	20.0000 U	19.0000 U	19.0000 U	17.0000 U
P Toxaphene	UG/KG	940.0000 U	180.0000 U	200.0000 U	190.0000 U	190.0000 U	170.0000 U
P alpha-BHC	UG/KG	9.4000 U	1.8000 U	2.0000 U	1.9000 U	1.9000 U	1.7000 U
P alpha-Chlordane	UG/KG	3.5000 J	1.9000	0.4100 J	2.7000	0.4600 J	0.5500 J
P beta-BHC	UG/KG	9.4000 U	1.8000 U	2.0000 U	1.9000 U	1.9000 U	1.7000 U
P delta-BHC	UG/KG	9.4000 U	1.8000 U	2.0000 U	1.9000 U	1.9000 U	1.7000 U
P gamma-BHC (Lindane)	UG/KG	9.4000 U	1.8000 U	2.0000 U	1.9000 U	1.9000 U	1.7000 U
P gamma-Chlordane	UG/KG	3.4000 J	1.7000 J	0.1900 J	0.7000 J	1.9000 U	1.7000 U
S 1,2,4-Trichlorobenzene	UG/KG		340.0000 U	380.0000 U	370.0000,U	370.0000 U	340.0000 U
S 1,2-Dichlorobenzene	UG/KG		340.0000 U	380.0000 U	370.0000 U	'370.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG		340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG		340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	.	830.0000 U	910.0000 U	900.0000 U	900.0000 U	810.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S51004DL	33S52002	33852004	33853002	33S53004	33854002
	LAB ID:	546889 D50	546880	546883	546755	546776	546779
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Carbazole	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Chrysene	UG/KG	.	39.0000 J	380.0000 U	170.0000 J	370.0000 U	340.0000 U
S Di-n-butylphthalate	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Di-n-octylphthalate	UG/KG	.	340.0000 U	380.0000 U	53.0000 U	370.0000 U	340.0000 U
S Dibenz(a,h)anthracene	UG/KG	.	340.0000 U	380.0000 U	47.0000 J	370.0000 U	340.0000 U
S Dibenzofuran	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Diethylphthalate	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Dimethylphthalate	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Fluoranthene	UG/KG	.	84.0000 J	380.0000 U	52.0000 J	370.0000 U	340.0000 U
S Fluorene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Hexachloroethane	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	.	49.0000 J	380.0000 U	130.0000 J	370.0000 U	340.0000 U
S Ieophorone	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Naphthalene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Nitrobenzene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	.	830.0000 U	910.0000 U	900.0000 U	900.0000 U	810.0000 U
S Phenanthrene	UG/KG	.	62.0000 J	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Phenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Pyrene	UG/KG	.	61.0000 J	380.0000 U	84.0000 J	370.0000 U	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	.	280.0000 U	380.0000 U	56.0000 U	46.0000 U	57.0000 U
S bis(2-chloroethyl)ether	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S51004DL	33852002	33S52004	33853002	33853004	33554002
	LAB ID:	546889 D50	546880	546883	546755	546776	546779
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2,4-Dichlorophenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2,4-Dimethylphenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2,4-Dinitrophenol	UG/KG	.	830.0000 U	910.0000 U	900.0000 U	900.0000 U	810.0000 U
S 2,4-Dinitrotoluene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2,6-Dinitrotoluene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2-Chloronaphthalene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2-Chlorophenol	UG/KG	.	340.0000 U	380.0000 U	370,0000 U	370.0000 U	340.0000 U
S 2-Methylnaphthalene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2-Methylphenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 2-Nitroaniline	UG/KG	.	830.0000 U	910.0000 U	900.0000 U	900.0000 U	810.0000 U
S 2-Nitrophenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 3-Nitroaniline	UG/KG	.	830.0000 U	910.0000 U	900.0000 U	900.0000 U	810.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	.	830.0000 U	910.0000 U	900.0000 U	900.0000 U	810.0000 U
S 4-Bromophenyl-phenylether	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 4-Chloro-3-methylphenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 4-Chloroaniline	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 4-Methylphenol	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S 4-Nitroaniline	UG/KG	.	830.0000 U	910.0000 U	900.0000 U	900.0000 U	810.0000 U
S 4-Nitrophenol	UG/KG	.	830.0000 U	910.0000 U	48.0000 J	900.0000 U	810.0000 U
S Acenaphthene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Acenaphthylene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Anthracene	UG/KG	.	340.0000 U	380.0000 U	370.0000 U	370.0000 U	340.0000 U
S Benzo(a)anthracene	UG/KG	.	38.0000 J	380.0000 U	63.0000'5	370.0000 U	340.0000 U
S Beneo(a)pyrene	UG/KG	.	44.0000 J	380.0000 U	120.0000 J	370.0000 U	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG	.	68.0000 J	380.0000 U	170.0000 J	370.0000 U	340.0000 U
S Benzo(g,h,i)perylene	UG/KG	.	340.0000 U	380.0000 U	71.0000 J	370.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S51004DL	33852002	33852004	33553002	33553004	33554002
	LAB ID:	546889 D50	546880	546883	546755	546776	546779
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	.	10.0000 U	11.0000 U	11,0000 U	11.0000 U	10.0000 U
V 2-Butanone	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 2-Hexanone	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Acetone	UG/KG	.	130.0000 U	200.0000 U	26.0000 U	11.0000 U	10.0000 U
V Benzene	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Bromoform	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Bromomethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Chlorobenzene	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Chloroethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Chloroform	UG/KG	.	1.0000 J	1.0000 J	11.0000 U	11.0000 U	10.0000 U
V Chloromethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Ethylbenzene	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Methylene Chloride	UG/KG	.	8.0000 U	29.0000 U	32.0000 U	12.0000 U	16.0000 U
V Styrene	UG/KG	.	10.0000 U	11,0000 U	11.0000 U	11.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Toluene	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Trichloroethene	UG/KG	.	10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33854004	33855002	33855004	33856002	33857001	33S57001DL
ENSAFE ID:		33854004	33855002	33855004	33856002	33857001	33S57001DL
LAB ID:		546787	546791	546799	546803	547126	547126 D51
DATE RCVD:		04/21/93	04/21/93	04/21/93	04/21/93	04/22/93	04/22/93
I Aluminum	MG/KG	37.0000 J	651.0000 J	64.5000 J	1330.0000 J	1830.0000	
I Antimony	MG/KG	10.2000 U	9.6000 U	10.6000 U	10.9000 U	11.9000 U	
I Arsenic	MG/KG	0.6600 U	0.6200 U	0.6900 U	0.7100 U	1.2000 J	
I Barium	MG/KG	1.4000 J	15.8000 J	1.1000 U	1.7000 J	43.9000 J	
I Beryllium	MG/KG	0.4400 U	0.4200 U	0.4600 U	0.4800 U	0.5200 U	
I Cadmium	MG/KG	1.1000 U	4.5000	1.1000 U	1.2000 U	241.0000	
I Calcium	MG/KG	59.2000 J	1370.0000	51.0000 J	108.0000 J	3050.0000	
I Chromium	MG/KG	11.7000	21.1000	10.0000	8.2000	4110.0000	
I Cobalt	MG/KG	2.0000 U	1.9000 U	2.1000 U	2.1000 U	2.3000 U	
I Copper	MG/KG	5.5000 U	5.2000 U	5.7000 U	5.9000 U	403.0000	
I Cyanide	MG/KG	0.5500 U	0.5200 U	0.5700 U	0.5900 U	52.1000	
I Iron	MG/KG	29.2000	528.0000	40.0000	824.0000	19300.0000	
I Lead	MG/KG	1.2000	2.5000	1.0000	12.8000	756.0000	
I Magnesium	MG/KG	105.0000 U	327.0000 J	109.0000 U	113.0000 U	167.0000 J	
I Manganese	MG/KG	1.9000 J	18.1000	1.8000 J	3.5000 J	88.8000 J	
I Mercury	MG/KG	0.1100 U	0.1000 U	0.1100 U	0.1200 U	2.6000	
I Nickel	MG/KG	6.9000 U	6.4000 U	7.1000 U	7.4000 U	15.4000	
I Potassium	MG/KG	494.0000 U	464.0000 U	513.0000 U	532.0000 U	579.0000 U	
I Selenium	MG/KG	0.6600 U	0.6200 U	0.6900 U	0.7100 U	0.7800 U	
I Silver	MG/KG	2.2000 U	2.1000 U	2.3000 U	2.4000 U	5.4000	
I Sodium	MG/KG	76.4000 J	97.0000 J	71.1000 J	84.3000 J	144.0000 J	
I Thallium	MG/KG	0.8800 U	0.8300 U	0.9200 U	0.9500 U	1.0000 U	
I Vanadium	MG/KG	1.5000 U	1.5000 U	1.6000 U	1.7000 U	8.5000 J	
I Zinc	MG/KG	3.9000 J	26.2000 J	3.8000 J	8.3000 J	656.0000	
P 4,4'-DDD	UG/KG	3.6000 U	3.4000 U	3.7000 U	3.9000 U	89.0000 J	79.0000 J
P 4,4'-DDE	UG/KG	0.7200 J	1.7000 J	1.1000 J	9.6000	420.0000 J	490.0000
P 4,4'-DDT	UG/KG	3.6000 U	3.4000 U	3.7000 U	10.0000	520.0000 J	580.0000
P Aldrin	UG/KG	1.9000 U	0.4000 J	1.9000 U	0.1200 J	11.0000 U	110.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33554004	33S55002	33855004	33856002	33S57001	33S57001DL
ENSAFE ID:		33554004	33S55002	33855004	33856002	33S57001	33S57001DL
LAB ID:		546787	546791	546799	546803	547126 D50	547126 D51
DATE RCVD:		04/21/93	04/21/93	04/21/93	04/21/93	04/22/93	04/22/93
P Aroclor 1016	UG/KG	36.0000 U	34.0000 U	37.0000 U	39.0000 U	210.0000 U	2100.0000 U
P Aroclor 1221	UG/KG	74.0000 U	69.0000 U	76.0000 U	78.0000 U	430.0000 U	4300.0000 U
P Aroclor 1232	UG/KG	36.0000 U	34.0000 U	37.0000 U	39.0000 U	210.0000 U	2100.0000 U
P Aroclor 1242	UG/KG	36.0000 U	34.0000 U	37.0000 U	39.0000 U	210.0000 U	2100.0000 U
P Aroclor 1248	UG/KG	36.0000 U	34.0000 U	37.0000 U	39.0000 U	210.0000 U	2100.0000 U
P Aroclor 1254	UG/KG	36.0000 U	34.0000 U	37.0000 U	39.0000 U	210.0000 U	2100.0000 U
P Aroclor 1260	UG/KG	62.0000	100.0000	65.0000	39.0000 U	210.0000 U	2100.0000 U
P Dieldrin	UG/KG	0.4100 J	0.5200 J	0.2300 J	3.9000 U	21.0000 U	210.0000 U
P Endoeulfan I	UG/KG	1.9000 U	1.8000 U	1.9000 U	0.1800 J	1.3000 J	110.0000 U
P Endosulfan II	UG/KG	3.6000 U	3.4000 U	3.7000 U	3.9000 U	21.0000 U	210.0000 U
P Endosulfan sulfate	UG/KG	3.6000 U	3.4000 U	3.7000 U	0.6200 J	21.0000 U	15.0000 U
P Endrin	UG/KG	3.6000 U	3.4000 U	3.7000 U	3.9000 U	21.0000 U	210.0000 U
P Endrin aldehyde	UG/KG	3.6000 U	3.4000 U	3.7000 U	0.4400 J	21.0000 U	15.0000 U
P Endrin ketone	UG/KG	3.6000 U	3.4000 U	3.7000 U	0.2200 J	21.0000 U	210.0000 U
P Heptachlor	UG/KG	1.9000 U	1.8000 U	1.9000 U	2.0000 U	11.0000 U	110.0000 U
P Heptachlor epoxide	UG/KG	1.9000 U	1.8000 U	1.9000 U	2.0000 U	11.0000 U	110.0000 U
P Methoxychlor	UG/KG	19.0000 U	18.0000 U	19.0000 U	20.0000 U	23.0000 J	1100.0000 U
P Toxaphene	UG/KG	190.0000 U	180.0000 U	190.0000 U	200.0000 U	1100.0000 U	11000.0000 U
P alpha-BHC	UG/KG	1.9000 U	1.8000 U	1.9000 U	2.0000 U	11.0000 U	110.0000 U
P alpha-Chlordane	UG/KG	0.4500 J	0.2300 J	0.1700 J	2.0000 U	11.0000 U	110.0000 U
P beta-BHC	UG/KG	1.9000 U	1.8000 U	1.9000 U	2.0000 U	11.0000 U	110.0000 U
P delta-BHC	UG/KG	1.9000 U	1.8000 U	1.9000 U	2.0000 U	11.0000 U	110.0000 U
P gamma-BHC (Lindane)	UG/KG	1.9000 U	1.8000 U	1.9000 U	2.0000 U	11.0000 U	110.0000 U
P gamma-Chlordane	UG/KG	0.5900 J	1.8000 U	1.9000 U	2.0000 U	0.9000 J	110.0000 U
S 1,2,4-Trichlorobenzene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	
S 1,2-Dichlorobenzene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	
S 1,3-Dichlorobenzene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	
S 1,4-Dichlorobenzene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	
S 2,4,5-Trichlorophenol	UG/KG	880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	

NAS PENSACOLA OU-11 SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S54004	33S55002	33S55004	33S56002	33S57001	33S57001DL
	LAB ID:	546787	546791	546799	546803	547126	547126 D51
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93	04/22/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2,4-Dichlorophenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2,4-Dimethylphenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2,4-Dinitrophenol	UG/KG	880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	.
S 2,4-Dinitrotoluene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2,6-Dinitrotoluene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2-Chloronaphthalene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2-Chlorophenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2-Methylnaphthalene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2-Methylphenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 2-Nitroaniline	UG/KG	880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	.
S 2-Nitrophenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 3,3'-Dichlorobenzidine	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 3-Nitroaniline	UG/KG	880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/KG	880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	.
S 4-Bromophenyl-phenylether	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 4-Chloro-3-methylphenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 4-Chloroaniline	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 4-Chlorophenyl-phenylether	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 4-Methylphenol	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S 4-Nitroaniline	UG/KG	880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	.
S 4-Nitrophenol	UG/KG	880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	.
S Acenaphthene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S Acenaphthylene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	43.0000 U	.
S Anthracene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	43.0000 U	.
S Benzo(a)anthracene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	220.0000 U	.
S Benzo(a)pyrene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	200.0000 U	.
S Benzo(b,k)fluoranthene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	480.0000 U	.
S Benzo(g,h,i)perylene	UG/KG	360.0000 U	340.0000 U	380.0000 U	390.0000 U	530.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME		ENSAFE ID:	33854004	33555002	33855004	33856002	33S57001	33S57001DL
		LAB ID:	546787	546791	546799	546803	547126	547126 D51
		DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93	04/22/93
		WITS						
Butylbenzylphthalate	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Carbazole	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Chrysene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	120.0000 J	.
Di-n-butylphthalate	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	99.0000 J	.
Di-n-octylphthalate	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Dibenz(a,h)anthracene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	87.0000 J	.
Dibenzofuran	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Diethylphthalate	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Dimethylphthalate	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Fluoranthene	UG/KG		360.0000 U	340.0000 U	380.0000 U	44.0000 J	200.0000 J	.
Fluorene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Hexachlorobenzene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Hexachlorobutadiene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Hexachlorocyclopentadiene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Hexachloroethane	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Indeno(1,2,3-cd)pyrene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	390.0000 J	.
Isophorone	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
N-Nitroso-di-n-propylamine	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
N-Nitrosodiphenylamine	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Naphthalene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
Nitrobenzene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S Pentachlorophenol	UG/KG		880.0000 U	820.0000 U	910.0000 U	940.0000 U	1000.0000 U	.
S Phenanthrene	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	49.0000 J	.
S Phenol	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S Pyrene	UG/KG		360.0000 U	340.0000 U	380.0000 U	45.0000 J	140.0000 J	.
S bis(2-Chloroethoxy)methane	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000' U	420.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/KG		42.0000 U	70.0000 U	380.0000 U	47.0000 U	48.0000 U	.
S bis(2-chloroethyl)ether	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.
S bis(2-chloroisopropyl)ether	UG/KG		360.0000 U	340.0000 U	380.0000 U	390.0000 U	420.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33854004	33555002	33855004	33856002	33S57001	33S57001DL
	LAB ID:	546787	546791	546799	546803	547126	547126 D51
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93	04/22/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 1,1-Dichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 1,1-Dichloroethene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 1,2-Dichloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 1,2-Dichloropropane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 2-Butanone	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 2-Hexanone	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Acetone	UG/KG	44.0000 U	10.0000 U	26.0000 U	27.0000 U	9.0000 U	
V Benzene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Bromodichloromethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Bromoform	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Bromomethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Carbon Disulfide	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Carbon Tetrachloride	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Chlorobenzene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Chloroethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Chloroform	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	1.0000 J	
V Chloromethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Dibromochloromethane	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Ethylbenzene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Methylene Chloride	UG/KG	24.0000 U	13.0000 U	11.0000 U	12.0000,U	44.0000 U	
V Styrene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Tetrachloroethene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Toluene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Trichloroethene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	3.0000 J	

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S54004	33S55002	33S55004	33S56002	33S57001	33S57001DL
	LAB ID:	546787	546791	546799	546803	547126	547126 D51
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93	04/22/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V Xylene (total)	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	10.0000 U	11.0000 U	12.0000 U	13.0000 U	

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
ENSAFE ID:		33S57B02	33S57B02DL	33S57B04			
LAB ID:		547128	547128 D50	547122			
DATE RCM:		04/22/93	04/22/93	04/22/93			
I Aluminum	MG/KG	381.0000		170.0000	.	.	.
I Antimony	MG/KG	9.5000 U		9.8000 U	.	.	.
I Arsenic	MG/KG	0.6200 U		0.6400 U	.	.	.
I Barium	MG/KG	2.5000 J		2.1000 J	.	.	.
I Beryllium	MG/KG	0.4100 U		0.4300 U	.	.	.
I Cadmium	MG/KG	1.0000 U		1.1000 U	.	.	.
I Calcium	MG/KG	32.3000 U		167.0000 J	.	.	.
I Chromium	MG/KG	9.9000		2.6000	.	.	.
I Cobalt	MG/KG	1.9000 U		1.9000 U	.	.	.
I Copper	MG/KG	5.2000 U		5.3000 U	.	.	.
I Cyanide	MG/KG	1.3000		0.5300 U	.	.	.
I Iron	MG/KG	307.0000		184.0000	.	.	.
I Lead	MG/KG	35.0000		2.7000	.	.	.
I Magnesium	MG/KG	98.7000 U		101.0000 U	.	.	.
I Manganese	MG/KG	1.8000 J		1.7000 J	.	.	.
I Mercury	MG/KG	0.1000 U		0.1100 U	.	.	.
I Nickel	MG/KG	6.4000 U		6.6000 U	.	.	.
I Potassium	MG/KG	463.0000 U		477.0000 U	.	.	.
I Selenium	MG/KG	0.6200 U		0.6400 U	.	.	.
I Silver	MG/KG	2.1000 U		2.1000 U	.	.	.
I Sodium	MG/KG	95.4000 J		63.3000 J	.	.	.
I Thallium	MG/KG	0.8300 U		0.8500 U	.	.	.
I Vanadium	MG/KG	1.5000 U		1.5000 U	.	.	.
I Zinc	MG/KG	6.8000		4.0000 J	.	.	.
P 4,4'-DDD	UG/KG	2.0000 J	17.0000 U	0.2700 J	.	.	.
P 4,4'-DDE	UG/KG	64.0000 J	73.0000	9.8000	.	.	.
P 4,4'-DDT	UG/KG	28.0000 J	29.0000 J	4.0000 J	.	.	.
P Aldrin	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.

NAS I'ENSACOLA OU-10 *AND* SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57B02	33S57B02DL	33557804			
	LAB ID:	547128	547128 D50	547122			
	DATE RCVD:	04/22/93	04/22/93	04/22/93	.	.	.
NAME	UNITS						
P Aroclor 1016	UG/KG	34.0000 U	170.0000 U	35.0000 U	.	.	
P Aroclor 1221	UG/KG	68.0000 U	340.0000 U	71.0000 U	.	.	.
P Aroclor 1232	UG/KG	34.0000 U	170.0000 U	35.0000 U	.	.	.
P Aroclor 1242	UG/KG	34.0000 U	170.0000 U	35.0000 U	.	.	.
P Aroclor 1248	UG/KG	34.0000 U	170.0000 U	35.0000 U	.	.	
P Aroclor 1254	UG/KG	34.0000 U	170.0000 U	35.0000 U	.	.	
P Aroclor 1260	UG/KG	34.0000 U	170.0000 U	35.0000 U	.	.	.
P Dieldrin	UG/KG	3.4000 U	17.0000 U	3.5000 U	.	.	
P Endosulfan I	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
P Endosulfan II	UG/KG	3.4000 U	17.0000 U	3.5000 U	.	.	.
P Endosulfan sulfate	UG/KG	0.4700 J	4.1000 J	3.5000 U	.	.	
P Endrin	UG/KG	3.4000 U	17.0000 U	3.5000 U	.	.	.
P Endrin aldehyde	UG/KG	0.8400 J	1.6000 J	3.5000 U	.	.	.
P Endrin ketone	UG/KG	3.4000 U	17.0000 U	3.5000 U	.	.	.
P Heptachlor	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
P Heptachlor epoxide	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
P Methoxychlor	UG/KG	0.4100 J	87.0000 U	18.0000 U	.	.	.
P Toxaphene	UG/KG	170.0000 U	870.0000 U	180.0000 U	.	.	.
P alpha-BHC	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
P alpha-Chlordane	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
P beta-BHC	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
P delta-BHC	UG/KG	1.7000 U	8.7000 U	1.8000 U		.	.
P gamma-BHC (Lindane)	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
P gamma-Chlordane	UG/KG	1.7000 U	8.7000 U	1.8000 U	.	.	.
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U		350.0000 U	.	.	.
S 1,2-Dichlorobenzene	UG/KG	340.0000 U		350.0000 U	.	.	
S 1,3-Dichlorobenzene	UG/KG	340.0000 U		350.0000 U	.	.	.
S 1,4-Dichlorobenzene	UG/KG	340.0000 U		350.0000 U	.	.	.
S 2,4,5-Trichlorophenol	UG/KG	810.0000 U		840.0000 U	.	.	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57B02	33S57B02DL	33S57B04			
	LAB ID:	547128	547128 D50	547122			
	DATE RCVD:	04/22/93	04/22/93	04/22/93			
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 2,4-Dichlorophenol	UG/KG	340.0000 U	.	350.0000 U	.	.	
S 2,4-Dimethylphenol	UG/KG	340.0000 U	.	350.0000 U	.	.	
S 2,4-Dinitrophenol	UG/KG	810.0000 U	.	840.0000 U	.	.	
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	.	350.0000 U	.	.	
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	.	350.0000 U	.	.	
S 2-Chloronaphthalene	UG/KG	340.0000 U	.	350.0000 U	.	.	
S 2-Chlorophenol	UG/KG	340.0000 U	.	350.0000 U	.	.	
S 2-Methylnaphthalene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 2-Methylphenol	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 2-Nitroaniline	UG/KG	810.0000 U	.	840.0000 U	.	.	
S 2-Nitrophenol	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 3-Nitroaniline	UG/KG	810.0000 U	.	840.0000 U	.	.	.
S 4,6-Dinitro-2-methylphenol	UG/KG	810.0000 U	.	840.0000 U	.	.	.
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	.	350.0000 U	.	.	
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 4-Chloroaniline	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 4-Methylphenol	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S 4-Nitroaniline	UG/KG	810.0000 U	.	840.0000 U	.	.	.
S 4-Nitrophenol	UG/KG	810.0000 U	.	840.0000 U	.	.	.
S Acenaphthene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Acenaphthylene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Anthracene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Benzo(a)anthracene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Benzo(a)pyrene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	.	350.0000 U	.	.	.

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57B02	33S57B02DL	33857804			
	LAB ID:	547128	547128 D50	547122			
	DATE RCM:	04/22/93	04/22/93	04/22/93	.	.	.
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	340.0000 U	.	350.0000 U		.	.
S Carbazole	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Chrysene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Di-n-butylphthalate	UG/KG	340.0000 U	.	350.0000 U		.	.
S Di-n-octylphthalate	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Dibenz (a,h) anthracene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Dibenzofuran	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Diethylphthalate	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Dimethylphthalate	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Fluoranthene	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Fluorene	UG/KG	340.0000 U	.	350.0000 U		.	
S Hexachlorobenzene	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Hexachlorobutadiene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Hexachloroethane	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Indeno (1,2,3-cd) pyrene	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Isophorone	UG/KG	340.0000 U	.	350.0000 U	.	.	
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	.	350.0000 U	.	.	
S Naphthalene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Nitrobenzene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Pentachlorophenol	UG/KG	810.0000 U	.	110.0000 J	.	.	
S Phenanthrene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Phenol	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S Pyrene	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S bis (2-Chloroethoxy) methane	UG/KG	340.0000 U	.	350.0000 U	.	.	.
S bis (2-Ethlyhexy) phthalate	UG/KG	340.0000 U	.	98.0000 U	.		.
S bis (2-chloroethyl) ether	UG/KG	340.0000 U	.	350.0000 U	.		.
S bis (2-chloroisopropyl) ether	UG/KG	340.0000 U	.	350.0000 U	.	.	.

NAS PENSACOLA OU-IO AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57B02	33S57B02DL	33S57B04			
	LAB ID:	547128	547128 D50	547122			
	DATE RCVD:	04/22/93	04/22/93	04/22/93	.	.	.
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	.	11.0000 U	.	.	
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	.	11.0000 U	.	.	
V 1,1-Dichloroethane	UG/KG	10.0000 U	.	11.0000 U	.	.	
V 1,1-Dichloroethene	UG/KG	10.0000 U	.	11.0000 U	.	.	
V 1,2-Dichloroethane	UG/KG	10.0000 U	.	11.0000 U	.	.	
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V 1,2-Dichloropropane	UG/KG	10.0000 U	.	11.0000 U		.	.
V 2-Butanone	UG/KG	10.0000 U	.	11.0000 U	.	.	
V 2-Hexanone	UG/KG	10.0000 U	.	11.0000 U	.	.	
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	.	11.0000 U	.	.	
V Acetone	UG/KG	33.0000 U	.	38.0000 U	.	.	
V Benzene	UG/KG	10.0000 U	.	11.0000 U	.	.	
V Bromodichloromethane	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V Bromoform	UG/KG	10.0000 U	.	11.0000 U	.	.	
V Bromomethane	UG/KG	10.0000 U	.	11.0000 U		.	
V Carbon Disulfide	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V Carbon Tetrachloride	UG/KG	10.0000 U	.	11.0000 U	.	.	
V Chlorobenzene	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V Chloroethane	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V Chloroform	UG/KG	10.0000 U	.	1.0000 J	.	.	.
V Chloromethane	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V Dibromochloromethane	UG/KG	10.0000 U	.	11.0000 U	.		.
V Ethylbenzene	UG/KG	10.0000 U	.	11.0000 U	.	.	.
V Methylene Chloride	UG/KG	37.0000 U		20.0000 U	.		.
V Styrene	UG/KG	10.0000 U	.	11.0000 U	.		.
V Tetrachloroethene	UG/KG	10.0000 U		11.0000 U	.		.
V Toluene	UG/KG	10.0000 U		11.0000 U	.	.	.
V Trichloroethene	UG/KG	10.0000 U	.	11.0000 U	.	.	

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57B02	33S57B02DL	33S57B04
	LAB ID:	547128	547128 D50	547122
	DATE RCVD:	04/22/93	04/22/93	04/22/93
NAME	UNITS			
V Vinyl Chloride	UG/KG	10.0000 U		11.0000 U
V xylene (total)	UG/KG	10.0000 U		11.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U		11.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U		11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SOIL SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S51004DL	33S52002	33S52004	33S53002	33S53004	33S54002
	LAB ID:	546889 D50	546880	546883	546755	546776	546779
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93
NAME	UNITS						
V Vinyl Chloride	UG/KG		10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V Xylene (total)	UG/KG		10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG		10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG		10.0000 U	11.0000 U	11.0000 U	11.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

APPENDIX M

AUXILIARY DATA — SEDIMENT SAMPLE COLLECTION

Sample Location/ Identification	Collection lime	Water Depth (feet)	Substrate Physical Description
33M01001	1525	1.0	Heavily vegetated substrate area with a mixture of sandy and dark gray silty sediment
33M02001	1510	2.5	Soft substrate dark gray to black silty sandy sediments
33M03001	1445	2.0	Firm substrate, mostly sandy, medium to dark brown sediment with minor amounts of silt
33M04001	1415	2.5	Soft substrate, dark gray to black silty-sandy sediments

Sediment Sampling
Auxiliary Field Data

Collection Data: 1/13/93
Weather: Mostly cloudy, light to moderate easterly winds
Temperature - approximately 60°F - 65°F

Sample Location/ Identification	Collection Time	Physical Description
13M61001	1400	Dark gray, silty, clay with minor amounts of sand, wet and highly plastic at edge of standing water
13M62001	1405	Dark gray clayey sand, wet at edge of standing water
13M63001	1345	Thin layer (2-3mm) of dark gray, silty clay mud, dry and cracked, underlain by light tan to orange, fine to medium grained quartz sand
13M65001	1350	Thin layer (2-3mm) of dark gray silty clay mud, dry and cracked, underlain by light tan and clayey orange, fine to medium grained quartz sand
13M65001	1440	Dark gray silty clay with minor amount of fine grained sand, wet and highly plastic at edge of standing water
13M66001	1445	Dark gray silty clay with minor amount of fine grained sand, wet and highly plastic at edge of standing water
13M67001	1450	Thin layer (2-3mm) of dark gray silty clay mud underlain by liaht tan. fine arained quartz sand

APPENDIX N

DETECTED CONCENTRATIONS — SEDIMENT SAMPLES

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end'of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001	13162001	13M63001	13M64001	13M65001	13M66001
	LAB ID:	532122	532123	532124	532125	532126	532127
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
I Aluminum	MG/KG	8970.0000	5110.0000	2950.0000	1310.0000	26600.0000	4290.0000
I Antimony	MG/KG	15.4000 U	24.4000 U	11.7000 U	11.9000 U	16.5000 U	20.7000 U
I Arsenic	MG/KG	0.8100 U	3.2000 J	0.8000 J	0.6200 U	16.7000	2.9000 J
I Barium	MG/KG	10.6000 J	6.6000 J	3.4000 J	1.2000 U	33.9000 J	6.8000 J
I Beryllium	MG/KG	0.8100 U	1.3000 U	0.6200 U	0.6200 U	1.4000 J	1.1000 U
I Cadmium	MG/KG	1.4000	2.1000 U	1.0000 U	1.0000 U	5.9000	1.8000 U
I Calcium	MG/KG	1610.0000	871.0000 J	1190.0000	98.8000 J	7620.0000	3620.0000
I Chromium	MG/KG	58.1000	35.0000	11.7000	3.6000	159.0000	21.0000
I Cobalt	MG/KG	3.2000 U	5.1000 U	2.5000 U	2.5000 U	6.8000 J	4.4000 U
I Copper	MG/KG	17.3000	10.7000 U	5.1000 U	5.2000 U	35.7000	9.1000 U
I Cyanide	MG/KG	0.6800 U	1.1000 U	0.5100 U	0.5200 U	0.7200 U	0.9100 U
I Iron	MG/KG	8620.0000	5170.0000	2380.0000	733.0000	27000.0000	3390.0000
I Lead	MG/KG	28.5000	11.2000	4.7000	12.7000	108.0000	12.6000
I Magnesium	MG/KG	2180.0000	1420.0000 J	541.0000 J	115.0000 U	7920.0000	909.0000 J
I Manganese	MG/KG	46.4000	31.8000	10.5000	1.7000 J	143.0000	21.3000
I Mercury	MG/KG	0.1400 U	0.2100 U	0.1000 U	0.1000 U	0.3100	0.1800 U
I Nickel	MG/KG	10.3000 U	16.2000 U	7.8000 U	7.9000 U	20.3000	13.8000 U
I Potassium	MG/KG	1070.0000 U	1690.0000 U	814.0000 U	824.0000 U	1830.0000	1440.0000 U
I Selenium	MG/KG	0.8100 U	1.3000 U	0.8200 U	0.6200 U	5.8000	1.5000 U
I Silver	MG/KG	2.7000 U	4.3000 U	2.1000 U	2.1000 U	2.9000 U	3.6000 U
I Sodium	MG/KG	257.0000 J	459.0000 J	121.0000 J	124.0000 J	869.0000 J	246.0000 J
I Thallium	MG/KG	0.8100 U	1.3000 U	0.6200 U	0.6200 U	0.8700 U	1.1000 U
I Vanadium	MG/KG	23.5000	12.3000 J	5.5000 J	3.1000 U	71.9000	8.9000 J
I Zinc	MG/KG	51.3000	30.5000	10.8000	2.5000 J	178.0000	23.9000
P 4,4'-DDD	UG/KG	4.5000 U	1.2000 J	0.7200 J	2.2000 J	4.7000 U	5.2000 J
P 4,4'-DDE	UG/KG	4.5000 U	1.6000 J	10.0000	5.8000	4.7000 U	3.5000 J
P 4,4'-DDT	UG/KG	4.5000 U	7,0000 U	5.8000	14.0000	4.7000 U	5.8000 U
P Aldrin	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	3.0000 U
P Aroclor 1016	UG/KG	45.0000 U	70.0000 U	33.0000 U	34.0000 U	47.0000 U	58.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001	13M62001	13M63001	13M64001	13M65001	13M66001
	LAB ID:	532122	532123	532124	532125	532126	532127
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
P Aroclor 1221	UG/KG	91.0000 U	140.0000 U	67.0000 U	69.0000 U	96.0000 U	120.0000 U
P Aroclor 1232	UG/KG	45.0000 U	70.0000 U	33.0000 U	34.0000 U	47.0000 U	58.0000 U
P Aroclor 1242	UG/KG	45.0000 U	70.0000 U	33.0000 U	34.0000 U	47.0000 U	58.0000 U
P Aroclor 1248	UG/KG	45.0000 U	70.0000 U	33.0000 U	34.0000 U	47.0000 U	58.0000 U
P Aroclor 1254	UG/KG	14.0000 J	70.0000 U	33.0000 U	34.0000 U	47.0000 U	58.0000 U
P Aroclor 1260	UG/KG	45.0000 U	28.0000 J	33.0000 U	34.0000 U	47.0000 U	31.0000 J
P Dieldrin	UG/KG	4.5000 U	7.0000 U	3.3000 U	3.4000 U	4.7000 U	5.8000 U
P Endosulfan I	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	3.0000 U
P Endosulfan II	UG/KG	4.5000 U	7.0000 U	3.3000 U	3.4000 U	4.7000 U	5.8000 U
P Endosulfan sulfate	UG/KG	4.5000 U	7.0000 U	3.3000 U	3.4000 U	4.7000 U	5.8000 U
P Endrin	UG/KG	4.5000 U	7.0000 U	3.3000 U	3.4000 U	4.7000 U	5.8000 U
P Endrin aldehyde	UG/KG	4.5000 U	7.0000 U	0.4400 J	0.5700 J	4.7000 U	1.3000 J
P Endrin ketone	UG/KG	4.5000 U	7.0000 U	3.3000 U	3.4000 U	4.7000 U	5.8000 U
P Heptachlor	UG/KG	2.3000 U	3.6000 U	0.2000 U	0.2300 U	2.4000 U	3.0000 U
P Heptachlor epoxide	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	0.4700 J
P Methoxychlor	UG/KG	1.1000 U	36.0000 U	17.0000 U	17.0000 U	24.0000 U	30.0000 U
P Toxaphene	UG/KG	230.0000 U	360.0000 U	170.0000 U	170.0000 U	240.0000 U	300.0000 U
P alpha-BHC	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	3.0000 U
P alpha-Chlordane	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	3.0000 U
P beta-BHC	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	3.0000 U
P delta-BHC	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	3.0000 U
P gamma-BHC (Lindane	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	3.0000 U
P gamma-Chlordane	UG/KG	2.3000 U	3.6000 U	1.7000 U	1.7000 U	2.4000 U	0.3300 U
S 1,2,4-Trichlorobenzene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 1,2-Dichlorobenzene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 1,3-Dichlorobenzene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 1,4-Dichlorobenzene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2,4,5-Trichlorophenol	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S 2,4,6-Trichlorophenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13M61001	13M62001	13M63001	13M64001	13M65001	13M66001
ENSAFE ID:		13M61001	13M62001	13M63001	13M64001	13M65001	13M66001
LAB ID:		532122	532123	532124	532125	532126	532127
DATE RCVD:		01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
S 2,4-Dichlorophenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2,4-Dimethylphenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2,4-Dinitrophenol	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S 2,4-Dinitrotoluene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2,6-Dinitrotoluene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2-Chloronaphthalene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2-Chlorophenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2-Methylnaphthalene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2-Methylphenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 2-Nitroaniline	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S 2-Nitrophenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 3-Nitroaniline	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S 4-Bromophenyl-phenylether	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 4-Chloro-3-methylphenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 4-Chloroaniline	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 4-Methylphenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S 4-Nitroaniline	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S 4-Nitrophenol	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S Acenaphthene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Acenaphthylene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Anthracene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Benzo(a)anthracene	UG/KG	440.0000 U	79.0000 J	340.0000 U	340.0000 U	52.0000 J	180.0000 J
S Benzo(a)pyrene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	180.0000 J
S Benzo(b,k)fluoranthene	UG/KG	440.0000 U	170.0000 J	340.0000 U	340.0000 U	120.0000 J	420.0000 J
S Benzo(g,h,i)perylene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Butylbenzylphthalate	UG/KG	440.0000 U	140.0000 U	340.0000 U	40.0000 U	83.0000 U	78.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001	13M62001	13M63001	13M64001	13M65001	13M66001
	LAB ID:	532122	532123	532124	532125	532126	532127
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
S Carbazole	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Chrysene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	140.0000 J
S Di-n-butylphthalate	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Di-n-octylphthalate	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Dibenz(a,h)anthracene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	70.0000 J
S Dibenzofuran	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Diethylphthalate	UG/KG	440.0000 U	75.0000 U	340.0000 U	340.0000 U	52.0000 U	70.0000 U
S Dimethylphthalate	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Fluoranthene	UG/KG	440.0000 U	110.0000 J	340.0000 U	340.0000 U	77.0000 J	260.0000 J
S Fluorene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Hexachlorobenzene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Hexachlorobutadiene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Hexachlorocyclopentadiene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Hexachloroethane	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	220.0000 J
S Isophorone	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S N-Nitrosodiphenylamine	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Naphthalene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Nitrobenzene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Pentachlorophenol	UG/KG	1100.0000 U	1700.0000 U	820.0000 U	830.0000 U	1100.0000 U	1400.0000 U
S Phenanthrene	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	99.0000 J
S Phenol	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S Pyrene	UG/KG	440.0000 U	100.0000 J	340.0000 U	340.0000 U	71.0000 J	250.0000 J
S bis(2-Chloroethoxy)methane	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	440.0000 U	85.0000 U	340.0000 U	340.0000 U	82.0000 U	84.0000 U
S bis(2-chloroethyl)ether	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	440.0000 U	700.0000 U	340.0000 U	340.0000 U	460.0000 U	580.0000 U
V 1,1,1-Trichloroethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	13M61001 / 532122 / 01/15/93	13M62001 / 532123 / 01/15/93	13M63001 / 532124 / 01/15/93	13M64001 / 532125 / 01/15/93	13M65001 / 532126 / 01/15/93	13M66001 / 532127 / 01/15/93
1,1,2,2-Tetrachloroethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
1,1,2-Trichloroethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
1,1-Dichloroethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
1,1-Dichloroethene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
1,2-Dichloroethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
1,2-Dichloroethene (total)	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
1,2-Dichloropropane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
2-Butanone	UG/KG	14.0000 U	15.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
2-Hexanone	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
4-Methyl-2-Pentanone	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Acetone	UG/KG	25.0000 U	90.0000 U	20.0000 U	14.0000 U	23.0000 U	39.0000 U
Benzene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Bromodichloromethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Bromoform	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Bromomethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Carbon Disulfide	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Carbon Tetrachloride	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Chlorobenzene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Chloroethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Chloroform	UG/KG	3.0000 U	5.0000 U	1.0000 U	1.0000 U	3.0000 U	3.0000 U
Chloromethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Dibromochloromethane	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Ethylbenzene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Methylene Chloride	UG/KG	69.0000 U	150.0000 U	14.0000 U	23.0000 U	30.0000 U	33.0000 U
Styrene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Tetrachloroethene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Toluene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Trichloroethene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
Vinyl Chloride	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001	13M62001	13M63001	13M64001	13M65001	13M66001
	LAB ID:	532122	532123	532124	532125	532126	532127
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
V Xylene (total)	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
V cis-1,3-Dichloropropene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U
V trans-1,3-Dichloropropene	UG/KG	14.0000 U	21.0000 U	10.0000 U	10.0000 U	14.0000 U	18.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M67001	33M01001	33302001	33M03001	33M03001DL	33M04001
	LAB ID:	532128	533364	533370	533344	533344 D50	533359
	DATE RCVD:	01/15/93	01/22/93	01/22/93	01/22/93	01/22/93	01/22/93
NAME	UNITS						
I Aluminum	MG/KG	188.0000	2660.0000	2340.0000	1100.0000	.	4150.0000
I Antimony	MG/KG	12.0000 U	15.2000 U	15.1000 U	14.5000 U	.	14.8000 U
I Arsenic	MG/KG	0.6300 U	1.3000 J	0.8200 J	1.0000 U	.	6.2000
I Barium	MG/KG	1.3000 U	5.3000 J	8.0000 J	1.6000 J	.	15.1000 J
I Beryllium	MG/KG	0.6300 U	0.8000 U	0.7900 U	0.7600 U	.	0.7800 U
I Cadmium	MG/KG	1.1000 U	1.3000 U	2.8000	1.3000 U	.	34.6000
I Calcium	MG/KG	55.5000 J	1260.0000 J	250.0000 J	155.0000 J	.	2760.0000
I Chromium	MG/KG	2.1000 U	18.5000	79.0000	9.3000	.	1180.0000
I Cobalt	MG/KG	2.5000 U	3.2000 U	3.2000 U	3.1000 U	.	3.1000 U
I Copper	MG/KG	5.3000 U	6.6000 U	9.3000	6.4000 U	.	45.1000
I Cyanide	MG/KG	0.5300 U	0.6600 U	0.6600 U	0.6400 U	.	0.6500 U
I Iron	MG/KG	117.0000	1210.0000	1010.0000	710.0000	.	1210.0000
I Lead	MG/KG	1.3000	6.5000	5.3000	5.7000	.	161.0000
I Magnesium	MG/KG	116.0000 U	223.0000 J	242.0000 J	140.0000 U	.	143.0000 U
I Manganese	MG/KG	1.3000 U	16.1000	5.8000	4.1000	.	27.1000
I Mercury	MG/KG	0.1100 U	0.1300 U	0.1300 U	0.1300 U	.	0.1300 U
I Nickel	MG/KG	8.0000 U	10.1000 U	10.1000 U	9.7000 U	.	52.1000
I Potassium	MG/KG	832.0000 U	1050.0000 U	1050.0000 U	1010.0000 U	.	1030.0000 U
I Selenium	MG/KG	0.8400 U	1.1000 U	1.1000 U	1.0000 U	.	1.0006 U
I Silver	MG/KG	2.1000 U	2.7000 U	3.0000 U	2.7000 U	.	3.6000 U
I Sodium	MG/KG	98.4000 J	122.0000 J	482.0000 J	137.0000 J		156.0000 J
I Thallium	MG/KG	0.6300 U	0.5300 U	0.7900 U	0.7600 U		0.7800 U
I Vanadium	MG/KG	3.2000 U	4.0000 U	4.0000 U	3.8000 U		3.9000 U
I Zinc	MG/KG	1.6000 J	38.2000	22.3000	8.0000		115.0000
P 4,4'-DDD	UG/KG	4.1000 J	37.0000 J	36.0000	120.0000 J	140.0000 J	25.0000 J
P 4,4'-DDE	UG/KG	1.6000 J	21.0000 J	54.0000 J	96.0000 J	120.0000 J	18.0000 J
P 4,4'-DDT	UG/KG	3.5000 U	6.2000 J	8.8000 J	48.0000 J	57.0000 J	7.8000
P Aldrin	UG/KG	1.8000 U	2.3000 U	0.1200 J	0.0780 J	11.0000 U	2.2000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M67001	331101001	331102001	33M03001	33M03001DL	33M04001
	LAB ID:	532120	533364	533370	533344	533344 D50	533359
	DATE RCVD:	01/15/93	01/22/93	01/22/93	01/22/93	01/22/93	01/22/93
NAME	UNITS						
P Aroclor 1016	UG/KG	35.0000 U	44.0000 U	43.0000 U	42.0000 U	210.0000 U	43.0000 U
P Aroclor 1221	UG/KG	70.0000 U	89.0000 U	87.0000 U	85.0000 U	420.0000 U	86.0000 U
P Aroclor 1232	UG/KG	35.0000 U	44.0000 U	43.0000 U	42.0000 U	210.0000 U	43.0000 U
P Aroclor 1242	UG/KG	35.0000 U	44.0000 U	43.0000 U	42.0000 U	210.0000 U	43.0000 U
P Aroclor 1248	UG/KG	35.0000 U	44.0000 U	43.0000 U	42.0000 U	210.0000 U	43.0000 U
P Aroclor 1254	UG/KG	35.0000 U	44.0000 U	43.0000 U	42.0000 U	210.0000 U	43.0000 U
P Aroclor 1260	UG/KG	16.0000 J	49.0000 J	12.0000 J	42.0000 U	210.0000 U	43.0000 U
P Dieldrin	UG/KG	3.5000 U	4.4000 U	4.3000 U	4.2000 U	21.0000 U	0.3900 J
P Endosulfan I	UG/KG	1.8000 U	2.3000 U	0.2100 J	0.1500 J	11.0000 U	0.1300 J
P Endosulfan II	UG/KG	3.5000 U	4.4000 U	4.3000 U	4.2000 U	21.0000 U	4.3000 U
P Endosulfan sulfate	UG/KG	3.5000 U	4.4000 U	4.3000 U	4.2000 U	21.0000 U	4.3000 U
P Endrin	UG/KG	3.5000 U	4.4000 U	4.3000 U	2.3000 J	21.0000 U	0.4100 J
P Endrin aldehyde	UG/KG	3.5000 U	4.4000 U	4.3000 U	4.2000 U	21.0000 U	0.5300 J
P Endrin ketone	UG/KG	3.5000 U	4.4000 U	4.3000 U	4.2000 U	21.0000 U	4.3000 U
P Heptachlor	UG/KG	1.8000 U	0.1100 U	0.1200 U	0.2300 U	11.0000 U	0.0990 U
P Heptachlor epoxide	UG/KG	1.8000 U	0.8400 J	0.1000 J	2,1000 U	11.0000 U	2.2000 U
P Methoxychlor	UG/KG	18.0000 U	23.0000 U	22.0000 U	21.0000 U	110.0000 U	22.0000 U
P Toxaphene	UG/KG	180.0000 U	230.0000 U	220.0000 U	210.0000 U	1100.0000 U	220.0000 U
P alpha-BHC	UG/KG	1.8000 U	2.3000 U	2.2000 U	2.1000 U	11.0000 U	2.2000 U
P alpha-Chlordane	UG/KG	1.8000 U	0.9600 J	0.2600 J	0.2100 J	11.0000 U	0.1300 J
P beta-BHC	UG/KG	1.8000 U	2.3000 U	2.2000 U	2.1000 U	11.0000 U	2.2000 U
P delta-BHC	UG/KG	1.8000 U	0.8800 J	0.4200 J	0.3800 J	11.0000 U	2.2000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U	0.0740 J	0.1200 J	2.1000 U	11.0000 U	2.2000 U
P gamma-Chlordane	UG/KG	1.8000 U	0.8200 J	0.5200 J	2.1000 U	11.0000 U	0.2700 J
S 1,2,4-Trichlorobenzene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 1,2-Dichlorobenzene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 1,3-Dichlorobenzene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 1,4-Dichlorobenzene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2,4,5-Trichlorophenol	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U		1000.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significan digits.

	ENSAFE ID:	13M67001	33M01001	33102001	33M03001	33M03001DL	33M04001
	LAB ID:	532120	533364	533370	533344	533344 D50	533359
	DATE RCVD:	01/15/93	01/22/93	01/22/93	01/22/93	01/22/93	01/22/93
NAME	WITS						
S 2,4,6-Trichlorophenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2,4-Dichlorophenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2,4-Dimethylphenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2,4-Dinitrophenol	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U		1000.0000 U
S 2,4-Dinitrotoluene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2,6-Dinitrotoluene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2-Chloronaphthalene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2-Chlorophenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2-Methylnaphthalene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2-Methylphenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U		420.0000 U
S 2-Nitroaniline	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U		1000.0000 U
S 2-Nitrophenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S 3-Nitroaniline	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U	.	1000.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U	.	1000.0000 U
S 4-Brmophenyl-phenylether	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S 4-Chloro-3-methylphenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410,0000 U	.	420.0000 U
S 4-Chloroaniline	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S 4-Methylphenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S 4-Nitroaniline	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U	.	1000.0000 U
S 4-Nitrophenol	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U	.	1000.0000 U
S Acenaphthene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Acenaphthylene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Anthracene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Benzo(a)anthracene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Benzo(a)pyrene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Benzo(b,k)fluoranthene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Benzo(g,h,i)perylene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M67001	33M01001	33M02001	33M03001	33M03001DL	33M04001
	LAB ID:	532128	533364	533370	533344	533344 D50	533359
	DATE RCVD:	01/15/93	01/22/93	01/22/93	01/22/93	01/22/93	01/22/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	350.0000 U	140.0000 U	140.0000 U	180.0000 U		230.0000 U
S Carbazole	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Chrysene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Di-n-butylphthalate	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Di-n-octylphthalate	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Dibenz(a,h)anthracene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Dibenzofuran	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Diethylphthalate	UG/KG	36.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Dimethylphthalate	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Fluoranthene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	43.0000 J
S Fluorene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Hexachlorobenzene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Hexachlorobutadiene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Hexachlorocyclopentadiene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Hexachloroethane	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Isophorone	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S N-Nitrosodiphenylamine	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Naphthalene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Nitrobenzene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Pentachlorophenol	UG/KG	840.0000 U	1000.0000 U	1000.0000 U	1000.0000 U	.	1000.0000 U
S Phenanthrene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Phenol	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S Pyrene	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	350,0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	350.0000 U	59.0000 U	150.0000 U	410.0000 U	.	420.0000 U
S bis(2-chloroethyl)ether	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	350.0000 U	430.0000 U	430.0000 U	410.0000 U	.	420.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SEDIMENT SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M67001	33M01001	33102001	33M03001	33M03001DL	33M04001
	LAB ID:	532128	533364	533370	533344	533344 D50	533359
	DATE RCVD:	01/15/93	01/22/93	01/22/93	01/22/93	01/22/93	01/22/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 1,1-Dichloroethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 1,1-Dichloroethene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 1,2-Dichloroethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 1,2-Dichloroethene (total)	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 1,2-Dichloropropane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 2-Butanone	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 2-Hexanone	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Acetone	UG/KG	19.0000 U	34.0000 U	47.0000 U	48.0000 U	.	45.0000 U
V Benzene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Bramodichloromethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Bramoform	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Bromomethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Chlorobenzene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Chloroethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Chloroform	UG/KG	2.0000 U	2.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Chloromethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Dibramochloromethane	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Ethylbenzene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Methylene Chloride	UG/KG	18.0000 U	10.0000 U	13.0000 U	15.0000 U	.	11.0000 U
V Styrene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Toluene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U
V Trichloroethene	UG/KG	11.0000 U	13.0000 U	13.0000 U	13.0000 U	.	13.0000 U

APPENDIX O

DETECTED CONCENTRATIONS — SURFACE WATER SAMPLES

NAS PENSACOLA OU-10 AND SITE 13
SURFACE WATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W01002	33W02002	33W03002	33W04002		
	LAB ID:	533140	533144	533147	533149		
	DATE RCVD:	01/21/93	01/21/93	01/21/93	01/21/93		
NAME	UNITS						
I Aluminum	UG/L	1280.0000	953.0000	696.0000	737.0000	.	.
I Antimony	UG/L	57.0000 U	57.0000 U	57.0000 U	57.0000 U	.	.
I Arsenic	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	.	.
I Barium	UG/L	16.1000 J	13.4000 J	12.1000 J	11.9000 J	.	.
I Beryllium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	.	.
I Cadmium	UG/L	5.2000	5.0000 U	5.0000 U	5.0000 U	.	.
I Calcium	UG/L	12900.0000	14500.0000	13900.0000	13000.0000	.	.
I Chromium	UG/L	10.0000 U	10.8000	10.0000 U	10.0000 U	.	.
I Cobalt	UG/L	12.0000 U	12.0000 U	12.0000 U	12.0000 U	.	.
I Copper	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
I Cyanide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
I Iron	UG/L	5110.0000	1090.0000	808.0000	808.0000	.	.
I Lead	UG/L	2.4000 J	2.0000 U	2.0000 U	2.0000 U	.	.
I Magnesium	UG/L	2950.0000 J	5470.0000	5160.0000	4810.0000 J	.	.
I Manganese	UG/L	280.0000	38.6000	19.3000	11.3000 J	.	.
I Mercury	UG/L	0.2000 U	0.2000 U	0.2000 U	0.2000 U	.	.
I Nickel	UG/L	38.0000 U	38.0000 U	38.0000 U	38.0000 U	.	.
I Potassium	UG/L	3950.0000 U	3950.0000 U	3950.0000 U	3950.0000 U	.	.
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	.	.
I Silver	UG/L	10.0000 U	10,0000	10.0000 U	10.0000 U	.	.
I Sodium	UG/L	5050.0000	24500.0000	21700.0000	20300.0000	.	.
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	.	.
I Vanadium	UG/L	15.0000 U	15.0000 U	15.0000 U	15.0000 U	.	.
I Zinc	UG/L	37.1000	39.2000	38.9000	29.0000	.	.
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.1100 J	0.0410'J	.	.
P 4,4'-DDE	UG/L	0.0021 J	0.0140 J	0.0450 J	0.0110 J	.	.
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	0.0470 J	0.0080 J	.	.
P Aldrin	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
SURFACE WATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W01002	33W02002	33W03002	33W04002
	LAB ID:	533140	533144	533147	533149
	DATE RCVD:	01/21/93	01/21/93	01/21/93	01/21/93
NAME	UNITS				
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Dieldrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Endosulfan II	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Heptachlor	UG/L	0.0014 U	0.0500 U	0.0500 U	0.0500 U
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0013 J	0.0500 U
P Methoxychlor	UG/L	0.5000 U	0.5000 U	0.5000 U	0.5000 U
P Toxaphene	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U
P alpha-BHC	UG/L	0.0500 U	0.0043 J	0.0027 J	0.0500 U
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0013 J	0.0500 U
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	0.0034 J	0.0500 U	0.0500 U	0.0030 J
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS I'ENSACOLA OU-IO AND SITE 13
SURFACE WATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W01002	33W02002	33W03002	33W04002		
	LAB ID:	533140	533144	533147	533149		
	DATE RCVD:	01/21/93	01/21/93	01/21/93	01/21/93	.	.
NAME	UNITS						
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.	.
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
SURFACE WATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W01002	33W02002	33W03002	33W04002
	LAB ID:	533140	533144	533147	533149
	DATE RCVD:	01/21/93	01/21/93	01/21/93	01/21/93
NAME	UNITS				
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Butylbenzylphthalate	UG/L	4.0000 U	4.0000 U	4.0000 U	4.0000 U
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	10.0000 U	10.0000 U	1.0000 U	1.0000 U
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Isophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Phenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
SURFACE WATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W01002	33W02002	33W03002	33W04002		
	LAB ID:	533140	533144	533147	533149		
	DATE RCVD:	01/21/93	01/21/93	01/21/93	01/21/93	.	.
NAME	UNITS						
S bis(2-Ethlyhexy)phthalate	UG/L	6.0000 U	6.0000 U	7.0000 U	1.0000 U	.	.
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 2-Butanone	UG/L	69.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10,0000 U	.	.
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Methylene Chloride	UG/L	11.0000 U	3.0000 U	14.0000 U	17.0000 U	.	.

NAS YENSACOLA OU-10 AND SITE 13
SURFACE WATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W01002	33W02002	33W03002	33W04002		
	LAB ID:	533140	533144	533147	533149		
	DATE RCVD:	01/21/93	01/21/93	01/21/93	01/21/93		
NAME	UNITS						
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Toluene	UG/L	10.0000 U	2.0000 J	10.0000 U	10.0000 U		.
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V Xylene (total)	UG/L	1.0000 J	10.0000 U	2.0000 J	2.0000 J	.	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

APPENDIX P

GROUNDWATER ANALYTICAL SUMMARY — FEBRUARY 1993
ECOLOGY AND ENVIRONMENT, INC.

SUMMARY ANALYTICAL RESULTS FOR SHALLOW ZONE MONITORING WELLS
(all results in ug/L unless noted)

Compound	Sample Number							FPDWS/ FSDWS
	UG-1	GM-8	GM-9	GM-10	GM-62	GM-67	PCS-1	
Barium	--	130	--	90	73	--	33	2,000
Cadmium	--	6.2	--	--	--	--	--	5
Chromium	--	--	--	12	--	--	--	100
Iron	1,600[a]	360	1,500	400	260	29,000	2,900	300
Manganese	--	48	24	68	270	130	100	50
Sodium (mg/L)	5	3.8	2.8	11	8.1	190	4	150
Zinc	--	46	31	--	130	23	23	5,000
Chloride (mg/L)	8.8	4.8	3.1	24	6.2	250	2.7	250
Sulfate (mg/L)	18	62	19	22	68	18	14	250
Complexed Cyanide (mg/L)	--	--	--	--	--	--	0.035	200
Fluoride (mg/L)	--	0.22	0.16	0.17	0.16	0.37	0.18	2,000
Methylene Chloride	(L,B)	(L,B)	(L,B)	(L,B)	--	--	(L,B)	5
Chlorobenzene	--	--	(L)	--	--	--	--	100
Acetone	(1)	43b	38	(L,B)	23	46	11	
Carbon Disulfide	--	--	38A	--	--	--	7.8A	
1,4-Dichlorobenzene	--	--	(L)	--	--	(L)	--	15
Bis(2-ethylhexyl)phthalate	(L)	(L)	(L,B)	--	--	--	--	
Fluorene	--	--	--	--	--	--	(L)	
Gross Alpha (pCi/L)	1.4+/-1.1	6.8+/-2.8	--	--	3.7+/-1.9	1.0+/-6.0	--	15
Gross Beta (pCi/L)	--	5.2+/-3.2	4.8+/-10.4	--	3.8+/-2.3	19.2+/-7.1	--	
Radium-226 (pCi/L)	0.2+/-0.05	1.5+/-0.07	--	0.9+/-0.06	0.2+/-0.05	4.8+/-0.2	0.3+/-0.1	5[b]
Radium-228 (pCi/L)	1.5+/-0.5	1.0+/-0.5	1.3+/-1.5	--	1.1+/-0.5	--	--	5[b]
Turbidity (NTU)	18	3.6	6.0	1.8	1.3	200	16	1
Total coliform (CLY/100)	--	28	--	--	--	--	--	1

Compound	Sample Number				FPDWS/ FSDWS
	GM-11	GM-12R	GM-13R	GM-14	
Complexed Cyanide	0.12	--	--	--	200

[a]Bold number indicates concentration exceeds the applicable Florida Primary Drinking Water Standard (FPDWS) or Florida Secondary Drinking Water Standard.
[b]The sum of the values for radium-226 and radium-228 has an FPDWS of 5.0 pCi/L.

Key :

-- = Not detected.
A = Phenonenon of methodology with acid preservation.
B = Compound present in blank.
CLY/100 = Colonies per 100 milliliters.
L = Compound present below detection limit.
µg/L = Micrograms per liter.
mg/L = Milligrams per liter.
NTU = Nethlometric turbidity unit.
pCi/L = Picocuries per liter.

Source: Ecoloqy and Environment, Inc., 1993.

SUMMARY ANALYTICAL RESULTS FOR
INTERMEDIATE ZONE MONITORING WELLS
(all results in ug/L unless noted)

Compound	Sample Number				FPDWS/ FSDWS
	GH-64	GM-66	GM-69	PCI-1	
Arsenic	--	13	--	--	50
Bat ium	--	60	35	--	2,000
Cadmium	--	44	--	--	5
Chromium	--	410	--	--	100
Copper	--	71	--	--	1,000
Iron	240	390,000	230	100	300
Manganese	42	1,900	100	14	50
Sodium (mg/L)	36	140	660	340	150
Nickel	--	200	--	--	100
Vanadium	--	540	--	--	
Zinc	--	420	--	20	5,000
Chloride (mg/L)	36	570	830	440	250
Sulfate (mg/L)	28	9,000	200	100	250
Complexed Cyanide (mg/L)	--	--	--	0.65	200
Fluoride (mg/L)	0.48	--	0.26	0.29	2,000
Methylene Chloride	--	--	14(B)	(L,B)	5
Chloroethane	--	--	(L)	--	
Toluene	--	--	(L)	--	1,000
Chlorobenzene	--	(L)	130	22	100
Acetone	30	--	(L)	70	
Total 1,2-Dichloroethene	--	750	--	--	
Trichlorethene	--	4,600	--	--	3
Carbon Disulfide	--	--	(L,A)	44(A)	
1,3-Dichlorobenzene	--	16	580	(L)	
1,4-Dichlorobenzene	--	31	510	(L)	75
1,2-Dichlorobenzene	--	42	460	(L)	600
Naphthalene	--	39	--	--	
Acenaphthene	(L)	--	--	--	
2,4-Dimethylphenol	--	49	--	--	
2-Mothylphonol	--	(L)	--	--	
4-Mothylphonol	--	(L)	--	--	
2-Methylnaphthalene	--	(L)	--	--	

Compound	Sample Number GM-64	GM-66	GM-69	PCI-1	FPDWS/ FSDWS
Gross Alpha (pCi/L)	3.8+/-1.5	83.9+/-62.3	9.0+/-9.8	--	15
Gross Beta (pCi/L)	3.8+/-2.3	142+/-81.5	34+/-14.3	3.5+/-3.1	
Radium-226 (pCi/L)	0.2+/-0.05	0.7+/-0.06	1.6+/-0.07	0.7+/-0.1	5*
Radium-228 (pCi/L)	1.1+/-0.5	--	2.4+/-0.6	--	5*
Turbidity (NTU)	3.1	4.9	32	5.3	1
Total Coliform (CLY/100)	28	--	--	4	1

Key :

-- = Not detected.
A = Phenomenon of methodology with acid preservation.
B = Compound present in blank.
CLY/100 = Colonies per 100 milliliters.
L = Compound present below detection limit.
µg/L = Micrograms per liter.
mg/L = Milligrams per liter.
NTU = Nethlemetric turbidity unit.
pCi/L = Picocuries per liter.

Source: Ecology and Environment, Inc., 1993.

SUMMARY ANALYTICAL RESULTS FOR
DEEP ZONE MONITORING WELLS
(all results in ug/L unless noted)

Compound	Sample Number GM-63	GM-65	GM-68	PCD-1	FPDWS/ FSDWS
Barium	26	--	96	21	2,000
Iron	1,100	880	I,200	280	300
Manganese	51	60	54	22	50
Sodium (mg/L)	570	95	87	220	150
Zinc	72	61	61	24	5,000
Chloride (mg/L)	850	160	140	340	250
Sulfate (mg/L)	70	12	10	30	250
Methlene Chloride	(L,B)	--	--	L,B)	5
Acetone	84	1,900	750	200	
Carbon Disulfide	7.2(A)	--	--	8(A)	
Di-N-Butyl-Phthalate	(L)	--	--	--	
Bis(2-ethylhexyl)phthalate	(L,B)	--	--	--	
Gross Alpha (pCi/L)	6.7+/-9.4	--	2.5+/-1.8	--	15
Gross Beta (pCi/L)	14.8+/-14.2	7.9+/-4.8	8.7+/-3.6	8.4+/-5.2	
Radium-226 (pCi/L)	1.1+/-0.07	1.0+/-0.1	0.9+/-0.1	--	5*
Radium-228 (pCi/L)	I.7+/-0.5	--	--	--	5*
Turbidity (NTU)	5.5	11	8.2	2.3	1
Nitrogen, Nitrite-Nitrate	--	0.2	--	0.2	
Nitroqon, Nitrate	--	0.2	--	0.2	

Key:

-- = Not detected.
A = Phenomenon of methodology with acid preservation
B = compound present in blank.
CLY/100 = Colonies prr 100 milliliters.
L = Compound present below detection limit.
µg/L = Micrograms per liter.
mg/L = Milligrams per litor.
NTU = Nethlemetric turbidity unit.
pCi/L = Picocuries per liter.

PENSACOLA, OU 10, SITE 13
PENSACOLA REFERENCE SAMPLES

METAL											
	SAMPLE ID ------->	001-G-GI68-00		001-G-GI70-00		001-G-GS67-00		001-G-GS69-00			
	ORIGINAL ID ----->	01GI68		01GI70		01GS67		01GS69			
	LAB SAMPLE ID --->	622850		622852		622849		622851			
	ID FROM REPORT -->	MAC011		MAC012		MAC016		MAC017			
	SAMPLE DATE ----->	07/01/94		07/01/94		07/01/94		07/01/94			
	MATRIX ---------->	Water		Water		Water		Water			
	UNITS ---------->	UG/L		UG/L		UG/L		UG/L			
CAS #	Parameter	93820	VAL	93820	VAL	93820	VAL	93820	VAL		
CN	Cyanide	NR		NR		NR		NR			
AL	Aluminum	293.	U	218.	U	4240.		3270.			
SB	Antimony	30.2	U	30.2	U	30.2	U	30.2	U		
AS	Arsenic	2.8	U	2.8	U	2.8	U	2.8	U		
BA	Barium	13.5	U	9.5	U	11.	U	18.9	U		
BE	Beryllium	1.1	U	1.1	U	1.1	U	1.1	U		
CD	Cadmium	3.4	U	3.4	U	3.4	U	3.4	U		
CA	Calcium	5670.		5350.		17800.		6300.			
CR	Chromium	5.2	U	5.2	U	9.7	U	59.9			
CO	Cobalt	4.1	U	4.1	U	4.1	U	4.1	U		
CU	Copper	10.8	U	10.8	U	10.8	U	16.2			
FE	Iron	942.		53.3	U	677.		1770.			
PB	Lead	1.6	UJ	1.6	UJ	1.6	UJ	1.6	UJ		
MG	Magnesium	1330.	U	3030.		1590.	U	2510.	U		
MN	Manganese	8.9		3.1	U	6.7		26.7			
HG	Mercury	0.2	U	0.2	U	0.2	U	0.2	U		
NI	Nickel	39.9	U	39.9	U	39.9	U	39.9	U		
K	Potassium	2550.	U	3200.		13300.		6560.			
SE	Selenium	3.9	U	3.9	U	3.9	U	3.9	UJ		
AG	Silver	4.	U	4.	U	4.	U	4.	U		
NA	Sodium	8350.		9810.		10700.		7830.			
TL	Thallium	3.6	UJ	3.6	UJ	3.6	UJ	3.6	UJ		
V	Vanadium	7.5	U	7.5	U	7.9		7.5	U		
		7.8	U	7.5	U	17.5	U	290.			

*** Validation Complete ***

PENSACOLA, OU 10, SITE 13
PENSACOLA REFERENCE SAMPLES

		001-S-0067-01 01S6701 553146 01S6701 Soil MG/KG		001-C-0067-01 01GS67 562669 01GS67 Water UG/L		001-S-0067-03 01S6703 553153 01S6703 Soil MG/KG		001-S-0067-05 01S6705 553155 01S6705 Soil MG/KG		001-S-0067-07 01S6707 553156 01S6707 soil MG/KG		001-S-0067-09 01S6709 553157 01S6709 Soil MG/KG	
CAS #	Parameter	731446	VAL	314936	VAL	731446	VAL	731446	VAL	731446	VAL	731446	VAL
AG	Silver	2.	U	4.	U	2.	U	2.	U	2.1	U	2.1	U
AL	Aluminum	12400.		202000.		1800.		9050.		1320.		932.	
AS	Arsenic	2.4	J	4.7		0.61	UJ	1.7	J	0.62	UJ	0.62	UJ
BA	Barium	5.8		237.		1.6		10.1		1.8		1.8	
BE	Beryllium	0.4	U	4.1		0.41	U	0.4	U	0.42	U	0.42	U
CA	Calciun	421.		11300.		227.		6740.		113.		78.7	
CD	Cadmium	1.	U	3.	U	1.	U	1.	U	1.	U	1.	U
CN	Cyanide	0.51	U	10.	U	0.51	U	0.5	U	0.52	U	0.52	U
CO	Cobalt	1.8	U	15.4		1.8	U	1.8	U	1.9	U	1.9	U
CR	Chromium	11.8		255.	J	2.	U	8.		2.1	U	2.1	U
CU	Copper	5.1		103.		5.1	U	5.1		5.2	U	5.2	U
		9180.		41100.	J	1450.		4790.		1140.		866.	
		0.1	U	2.2		0.1	U	0.1	U	0.1	U	0.1	U
		452.	U	8450.		458.	U	450.	U	466.	U	466.	U
		96.3	U	7580.		97.4	U	365.		99.2	U	99.2	U
		63.7		303.		3.7		63.6		3.8		4.4	
		116.	U	9800.		87.5	U	112.	U	111.	U	135.	U
		6.3	U	55.4	J	6.3	U	6.2	U	6.5	U	6.5	U
		20.1		78.9		1.3		28.		1.2		0.53	
		9.3	U	35.	U	9.4	U	9.3	U	9.6	U	9.6	U
		0.61	U	30.	U	0.61	U	0.6	U	0.62	U	0.62	U
		0.81	U	3.	U	0.82	U	0.81	U	0.83	U	0.83	U
V	Vanadium	20.6		179.		2.6		12.5		2.		1.6	
ZN	Zinc	10.9		146.		3.	U	29.		2.8	U	3.3	

PENSACOLA, OU 10, SITE 13
PENSACOLA REFERENCE SAMPLES

METAL													
	SAMPLE ID ------->	001-S-0067-11		001-S-0067-13		001-S-0067-17		001-C-0067-17		001-S-0067-19		001-S-0067-21	
	ORIGINAL ID ----->	D1S6711		D1S6713		D1S6717		D1S6717D		D1S6719		D1S6721	
	LAB SAMPLE ID --->	553158		553160		553720		553168		553194		553169	
	ID FROM REPORT -->	D1S6711		D1S6713		0156717		D1S6717D		0156719		D1S6721	
	SAMPLE DATE ----->												
	MATRIX ---------->	soil		Soil		Soil		Soil		soil		Soil	
	UNITS ---------->	MG/KG		MG/KG		MG/KG		MG/KG		MG/KG		MG/KG	
CAS #	Parameter	731446	VAL	731446	VAL	314135	VAL	731446	VAL	314135	VAL	731446	VAL
AC	Silver	2.1	U	2.1	U	2.1	U	2.1	U	2.1	U	2.1	U
AL	Aluminum	574.		502.		473.		318.		179.		156.	
AS	Arsenic	0.62	UJ	0.62	UJ	0.7	UJ	1.3	J	0.62	UJ	0.62	UJ
BA	Barium	1.2		1.	U	1.	U	1.	U	1.	U	1.	U
BE	Beryllium	0.41	U	0.42	U	0.41	U	0.41	U	0.42	U	0.41	U
CA	Calcium	67.		50.6		47.7		37.3		32.5	U	35.7	
CD	Cadmium	1.	U	1.	U	1.	U	1.	U	1.	U	1.	U
CN	Cyanide	0.52	U	0.52	U	0.51	U	0.52	U	0.52	U	0.51	U
CO	Cobalt	1.9	U	1.9	U	1.9	U	1.9	U	1.9	U	1.9	U
CR	Chromium	2.1	U	2.1	U	2.1	U	2.1	U	3.7		2.1	U
CU	Copper	5.2	U	5.2	U	5.1	U	5.2	U	5.2	U	5.1	U
FE	Iron	525.		482.		483.		367.		449.		237.	
HG	Mercury	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
K	Potassium	464.	U	464.	U	460.	U	463.	U	466.	U	460.	U
MG	Magnesium	98.8	U	98.9	U	97.9	U	98.6	U	99.2	U	97.9	U
MN	Manganese	5.7		5.7		5.6		3.7		4.1		1.7	
NA	Sodium	98.	U	93.9	U	88.8	U	91.5	U	73.8	U	108.	U
NI	Nickel	6.4	U	6.4	U	6.4	U	6.4	U	6.5	U	6.4	U
PB	lead	0.41	U	0.44		1.1		0.61		0.6		0.41	U
SB	Antimony	9.5	U	9.6	U	9.5	R	9.5	U	9.6	R	9.5	U
SE	Selenium	0.62	U	0.62	U	0.62	U	0.62	U	0.62	U	0.62	U
TL	Thallium	0.83	U	0.83	U	0.82	U	0.83	U	0.83	U	0.82	U
V	Vanadium	1.5	U	1.5	U	1.4	U	1.4	U	1.5	U	1.4	U
ZN	Zinc	7.2		7.3		6.		6.7		29.		19.9	

*** Validation Complete ***

PENSACOLA, OU 10, SITE 13
PENSACOLA REFERENCE SAMPLES

METAL													
	SAMPLE ID ------->	001-C-0067-21		001-S-0067-23		001-S-0069-01		001-G-0069-01		001-S-0069-03		001-S-0069-05	
	ORIGINAL ID ----->	01S6721D		01S6723		01S6901		01GS69		01S6903		01S6905	
	LAB SAMPLE ID --->	553171		553173		552866		562667		552867		552868	
	ID FROM REPORT -->	01S6721D		01S6723		01S6901		01GS69		01S6903		01S6905	
	SAMPLE DATE ----->												
	MATRIX ---------->	Soil		Soil		Soil		Water		Soil		Soil	
	UNITS ---------->	MG/KG		MG/KG		MG/KG		UG/L		MG/KG		MG/KG	
CAS #	Parameter	731446	VAL	731446	VAL	731446	VAL	314936	VAL	731446	VAL	731446	VAL
AG	Silver	2.1	U	2.1	U	2.	U	4.	U	2.	U	2.1	U
AL	Aluminum	141.		95.2		1240.		51.6		1430.		1550.	
AS	Arsenic	0.63	UJ	0.62	UJ	0.61	UJ	4.	U	0.61	UJ	0.62	UJ
BA	Barium	1.	U	1.	U	5.9		2.	U	3.		2.4	
BE	Beryllium	0.42	U	0.41	U	0.41	U	1.	U	0.41	U	0.42	U
CA	Calciun	46.8		36.5		162.		115.	U	71.4		50.8	
CD	Cadmium	1.	U	1.	U	1.	U	3.	U	1.	U	1.	U
CN	Cyanide	0.52	U	0.52	U	0.51	U	11.7		0.51	U	0.52	U
CO	Cobalt	1.9	U	'1.9	U	1.8	U	9.	U	1.8	U	1.9	U
CR	Chromium	2.1	U	2.1	U	2.	U	8.	U	3.5		2.4	
CU	Copper	5.2	U	5.2	U	5.1	U	4.	U	5.1	U	5.2	U
FE	Iron	243.		225.		911.		43.6	J	1010.		963.	
HC	Mercury	0.1	U	0.1	U	0.1	U	1.1		0.1	U	0.1	U
K	Potassiun	467.	U	463.	U	453.	U	1270.	U	456.	U	465.	U
MG	Magnesiun	99.4	U	98.7	U	96.6	U	48.	U	97.1	U	99.	U
MN	Manganese	2.2		1.7		3.7		1.5		3.7		3.7	
NA	Sodium	101.	U	101.	U	101.	U	140.	U	99.6	U	109.	U
Nf	Nickel	6.5	U	6.4	U	6.3	U	18.	U	6.3	U	6.4	U
PB	Lead	0.42	U	0.41	U	6.7		2.	U	2.9		0.79	
SB	Antimony	9.6	U	9.5	U	9.3	U	35.	U	9.4	U	9.6	U
SE	Selenium	0.63	U	0.62	U	0.61	U	3.	U	0.61	U	0.62	U
TL	Thallium	0.84	U	0.83	U	0.81	U	3.	U	0.82	U	0.83	U
V	Vanadiun	1.5	U	1.5	U	2.5		6.	U	1.7		1.6	
ZN	Zinc	20.6		17.2		8.2		50.3		4.6		3.6	U

PENSACOLA, OU 10, SITE 13
PENSACOLA REFERENCE SAMPLES

METAL											
		SAMPLE ID ------->	001-S-0069-07		001-S-0069-09		001-S-0069-11		001-S-0069-13		
		ORIGINAL ID ----->	01S6907		01S6909		01S6911		01S6913		
		LAB SAMPLE ID --->	552869		552870		552871		552872		
		ID FROM REPORT -->	01S6907		01S6909		01S6911		01S6913		
		SAMPLE DATE ----->									
		MATRIX ---------->	Soil		Soil		Soil		Soil		
		UNITS ---------->	MG/KG		MG/KG		MG/KG		MG/KG		
CAS #	Parameter		731446	VAL	731446	VAL	731446	VAL	731446	VAL	
AC	Silver		2.1	U	2.	U	2.1	U	2.1	U	
AL	Aluminum		991.		820.		293.		695.		
AS	Arsenic		0.62	UJ	0.61	UJ	0.62	UJ	0.62	UJ	
BA	Barium		1.7		1.6		1.	U	1.8		
BE	Beryllium		0.41	U	0.41	U	0.41	U	0.42	U	
CA	Calcium		32.3	U	31.9	U	32.3	U	45.4		
CD	Cadmiun		1.	U	1.	U	1.	U	1.	U	
CN	Cyanide		0.52	U	0.51	U	0.52	U	0.52	U	
CO	Cobalt		1.9	U	1.8	U	1.9	U	1.9	U	
CR	Chromium		3.3		3.5		5.4		4.6		
CU	Copper		5.2	U	5.1	U	5.2	U	5.2	U	
FE	Iron		697.		595.		205.		497.		
HG	Mercury		0.1	U	0.1	U	0.1	U	0.1	U	
K	Potassiun		463.	U	457.	U	463.	U	466.	U	
MG	Magnesium		98.7	U	97.2	U	98.7	U	99.2	U	
MN	Manganese		4.6		3.7		3.4		5.7		
NA	Sodium		103.	U	118.	U	97.2		91.3	U	
NI	Nickel		6.4	U	6.3	U	6.4	U	6.5	U	
PB	Lead		0.56		0.49		0.41	U	0.59		
			9.5	U	9.4	U	9.5	U	9.6	U	
			0.62	U	0.61	U	0.62	U	0.62	U	
			0.83	U	0.82	U	0.83	U	0.83	U	
V	Vanadium		1.5	U	1.4	U	1.5	U	1.5	U	
ZN	Zinc		2.	U	2.4	U	1.6	U	3.	U	

*** Validation Complete ***

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key preeented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	33G11001F	33G12001	33G12001DL	33G12001F	33G12002	33G13001
	LAB ID:	534334	534604	534604	534626	547267	534307
	DATE RCVD:	01/30/93	02/03/93	02/03/93	02/03/93	04/23/93	01/30/93
NAME	UNITS						
S bis(2-Chloroethoxy)methane	UG/L	.	50.0000 U	100,0000 U	.	.	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	.	50.0000 U	100.0000 U	.	.	1.0000 U
S bis(2-chloroethyl)ether	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
V 1,1,1-Trichloroethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V 1,1,2-Trichloroethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V 1,1-Dichloroethane	UG/L	.	200.0000	.	.	.	4.0000 J
V 1,1-Dichloroethene	UG/L	.	20.0000 U		.	.	10.0000 U
V 1,2-Dichloroethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	.	10.0000 J		.	.	10.0000 U
V 1,2-Dichloropropane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V 2-Butanone	UG/L	.	160.0000 U	.	.	.	110.0000 U
V 2-Hexanone	UG/L	.	20.0000 U		.	.	10,0000 U
V 4-Methyl-2-Pentanone	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Acetone	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Benzene	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Bromodichloromethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Bromoform	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Bromomethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Carbon Disulfide	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Carbon Tetrachloride	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Chlorobenzene	UG/L	.	130.0000	.	.	.	13.0000
V Chloroethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Chloroform	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Chloromethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Dibromochloromethane	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Ethylbenzene	UG/L	.	20.0000 U	.	.	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G11001F	33612001	33612001DL	33612001F	33612002	33613001
	LAB ID:	534334	534604	534604	534626	547267	534307
	DATE RCVD:	01/30/93	02/03/93	02/03/93	02/03/93	04/23/93	01/30/93
NAME	WITS						
V Methylene Chloride	UG/L	.	20.0000 U	.	.	.	10.0000 U
V Styrene	UG/L		5.0000 J	.	.	.	2.0000 J
V Tetrachloroethene	UG/L		6.0000 J	.	.	.	10.0000 U
V Toluene	UG/L		4.0000 J	.	.	.	10.0000 U
V Trichloroethene	UG/L	.	2.0000 J	.	.	.	10.0000 U
V Vinyl Chloride	UG/L	.	15.0000 J	.	.	.	10.0000 U
V Xylene (total)	UG/L		4.0000 J		.	.	10.0000 U
V cis-1,3-Dichloropropene	UG/L		20.0000 U	.	.	.	10.0000 U
V trans-1,3-Dichloropropene	UG/L	.	20.0000 U	.	.	.	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G13001F	33G13002	33G14001	33G14001F	33G15001	33G15001F
	LAB ID:	534327	547269	534296	534326	533778	533793
	DATE RCVD:	01/30/93	04/23/93	01/30/93	01/30/93	01/27/93	01/27/93
NAME	UNITS						
I Aluminum	UG/L	202.0000	376.0000	2230.0000	196.0000 J	18300.0000	351.0000
I Antimony	UG/L	52.0000 U	36.0000 U	57.0000 U	52.0000 U	57.0000 U	52.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	3.0000 U	6.0000 J	18.7000	3.0000 U
I Barium	UG/L	50.0000 J	33.6000 J	74.4000 J	68.6000 J	51.5000 J	7.8000 J
I Beryllium	UG/L	1.0000 U	1.0000 U	3.0000 U	1.0000 U	3.0000 U	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	78300.0000	69200.0000	142000.000	141000.000	11500.0000	12900.0000
I Chromium	UG/L	6.0000 U	8.0000 U	10.0000 U	6.0000 U	35.6000	6.2000 J
I Cobalt	UG/L	9.0000 U	8.0000 U	12.0000 U	9.0000 U	12.0000 U	9.0000 U
I Copper	UG/L	4.0000 U	4.7000 J	25.0000 U	4.0000 U	25.0000 U	4.0000 U
I Cyanide	UG/L		2.0000	10.0000 U		10.0000 U	
I Iron	UG/L	1320.0000	2780.0000	11200.0000	9400.0000	21500.0000	265.0000
I Lead	UG/L	2.0000 U	2.0000 U	40.0000 U	2.0000 U	10.8000	2.0000 U
I Magnesium	UG/L	93800.0000	79500.0000	346000.000	349000.000	11000.0000	8790.0000
I Manganese	UG/L	70.7000	59.2000	240.0000	226.0000	85.6000	26.0000
I Mercury	UG/L	0.2000 U	0.2000 U	0.2000 U	0.2000 U	0.2000 U	0.2000 U
I Nickel	UG/L	13.0000 U	17.0000 U	38.0000 U	13.0000 U	38.0000 U	13.0000 U
I Potassium	UG/L	41400.0000	34900.0000	109000.000	108000.000	8750.0000	14100.0000
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	5.0000 U	7.0000 U	10.0000	5.0000 U	10.0000 U	5.0000 U
I sodium	UG/L	861000.000	719000.000	2940000.00	2880000.00	12900.0000	14700.0000
I Thallium	UG/L	3.0000 U	40.0000 U	60.0000 U	3.0000 U	3.0000 U	10.0000 U
I Vanadium	UG/L	6.7000 J	8.1000 J	15.0000 U	4.0000 U	65.1000	18.5000 J
I Zinc	UG/L	9.5000 J	9.7000 J	5.7000 J	16.4000 J	152.0000	7.8000 J
P 4,4'-DDD	UG/L	.		0.1000 U	.	0.1000 U	
P 4,4'-DDE	UG/L	.		0.1000 U	.	0.1000 U	
P 4,4'-DDT	UG/L	.		0.0095 J	.	0.1000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G11001F	33012001	33G12001DL	33G12001F	33G12002	33G13001
	LAB ID:	534334	534604	534604	534626	547267	534307
	DATE RCVD:	01/30/93	02/03/93	02/03/93	02/03/93	04/23/93	01/30/93
NAME	UNITS						
I Aluminum	UG/L	50.0000 U	840,0000	.	50.0000 U	498.0000	2000.0000
I Antimony	UG/L	30.0000 U	29.0000 U	.	30.0000 U	36.0000 U	57.0000 U
I Arsenic	UG/L	7.7000 J	4.0000 J	.	3.0000 U	3.0000 U	3.0000 U
I Barium	UG/L	76.6000 J	87.2000 J	.	81.1000 J	47.8000 J	54.4000 J
I Beryllium	UG/L	1.0000 U	3.0000 U	.	1.0000 U	1.0000 U	3.0000 U
I Cadmium	UG/L	5.0000 U	5.9000 U	.	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	49000.0000	169000.000	.	166000.000	130000.000	74900.0000
I Chromium	UG/L	6.0000 U	10.0000 U		6.0000 U	8.0000 U	10.0000 U
I Cobalt	UG/L	9.0000 U	12.0000 U	.	9.0000 U	8.0000 U	12.0000 U
I Copper	UG/L	4.0000 U	25.0000 U	.	4.0000 U	4.0000 U	25.0000 U
I Cyanide	UG/L		10.0000 U	.		2.0000	10.0000 U
I Iron	UG/L	16800.0000	4500.0000	.	4180.0000	5990.0000	2380.0000
I Lead	UG/L	2.0000 U	20.0000 U	.	2.0000 U	100.0000 U	2.0000 U
I Magnesium	UG/L	41000.0000	319000.000	.	319000.000	265000.000	86400.0000
I Manganese	UG/L	264.0000	501.0000	.	460.0000	205.0000	73.4000
I Mercury	UG/L	0.2600	0.2000 U	.	0.2000 U	0.2000 U	0.2000 U
I Nickel	UG/L	13.0000 U	38.0000 U	.	13.0000 U	17.0000 U	38.0000 U
I Potassium	UG/L	20800.0000	107000.000	.	104000.000	85700.0000	39300.0000
I Selenium	UG/L	3.0000 U	3.0000 U	.	3.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	5.0000 U	10.0000 U	.	5.0000 U	7.0000 U	10.0000 U
I Sodium	UG/L	363000.000	1420000.00	.	2600000.00	2170000.00	814000.000
I Thallium	UG/L	3.0000 U	30.0000 U	.	3.0000 U	400.0000 U	3.0000 U
I Vanadium	UG/L	4.0000 U	15.0000 U	.	4.0000 U	6.0000 U	15.0000 U
I Zinc	UG/L	39.4000	34.6000	.	5.8000 J	20.2000	19.2000 J
P 4,4'-DDD	UG/L	.	0.1000 U	.			0.1000 U
P 4,4'-DDE	UG/L	.	0.1000 U	.			0.1000 U
P 4,4'-DDT	UG/L	.	0.1000 U	.			0.1000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	33G11001F	33G12001	33G12001DL	33G12001F	33612002	33G13001
	LAB ID:	534334	534604	534604	534626	547267	534307
	DATE RCVD:	01/30/93	02/03/93	02/03/93	02/03/93	04/23/93	01/30/93
P Aldrin	UG/L	.	0.0500 U	.	.	.	0.0500 U
P Aroclor 1016	UG/L		1.0000 U	.	.	.	1.0000 U
P Aroclor 1221	UG/L	.	2.0000 U	.	.	.	2.0000 U
P Aroclor 1232	UG/L	.	1.0000 U	.	.	.	4.0000 U
P Aroclor 1242	UG/L	.	1.0000 U	.	.	.	1.0000 U
P Aroclor 1248	UG/L	.	1.0000 U	.	.	.	1.0000 U
P Aroclor 1254	UG/L	.	1.0000 U	.	.	.	1.0000 U
P Aroclor 1260	UG/L	.	1.0000 U	.	.	.	1.0000 U
P Dieldrin	UG/L	.	0.1000 U	.	.	.	0.1000 U
P Endosulfan I	UG/L	.	0.0500 U	.	.	.	0.0500 U
P Endosulfan II	UG/L	.	0.1000 U	.	.	.	0.1000 U
P Endosulfan sulfate	UG/L	.	0.1000 U	.	.	.	0.1000 U
P Endrin	UG/L	.	0.1000 U	.	.	.	0.1000 U
P Endrin aldehyde	UG/L	.	0.0300 J	.	.	.	0.1000 U
P Endrin ketone	UG/L	.	0.1000 U	.	.	.	0.0022 J
P Heptachlor	UG/L	.	0.0500 U	.	.	.	0.0031 U
P Heptachlor epoxide	UG/L	.	0.0500 U	.	.	.	0.0500 U
P Methoxychlor	UG/L	.	0.5000 U	.	.	.	0.5000 U
P Toxaphene	UG/L	.	5.0000 U	.	.	.	5.0000 U
P alpha-BHC	UG/L	.	0.0500 U	.	.	.	0.0500 U
P alpha-Chlordane	UG/L	.	0.0500 U	.	.	.	0.0500 U
P beta-BHC	UG/L	.	0.0500 U	.	.	.	0.0500 U
P delta-BHC	UG/L	.	0.0500 U	.	.	.	0.0500 U
P gamma-BHC (Lindane)	UG/L	.	0.0500 U	.	.	.	0.0500 U
P gma-Chlordane	UG/L	.	0.0500 U		.	.	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 1,2-Dichlorobenzene	UG/L	.	510.0000 J	770.0000 J	.	.	2.0000 J
S 1,3-Dichlorobenzene	UG/L	.	290.0000	460.0000 J	.	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G11001F	33G12001	33G12001DL	33G12001F	33G12002	33G13001
	LAB ID:	534334	534604	534604	534626	547267	534307
	DATE RCVD:	01/30/93	02/03/93	02/03/93	02/03/93	04/23/93	01/30/93
NAME	UNITS						
S 1,4-Dichlorobenzene	UG/L	.	470.0000 J	700.0000 J	.	.	1.0000 J
S 2,4,5-Trichlorophenol	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2,4-Dichlorophenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2,4-Dimethylphenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2,4-Dinitrophenol	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S 2,4-Dinitrotoluene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2,6-Dinitrotoluene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2-Chloronaphthalene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2-Chlorophenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2-Methylnaphthalene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2-Methylphenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 2-Nitroaniline	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S 2-Nitrophenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 3-Nitroaniline	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S 4-Brainophenyl-phenylether	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 4-Chloroaniline	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 4-Methylphenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S 4-Nitroaniline	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S 4-Nitrophenol	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S Acenaphthene	UG/L	.	15.0000 J	26.0000 J	.	.	10.0000 U
S Acenaphthylene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Anthracene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Benzo(a)anthracene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G11001F	33012001	33G12001DL	33G12001F	33012002	33G13001
	LAB ID:	534334	534604	534604	534626	547267	534307
	DATE RCVD:	01/30/93	02/03/93	02/03/93	02/03/93	04/23/93	01/30/93
NAME	UNITS						
S Benzo(a)pyrene	UG/L		50.0000 U	100.0000 U	.	.	10.0000 U
S Benzo(b,k)fluoranthene	UG/L		50.0000 U	100.0000 U	.	.	10.0000 U
S Benzo(g,h,i)perylene	UG/L		50.0000 U	100.0000 U	.	.	10.0000 U
S Butylbenzylphthalate	UG/L		50.0000 U	100.0000 U	.	.	10.0000 U
S Carbazole	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Chrysene	UG/L		50.0000 U	100.0000 U	.	.	10.0000 U
S Di-n-butylphthalate	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Di-n-octylphthalate	UG/L		50.0000 U	100.0000 U	.	.	10.0000 U
S Dibenz(a,h)anthracene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Dibenzofuran	UG/L		5.0000 J	100.0000 U	.	.	10.0000 U
S Diethylphthalate	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Dimethylphthalate	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Fluoranthene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Fluorene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Hexachlorobenzene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Hexachlorobutadiene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Hexachlorocyclopentadiene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Hexachloroethane	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Isophorone	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S N-Nitrosodiphenylamine	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Naphthalene	UG/L		22.0000 J	38.0000 J	.	.	10.0000 U
S Nitrobenzene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Pentachlorophenol	UG/L	.	120.0000 U	250.0000 U	.	.	25.0000 U
S Phenanthrene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Phenol	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U
S Pyrene	UG/L	.	50.0000 U	100.0000 U	.	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID:	33G13001F	33G13002	33014001	33G14001F	33015001	33G15001F
	LAB ID:	534327	547269	534296	534326	533778	533793
	DATE RCVD:	01/30/93	04/23/93	01/30/93	01/30/93	01/27/93	01/27/93
	UNITS						
P Aldrin	UG/L	.		0.0500 U	.	0.0500 U	.
P Aroclor 1016	UG/L		.	1.0000 U	.	1.0000 U	.
P Aroclor 1221	UG/L	.	.	2.0000 U	.	2.0000 U	.
P Aroclor 1232	UG/L	.		1.0000 U	.	1.0000 U	.
P Aroclor 1242	UG/L	.	.	1.0000 U	.	1.0000 U	.
P Aroclor 1248	UG/L	.	.	1.0000 U	.	1.0000 U	.
P Aroclor 1254	UG/L	.		1.0000 U	.	1.0000 U	.
P Aroclor 1260	UG/L	.	.	1.0000 U	.	1.0000 U	.
P Dieldrin	UG/L	.	.	0.1000 U	.	0.1000 U	.
P Endosulfan I	UG/L	.	.	0.0500 U	.	0.0500 U	.
P Endosulfan II	UG/L	.	.	0.1000 U	.	0.1000 U	.
P Endosulfan sulfate	UG/L	.	.	0.1000 U	.	0.1000 U	.
P Endrin	UG/L	.	.	0.1000 U	.	0.1000 U	.
P Endrin aldehyde	UG/L	.	.	0.1000 U	.	0.1000 U	.
P Endrin ketone	UG/L	.	.	0.1000 U	.	0.1000 U	.
P Heptachlor	UG/L	.	.	0.0030 U	.	0.0500 U	.
P Heptachlor epoxide	UG/L	.	.	0.0500 U	.	0.0500 U	.
P Methoxychlor	UG/L	.	.	0.5000 U	.	0.5000 U	.
P Toxaphene	UG/L	.	.	5.0000 U	.	5.0000 U	.
P alpha-BHC	UG/L	.	.	0.0500 U	.	0.0500 U	.
P alpha-Chlordane	UG/L	.	.	0.0500 U	.	0.0500 U	.
P beta-BHC	UG/L	.	.	0.0500 U	.	0.0500 U	.
P delta-BHC	UG/L	.	.	0.0500 U	.	0.0500 U	.
P gamma-EHC (Lindane)	UG/L	.	.	0.0500 U	.	0.0500 U	.
P gamma-Chlordane	UG/L	.	.	0.0500 U	.	0.0500 U	.
S 1,2,4-Trichlorobenzene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 1,2-Dichlorobenzene	UG/L	.	.	4.0000 J	.	10.0000 U	.
S 1,3-Dichlorobenzene	UG/L	.	.	10.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 ANI) SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G13001F	33G13002	33G14001	33G14001F	33G15001	330150011
	LAB ID:	534327	547269	534296	534326	533778	533793
	DATE RCVD:	01/30/93	04/23/93	01/30/93	01/30/93	01/27/93	01/27/93
NAME	UNITS						
S 1,4-Dichlorobenzene	UG/L	.	.	1.0000 J	.	10.0000 U	.
S 2,4,5-Trichlorophenol	UG/L	.	.	25.0000 U	.	25.0000 U	.
S 2,4,6-Trichlorophenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2,4-Dichlorophenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2,4-Dimethylphenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2,4-Dinitrophenol	UG/L	.	.	25.0000 U	.	25.0000 U	.
S 2,4-Dinitrotoluene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2,6-Dinitrotoluene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2-Chloronaphthalene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2-Chlorophenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2-Methylnaphthalene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2-Methylphenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 2-Nitroaniline	UG/L	.	.	25.0000 U	.	25.0000 U	.
S 2-Nitrophenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 3-Nitroaniline	UG/L	.	.	25.0000 U	.	25.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	.	.	25.0000 U	.	25.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 4-Chloroaniline	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 4-Methylphenol	UG/L	.	.	10.0000 U	.	10.0000 U	.
S 4-Nitroaniline	UG/L	.	.	25.0000 U	.	25.0000 U	.
S 4-Nitrophenol	UG/L	.	.	25.0000 U	.	25.0000 U	.
S Acenaphthene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Acenaphthylene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Anthracene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Bento(a)anthracene	UG/L	.	.	10.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33G13001F / 534327 / 01/30/93	33G13002 / 547269 / 04/23/93	33G14001 / 534296 / 01/30/93	33G14001F / 534326 / 01/30/93	33G15001 / 533778 / 01/27/93	33G15001F / 533793 / 01/27/93
S Benzo(a)pyrene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Benzo(b,k)fluoranthene	UG/L		.	10.0000 U	.	10.0000 U	.
S Benzo(g,h,i)perylene	UG/L		.	10.0000 U	.	10.0000 U	.
S Butylbenzylphthalate	UG/L		.	10.0000 U	.	3.0000 U	.
S Carbazole	UG/L		.	10.0000 U	.	10.0000 U	.
S Chrysene	UG/L		.	10.0000 U	.	10.0000 U	.
S Di-n-butylphthalate	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Di-n-octylphthalate	UG/L		.	10.0000 U	.	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L		.	10.0000 U	.	10.0000 U	.
S Dibenzofuran	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Diethylphthalate	UG/L	.	.	10.0000 U	.	1.0000 U	.
S Dimethylphthalate	UG/L		.	10.0000 U	.	10.0000 U	.
S Fluoranthene	UG/L	.	.	10.0000 U	.	10.0000 U	
S Fluorene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Hexachlorobenzene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Hexachlorobutadiene	UG/L		.	10.0000 U	.	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Hexachloroethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Ieophorone	UG/L	.	.	10.0000 U	.	10.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	.	.	10.0000 U	.	10.0000 U	.
S N-Nitrosodiphenylamine	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Naphthalene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Nitrobenzene	UG/L	.	.	10.0000 U	.	10.0000 U	.
S Pentachlorophenol	UG/L	.	.	25.0000 U	.	25.0000 U	.
S Phenanthrene	UG/L		.	10.0000 U	.	10.0000 U	.
S Phenol	UG/L	.	.	10.0000 U	.	10.0000	.
S Pyrene	UG/L	.	.	10.0000 U	.	10.0000 U	.

NAS I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33G13001F	33G13002	33G14001	33G14001F	33G15001	33G15001F
ENSAFE ID:		33G13001F	33G13002	33G14001	33G14001F	33G15001	33G15001F
LAB ID:		534327	547269	534296	534326	533778	533793
DATE RCVD:		01/30/93	04/23/93	01/30/93	01/30/93	01/27/93	01/27/93
S bis(2-Chloroethoxy)methane	UG/L	.	.	10.0000 U	.	10.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	.	.	10.0000 U	.	3.0000 U	.
S bis(2-chloroethyl)ether	UG/L	.	.	10.0000 U	.	10.0000 U	
S bis(2-chloroisopropyl)ether	UG/L	.	.	10.0000 U	.	10.0000 U	.
V 1,1,1-Trichloroethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V 1,1,2,2-Tetrachloroethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V 1,1,2-Trichloroethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V 1,1-Dichloroethane	UG/L	.	.	2.0000 J	.	10.0000 U	.
V 1,1-Dichloroethene	UG/L		.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethene (total)	UG/L	.	.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloropropane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V 2-Butanone	UG/L	.	.	10.0000 U	.	92.0000 U	.
V 2-Hexanone	UG/L	.		10.0000 U	.	10.0000 U	.
V 4-Methyl-2-Pentanone	UG/L	.	.	10.0000 U	.	10.0000 U	.
V Acetone	UG/L	.	.	10.0000 U	.	28.0000 U	.
V Benzene	UG/L	.		10.0000 U	.	10.0000 U	.
V Bromodichloromethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V Bromoform	UG/L	.	.	10.0000 U	.	10.0000 U	.
V Bromomethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V Carbon Disulfide	UG/L	.	.	10.0000 U	.	4.0000 J	.
V Carbon Tetrachloride	UG/L	.	.	10.0000 U	.	10.0000 U	.
V Chlorobenzene	UG/L	.	.	10.0000	.	10.0000 U	.
V Chloroethane	UG/L	.	.	10.0000 U	.	10.0000 U	
V Chloroform	UG/L	.	.	10.0000 U	.	10.0000 U	.
V Chloromethane	UG/L	.	.	10.0000 U	.	10.0000 U	.
V Dibromochloromethane	UG/L	.	.	10.0000 U	.	10.0000 U	
V Ethylbenzene	UG/L	.	.	10.0000 U	.	10.0000 U	

NAS YENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G13001F	33G13002	33G14001	33G14001F	33G15001	33G15001F
	LAB ID:	534327	547269	534296	534326	533778	533793
	DATE RCVD:	01/30/93	04/23/93	01/30/93	01/30/93	01/27/93	01/27/93
NAME	UNITS						
V Methylene Chloride	UG/L	.		10.0000 U		16.0000 U	.
V Styrene	UG/L	.		2.0000 J		10.0000 U	.
V Tetrachloroethene	UG/L			10.0000 U		190.0000	.
V Toluene	UG/L		.	10.0000 U		10.0000 U	.
V Trichloroethene	UG/L	.	.	10.0000 U		5.0000 J	.
V Vinyl Chloride	UG/L		.	10.0000 U		10.0000 U	.
V Xylene (total)	UG/L	.	.	10.0000 U		4.0000 J	.
V cis-1,3-Dichloropropene	UG/L	.	.	10.0000 U		10.0000 U	.
V trans-1,3-Dichloropropene	UG/L	.	.	10.0000 U		10.0000 U	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33016001	33G16001DL	33G16OOlP	33G17001	33G17OOlP	33018001
ENSAFE ID:		33016001	33G16001DL	33G16OOlP	33G17001	33G17OOlP	33018001
LAB ID:		534620	534646	534627	534602	534624	534599
DATE RCVD:		02/03/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
I Aluminum	UG/L	11400.0000	.	236.0000	925.0000	309.0000	11800.0000
I Antimony	UG/L	29.0000 U	.	30.0000 U	29.0000 U	30.0000 U	29.0000 U
I Arsenic	UG/L	11.9000	.	3.0000 U	13.8000	5.0000 J	4.8000 J
I Barium	UG/L	33.6000 J	.	30.8000 J	93.5000 J	88.2000 J	44.8000 J
I Beryllium	UG/L	3.0000 U	.	1.0000 U	3.0000 U	1.0000 U	3.0000 U
I Cadmium	UG/L	5.0000 U	.	5.0000 U	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	124000.000	.	130000.000	54300.0000	55600.0000	33500.0000
I Chromium	UG/L	12.3000	.	*6.0000* U	10.0000 U	6.0000 U	27.0000
I Cobalt	UG/L	12.0000 U	.	9.0000 U	12.0000 U	9.0000 U	12.0000 U
I Copper	UG/L	25.0000 U	.	4.0000 U	25.0000 U	4.0000 U	25.0000 U
I Cyanide	UG/L	10.0000 U	.		10.0000 U		10.0000 U
I Iron	UG/L	5000,0000	.	78.5000 J	786.0000	78.4000 J	9950.0000
I Lead	UG/L	*2.0000* U	.	2.0000 U	2.0000 U	2.0000 U	4.6000
I Magnesium	UG/L	32700.0000	.	34100.0000	11800.0000	11800.0000	10300.0000
I Manganese	UG/L	176.0000	.	187.0000	366.0000	348.0000	86.8000
I Mercury	UG/L	1.4000	.	0.2000 U	0.4300	0.2000 U	0.4100
I Nickel	UG/L	38.0000 U	.	13.0000 U	38.0000 U	13.0000 U	38.0000 U
I Potassium	UG/L	23000.0000	.	29400.0000	76100.0000	90900.0000	10500.0000
I Selenium	UG/L	3.0000 U	.	30.0000 U	13.9000	12.5000	3.0000 U
I Silver	UG/L	10.0000 U	.	*5.0000* U	10.0000 U	*5.0000* U	10.0000 U
I Sodium	UG/L	209000.000	.	229000.000	286000.000	291000.000	187000.000
I Thallium	UG/L	3.0000 U	.	3.0000 U	3.0000 U	3.0000 U	15.0000 U
I Vanadium	UG/L	15.0000 U	.	4.0000 U	23.6000 J	24.3000 J	22.6000 J
I Zinc	UG/L	66.7000	.	5.0000 U	44.0000	5.0000 U	95.6000
P 4,4'-DDD	UG/L	0.1000 U	.	.	0.1000 U		0.1000 U
P 4,4'-DDE	UG/L	0.0420 J	.		0.1000 U		0.1000 U
P 4,4'-DDT	UG/L	0.1000 U	.	.	0.1000 U		0.1000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G16001	33G16001DL	33G16001F	33G17001	33G17001F	33G18001
	LAB ID:	534646 R1	534646	534627	534602	534624	534599
	DATE RCVD:	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
NAME	UNITS						
P Aldrin	UG/L	0.0500 U	.	.	0.0500 U	.	0.0500 U
P Aroclor 1016	UG/L	1.0000 U	.	.	1.0000 U	.	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	.	.	2.0000 U	.	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	.	.	1.0000 U	.	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	.	.	1.0000 U		1.0000 U
P Aroclor 1248	UG/L	1.0000 U		.	1.0000 U		1.0000 U
P Aroclor 1254	UG/L	1.0000 U		.	1.0000 U	.	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	.	.	1.0000 U	.	1.0000 U
P Dieldrin	UG/L	0.0420 J		.	0.1000 U	.	0.1000 U
P Endosulfan I	UG/L	0.0500 U	.	.	0.0500 U	.	0.0500 U
P Endosulfan II	UG/L	0.1900 J		.	0.1000 U	.	0.1000 U
P Endoeulfan sulfate	UG/L	0.1000 U	.	.	0.1000 U		0.1000 U
P Endrin	UG/L	0.1000 U	.	.	0.1000 U	.	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U		.	0.1000 U	.	0.1000 U
P Endrin ketone	UG/L	0.1000 U	.	.	0.1000 U	.	0.1000 U
P Heptachlor	UG/L	0.0500 U		.	0.0500 U	.	0.0500 U
P Heptachlor epoxide	UG/L	0.0500 U		.	0.0500 U	.	0.0500 U
P Methoxychlor	UG/L	0.5000 U		.	0.5000 U	.	0.5000 U
P Toxaphene	UG/L	5.0000 U	.	.	5.0000 U	.	5.0000 U
P alpha-BHC	UG/L	0.0500 U		.	0.0500 U	.	0.0500 U
P alpha-Chlordane	UG/L	0.0500 U	.	.	0.0500 U	.	0.0500 U
P beta-BHC	UG/L	0.0500 U	.	.	0.0500 U	.	0.0500 U
P delta-BHC	UG/L	0.0500 U		.	0.0500 U	.	0.0500 U
P gamma-BHC (Lindane)	UG/L	0.0210 J		.	0.0500 U	.	0.0500 U
P gamma-Chlordane	UG/L	0.0500 U		.	0.0500 U	.	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 1,2-Dichlorobenzene	UG/L	850.0000 J	930.0000 J	.	110.0000	.	10.0000 U
S 1,3-Dichlorobenzene	UG/L	180.0000	210.0000 J	.	12.0000 J		10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33616001	33G16001DL	33G16001F	33G17001	336170013	33G18001
	LAB ID:	534646	534646	534627	534602	534624	534599
	DATE RCVD:	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
NAME	UNITS						
S 1,4-Dichlorobenzene	UG/L	310.0000	350.0000 J	.	30.0000	.	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	250.0000 U	380.0000 U		50.0000 U	.	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 2,4-Dichlorophenol	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 2,4-Dimethylphenol	UG/L	100.0000 U	150.0000 U	.	20.0000 U		10.0000 U
S 2,4-Dinitrophenol	UG/L	250.0000 U	380.0000 U	.	50.0000 U		25.0000 U
S 2,4-Dinitrotoluene	UG/L	100.0000 U	150.0000 U	.	20.0000 U		10.0000 U
S 2,6-Dinitrotoluene	UG/L	100.0000 U	150.0000 U		20.0000 U	.	10.0000 U
S 2-Chloronaphthalene	UG/L	100.0000 U	150.0000 U		20.0000 U		10.0000 U
S 2-Chlorophenol	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 2-Methylnaphthalene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 2-Methylphenol	UG/L	13.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 2-Nitroaniline	UG/L	250.0000 U	380.0000 U	.	50.0000 U	.	25.0000 U
S 2-Nitrophenol	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	100.0000 U	150.0000 U		20.0000 U		10.0000 U
S 3-Nitroaniline	UG/L	250.0000 U	380.0000 U		50.0000 U	.	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	250.0000 U	380.0000 U	.	50.0000 U	.	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 4-Chloroaniline	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 4-Methylphenol	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S 4-Nitroaniline	UG/L	250.0000 U	380.0000 U	.	50.0000 U	.	25.0000 U
S 4-Nitrophenol	UG/L	250.0000 U	380.0000 U	.	50.0000 U		25.0000 U
S Acenaphthene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Acenaphthylene	UG/L	100.0000 U	150.0000 U	.	20.0000 U		10.0000 U
S Anthracene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Benzo(a)anthracene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G16001	33G16001DL	33G16001F	33G17001	33G17001F	33G18001
	LAB ID:	534646	534646	534627	534602	534624	534599
	DATE RCVD:	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
NAME	UNITS						
S Benzo(a)pyrene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Benzo(g,h,i)perylene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Butylbenzylphthalate	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Carbazole	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Chrysene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Di-n-butylphthalate	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Di-n-octylphthalate	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Dibenz(a,h)anthracene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Dibenzofuran	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Diethylphthalate	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Dimethylphthalate	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Fluoranthene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Fluorene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Hexachlorobenzene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Hexachlorobutadiene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Hexachlorocyclopentadiene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Hexachloroethane	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Ieophorone	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S N-Nitrosodiphenylamine	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Naphthalene	UG/L	16.0000 J	26.0000 J	.	20.0000 U	.	10.0000 U
S Nitrobenzene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Pentachlorophenol	UG/L	250.0000 U	380.0000 U	.	50.0000 U	.	25.0000 U
S Phenanthrene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U
S Phenol	UG/L	100.0000 U	150.0000 U	.	2.0000 J	.	10.0000 U
S Pyrene	UG/L	100.0000 U	150.0000 U	.	20.0000 U	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G16001	33G16001DL	330160013	33G17001	33G17001F	33G18001
	LAB ID:	534646	534646	534627	534602	534624	534599
	DATE RCVD:	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
NAME	UNITS						
S bis(2-Chloroethoxy)methane	UG/L	100.0000 U	150.0000 U	.	20.0000 U		10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	100.0000 U	150.0000 U		20.0000 U		1.0000 U
S bis(2-chloroethyl)ether	UG/L	100.0000 U	150.0000 U		20.0000 U	.	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	100.0000 U	150.0000 U		20.0000 U	.	10.0000 U
V 1,1,1-Trichloroethane	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	20.0000 U			10.0000 U	.	10.0000 U
V 1,1,2-Trichloroethane	UG/L	20.0000 U			10.0000 U	.	10.0000 U
V 1,1-Dichloroethane	UG/L	54.0000		.	160.0000	.	10.0000 U
V 1,1-Dichloroethene	UG/L	20.0000 U			10.0000 U	.	10.0000 U
V 1,2-Dichloroethane	UG/L	20.0000 U			10.0000 U	.	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	20.0000 U			3.0000 J		10.0.000 U
V 1,2-Dichloropropane	UG/L	20.0000 U			10.0000 U	.	10.0000 U
V 2-Butanone	UG/L	95.0000 U		.	180.0000 U	.	110.0000 U
V 2-Hexanone	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V Acetone	UG/L	20.0000 U		.	30.0000	.	40.0000 U
V Benzene	UG/L	3.0000 J		.	10.0000 U	.	10.0000 U
V Bromodichloromethane	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V Bromoform	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V Bromomethane	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V Carbon Disulfide	UG/L	7.0000 J		.	3.0000 J	.	10.0000 U
V Carbon Tetrachloride	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V Chlorobenzene	UG/L	340.0000		.	37.0000	.	10.0000 U
V Chloroethane	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V Chloroform	UG/L	20.0000 U		.	10.0000 U	.	10.0000 U
V Chloromethane	UG/L	20.0000 U			10.0000 U	.	10.0000 U
V Dibromochloromethane	UG/L	20.0000 U		.	10.0000 U		10.0000 U
V Ethylbenzene	UG/L	4.0000 J		.	10.0000 U	.	10.0000 U

NAS YENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33016001	33G16001DL	33G16001F	33017001	33G17001F	33018001
	LAB ID:	534646	534646	534627	534785	534624	534599
	DATE RCVD:	02/03/93	02/03/93	02/03/93	02/04/93	02/03/93	02/03/93
NAME	UNITS						
V Methylene Chloride	UG/L	20.0000 U	.	.	10.0000 U	.	6.0000 U
V Styrene	UG/L	20.0000 U	.	.	10.0000 U	.	10.0000 U
V Tetrachloroethene	UG/L	20.0000 U	.	.	10.0000 U	.	10.0000 U
V Toluene	UG/L	10.0000 J	.	.	10.0000 U	.	10.0000 U
V Trichloroethene	UG/L	20.0000 U	.	.	10.0000 U	.	10.0000 U
V Vinyl Chloride	UG/L	20.0000 U	.	.	10.0000 U	.	10.0000 U
V Xylene (total)	UG/L	10.0000 J	.	.	10.0000 U	.	10.0000 U
V cis-1,3-Dichloropropene	UG/L	20.0000 U	.	.	10.0000 U	.	10.0000 U
V trans-1,3-Dichloropropene	UG/L	20.0000 U	.	.	10.0000 U	.	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33618001F	33G20001	33020001DL	33G20001F	33081001	33G81001F
ENSAFE ID:		33618001F	33G20001	33020001DL	33G20001F	33081001	33G81001F
LAB ID:		534621	547566	547566	547570	533777	533792
DATE RCVD:		02/03/93	04/27/93	04/27/93	04/27/93	01/27/93	01/27/93
I Aluminum	UG/L	56.4000 J	2630.0000	.	54.0000 J	3110.0000	2040.0000
I Antimony	UG/L	38.7000 J	36.0000 U	.	36.0000 U	57.0000 U	52.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	.	3.0000 U	3.0000 U	3.3000 J
I Barium	UG/L	25.8000 J	19.7000 J	.	17.7000 J	7.7000 J	4.6000 J
I Beryllium	UG/L	1.0000 U	1.0000 U	.	1.0000 U	3.0000 U	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	.	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	31800.0000	60600.0000	.	64300.0000	2700.0000 J	2820.0000 J
I Chromium	UG/L	6.0000 U	10.4000	.	8.0000 U	11.9000	6.0000 U
I Cobalt	UG/L	9.0000 U	8.0000 U	.	8.0000 U	12.0000 U	9.0000 U
I Copper	UG/L	4.0000 U	4.7000 J	.	4.0000 U	25.0000 U	4.0000 U
I Cyanide	UG/L		10.0000 U	.	2.0000	10.0000 U	
I Iron	UG/L	791.0000	4770.0000	.	1340.0000	2910.0000	1760.0000
I Lead	UG/L	2.0000 U	2.0000 U	.	2.0000 U	2.0000 U	2.0000 U
I Magnesium	UG/L	8410.0000	33600.0000	.	36100.0000	5320.0000	5620.0000
I Manganese	UG/L	47.1000	45.2000	.	42.7000	14.6000 J	6.2000 J
I Mercury	UG/L	0.2000 U	0.2000 U	.	0.2000 U	0.2000 U	0.2000 U
I Nickel	UG/L	13.0000 U	17.0000 U	.	17.0000 U	38.0000 U	13.0000 U
I Potassium	UG/L	8080.0000	22600.0000	.	24900.0000	3950.0000 U	1440.0000 U
I Selenium	UG/L	3.0000 U	3.0000 U	.	3.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	5.0000 U	7.0000 U	.	7.0000 U	7.0000	5.0000 U
I Sodium	UG/L	196000.000	275000.000	.	300000.000	7220.0000	7880.0000
I Thallium	UG/L	3.0000 U	40.0000 U	.	40.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	4.0000 U	17.6000 J	.	10.4000 J	15.0000 U	8.4000 J
I Zinc	UG/L	8.6000 J	9.7000 J	.	4.0000 U	57.2000	7.1000 J
P 4,4'-DDD	UG/L		0.1000 U	.		0.1000 U	
P 4,4'-DDE	UG/L		0.1000 U	.		0.1000 U	
P 4,4'-DDT	UG/L		0.1000 U	.		0.1000 U	

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G18001F	33G20001	33G20001DL	33G20001F	33G81001	33G81001F
	LAB ID:	534621	547566	547566	547570	533777	533792
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	01/27/93	01/27/93
NAME	UNITS						
P Aldrin	UG/L	.	0.0500 U	.	.	0.0500 U	.
P Aroclor 1016	UG/L	.	1.0000 U		.	1.0000 U	.
P Aroclor 1221	UG/L	.	2.0000 U		.	2.0000 U	.
P Aroclor 1232	UG/L	.	1.0000 U		.	1.0000 U	.
P Aroclor 1242	UG/L	.	1.0000 U		.	1.0000 U	.
P Aroclor 1240	UG/L	.	1.0000 U		.	1.0000 U	.
P Aroclor 1254	UG/L	.	1.0000 U		.	1.0000 U	.
P Aroclor 1260	UG/L	.	1.0000 U		.	1.0000 U	.
P Dieldrin	UG/L	.	0.1000 U		.	0.1000 U	.
P Endosulfan I	UG/L	.	0.0500 U		.	0.0500 U	.
P Endoeulfan II	UG/L	.	0.1000 U		.	0.1000 U	.
P Endosulfan sulfate	UG/L	.	0.1000 U		.	0.1000 U	.
P Endrin	UG/L	.	0.1000 U		.	0.1000 U	.
P Endrin aldehyde	UG/L	.	0.1000 U		.	0.1000 U	.
P Endrin ketone	UG/L	.	0.1000 U		.	0.1000 U	.
P Heptachlor	UG/L	.	0.0500 U		.	0.0011 J	.
P Heptachlor epoxide	UG/L	.	0.0500 U		.	0.0500 U	.
P Methoxychlor	UG/L	.	0.5000 U		.	0.5000 U	.
P Toxaphene	UG/L	.	5.0000 U		.	5.0000 U	.
P alpha-BHC	UG/L	.	0.0500 U		.	0.0500 U	.
P alpha-Chlordane	UG/L	.	0.0500 U		.	0.0500 U	.
P beta-BHC	UG/L	.	0.0500 U		.	0.0500 U	.
P de1ta-BHC	UG/L	.	0.0500 U		.	0.0500 U	.
P gamma-BHC (Lindane)	UG/L	.	0.0500 U		.	0.0500 U	.
P gamma-Chlordane	UG/L	.	0.0500 U		.	0.0500 U	.
S 1,2,4-Trichlorobenzene	UG/L	.	1.0000 J	300.0000 U	.	10.0000 U	.
S 1,2-Dichlorobenzene	UG/L	.	570.0000 J	1200.0000	.	10.0000 U	.
S 1,3-Dichlorobenzene	UG/L	.	400.0000 J	700.0000	.	10.0000 U	.

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33618001F	33620001	33G20001DL	33G20001F	33081001	33G81001F
	LAB ID:	534621	547566	547566	547570	533777	533792
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	01/27/93	01/27/93
NAME	UNITS						
S 1,4-Dichlorobenzene	UG/L	.	360.0000 J	670.0000	.	10.0000 U	
S 2,4,5-Trichlorophenol	UG/L	.	1.0000 J	750.0000 U	.	25.0000 U	
S 2,4,6-Trichlorophenol	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 2,4-Dichlorophenol	UG/L	.	12.0000	300.0000 U	.	10.0000 U	
S 2,4-Dimethylphenol	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 2,4-Dinitrophenol	UG/L	.	25.0000 U	750.0000 U	.	25.0000 U	
S 2,4-Dinitrotoluene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 2,6-Dinitrotoluene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 2-Chloronaphthalene	UG/L	.	2.0000 J	300.0000 U	.	10.0000 U	
S 2-Chlorophenol	UG/L	.	2.0000 J	300.0000 U	.	10.0000 U	
S 2-Methylnaphthalene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 2-Methylphenol	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 2-Nitroaniline	UG/L	.	25.0000 U	750.0000 U	.	25.0000 U	
S 2-Nitrophenol	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 3,3'-Dichlorobenzidine	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 3-Nitroaniline	UG/L	.	25.0000 U	750.0000 U	.	25.0000 U	
S 4,6-Dinitro-2-methylphenol	UG/L	.	25.0000 U	750.0000 U	.	25.0000 U	
S 4-Bromophenyl-phenylethar	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 4-Chloro-3-methylphenol	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 4-Chloroaniline	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 4-Chlorophenyl-phenylether	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 4-Methylphenol	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S 4-Nitroaniline	UG/L	.	25.0000 U	750.0000 U	.	25.0000 U	
S 4-Nitrophenol	UG/L	.	25.0000 U	750.0000 U	.	25.0000 U	
S Acenaphthene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S Acenaphthylene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S Anthracene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	
S Benzo(a)anthracene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G18001F	33G20001	33G20001DL	33G20001F	33G81001	33G81001F
	LAB ID:	534621	547566	547566	547570	533777	533792
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	01/27/93	01/27/93
NAME	UNITS						
S Benzo(a)pyrene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Benzo(b,k)fluoranthene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Benzo(g,h,i)perylene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Butylbenzylphthalate	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Carbazole	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Chrysene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Di-n-butylphthalate	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Di-n-octylphthalate	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Dibenzofuran	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Diethylphthalate	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Dimethylphthalate	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Fluoranthene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Fluorene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Hexachlorobenzene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Hexachlorobutadiene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Hexachloroethane	UG/L	.	3.0000 J	300.0000 U	.	10.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Ieophorone	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S N-Nitroeodiphenylamine	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Naphthalene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Nitrobenzene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Pentachlorophenol	UG/L	.	25.0000 U	750.0000 U	.	25.0000 U	.
S Phenanthrene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Phenol	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S Pyrene	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G18001F	33G20001	33G20001DL	33G20001F	33G81001	33G81001F
	LAB ID:	534621	547566	547566	547570	533777	533792
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	01/27/93	01/27/93
NAME	UNITS						
S bis(2-Chloroethoxy)methane	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	.	4.0000 U	300.0000 U	.	2.0000 U	.
S bis(2-chloroethyl)ether	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L	.	10.0000 U	300.0000 U	.	10.0000 U	.
V 1,1,1-Trichloroethane	UG/L	.	33.0000 U		.	10.0000 U	.
V 1,1,2,2-Tetrachloroethane	UG/L	.	33.0000 U		.	10.0000 U	.
V 1,1,2-Trichloroethane	UG/L	.	33.0000 U		.	10.0000 U	.
V 1,1-Dichloroethane	UG/L	.	190.0000		.	10.0000 U	.
V 1,1-Dichloroethene	UG/L	.	33.0000 U		.	10.0000 U	.
V 1,2-Dichloroethane	UG/L	.	33.0000 U		.	10.0000 U	.
V 1,2-Dichloroethene (total)	UG/L	.	12.0000 J		.	10.0000 U	.
V 1,2-Dichloropropane	UG/L	.	33.0000 U		.	10.0000 U	.
V 2-Butanone	UG/L	.	33.0000 U		.	10.0000 U	.
V 2-Hexanone	UG/L	.	33.0000 U		.	10.0000 U	.
V 4-Methyl-2-Pentanone	UG/L	.	33.0000 U		.	10.0000 U	.
V Acetone	UG/L	.	33.0000 U		.	10.0000 U	.
V Benzene	UG/L	.	33.0000 U		.	10.0000 U	.
V Bromodichloromethane	UG/L	.	33.0000 U		.	10.0000 U	.
V Bromoform	UG/L	.	33.0000 U		.	10.0000 U	.
V Bromomethane	UG/L	.	33.0000 U		.	10.0000 U	.
V Carbon Disulfide	UG/L	.	33.0000 U		.	10.0000 U	.
V Carbon Tetrachloride	UG/L	.	33.0000 U		.	10.0000 U	.
V Chlorobenzene	UG/L	.	200.0000		.	10.0000 U	.
V Chloroethane	UG/L	.	33.0000 U		.	10.0000 U	.
V Chloroform	UG/L	.	33.0000 U		.	10.0000 U	.
V Chloromethane	UG/L	.	33.0000 U		.	10.0000 U	.
V Dibromochloromethane	UG/L	.	33.0000 U		.	10.0000 U	.
V Ethylbenzene	UG/L	.	33.0000 U		.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G18001F	33G20001	33G20001DL	33G20001F	33G81001	33G81001F
	LAB ID:	534621	547566	547566	547570	533777	533792
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	01/27/93	01/27/93
NAME	UNITS						
V Methylene Chloride	UG/L	.	26.0000 U	.	.	5.0000 U	.
V Styrene	UG/L	.	33.0000 U	.	.	10.0000 U	.
V Tetrachloroethene	UG/L	.	6.0000 J	.	.	10.0000 U	.
V Toluene	UG/L	.	6.0000 J	.	.	10.0000 U	.
V Trichloroethene	UG/L	.	7.0000 J	.	.	10.0000 U	.
V Vinyl Chloride	UG/L	.	33.0000 U	.	.	10.0000 U	.
V Xylene (total)	UG/L	.	6.0000 J	.	.	10.0000 U	.
V cis-1,3-Dichloropropene	UG/L	.	33.0000 U	.	.	10.0000 U	.
V trans-1,3-Dichloropropene	UG/L	.	33.0000 U	.	.	10.0000 U	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA ou-io AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
ENSAFE ID:		33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
LAB ID:		534601	534623	547145	547146	534123	534188
DATE RCVD:		02/03/93	02/03/93	04/22/93	04/22/93	01/29/93	01/29/93
I Aluminum	UG/L	1100.0000	123.0000 J	887000.000 J	120.0000 J	1010.0000	242.0000
I Antimony	UG/L	29.0000 U	30.0000 U	41.2000 J	36.0000 U	57.0000 U	52.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	30.0000 U	3.0000 U	6.3860 J	6.2000 J
I Barium	UG/L	19.3000 J	9.8000 J	29.2000 J	28.0000 J	17.9000 J	15.8000 J
I Beryllium	UG/L	3.0000 U	1.0000 U	78.6000	1.0000 U	3.0000 U	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.4000	5.0096 U	20.2000	5.0000 U
I Calcium	UG/L	28800.0000	30800.0000	210000.000	49400.0000	25300.0000	27500.0000
I Chromium	UG/L	17.0000	6.0000 U	316.0000	8.0000 U	12.7000	7.3000 J
I Cobalt	UG/L	12.0000 U	9.0000 U	26.9000 J	8.0000 U	12.0000 U	9.0000 U
I Copper	UG/L	25.0000 U	4.0000 U	4.0000 U	4.0000 U	25.0000 U	4.0000 U
I Cyanide	UG/L	10.0000 U		2.0000	2.0000	10.0000 U	
I Iron	UG/L	2520.0000	1540.0000	399000.000	229.0000	2760.0000	1110.0000
I Lead	UG/L	2.0000 U	2.0000 U	100.0000 U	2.0000 U	2.2000 J	2.0000 U
I Magnesium	UG/L	3820.0000 J	3330.0000 J	237000.000	39700.0000	3030.0000 J	3280.0000 J
I Manganese	UG/L	76.0000	71.6000	1560.0000	86.8000	36.4000	37.7000
I Mercury	UG/L	1.5000	0.2000 U	0.2000 U	0.2000 U	0.3700	0.2000 U
I Nickel	UG/L	38.0000 U	13.0000 U	104.0000	17.0000 U	38.0000 U	13.0000 U
I Potassium	UG/L	3950.0000 U	2480.0000 J	97300.0000	27500.0000	3950.0000 U	1710.0000 J
I Selenium	UG/L	3.0000 U	3.0000 U	30.0000 U	3.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	5.0000 U	5.0000 U	7.0000 U	7.0000 U	10.0000 U	5.0000 U
I Sodium	UG/L	15600.0000	17300.0000	593000.000	599000.000	2880.0000 J	3260.0000 J
I Thallium	UG/L	3.0000 U	3.0000 U	200.0000 U	40.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	15.0000 U	4.0000 U	521.0000	10.2000 J	15.0000 U	4.0000 U
I Zinc	UG/L	53.5000	5.0000 U	388.0000	7.1000 J	208.0000	193.0000
P 4,4'-DDD	UG/L	0.1000 U				0.1000 U	
P 4,4'-DDE	UG/L	0.1000 U				0.1000 U	
P 4,4'-DDT	UG/L	0.1000 U				0.1000 U	

NAS I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
	LAB ID:	534601	534623	547145	547146	534227	534188
	DATE RCVD:	02/03/93	02/03/93	04/22/93	04/22/93	01/29/93	01/29/93
NAME	UNITS						
P Aldrin	UG/L	0.0500 U	.	.	.	0.0500 U	.
P Aroclor 1016	UG/L	1.0000 U	.	.	.	1.0000 U	.
P Aroclor 1221	UG/L	2.0000 U	.	.	.	2.0000 U	.
P Aroclor 1232	UG/L	1.0000 U	.	.	.	1.0000 U	.
P Aroclor 1242	UG/L	1.0000 U	.	.	.	1.0000 U	.
P Aroclor 1240	UG/L	1.0000 U	.	.	.	1.0000 U	.
P Aroclor 1254	UG/L	1.0000 U	.	.	.	1.0000 U	.
P Aroclor 1260	UG/L	1.0000 U	.	.	.	1.0000 U	.
P Dieldrin	UG/L	0.1000 U	.	.	.	0.1000 U	.
P Endosulfan I	UG/L	0.0500 U	.	.	.	0.0500 U	.
P Endosulfan II	UG/L	0.1000 U	.	.	.	0.1000 U	.
P Endosulfan sulfate	UG/L	0.1000 U	.	.	.	0.1000 U	.
P Endrin	UG/L	0.1000 U	.	.	.	0.1000 U	.
P Endrin aldehyde	UG/L	0.1000 U	.	.	.	0.1000 U	.
P Endrin ketone	UG/L	0.1000 U	.	.	.	0.1000 U	.
P Heptachlor	UG/L	0.0500 U	.	.	.	0.0013 U	.
P Heptachlor epoxide	UG/L	0.0500 U	.	.	.	0.0500 U	.
P Methoxychlor	UG/L	0.5000 U	.	.	.	0.5000 U	.
P Toxaphene	UG/L	5.0000 U	.	.	.	5.0000 U	.
P alpha-BHC	UG/L	0.0500 U	.	.	.	0.0500 U	.
P alpha-Chlordane	UG/L	0.0500 U	.	.	.	0.0500 U	.
P beta-BHC	UG/L	0.0500 U	.	.	.	0.0500 U	.
P delta-BHC	UG/L	0.0500 U	.	.	.	0.0500 U	.
P gamma-BHC (Lindane)	UG/L	0.0500 U	.	.	.	0.0500 U	.
P gamma-Chlordane	UG/L	0.0500 U	.	.	.	0.0013 J	.
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	.	.	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

NAME	UNITS	33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
ENSAFE ID:		33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
LAB ID:		534601	534623	547145	547146	534227	534188
DATE RCVD:		02/03/93	02/03/93	04/22/93	04/22/93	01/29/93	01/29/93
S 1,4-Dichlorobenzene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	.	.	.	25.0000 U	.
S 2,4,6-Trichlorophenol	UG/L	10.0000 U			.	10.0000 U	.
S 2,4-Dichlorophenol	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 U	.		.	10.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	.	.	.	25.0000 U	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U	.	.	.	25.0000 U	.
S 2-Nitrophenol	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U	.	.	.	25.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	.	.	.	25.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 4-Chloroaniline	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 4-Methylphenol	UG/L	10.0000 U	.	.	.	10.0000 U	.
S 4-Nitroaniline	UG/L	25.0000 U	.	.	.	25.0000 U	.
S 4-Nitrophenol	UG/L	25.0000 U	.	.	.	25.0000 U	.
S Acenaphthene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Acenaphthylene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Anthracene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Beneo(a)anthracene	UG/L	10.0000 U	.	.	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
ENSAFE ID:		33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
LAB ID:		534601	534623	547145	547146	534227	534188
DATE RCVD:		02/03/93	02/03/93	04/22/93	04/22/93	01/29/93	01/29/93
S Benzo(a)pyrene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Benzo(g,h,i)perylene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Eutylbenzylphthalate	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Carbazole	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Chrysene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Di-n-butylphthalate	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Di-n-octylphthalate	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Dibenzofuran	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Diethylphthalate	UG/L	10.0000 U	.	.	.	1.0000 U	.
S Dimethylphthalate	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Fluoranthene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Fluorene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Hexachlorobenzene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Hexachlorobutadiene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Hexachloroethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Isophorone	UG/L	10.0000 U	.	.	.	10.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	.	.	.	10.0000 U	.
S N-Nitrosodiphenylamine	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Naphthalene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Nitrobenzene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Pentachlorophenol	UG/L	25.0000 U	.	.	.	25.0000 U	.
S Phenanthrene	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Phenol	UG/L	10.0000 U	.	.	.	10.0000 U	.
S Pyrene	UG/L	10.0000 U	.	.	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
	LAB ID:	534601	534623	547145	547146	534227	534188
	DATE RCVD:	02/03/93	02/03/93	04/22/93	04/22/93	01/29/93	01/29/93
NAME	UNITS						
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	.	.	.	10.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	1.0000 U	.	.	.	2.0000 U	.
S bis(2-chloroethyl)ether	UG/L	10.0000 U	.	.	.	10.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,1,1-Trichloroethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,1,2-Trichloroethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,1-Dichloroethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,1-Dichloroethene	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,2-Dichloroethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 1,2-Dichloropropane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 2-Butanone	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 2-Hexanone	UG/L	10.0000 U	.	.	.	10.0000 U	.
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Acetone	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Benzene	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Bromodichloromethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Bromoform	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Bromomethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Carbon Disulfide	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Carbon Tetrachloride	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Chlorobenzene	UG/L	10.0000 U	.	.	.	1.0000 J	.
V Chloroethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Chloroform	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Chloromethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Dibromochloromethane	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Ethylbenzene	UG/L	10.0000 U	.	.	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM131	33GGM131F	33GGM661	33GGM691	33GGM711	33GGM711F
	LAB ID:	534601	534623	547145	547146	534227	534188
	DATE RCVD:	02/03/93	02/03/93	04/22/93	04/22/93	01/29/93	01/29/93
NAME	UNITS						
V Methylene Chloride	UG/L	3.0000 U	.	.	.	1.0000 U	.
V Styrene	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Tetrachloroethene	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Toluene	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Trichloroethene	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Vinyl Chloride	UG/L	10.0000 U	.	.	.	10.0000 U	.
V Xylene (total)	UG/L	10.0000 U	.	.	.	10.0000 U	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	.	.	.	10.0000 U	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	.	.	.	10.0000 U	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID:	33GGM721	33GGM721F	33GGM761	33GGM761F	33GGM771	33GGM771F
	LAB ID:	534218	534193	534197	534185	534215	534190
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
	UNITS						
I Aluminum	UG/L	359.0000	251.0000	500.0000	221.0000	3830.0600	244.0000
I Antimony	UG/L	57.0000 U	52.0000 U	57.0000 U	52.0000 U	57.0000 U	52.0000 U
I Arsenic	UG/L	16.2000	11.5000	3.0000 U	4.1000 J	3,0000 U	3.0000 U
I Barium	UG/L	27.1000 J	28.6000 J	11.5000 J	12.1000 J	17.1000 J	13.1000 J
I Beryllium	UG/L	3.0000 U	1.0000 U	3.0000 U	1.0000 U	3.0000 U	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0006 U	5.0000 U
I Calcium	UG/L	31000.0000	35100.0000	23900.0000	24000.0000	43000.0000	48000.0000
I Chromium	UG/L	10.0000 U	6.0000 U	10.0000 U	6.0000 U	10.0000 U	6.0000 U
I Cobalt	UG/L	12.0000 U	9.0000 U	12.0000 U	9.0000 U	12.0000 U	9.0000 U
I Copper	UG/L	25.0000 U	4.0000 U	25.0000 U	4.0000 U	25.0000 U	4.0000 U
I Cyanide	UG/L	10.0000 U		10.0000 U		10.0000 U	
I Iron	UG/L	2270.0000	1990.0000	467.0000	13.0000 U	2930.0000	113.0000
I Lead	UG/L	2.1000 J	2.0000 U	2.0000 U	2.0000 U	5.4000	2.0000 U
I Magnesium	UG/L	3690.0000 J	3470.0000 J	9050.0000	9030.0000	11800.0000	12500.0000
I Manganese	UG/L	149.0000	120.0000	157.0000	164.0000	75.00	79.5000
I Mercury	UG/L	0.2200	0.2000 U	0.2900	0.2000 U	0.2000 U	0.2000 U
I Nickel	UG/L	38.0000 U	13.0000 U	38.0000 U	13.0000 U	38.0000 U	13.0000 U
I Potassium	UG/L	3950.0000 U	1440.0000 U	6440.0000	5320.0000	6220.0000	6120.0000
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	10.0000 U	5.0000 U	10.0000 U	5.0000 U	10.0000 U	5.0000
I Sodium	UG/L	37000.0000	45400.0000	71800.0000	74600.0000	32300.0000	40500.0000
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	15.0000 U	4.0000 U	15.0000 U	4.0000 U	15.0000 U	4.0000 U
I Zinc	UG/L	76.2000	52.3000	3.0000 U	5.0000 U	15.4000 J	5.0000 U
P 4,4'-DDD	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDE	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDT	UG/L	0.1000 U		0.1000 U		0.1000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33GGM721	33GGM721F	33GGM761	33GGM761F	33GGM771	33GGM771F
ENSAFE ID:		33GGM721	33GGM721F	33GGM761	33GGM761F	33GGM771	33GGM771F
LAB ID:		534232	534193	534224	534185	534229	534190
DATE RCVD:		01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
P Aldrin	UG/L	0.0500 U		0.0500 U	.	0.0500 U	
P Aroclor 1016	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	
P Aroclor 1221	UG/L	2.0000 U		2.0000 U		2.0000 U	.
P Aroclor 1232	UG/L	1.0000 U		1.0000 U		1.0000 U	.
P Aroclor 1242	UG/L	1.0000 U		1.0000 U	.	1.0000 U	.
P Aroclor 1248	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1254	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1260	UG/L	1.0000 U	.	1.0000 U		1.0000 U	.
P Dieldrin	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endosulfan I	UG/L	0.0021 J	.	0.0500 U	.	0.0500 U	.
P Endosulfan II	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endosulfan sulfate	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endrin	UG/L	0.1000 U	.	0.1000 U		0.1000 U	.
P Endrin aldehyde	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	
P Endrin ketone	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Heptachlor	UG/L	0.0500 U	.	0.0500 U		0.0500 U	
P Heptachlor epoxide	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P Methoxychlor	UG/L	0.5000 U	.	0.5000 U	.	0.5000 U	.
P Toxaphene	UG/L	5.0000 U	.	5.0000 U	.	5.0000 U	.
P alpha-BHC	UG/L	0.0500 U	.	0.0500 U		0.0500 U	.
P alpha-Chlordane	UG/L	0.0500 U	.	0.0500 U		0.0500 U	.
P beta-BHC	UG/L	0.0500 U	.	0.0500 U		0.0500 U	
P delta-BHC	UG/L	0.0500 U	.	0.0500 U		0.0500 U	.
P gamma-BHC (Lindane)	UG/L	0.0500 U	.	0.0220 J	.	0.0500 U	.
P gamma-Chlordane	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	.	20.0000 U		10.0000 U	.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	.	20.0000 U		1.0000 J	.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33GGM721 / 534232 / 01/29/93	33GGM721F / 534193 / 01/29/93	33GGM761 / 534224 / 01/29/93	33GGM761F / 534185 / 01/29/93	33GGM771 / 534229 / 01/29/93	33GGM771F / 534190 / 01/29/93
S 1,4-Dichlorobenzene	UG/L	10.0000 U	.	20.0000 U	.	1.0000 J	.
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2,4-Dichlorophenol	UG/L	10.0000 U	.	20.0000 U		10.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S 2-Nitrophenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 4-Chloroaniline	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 4-Methylphenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S 4-Nitroaniline	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S 4-Nitrophenol	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S Acenaphthene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Acenaphthylene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Anthracene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Benzo(a)anthracene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM721	33GGM721F	33GGM761	33GGM761F	33GGM771	33GGM771F
	LAB ID:	534232	534193	534224	534185	534229	534190
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29 /93	01/29/93	01/29 /93
NAME	UNITS						
S Benzo(a)pyrene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 u	.
S Benzo(b,k)fluoranthene	UG/L	10.0000 u	.	20.0000 U	.	10.0000 U	.
S Benzo(g,h,i)perylene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Butylbenzylphthalate	UG/L	10.0000 U	.	3.0000 U	.	10.0000 U	.
S Carbazole	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Chrysene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Di-n-butylphthalate	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Di-n-octylphthalate	UG/L	10.0000 U	.	20.0000 u	.	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.
S Dibenzofuran	UG/L	10.0000 U	.	20.0000 u	.	10.0000 U	.
S Diethylphthalate	UG/L	10.0000 U	.	20.0000 u	.	10.0000 U	.
S Dimethylphthalate	UG/L	10.0000 U	.	20.0000 u	.	10.0000 u	.
S Fluoranthene	UG/L	10.0000 U	.	20.0000 u	.	10.0000 U	.
S Fluorene	UG/L	10.0000 u	.	20.0000 u	.	10.0000 u	.
S Hexachlorobenzene	UG/L	10.0000 U	.	20.0000 u	.	10.0000 u	.
S Hexachlorobutadiene	UG/L	10.0000 U	.	20.0000 u	.	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	10.0000 u	.	20.0000 u	.	10.0000 U	.
S Hexachloroethane	UG/L	10.0000 u	.	20.0000 u	.	10.0000 u	.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 u	.	20.0000 u	.	10.0000 u	.
S Ieophorone	UG/L	10.0000 u	.	20.0000 u	.	10.0000 u	.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	.	20.0000 u	.	10.0000 u	.
S N-Nitrosodiphenylamine	UG/L	10.0000 u	.	20.0000 U	.	10.0000 u	.
S Naphthalene	UG/L	10.0000 U	.	20.0000 u	.	10.0000 U	.
S Nitrobenzene	UG/L	10.0000 U	.	20.0000 u	.	10.0000 u	.
S Pentachlorophenol	UG/L	25.0000 U	.	50.0000 U	.	25.0000 U	.
S Phenanthrene	UG/L	10.0000 u	.	20.0000 u	.	10.000'0 u	.
S Phenol	UG/L	10.0000 U	.	20.0000 U	.	10.0000 u	.
S Pyrene	UG/L	10.0000 U	.	20.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM721	33GGM721F	33GGM761	33GGM761F	33GGM771	33GGM771F
	LAB ID:	534232	534193	534224	534185	534229	534190
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U		20.0000 U	.	10.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	2.0000 U		88.0000	.	2.0000 U	.
S bis(2-chloroethyl)ether	UG/L	10.0000 U		20.0000 U	.	10.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U		20.0000 U	.	10.0000 U	.
V 1,1,1-Trichloroethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 1,1,2-Trichloroethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 1,1-Dichloroethane	UG/L	10.0000 U		10.0000 U	.	2.0000 J	.
V 1,1-Dichloroethene	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethene (total)	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 1,2-Dichloropropane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 2-Butanone	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 2-Hexanone	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V 4-Methyl-2-Pentanone	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Acetone	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Benzene	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Bromodichloromethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Bromoform	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Bromomethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Carbon Disulfide	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Carbon Tetrachloride	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Chlorobenzene	UG/L	10.0000 U		4.0000 J	.	6.0000 J	.
V Chloroethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Chloroform	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Chloromethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Dibromochloromethane	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.
V Ethylbenzene	UG/L	10.0000 U		10.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM721	33GGM721F	33GGM761	33GGM761F	33GGM771	33GGM771F
	LAB ID:	534232	534193	534224	534185	534229	534190
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
V Methylene Chloride	UG/L	10.0000 U	.	1.0000 U	.	1.0000 U	.
V Styrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Tetrachloroethene	UG/L	10.0000 U	.	10.0000 U	.	L0.0000 U	.
V Toluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Trichloroethene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Vinyl Chloride	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Xylene (total)	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM781	33GGM781F	33GR3001	33GR3001DL	33GR3001F
	LAB ID:	534216	534191	534779	534779	534781
	DATE RCVD:	01/29/93	01/29/93	02/04/93	02/03/93	02/04/93
NAME	UNITS					
I Aluminum	UG/L	580.0000	277.0000	4300.0000	.	444.0000
I Antimony	UG/L	57.6000 U	52.0000 U	29.0000 U	.	30.0000 U
I Arsenic	UG/L	3.0000 U	6.5000 J	3.0000 U	.	3.0000 U
I Barium	UG/L	27.8000 J	28.0000 J	8.5000 J	.	3.3000 J
I Beryllium	UG/L	3.0000 U	1.0000 U	3.0000 U	.	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	.	5.0000 U
I Calcium	UG/L	21000.0000	21700.0000	18800.0000	.	19300.0000
I Chromium	UG/L	10.0000 U	6.0000 U	16.2000	.	6.2000 J
I Cobalt	UG/L	12.0000 U	9.0000 U	12.0000 U	.	9.0000 U
I Copper	UG/L	25.0000 U	4.0000 U	25.0000 U	.	4.0000 U
I Cyanide	UG/L	10.0000 U		10.0000 U	.	
I Iron	UG/L	543.0000	165.0000	3330.0000	.	1510.0000
I Lead	UG/L	2.0000 U	2.0000 U	2.0000 U	.	2.0000 U
I Magnesium	UG/L	2640.0000 J	2820.0000 J	13600.0000	.	14100.0000
I Manganese	UG/L	8.2000 J	10.3000 J	21.7000	.	14.1000 J
I Mercury	UG/L	0.2000 U	0.2000 U	0.5000	.	0.2000 U
I Nickel	UG/L	38.0000 U	13.0000 U	38.0000 U	.	13.0000 U
I Potassium	UG/L	3950.0000 U	1440.0000 U	8050.0000	.	7230.0000
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	.	3.0000 U
I Silver	UG/L	10.0000 U	5.0000 U	10.0000 U	.	5.0000 U
I Sodium	UG/L	7740.0000	8550.0000	87600.0000	.	96300.0000
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	.	3.0000 U
I Vanadium	UG/L	15.0000 U	4.0000 U	29.0000 J	.	18.7000 J
I Zinc	UG/L	4.4000 J	5.0000 U	57.9000	.	5.0000 U
P 4,4'-DDD	UG/L	0.1000 U		0.1000 U	.	
P 4,4'-DDE	UG/L	0.1000 U		0.1000 U	.	
P 4,4'-DDT	UG/L	0.1000 U		0.1000 U	.	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM781	33GGM781F	33GR3001	33GR3001DL	33GR3001F
	LAB ID:	534230	534191	534779	534779	534781
	DATE RCVD:	01/29/93	01/29/93	02/04/93	02/03/93	02/04/93
NAME	UNITS					
P Aldrin	UG/L	0.0500 U	.	0.0500 U		. .
P Aroclor 1016	UG/L	1.0000 U	.	1.0000 U		. .
P Aroclor 1221	UG/L	2.0000 U	.	2.0000 U		. .
P Aroclor 1232	UG/L	1.0000 U	.	1.0000 U		. .
P Aroclor 1242	UG/L	1.0000 U	.	1.0000 U		. .
P Aroclor 1248	UG/L	1.0000 U	.	1.0000 U		. .
P Aroclor 1254	UG/L	1.0000 U	.	1.0000 U		. .
P Aroclor 1260	UG/L	1.0000 U	.	1.0000 U		. .
P Dieldrin	UG/L	0.1000 U	.	0.1000 U		. .
P Endoeulfan I	UG/L	0.0500 U	.	0.0500 U		. .
P Endoeulfan II	UG/L	0.1000 U	.	0.1000 U		. .
P Endosulfan sulfate	UG/L	0.1000 U	.	0.0370 J		. .
P Endrin	UG/L	0.1000 U	.	0.1000 U		. .
P Endrin aldehyde	UG/L	0.1000 U	.	0.1000 U		. .
P Endrin ketone	UG/L	0.1000 U	.	0.1000 U		. .
P Heptachlor	UG/L	0.0023 U	.	0.0500 U		. .
P Heptachlor epoxide	UG/L	0.0500 U	.	0.0500 U		. .
P Methoxychlor	UG/L	0.5000 U	.	0.5000 U		. .
P Toxaphene	UG/L	5.0000 U	.	5.0000 U		. .
P alpha-BHC	UG/L	0.0500 U	.	0.0500 U		. .
P alpha-Chlordane	UG/L	*0.0500* U	.	0.0500 U		. .
P beta-BHC	UG/L	0.0500 U	.	0.0500 U		. .
P delta-BHC	UG/L	0.0500 U	.	0.0500 U		. .
P gamma-BHC (Lindane)	UG/L	0.0500 U	.	0.0500 U		. .
P gamma-Chlordane	UG/L	0.0500 U	.	0.0500 U		. .
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	. .
S 1,2-Dichlorobenzene	UG/L	10.0000 U	.	72.0000	66.0000 J	. .
S 1,3-Dichlorobenzene	UG/L	10.0000 U	.	120.0000 J	110.0000 J	. .

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM781	33GGM781F	33GR3001	33GR3001DL	33GR3001F
	LAB ID:	534230	534191	534779	534779	534781
	DATE RCVD:	01/29/93	01/29/93	02/03/93	02/03/93	02/04/93
NAME	UNITS					
S 1,4-Dichlorobenzene	UG/L	10.0000 U	.	90.0000 J	85.0000 J	.
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2,4-Dichlorophenol	UG/L	10.0000 U	.	2.0000 J	30.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S 2-Nitrophenol	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	.	2.0000 J	30.0000 U	.
S 4-Chloroaniline	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 4-Methylphenol	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S 4-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S 4-Nitrophenol	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S Acenaphthene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Acenaphthylene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Anthracene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Benzo(a)anthracene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33GGM781	33GGM781F	33GR3001	33GR3001DL	33GR3001F
ENSAFE ID:		33GGM781	33GGM781F	33GR3001	33GR3001DL	33GR3001F
LAB ID:		534230	534191	534779	534779	534781
DATE RCVD:		01/29/93	01/29/93	02/03/93	02/03/93	02/04/93
S Benzo(a)pyrene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Benzo(g,h,i)perylene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Butylbenzylphthalate	UG/L	10.0000 U	.	1.0000 U	30.0000 U	.
S Carbazole	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Chrysene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	
S Di-n-butylphthalate	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Di-n-octylphthalate	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Dibenz(a,h)anthracene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Dibenzofuran	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Diethylphthalate	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Dimethylphthalate	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Fluoranthene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Fluorene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Hexachlorobenzene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Hexachlorobutadiene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	
S Hexachlorocyclopentadiene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Hexachloroethane	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Isophorone	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S N-Nitrosodiphenylamine	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Naphthalene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Nitrobenzene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	
S Pentachlorophenol	UG/L	25.0000 U	.	25.0000 U	75.0000 U	.
S Phenanthrene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Phenol	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S Pyrene	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM781	33GGM781F	33GR3001	33GR3001DL	33GR3001F
	LAB ID:	534230	534191	534779	534779	534781
	DATE RCVD:	01/29/93	01/29/93	02/03/93	02/03/93	02/04/93
NAME	UNITS					
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	2.0000 U	.	4.0000 u	3.0000 U	.
S bis(2-chloroethyl)ether	UG/L	10.0000 U	.	10,0000 U	30.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	.	10.0000 U	30.0000 U	.
V 1,1,1-Trichloroethane	UG/L	10.0000 U	.	10.0000 U	.	
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	.	10.0000 U	.	
V 1,1,2-Trichloroethane	UG/L	10.0000 U	.	10.0000 U	.	
V 1,1-Dichloroethane	UG/L	10.0000 U	.	9.0000 J	.	
V 1,1-Dichloroethene	UG/L	10.0000 U	.	10.0000 u	.	
V 1,2-Dichloroethane	UG/L	10.0000 U	.	10.0000 U	.	
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	.	10.0000 U	.	. .
V 1,2-Dichloropropane	UG/L	10.0000 U	.	10.0000 U	.	
V 2-Butanone	UG/L	10.0000 U	.	10.0000 U	.	
V 2-Hexanone	UG/L	10.0000 U	.	10.0000 U	.	
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	.	10.0000 U	.	
V Acetone	UG/L	10.0000 U	.	10.0000 U	.	
V Benzene	UG/L	10.0000 U	.	10.0000 U	.	
V Bromodichloromethane	UG/L	10.0000 U	.	10.0000 U	.	
V Bromoform	UG/L	10.0000 U	.	10.0000 U	.	
V Bromomethane	UG/L	10.0000 U	.	10.0000 U	.	
V Carbon Disulfide	UG/L	10.0000 U	.	10.0000 U	.	
V Carbon Tetrachloride	UG/L	10.0000 U	.	10.0000 U	.	
V Chlorobenzene	UG/L	10.0000 U	.	31.0000	.	.
V Chloroethane	UG/L	10.0000 U	.	4.0000 J	.	.
V Chloroform	UG/L	10.0000 U	.	10.0000 U	.	.
V Chloromethane	UG/L	10.0000 U	.	10.0000 U	.	.
V Dibromochloromethane	UG/L	10.0000 U	.	10.0000 U		.
V Ethylbenzene	UG/L	10.0000 U	.	10.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GGM781	33GGM781F	33GR3001	33GR3001DL	33GR3001F
	LAB ID:	534230	534191	534779	534779	534781
	DATE RCVD:	01/29/93	01/29/93	02/04/93	02/03/93	02/04/93
NAME	UNITS					
V Methylene Chloride	UG/L	10.0000 U		10.0000 U	.	.
V Styrene	UG/L	10.0000 U	.	10.0000 U	.	.
V Tetrachloroethene	UG/L	10.0000 U	.	10.0000 U	.	.
V Toluene	UG/L	10.0000 U	.	10.0000 U	.	.
V Trichloroethene	UG/L	10.0000 U	.	10.0000 U	.	.
V Vinyl Chloride	UG/L	10.0000 U	.	10.0000 U	.	.
V Xylene (total)	UG/L	10.0000 U	.	10.0000 U	.	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G06001	13G06001F	13607001	13G07001F	13G19001	13G19001F
	LAB ID:	534316	534329	533766	533788	533781	533794
	DATE RCVD:	01/30/93	01/30/93	01/27/93	01/27/93	01/27/93	01/27/93
NAME	UNITS						
I Aluminum	UG/L	2920.0000	257.0000	7740.0000	208.0000	1890.0000	172.0000 J
I Antimony	UG/L	57.0000 U	52.0000 U	57.0000 U	52.0000 U	57.0000 U	52.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	3.1000 J	3.0000 U	3.0000 U	4.5000 J
I Barium	UG/L	17.8000 J	11.2000 J	25.2000 J	5.9000 J	29.9000 J	28.4000 J
I Beryllium	UG/L	3.0000 U	1.0000 U	3.0000 U	1.0000 U	3.0000 U	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	9720.0000	10600.0000	11100.0000	10000.0000	48800.0000	63500.0000
I Chromium	UG/L	10.0000 U	6.0000 U	28.7000	6.0000 U	14.7000	6.0000 U
I Cobalt	UG/L	12.0000 U	9.0000 U	12.0000 U	9.0000 U	12.0000 U	9.0000 U
I Copper	UG/L	25.0000 U	4.0000 U	53.7000	4.0000 U	25.0000 U	4.0000 U
I Cyanide	UG/L	10.0000 U		10.0000 U		10.0000 U	
I Iron	UG/L	3590.0000	165.0000	7470.0000	109.0000	1600.0000	30.4000 J
I Lead	UG/L	5.9000	2.0000 U	18.2000	2.0000 U	2.0000 U	2.0000 U
I Magnesium	UG/L	1510.0000 J	1210.0000 J	981.0000 J	505.0000 J	130000.000	167000.000
I Manganese	UG/L	32.9000	23.8000	29.2000	7.4000 J	30.7000	39.8000
I Mercury	UG/L	0.2100	0.2000 U	0.2000 U	0.2000 U	0.2000 U	0.2000 U
I Nickel	UG/L	38.0000 U	13.0000 U	38.6000 U	13.0000 U	38.0600 U	13.0000 U
I Potassium	UG/L	3950.0000 U	1440.0000 U	3950.0000 U	1440.0000 U	42100.0000	57200.0000
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	10.0000 U	5.0000 U	10.0000 U	5.0000 U	10.0000 U	5.0000 U
I Sodium	UG/L	1550.0000 J	1830.0000 J	2290.0000 J	2240.0000 J	1050000.00	1370000.00
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	15.0000 U	4.0000 U	15.0000 U	4.0000 U	15.0000 U	4.0000 U
I Zinc	UG/L	55.7000	5.0000 J	73.2000	21.0000	87.1000	5.0000 U
P 4,4'-DDD	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDE	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDT	UG/L	0.1000 U		0.1000 U		0.1000 U	
P Aldrin	UG/L	0.0500 U		0.0500 U		0.0500 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G06001	13G06001F	13607001	13607001F	13619001	13G19001F
	LAB ID:	534316	534329	533766	533788	533781	533794
	DATE RCVD:	01/30/93	01/30/93	01/27/93	01/27/93	01/27/93	01/27/93
NAME	UNITS						
P Aroclor 1016	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	
P Aroclor 1221	UG/L	2.0000 U	.	2.0000 U	.	2.0000 U	
P Aroclor 1232	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	
P Aroclor 1242	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	
P Aroclor 1248	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	
P Aroclor 1254	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	
P Aroclor 1260	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	
P Die1drin	UG/L	0.1000 U	.	0.1000 U	.	0.0027 J	
P Endosulfan I	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
P Endosulfan II	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	
P Endoaulfan sulfate	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	
P Endrin	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	
P Endrin aldehyde	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	
P Endrin ketone	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	
P Heptachlor	UG/L	0.0016 U	.	0.0016 U	.	0.0500 U	
P Heptachlor epoxide	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
P Methoxychlor	UG/L	0.5000 U	.	0.5000 U	.	0.5000 U	
P Toxaphene	UG/L	5.0000 U	.	5.0000 U	.	5.0000 U	
P alpha-BHC	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
P alpha-Chlordane	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
P beta-BHC	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
P delta-BHC	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
P gamma-BHC (Lindane)	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
P gamma-Chlordane	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S 1,2-Dichlorobenzene	UG/L	10,0000 U	.	10.0000 U	.	10.0000 U	
S 1,3-Dichlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S 1,4-Dichlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	

NAS I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	13G06001	13G06001F	13G07001	13G07001F	13G19001	13G19001F
	LAB ID:	534316	534329	533766	533788	533781	533794
	DATE RCVD:	01/30/93	01/30/93	01/27/93	01/27/93	01/27/93	01/27/93
S 2,4,5-Trichlorophenol	UG/L	25.0000 U		25.0000 U		25.0000 U	.
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,4-Dichlorophenol	UG/L	10.0000 U	.	10,0000 U	.	10.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 U	.	10.0000 U		10.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	.	25.0000 U		25.0000 U	
S 2,4-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 2-Nitrophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Chloroaniline	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S 4-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 4-Nitrophenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S Acenaphthene	UG/L	10.0000 U		10.0000 U	.	1.0000 J	.
S Acenaphthylene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Anthracene	UG/L	10.0000 U		10.0000 U	.	10.0000 U	
S Benzo(a)anthracene	UG/L	10.0000 U		10.0000 U	.	10.0000 U	
S Benzo(a)pyrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G06001	13G06001F	13G07001	13G07001F	13G19001	13G19001F
	LAB ID:	534316	534329	533766	533788	533781	533794
	DATE RCVD:	01/30/93	01/30/93	01/27/93	01/27/93	01/27/93	01/27/93
NAME	UNITS						
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Benzo(g,h,i)perylene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Butylbenzylphthalate	UG/L	10.0000 U	.	2.0000 U	.	7.0000 U	.
S Carbazole	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Chrysene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Di-n-butylphthalate	UG/L	10.0000 U	.	10.0000 U	.	1.0000 U	.
S Di-n-octylphthalate	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Dibenzofuran	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Diethylphthalate	UG/L	10.0000 U	.	10.0000 U	.	1.0000 U	.
S Dimethylphthalate	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Fluoranthene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Fluorene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Hexachlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Hexachlorobutadiene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Hexachloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Isophorone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S N-Nitrosodiphenylamine	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Naphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Nitrobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Pentachlorophenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S Phenanthrene	UG/L	10.0000 U	.	10.0000 U	.	'10.0000 U	.
S Phenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Pyrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.

NAS I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME		ENSAFE ID:	13G06001	13606001F	13G07001	13G07001F	13G19001	13G19001F
		LAB ID:	534316	534329	533766	533780	533781	533794
		DATE RCVD:	01/30/93	01/30/93	01/27/93	01/27/93	01/27/93	01/27/93
	WITS							
S bis(2-Ethlyhexy)phthalate	UG/L		2.0000 U	.	2.0000 U	.	4.0000 U	.
S bis(2-chloroethyl)ether	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1,1-Trichloroethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1,2,2-Tetrachloroethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1,2-Trichloroethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1-Dichloroethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1-Dichloroethene	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethene (total)	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloropropane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 2-Butanone	UG/L		10.0000 U	.	10.0000 U	.	7.0000 U	.
V 2-Hexanone	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V 4-Methyl-2-Pentanone	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Acetone	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Benzene	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Bromodichloromethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Bromoform	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Bromomethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Carbon Disulfide	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Carbon Tetrachloride	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Chlorobenzene	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Chloroethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Chloroform	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Chloromethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Dibromochloromethane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Ethylbenzene	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U	.
V Methylene Chloride	UG/L		10.0000 U	.	3.0000 U	.	8.0000 U	.

NAS I'ENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G06001	13G06001F	13G07001	13G07001F	13G19001	13G19001F
	LAB ID:	534316	534329	533766	533788	533781	533794
	DATE RCVD:	01/30/93	01/30/93	01/27/93	01/27/93	01/27/93	01/27/93
NAME	UNITS						
V Styrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Tetrachloroethene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Toluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Trichloroethene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Vinyl Chloride	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Xylene (total)	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
ENSAFE ID:		33G01001	33G01001F	33602001	33G02001F	33G03001	33603001F
LAB ID:		534214	534189	534219	534194	534204	534186
DATE RCVD:		01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
I Aluminum	UG/L	1630.0000	201.0000	2170.0000	265.0000	33600.0000	1170.0000
I Antimony	UG/L	57.0000 U	52.0000 U	57.0000 U	52.0000 U	57.0000 U	52.0000 U
I Arsenic	UG/L	5.6000 J	5.2000 J	4.1000 J	3.0000 U	13.2000	3.0000 U
I Barium	UG/L	23.7000 J	23.9000 J	26.2000 J	22.7000 J	82.4000 J	13.4000 J
I Beryllium	UG/L	3.0000 U	1.0000 U	3.0000 U	1.0000 U	3.0000 U	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	71400.0000	82200.0000	30700.0000	31900.0000	13500.0000	11000.0000
I Chromium	UG/L	31.6000	6.0000 U	10.7000	6.0000 U	75.7000	13.5000
I Cobalt	UG/L	12.0000 U	9.0000 U	12.0000 U	9.0000 U	12.0000 U	9.0000 U
I Copper	UG/L	25.0000 U	4.0000 U	25.0000 U	4.0000 U	30.9000	4.0000 U
I Cyanide	UG/L	10.0000 U		10.0000 U		10.0000 U	
I Iron	UG/L	3350.0000	1950.0000	5730.0000	3540.0000	26000.0000	2230.0000
I Lead	UG/L	2.3000 J	2.0000 U	11.4000	2.0000 U	14.3000	2.0000 U
I Magnesium	UG/L	4160.0000 J	4590.0000 J	1930.0000 J	1950.0000 J	8740.0000	4440.0000 J
I Manganese	UG/L	111.0000	121.0000	89.9000	88.3000	108.0000	48.2000
I Mercury	UG/L	0.2400	0.2000 U	0.$700	0.2000 U	1.6000	0.2000 U
I Nickel	UG/L	38.0000 U	13.0000 U	38.0000 U	13.0000 U	38.0000 U	13.0000 U
I Potassium	UG/L	3950.0000 U	3910.0000 J	3950.0000 U	1440.0000 U	3950.0000 U	1440.0000 U
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	15.0000 U	3.0000 U
I Silver	UG/L	10.0000 U	5.0000 U	10.0000 U	5.0000 U	10.0000 U	5.0000 U
I Sodium	UG/L	11600.0000	12500.0000	2780.0000 J	3100.0000 J	15000.0000	17100.0000
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	15.0000 U	4.0000 U	15.0000 U	4.0000 U	76.0000	6.2000 J
I Zinc	UG/L	19.6000 J	9.7000 J	40.0000	7.6000 J	97.5000	7.8000 J
P 4,4'-DDD	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDE	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4.4'-DDT	UG/L	0.1000 U		0.1000 U		0.1000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G01001	33G01001F	33G02001	33G02001F	33G03001	33G03001F
	LAB ID:	534228	534189	534233	534194	534225	534186
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
P Aldrin	UG/L	0.0500 U		0.0500 U	.	0.0500 U	.
P Aroclor 1016	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1221	UG/L	2.0000 U	.	2.0000 U	.	2.0000 U	.
P Aroclor 1232	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1242	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1248	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1254	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1260	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Dieldrin	UG/L	0.0053 J	.	0.1000 U	.	0.1000 U	.
P Endosulfan I	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P Endosulfan II	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endosulfan sulfate	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endrin	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endrin aldehyde	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endrin ketone	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Heptachlor	UG/L	0.0027 U	.	0.0030 U	.	0.0029 U	.
P Heptachlor epoxide	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P Methoxychlor	UG/L	*0.5000* U	.	0.5000 U	.	*0.5000* U	.
P Toxaphene	UG/L	*5.0000* U	.	5.0000 U	.	*5.0000* U	.
P alpha-BHC	UG/L	0.0500 U	.	0.0500 U	.	*0.0500* U	.
P alpha-Chlordane	UG/L	0.0500 U	.	*0.0500* U	.	*0.0500* U	.
P beta-BHC	UG/L	0.0500 U	.	0.0500 U	.	*0.0500* U	.
P delta-BHC	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P gama-BHC (Lindane)	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P gamma-Chlordane	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G01001	33G01001	33G02001	33G02001F	33G03001	33G03001
	LAB ID:	534228	534189	534233	534194	534223	534186
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
S 1,4-Dichlorobenzene	UG/L	2.0000 J	.	10.0000 U	.	10.0000 U	.
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,4-Dichlorophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 2-Nitrophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Chloroaniline	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 4-Nitrophenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S Acenaphthene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Acenaphthylene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Anthracene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S Benzo(a)anthracene	UG/L	10.0000 U	.	10.0000 U		10.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G01001	33G01001F	33002001	330020018	33G03001	330030011
	LAB ID:	534228	534189	534233	534194	534225	534186
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
S Benzo(a)pyrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Benzo(g,h,i)perylene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Butylbenzylphthalate	UG/L	1.0000 U	.	10.0000 U	.	10.0000 U	
S Carbazole	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Chrysene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Di-n-butylphthalate	UG/L	1.0000 U	.	10.0000 U	.	10.0000 U	
S Di-n-octylphthalate	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Dibenz(a,h)anthracene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Dibenzofuran	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Diethylphthalate	UG/L	1.0000 U	.	10.0000 U	.	10.0000 U	
S Dimethylphthalate	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Fluoranthene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Fluorene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Hexachlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Hexachlorobutadiene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Hexachlorocyclopentadiene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Hexachloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Ieophorone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S N-Nitroeo-di-n-propylamine	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S N-Nitrosodiphenylamine	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Naphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Nitrobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Pentachlorophenol	UG/L	25.0000 U	.	25.0000 U	.	'25.0000 U	
S Phenanthrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Phenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S Pyrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	

NAY I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33G01001 / 534220 / 01/29/93	33G01001F / 534109 / 01/29/93	33602001 / 534233 / 01/29/93	33602001F / 534194 / 01/29/93	33G03001 / 534225 / 01/29/93	33603001F / 534186 / 01/29/93
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	3.0000 U	.	4.0000 U		2.0000 U	.
S bis(2-chloroethyl)ether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1,1-Trichloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1,2-Trichloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1-Dichloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,1-Dichloroethene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 1,2-Dichloropropane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 2-Butanone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 2-Hexanone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Acetone	UG/L	10.0000 U	.	160.0000 U	.	10.0000 U	.
V Benzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Bromodichloromethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Bromoform	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Bromomethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Carbon Disulfide	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Carbon Tetrachloride	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Chlorobenzene	UG/L	18.0000	.	7.0000 J	.	10.0000 U	.
V Chloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Chloroform	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Chloromethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Dibromochloromethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Ethylbenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G01001	33G01001F	33G02001	33G02001F	33G03001	33603001F
	LAB ID:	534228	534189	534233	534194	534225	534186
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
V Methylene Chloride	UG/L	2.0000 U	.	1.0000 U	.	2.0000 U	.
V Styrene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Tetrachloroethene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Toluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Trichloroethene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Vinyl Chloride	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V Xylene (total)	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G04001	330040018	33G05001	336050018	33008001	33G080018
	LAB ID:	534221	534195	534222	534196	534317	534330
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/30/93	01/30/93
NAME	UNITS						
I Aluminum	UG/L	4380.0000	264.0000	1690.0000	249.0000	3640.0000	72.0000 J
I Antimony	UG/L	57.0000 U	52.0000 U	57.0000 U	52.0000 U	57.0000 U	30.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	7.1000 J	3.0000 U
I Barium	UG/L	35.9000 J	34.3000 J	12.3000 J	9.8000 J	38.2000 J	33.3000 J
I Beryllium	UG/L	3.0000 U	1.0000 U	3.0000 U	1.0000 U	3.0000 U	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	33000.0000	35000.0000	24300.0000	26300.0000	56800.0000	62100.0000
I Chromium	UG/L	11.6000	6.0000 U	10.0000 U	6.0000 U	10.0000 U	6.0000 U
I Cobalt	UG/L	12.0000 U	9.0000 U	12.0000 U	9.0000 U	12.0000 U	9.0000 U
I Copper	UG/L	25.0000 U	4.0000 U	25.0000 U	4.0000 U	25.0000 U	4.0000 U
I Cyanide	UG/L	10.0000 U		10.0000 U		10.0000 U	
I Iron	UG/L	4120.0000	964.0000	2910.0000	1080.0000	2390.0000	26.6000 J
I Lead	UG/L	5.8000	2.0000 U	4.9000	2.0000 U	2.0000 U	2.0000 U
I Magnesium	UG/L	3580.0000 J	3580.0000 J	7800.0000	9190.0000	47000.0000	52900.0000
I Manganese	UG/L	51.2000	36.8000	18.8000	23.6000	93.7000	91.0000
I Mercury	UG/L	0.2600	0.2000 U	0.2000 U	0.2000 U	0.3800	0.2000 U
I Nickel	UG/L	38.0000, U	13.0000 U	38.0000 U	13.0000 U	38.0000 U	13.0000 U
I Potassium	UG/L	5470.0000	4090.0000 J	3950.0000 U	2290.0000 J	20700.0000	22400.0000
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	30.0000 U
I Silver	UG/L	10.0000 U	5.0000 U	10.0000 U	5.0000 U	10.0000 U	5.0000 U
I Sodium	UG/L	37400.0000	44700.0000	17800.0000	20500.0000	510000.000	592000.000
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	15.0000 U	4.0000 U	15.0000 U	4.0000 U	19.7000 J	9.1000 J
I Zinc	UG/L	18.7000 J	5.7000 J	25.6000	5.0000 U	25.4000	5.0000 U
P 4,4'-DDD	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDE	UG/L	0.1000 U		0.0120 J		0.1000 U	
P 4,4'-DDT	UG/L	0.1000 U		0.1000 U		0.1000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33G04001	33G04001F	33G05001	33G05001F	33G08001	33G08001F
ENSAFE ID:		33G04001	33G04001F	33G05001	33G05001F	33G08001	33G08001F
LAB ID:		534235	534195	534236	534196	534317	534330
DATE RCVD:		01/29/93	01/29/93	01/29/93	01/29/93	01/30/93	01/30/93
P Aldrin	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P Aroclor 1016	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1221	UG/L	2.0000 U	.	2.0000 U	.	2.0000 U	.
P Aroclor 1232	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1242	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1249	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1254	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Aroclor 1260	UG/L	1.0000 U	.	1.0000 U	.	1.0000 U	.
P Dieldrin	UG/L	0.0037 J	.	0.1000 U	.	0.1000 U	.
P Endosulfan I	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P Endosulfan II	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endosulfan sulfate	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endrin	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endrin aldehyde	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Endrin ketone	UG/L	0.1000 U	.	0.1000 U	.	0.1000 U	.
P Heptachlor	UG/L	0.0015 U	.	0.0500 U	.	0.0500 U	.
P Heptachlor epoxide	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P Methoxychlor	UG/L	0.5000 U	.	0.5000 U	.	0.5000 U	.
P Toxaphene	UG/L	5.0000 U	.	5.0000 U	.	5.0000 U	.
P alpha-BHC	UG/L	0.0500 U	.	0.0500 U	.	0.0052 J	.
P alpha-Chlordane	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P beta-BHC	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P delta-BHC	UG/L	0.0500 U	.	0.0015 J	.	0.0500 U	.
P gamma-BHC (Lindane)	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
P gamma-Chlordane	UG/L	0.0500 U	.	0.0500 U	.	0.0500 U	.
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	.	7.0000 J	.	32.0000	.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	.	30.0000	.	1.0000 J	.

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key preeented at end of Appendix. Data not shown in significant digita.

	ENSAFE ID:	33G04001	33G04001F	33G05001	330050018	33G08001	330080018
	LAB ID:	534235	534195	534236	534196	534317	534330
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/30/93	01/30/93
NAME	UNITS						
S 1,4-Dichlorobenzene	UG/L	10.0000 U	.	40.0000	.	4.0000 J	
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	.	25.0000 U		25.0000 U	
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	.	10.0000 U		10.0000 U	
S 2,4-Dichlorophenol	UG/L	10.0000 U	.	10.0000 U		10.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	
S 2-Nitrophenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S 4-Chloroaniline	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S 4-Methylphenol	UG/L	10.0000 U		10.0000 U	.	10.0000 U	
S 4-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	.	25.0000 U	
S 4-Nitrophenol	UG/L	25.0000 U		25.0000 U	.	25.0000 U	.
S Acenaphthene	UG/L	10.0000 U		19.0000	.	10.0000 U	
S Acenaphthylene	UG/L	10.0000 U	.	10.0000 U		10.0000 U	
S Anthracene	UG/L	10.0000 U		10.0000 U		10.0000 U	.
S Benzo(a)anthracene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	

NAS I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key pteeented at end of Appendix. Data not shown in significant digits.

NAME		ENSAFE ID:	33G04001	33G04001F	33G05001	330050013	33G08001	33G08001F
		LAB ID:	534235	534195	534236	534196	534317	534330
		DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/30/93	01/30/93
	WITS							
S Benzo(a)pyrene	UG/L	10.0000 U			10.0000 U		10.0000 U	.
S Benzo(b,k)fluoranthene	UG/L	10.0000 U			10.0000 U	.	10.0000 U	.
S Benzo(g,h,i)perylene	UG/L	10.0000 U			10.0000 U	.	10.0000 U	.
S Butylbenzylphthalate	UG/L	12.0000 U			16.0000 U		10.0000 U	.
S Carbazole	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Chrysene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Di-n-butylphthalate	UG/L	10.0000 U			10.0000 U	.	10.0000 U	.
S Di-n-octylphthalate	UG/L	10.0000 U			10.0000 U	.	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L	10.0000 U			10.0000 U	.	10.0000 U	
S Dibenzofuran	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Diethylphthalate	UG/L	1.0000 U		.	2.0000 U	.	10.0000 U	.
S Dimethylphthalate	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Fluoranthene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Fluorene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Hexachlorobenzene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Hexachlorobutadiene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Hexachloroethane	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Isophorone	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S N-Nitrosodiphenylamine	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Naphthalene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Nitrobenzene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Pentachlorophenol	UG/L	25.0000 U		.	25.0000 U	.	25.0000 U	.
S Phenanthrene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Phenol	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.
S Pyrene	UG/L	10.0000 U		.	10.0000 U	.	10.0000 U	.

NAS I'ENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33004001	33G04001F	33G05001	330050018	33G08001	330080018
	LAB ID:	534235	534195	534236	534196	534317	534330
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/30/93	01/30/93
NAME	UNITS						
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S bis(2-Ethlyhexy)phthalate	UG/L	3.0000 U	.	4.0000 U	.	10.0000 U	
S bis(2-chloroethyl)ether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 1,1,1-Trichloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 1,1,2-Trichloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 1,1-Dichloroethane	UG/L	10.0000 U	.	10.0000 U	.	46.0000	
V 1,1-Dichloroethene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 1,2-Dichloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	.	10.0000 U	.	5.0000 J	
V 1,2-Dichloropropane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 2-Butanone	UG/L	10.0000 U	.	10.0000 U	.	370.0000 U	
V 2-Hexanone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Acetone	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Benzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Bromodichloromethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Bromoform	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Bromomethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Carbon Disulfide	UG/L	10.0000 U	.	10.0000 U	.	5.0000 J	
V Carbon Tetrachloride	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Chlorobenzene	UG/L	10.0000 U	.	45.0000	.	28.0000	
V Chloroethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Chloroform	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Chloromethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Dibromochloromethane	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	
V Ethylbenzene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33604001	33604001F	33605001	336050013	33608001	33G08001F
	LAB ID:	534235	534195	534236	534196	534317	534330
	DATE RCVD:	01/29/93	01/29/93	01/29/93	01/29/93	01/30/93	01/30/93
NAME	UNITS						
V Methylene Chloride	UG/L	10.0000 U		10.0000 U		10.0000 U	.
V Styrene	UG/L	10.0000 U	.	10.0000 U	.	2.0000 J	.
v Tetrachloroethene	UG/L	10.0000 U		10.0000 U		10.0000 U	.
V Toluene	UG/L	10.0000 U		1.0000 J		10.0000 U	.
V Trichloroethene	UG/L	10.0000 U	.	10.0000 U	.	3.0000 J	
V Vinyl Chloride	UG/L	10.0000 U		10.0000 U		10.0000 U	.
V Xylene (total)	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	.	10.0000 U	.	10.0000 U	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G08001RE	33G09001	33609001F	33610001	33G10001F	33G11001
	LAB ID:	534317	534313	534328	534600	534622	534318
	DATE RCVD:	01/30/93	01/30/93	01/30/93	02/03/93	02/03/93	01/30/93
NAME	UNITS						
I Aluminum	UG/L	14600.0000	350.0000	5320.0000	5320.0000	711.0000	
I Antimony	UG/L	57.0000 U	52.0000 U	29.0000 U	*30.0000* U	57.0000 U	
I Arsenic	UG/L	7.6000 J	3.0000 U	5.5900 J	5.9000 J	3.0000 U	
I Barium	UG/L	49.6000 J	24.4000 J	35.7000 J	36.0000 J	75.0000 J	
I Beryllium	UG/L	3.0000 U	1.0000 U	3.0000 U	1.0000 U	3.0000 U	
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	
I Calcium	UG/L	15700.0000	15300.0000	17600.0000	19400.0000	47200.0000	
I Chromium	UG/L	27.2000	6.2000 J	10.0000 U	6.0000 U	10.0000 U	
I Cobalt	UG/L	12.0000 U	9.0000 U	12.0000 U	9.0000 U	12.0000 U	
I Copper	UG/L	25.0000 U	4.0000 U	25.0000 U	4.0000 U	25.0000 U	
I Cyanide	UG/L	10.0000 U		10.0000 U		10.0000 U	
I Iron	UG/L	8440.0000	2380.0000	4990.0000	3700.0000	16200.0000	
I Lead	UG/L	6.8000	2.0000 U	4.1000	2.2000 J	2.0000 U	
I Magnesium	UG/L	6040.0000	5320.0000	9900.0000	10600.0000	36600.0000	
I Manganese	UG/L	111.0000	104.0000	95.5000	99.5000	259.0000	
I Mercury	UG/L	*0.3960*	0.2000 U	0.6600	0.2000 U	1.0000	
I Nickel	UG/L	38.0000 U	13.0000 U	38.0000 U	13.0000 U	*38.0000* U	
I Potaseium	UG/L	5050.0000	3640.0000 J	5520.0000	6050.0000	19400.0000	
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	*3.0000* U	
I Silver	UG/L	10.0000 U	5.0000 U	10.0000 U	5.0000 U	10.0000 U	
I Sodium	UG/L	95800.0000	110000.000	107000.000	122000.000	321000.000	
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U	
I Vanadium	UG/L	67.2000	29.4000 J	15.9000 J	14.5000 J	15.0000 U	
I Zinc	UG/L	26.8000	8.0000 J	46.7000	10.8000 J	23.5000	
P 4,4'-DDD	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDE	UG/L	0.1000 U		0.1000 U		0.1000 U	
P 4,4'-DDT	UG/L	0.1000 U		0.1000 U		0.0029 J	

NAS I'ENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G08001RE	33G09001	33G09001F	33G10001	33G10001F	33G11001
	LAB ID:	534317	534313	534328	534600	534622	534318
	DATE RCVD:	01/30/93	01/30/93	01/30/93	02/03/93	02/03/93	01/30/93
NAME	UNITS						
P Aldrin	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P Aroclor 1016	UG/L	.	1.0000 U	.	1.0000 U	.	1.0000 U
P Aroclor 1221	UG/L	.	2.0000 U	.	2.0000 U	.	2.0000 U
P Aroclor 1232	UG/L	.	1.0000 U	.	1.0000 U	.	1.0000 U
P Aroclor 1242	UG/L	.	1.0000 U	.	1.0000 U	.	1.0000 U
P Aroclor 1248	UG/L	.	1.0000 U	.	1.0000 U	.	1.0000 U
P Aroclor 1254	UG/L	.	1.0000 U	.	1.0000 U	.	1.0000 U
P Aroclor 1260	UG/L	.	1.0000 U	.	1.0000 U	.	1.0000 U
P Dieldrin	UG/L	.	0.1000 U	.	0.1000 U	.	0.1000 U
P Endosulfan I	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P Endosulfan II	UG/L	.	0.1000 U	.	0.1000 U	.	0.1000 U
P Endosulfan sulfate	UG/L	.	0.1000 U	.	0.1000 U	.	0.1000 U
P Endrin	UG/L	.	0.1000 U	.	0.1000 U	.	0.1000 U
P Endrin aldehyde	UG/L	.	0.1000 U	.	0.1000 U	.	0.1000 U
P Endrin ketone	UG/L	.	0.1000 U	.	0.1000 U	.	0.1000 U
P Heptachlor	UG/L	.	0.0021 U	.	0.0090 U	.	0.0500 U
P Heptachlor epoxide	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P Methoxychlor	UG/L	.	0.5000 U	.	0.5000 U	.	0.5000 U
P Toxaphene	UG/L	.	5.0000 U	.	5.0000 U	.	5.0000 U
P alpha-BHC	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P alpha-Chlordane	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P beta-BHC	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P delta-BHC	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P gamma-BHC (Lindane)	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
P gamma-Chlordane	UG/L	.	0.0500 U	.	0.0500 U	.	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 1,2-Dichlorobenzene	UG/L	.	10.0000 U	.	10.0000 U	.	3.0000 J
S 1,3-Dichlorobenzene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U

NAS YENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G08001RE	33G09001	33G09001F	33610001	33G10001F	33G11001
	LAB ID:	534317	534313	534328	534600	534622	534318
	DATE RCVD:	01/30/93	01/30/93	01/30/93	02/03/93	02/03/93	01/30/93
NAME	UNITS						
S 1,4-Dichlorobenzene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	.	25.0000 U	.	25.0000 U	.	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2,4-Dichlorophenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2,4-Dimethylphenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2,4-Dinitrophenol	UG/L	.	25.0000 U	.	25.0000 U	.	25.0000 U
S 2,4-Dinitrotoluene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2,6-Dinitrotoluene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2-Chloronaphthalene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2-Chlorophenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2-Methylnaphthalene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2-Methylphenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 2-Nitroaniline	UG/L	.	25.0000 U	.	25.0000 U	.	25.0000 U
S 2-Nitrophenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 3-Nitroaniline	UG/L	.	25.0000 U	.	25.0000 U	.	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	.	25.0000 U	.	25.0000 U	.	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 4-Chloroaniline	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 4-Methylphenol	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S 4-Nitroaniline	UG/L	.	25.0000 U	.	25.0000 U	.	25.0000 U
S 4-Nitrophenol	UG/L	.	25.0000 U	.	25.0000 U	.	25.0000 U
S Acenaphthene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S Acenaphthylene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S Anthracene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U
S Benzo(a)anthracene	UG/L	.	10.0000 U	.	10.0000 U	.	10.0000 U

NAS I'ENSACOLA OU-10 *AND* SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G08001RE	33G09001	33G09001F	33G10001	33G10001F	33G11001
	LAB ID:	534317	534313	534328	534600	534622	534318
	DATE RCVD:	01/30/93	01/30/93	01/30/93	02/03/93	02/03/93	01/30/93
NAME	UNITS						
S Benzo(a)pyrene	UG/L		10.0000 U	.	10.0000 U		10.0000 U
S Benzo(b,k)fluoranthene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Benzo(g,h,i)perylene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Butylbenzylphthalate	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Carbazole	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Chrysene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Di-n-butylphthalate	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Di-n-octylphthalate	UG/L		10.0000 U	.	10.0000 U		10.0000 u
S Dibenz(a,h)anthracene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Dibenzofuran	UG/L		10.0000 U	.	10.0000 U		10.0000 U
S Diethylphthalate	UG/L		10.0000 U	.	10.0000 U		10.0000 U
S Dimethylphthalate	UG/L		10.0000 U		10.0000 U		10.0000 U
S Fluoranthene	UG/L	.	10.0000 U	.	10.0000 u		10.0000 u
S Fluorene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Hexachlorobenzene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Hexachlorobutadiene	UG/L		10.0000 U	.	10.0000 U		10.0000 U
S Hexachlorocyclopentadiene	UG/L		10.0000 U		10.0000 U		10.0000 U
S Hexachloroethane	UG/L		10.0000 U		10.0000 U		10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Isophorone	UG/L		10.0000 U		10.0000 U		10.0000 U
S N-Nitroso-di-n-propylamine	UG/L		10.0000 U		10.0000 U		10.0000 U
S N-Nitrosodiphenylamine	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Naphthalene	UG/L		10.0000 U		10.0000 U		10.0000 U
S Nitrobenzene	UG/L		10.0000 U		10.0000 U		10.0000 U
S Pentachlorophenol	UG/L		25.0000 U		25.0000 U		25.0000 U
S Phenanthrene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U
S Phenol	UG/L		10.0000 U		10.0000 U		10.0000 U
S Pyrene	UG/L	.	10.0000 U	.	10.0000 U		10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G08001RE	33009001	33G09001F	33010001	33G10001F	33011001
	LAB ID:	534317	534313	534328	534600	534622	534318
	DATE RCVD:	01/30/93	01/30/93	01/30/93	02/03/93	02/03/93	01/30/93
NAME	UNITS						
S bis(2-Chloroethoxy)methane	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L		10.0000 U		3.0000 U	.	3.0000 U
S bis(2-chloroethyl)ether	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L		10.0000 U	.	10.0000 U	.	10.0000 U
V 1,1,1-Trichloroethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V 1,1,2-Trichloroethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V 1,1-Dichloroethane	UG/L	57.0000 J	10.0000 U	.	10.0000 U	.	2.0000 J
V 1,1-Dichloroethene	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V 1,2-Dichloroethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U		10.0000 U
V 1,2-Dichloroethene (total)	UG/L	5.0000 J	10.0000 U	.	10.0000 U		10.0000 U
V 1,2-Dichloropropane	UG/L	25.0000 U	10.0000 U	.	10.0000 U		10.0000 U
V 2-Butanone	UG/L	270.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V 2-Hexanone	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Acetone	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Benzene	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Bromodichloromethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Bromoform	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Bromomethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Carbon Disulfide	UG/L	6.0000 J	10.0000 U	.	10.0000 U	.	10.0000 U
V Carbon Tetrachloride	UG/L	25.0000 U	10.0000 U		10.0000 U	.	10.0000 U
V Chlorobenzene	UG/L	29.0000 J	10.0000 U	.	10.0000 U	.	29.0000
V Chloroethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Chloroform	UG/L	25.0000 U	10.0000 U	.	10.0000 U		10.0000 U
V Chloromethane	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Dibromochloromethane	UG/L	25.0000 U	10.0000 U		10.0000 U		10.0000 U
V Ethylbenzene	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
GROUNDWATER SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G08001RE	33G09001	33G09001F	33G10001	33G10001F	33G11001
	LAB ID:	534317	534313	534328	534600	534622	534318
	DATE RCVD:	01/30/93	01/30/93	01/30/93	02/03/93	02/03/93	01/30/93
NAME	UNITS						
V Methylene Chloride	UG/L	25.0000 U	1.0000 U	.	2.0000 U	.	10.0000 U
V Styrene	UG/L	25.0000 U	2.0000 J	.	10.0000 U	.	3.0000 J
V Tetrachloroethene	UG/L	3.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Toluene	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Trichloroethene	UG/L	3.0000 J	10.0000 U	.	10.0000 U	.	10.0000 U
V Vinyl Chloride	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V Xylene (total)	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V cis-1,3-Dichloropropene	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U
V trans-1,3-Dichloropropene	UG/L	25.0000 U	10.0000 U	.	10.0000 U	.	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

APPENDIX R

SOIL *AND* GROUNDWATER ANALYTICAL RESULTS — MARCH 1994
UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84219 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S06

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0905 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS	UG/KG	ANALYTICAL RESULTS
35u	CHLOROMETHANE	35u	CIS-1.3-DICHLOROPROPENE
35u	VINYL CHLORIDE	87U	METHYL ISOBUTYL KETONE
35u	BROMOMETHANE	35u	TOLUENE
35U	CHLOROETHANE	35u	TRANS-1.3-DICHLOROPROPENE
35u	TRICHLOROFLUOROMETHANE	35U	1.1.2-TRICHLOROETHANE
35u	1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)	35u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
350U	ACETONE	35u	1.3-DICHLOROPROPANE
87U	CARBON DISULFIDE	87U	METHYL BUTYL KETONE
35u	METHYLENE CHLORIDE	35u	DIBROMOCHLOROMETHANE
35u	TRANS-1.2-DICHLOROETHENE	35u	CHLOROBENZENE
35u	1.1-DICHLOROETHANE	35u	1.1.1.2-TETRACHLOROETHANE
35u	CIS-1,2-DICHLOROETHENE	35u	ETHYL BENZENE
35u	2.2-DICHLOROPROPANE	35u	(M- AND/OR P-)XYLENE
350U	METHYL ETHYL KETONE	35u	O-XYLENE
35U	BROMOCHLOROMETHANE	35U	STYRENE
35U	CHLOROFORM	35u	BROMOFORM
35U	1.1.1-TRICHLOROETHANE	35u	BROMOBENZENE
35U	1.1-DICHLOROPROPENE	35u	1.1.2.2-TETRACHLOROETHANE
35u	CARBON TETRACHLORIDE	35u	1.2.3-TRICHLOROPROPANE
35U	1.2-DICHLOROETHANE	35u	O-CHLOROTOLUENE
35U	BENZENE	35u	P-CHLOROTOLUENE
35U	TRICHLOROETHENE(TRICHLOROETHYLENE)	35u	1.3-DICHLOROBENZENE
35U	1.2-DICHLOROPROPANE	35U	1.4-DICHLOROBENZENE
35U	DIBROMOMETHANE	35U	1.2-DICHLOROBENZENE
35U	BROMODICHLOROMETHANE	4.6	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS... MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS. GA.

03/30/94

ENTRACTABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84219 SAMPLE TYPE: SOIL
SWRCE: NAS PENSACOLA
STATION ID: 32S06

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0905 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1400U	(3-AND/OR 4-)METHYLPHENOL
1400U	1,2,4-TRICHLOROBENZENE
1400U	2,2'-CHLOROISOPROPYLETHER
1400U	2,3,4,6-TETRACHLOROPHENOL
1400U	2.4.5-TRICHLOROPHENOL
1400U	2.4.6-TRJCHLOROPHEWL
1400U	2,4-DICHLOROPHENOL
1400U	2.4-DIMETHYLPHENOL
2800U	2.4-DINITROPHENOL
1400U	2,4-DINITROTOLUENE
1400U	2.6-DINITROTOLUENE
1400U	2-CHLORONAPHTHALENE
1400U	2-CHLOROPHENOL
2800U	2-METHYL-4.6-DINITROPHENOL
1400U	2-METHVLNAPHTHALENE
1400U	2-METHYLPHENOL
1400U	2-NITROANILINE
1400U	2-NITROPHENOL
1400U	3,3'-DICHLOROBENZIDINE
1400U	3-NITROANILINE
1400U	4-BROMOPHENYL PHENYL ETHER
1400U	4-CHLORO-3-METHYLPHENOL
1400U	4-CHLOROANILINE
1400U	4-CHLOROPHENYL PHENYL ETHER
1400U	4-NITROANIL JN€
2800U	4-NITROPHENOL
1400U	ACENAPHTHENE
1400U	ACENAPHTHYLENE
1400U	ANTHRACENE
1400U	BENZO(A)ANTHRACENE
1400U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1400U	BENZO(GHI)PERYLENE
1400U	BENZO-A-PYRENE
1400U	BENZYL BUTYL PHTHALATE
1400U	BIS(2-CHLOROETHOXY) METHANE
1400U	BIS(2-CHLOROETHYL) ETHER
1400U	BIS(2-ETHYLHEXYL) PHTHALATE
1400U	CARBAZOLE
1400U	CHRYSENE
1400U	DI-N-BUTYLPHTHALATE
1400U	DI-N-OCTYLPHTHALATE
1400U	DIBENZO(A,H)ANTHRACENE
1400U	DIBENZOFURAN
1400U	DIETHYL PHTHALATE
1400U	DIMETHYL PHTHALATE
1400U	FLUORANTHENE
1400U	FLUORENE
1400U	HEXACHLOROBENZENE (HCB)
1400U	HEXACHLOROBUTADIENE
1400U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1400U	HEXACHLOROETHANE
1400U	INDENO (1,2,3-CD) PYRENE
1400U	ISOPHORONE
1400U	N-NITROSODI-N-PROPYLAMINE
1400U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1400U	NAPHTHALENE
1400U	NITROBENZENE
2800U	PENTACHLOROPHENOL
1400U	PHENANTHRENE
1400U	PHENOL
1400U	PYRENE
4.6	PERCENT MOISTURE

REMARKS

••*REMARKS* ••

FOOTNOTES

*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

04/05/94

*PESTICIDES/PCB'S DATA REPORT

PROJECT NO, 94-0323 SAMPLE NO. 84219 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S06

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0905 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN 1 (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4.4'-OD0 (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200u	CHLORDANE (TECH. MIXTURE) /1
500u	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	FCO-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXVCHLOR
50U	ENDRIN KETONE
4.5	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES

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*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN *CONFIRMED BY GC/MS
*U-MATERIAL WAS ANALYZED FOR BUT MOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT
1. WHEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84219 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32506

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PCNSACOLA ST: FL
COLLECTION START: 03/08/94 0905 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.21U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS. GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84219 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32S06 COLLECTION START: 03/08/94 0905 STOP: 00/00/00

ANALYTICAL RESULTS

MG/KG	
1 OU	SILVER
3.OU	ARSENIC
NA	BORON
9.6	BARIUM
0.50U	BERYLLIUM
1.4	CADMIUM
1.0U	COBALT
50	CHROMIUM
5.2	COOPER
1.0U	MOLYBDENUM
2.OU	NICKFL
28	LEAD
3.OU	ANTIMONY
4.OU	SELENIUM
4.OU	TIN
34	STRONTIUM
5.OU	TELLURIUM
39	TITANIUM
10U	THALLIUM
3.2	VANADIUM
1.OU	YTTRIUM
23	ZINC
NA	21RCONIUM
0.05U	MERCURY
2300	ALUMINUM
38	MANGANESE

ANALYTICAL RESULTS

MG/KG	
10000	CALCIUM
150	MAGNESIUM
1500	IRON
100	SODIUM
200U	POTASSIUM
4	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALY, MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84220 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32524

PROC ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0925 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
42U	CHLOROMETHANE
42U	VINYL CHLORIDE
42U	BROMOMETHANE
42U	CHLOROETHANE
42U	TRICHLOROFLUOROMETHANE
42U	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
420U	ACETONE
100U	CARBON DISULFIDE
42U	METHYLENE CHLORIDE
42U	TRANS-1,2-DICHLOROETHENE
42U	1,1-DICHLOROETHANE
42U	CIS-1,2-DICHLOROETHENE
42U	2,2-DICHLOROPROPANE
420U	METHYL ETHYL KETONE
42U	BROMOCHLOROMETHANE
42U	CHLOROFORM
42U	1,1,1-TRICHLOROETHANE
42U	1,1-DICHLOROPROPENE
42U	CARBON TETRACHLORIDE
42U	1,2-DICHLOROETHANE
42U	BENZENE
42U	TRICHLOROETHENE(TRICHLOROETHYLENE)
42U	1,2-DICHLOROPROPANE
42U	DIBROMOMETHANE
42U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
42u	CIS-1,3-DICHLOROPROPENE
100U	METHYL ISOBUTYL KETONE
42U	TOLUENE
42U	TRANS-1,3-DICHLOROPROPENE
42U	1,1,2-TRICHLOROETHANE
42U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
42U	1,3-DICHLOROPROPANE
100U	METHYL BUTYL KETONE
42U	DIBROMOCHLOROMETHANE
42U	CHLOROBENZENE
42U	1,1,1,2-TETRACHLOROETHANE
42U	ETHYL BENZENE
42U	(M- AND/OR P-)XYLENE
5.8J	O-XYLENE
42U	STYRENE
42U	BROMOFORM
42U	BROMOBENZENE
42U	1,1,2,2-TETRACHLOROETHANE
42U	1,2,3-TRICHLOROPROPANE
42U	O-CHLOROTOLUENE
42U	P-CHLOROTOLUENE
42U	1,3-DICHLOROBENZENE
42U	1,4-DICHLOROBENZENE
42U	1,2-DICHLOROBENZENE
7.6	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYS. MANAGEMENT SYSTEM
€PA-REGION I V ESD. ATHENS. GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84220 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32524

PROG LLEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLCCTION START: 03/06/94 0925 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1400U	(3-AND/OR 4-)METHYLPHENOL
1400U	1,2,4-TRICHLOROBENZENE
1400U	2,2'-CHLOROISOPROPYLETHER
1400U	2,3,4,6-TETRACHLOROPHENOL
1400U	2,4,5-TRICHLOROPHENOL
1400U	2,4,6-TRICHLOROPHENOL
1400U	2,4-DICHLOROPHENOL
1400U	2,4-DIMETHYLPHENOL
2900U	2,4-DINITROPHENOL
1400U	2,4-DINITROTOLUENE
1400U	2,6-DINITROTOLUENE
1400U	2-CHLORONAPHTHALENE
1400U	2-CHLOROPHENOL
2900U	2-METHYL-4,6-DINITROPHENOL
1400U	2-METHYLNAPHTHALENE
1400U	2-METHYLPHENOL
1400U	2-NITROANILINE
1400U	2-NITROPHENOL
1400U	3,3'-DICHLOROBENZIDINE
1400U	3-NITROANILINE
1400U	4-BROMOPHENYL PHENYL ETHER
1400U	4-CHLORO-3-METHYLPHENOL
1400U	4-CHLOROANILINE
1400U	4-CHLOROPHENYL PHENYL ETHER
1400U	4-NITROANILINE
2900U	4-NITROPHENOL
1400U	ACENAPHTHENE
1400U	ACENAPHTHYLENE
1400U	ANTHRACENE
1400U	BENZO(A)ANTHRACENE
1400U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1400U	BENZO(GHI)PERYLENE
1400U	BENZO-A-PYRENE
1400U	BENZYL BUTYL PHTHALATE
1400U	BIS(2-CHLOROETHOXY) METHANE
1400U	BIS(2-CHLOROETHYL) ETHER
1400U	BIS(2-ETHYLHEXYL) PHTHALATE
1400U	CARBAZOLE
1400U	CHRYSENE
1400U	DI-N-BUTYLPHTHALATE
1400U	DI-N-OCTYLPHTHALATE
1400U	DIBENZO(A,H)ANTHRACENE
1400U	DIBENZOFURAN
1400U	DIETHYL PHTHALATE
1400U	DIMETHYL PHTHALATE
1400U	FLUORANTHENE
1400U	FLUORENE
1400U	HEXACHLOROBENZENE (HCB)
1400U	HEXACHLOROBUTADIENE
1400U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1400U	HEXACHLOROETHANE
1400U	INDENO (1,2,3-CD) PYRENE
1400U	ISOPHORONE
1400U	N-NITROSODI-N-PROPYLAMINE
1400U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1400U	NAPHTHALENE
1400U	NITROBENZENE
2900U	PENTACHLOROPHENOL
1400U	PHENANTHRENE
1400U	PHENOL
1400U	PYRENE
7.0	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANAL), MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, CA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84220 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S24

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0925 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCD-1232 (AROCLOR 1232)
500U	PCB-1218 (AROCLOR 1218)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
7.8	PERCENT MOISTURE

@+@REMARKS***

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO OE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR OUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT C-CONFIRMED BY GC/MS
1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84220 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S24

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0925 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.29	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84220 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION IO: 32524

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0925 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
3.5	SILVER
3.0U	ARSENIC
NA	BORON
37	BARIUM
0.50U	BERYLLIUM
12	CADMIUM
1.1	COBALT
580	CHROMIUM
27	COPPER
1.0U	MOLYBDENUM
46	NICKEL
39	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
7.0U	TIN
47	STRONTIUM
5.0U	TELLURIUM
47	TITANIUM
10U	THALLIUM
4.9	VANADIUM
1 6	YTTRIUM
60	ZINC
NA	ZIRCONIUM
0.12	MERCURY
4500	ALUMINUM
82	MANGANESE

MG/KG	ANALYTICAL RESULTS
12000	CALCIUM
620	MAGNESIUM
1900	IRON
110	SODIUM
200U	POTASSIUM
8	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84221 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S53

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0940 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
40U	CHLOROMCTHANE
40U	VINYL CHLORIDE
40U	BROMOMETHANE
40U	CHLOROETHANE
40U	TRICHLOROFLUOROMETHANE
40U	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
400U	ACETONE
100U	CARBON DISULFIDE
40U	METHYLENE CHLORIDE
40U	TRANS-1,2-DICHLOROETHENE
40U	1,1-DICHLOROETHANE
40U	CIS-1,2-DICHLOROETHENE
40U	2,2-DICHLOROPROPANE
4wu	METHYL ETHYL KETONE
40U	BROMOCHLOROMETHANE
40U	CHLOROFORM
40U	1,1,1-TRICHLOROETHANE
40U	1,1-DICHLOROPROPENE
40U	CARBON TETRACHLORIDE
40U	1,2-DICHLOROETHANE
40U	BENZENE
40U	TRICHLOROETHENE(TRICHLOROETHYLENE
40U	1,2-DICHLOROPROPANE
40U	DIBROMOMETHANE
40U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
40U	CIS-1,3-DICHLOROPROPENE
100U	METHYL ISOBUTYL KETONE
40U	TOLUENE
40U	TRANS-1,3-DICHLOROPROPENE
40U	1,1,2-TRICHLOROETHANE
40U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
40U	1.3-DICHLOROPROPANE
100U	METHYL BUTYL KETONE
40U	DIBROMOCHLOROMETHANE
40U	CHLOROBENZENE
40U	1.1.1.2-TETRACHLOROETHANE
40U	ETHYL BENZENE
40U	(M- AND/OR P-)XYLENE
40U	O-XYLENE
40U	STYRENE
40U	BROMOFORM
40U	BROMOBENZENE
40U	1.1.2.2-TETRACHLOROETHANE
40U	1.2,3-TRICHLOROPROPANE
40U	O-CHLOROTOLUENE
40U	P-CHLOROTOLUENE
40U	1.3-DICHLOROBENZENE
40U	1.4-DICHLOROBENZENE
40U	1.2-DICHLOROBENZENE
4 8	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS. MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, EAT

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84221 SAMPLE TYPE; SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32553

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0940 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1300U	(3-AND/OR 4-)METHYLPHENOL
1300U	1,2,4-TRICHLOROBENZEPIE
1300U	2,2'-CHLOROISOPROPYLETHER
1300U	2,3,4,6-TETRACHLOROPHENOL
1300U	2,4,5-TRICHLOROPHENOL
1300U	2,4,6-TRICHLOROPHENOL
1300U	2,4-DICHLOROPHENOL
1300U	2,4-DIMETHYLPHENOL
2700U	2.4-DINITROPHENOL
1300U	2,4-DINITROTOLUENE
1300U	2,6-DINITROTOLUENE
1300U	2-CHLORONAPHTHALENE
1300U	2-CHLOROPHENOL
2700U	2-METHYL-4,6-DINITROPHENOL
1300U	2-METHYLNAPHTHALENE
1300U	2-METHYLPHENOL
1300U	2-NITROANILINE
1300U	2-NITROPHENOL
1300U	3,3'-DICHLOROBENZIDINE
1300U	3-NITROANILINE
1300U	4-BROMOPHENYL PHENYL ETHER
1300U	4-CHLORO-3-METHYLPHENOL
1300U	4-CHLOROANILINE
1300U	4-CHLOROPHENYL PHENYL ETHER
1300U	4-NITROANILINE
2700U	4-NITROPHENOL
1300U	ACENAPHTHENE
1300U	ACENAPHTHYLENE
1300U	ANTHRACENE
1300U	BENZO(A)ANTHRACENE
1300U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1300U	BENZO(GHI)PERYLENE
1300U	BENZO-A-PYRENE
1300U	BENZYL BUTYL PHTHALATE
1300U	BIS(2-CHLOROETHOXY) METHANE
1300U	BIS(2-CHLOROETHYL) ETHER
1300U	BIS(2-ETHYLHEXYL) PHTHALATE
1300U	CARBAZOLE
1300U	CHRYSENE
1300U	DI-N-BUTYLPHTHALATE
1300U	DI-N-OCTYLPHTHALATE
1300U	DIBENZO(A,H)ANTHRACENE
1300U	DIBENZOFURAN
1300U	DIETHYL PHTHALATE
1300U	DIMETHYL PHTHALATE
1300U	FLUORANTHENE
1300U	FLUORENE
1300U	HEXACHLOROBENZENE (HCB)
1300U	HEXACHLOROBUTADIENE
1300U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1300U	HEXACHLOROETHANE
1300U	INDENO (1,2,3-CD) PYRENE
1300U	ISOPHORONE
1300U	N-NITROSODI-N-PROPYLAMINE
1300U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1300U	NAPHTHALENE
1300U	NITROBENZENE
2700U	PENTACHLOROPHENOL
1300U	PHENANTHRENE
1300U	PHENOL
1300U	PYRENE
4.8	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALY_ _ MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84221 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S53

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0940 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4.4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCO-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
4.8	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYS.. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84221 SAMPLE TYPE: SOIL
SOURCC: NAS PENSACOLA
STATION ID: 32S53

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0940 STOP: 00/00/00

RESULTS UNITS PARAMCTCR
0.21U MG/KG CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS ...NAGEMENT SYSTEM
€PA-REGION I V ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84221 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S53

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0940 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1.4	SILVER
3.0U	ARSENIC
NA	BORON
5.9	BARIUM
0.50U	BERYLLIUM
2.0	CADMIUM
1.0U	COBALT
33	CHROMIUM
4.0	COPPER
1.0U	MOLYBDENUM
2.0U	NICKEL
14	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
5.5U	TIN
1.2	STRONTIUM
5.0U	TELLURIUM
23	TITANIUM
10U	THALLIUM
1.0U	VANADIUM
1.0U	YTTRIUM
7.5	ZINC
NA	ZIRCONIUM
0.05U	MERCURY
390	ALUMINUM
3.3	MANGANESE

MG/KG	ANALYTICAL RESULTS
130	CALCIUM
21	MAGNCSIUM
360	IRON
100U	SODIUM
200U	POTASSIUM
5	PERCENT MOISTURE

REMARKS

*****REMARKS*****

*****FOOTNOTES*****

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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84222 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32554

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0955 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
41U	CHLOROMETHANE
41U	VINYL CHLORIDE
41U	BROMOMETHANE
41U	CHLOROETHANE
41U	TRICHLOROFLUOROMETHANE
41U	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
410U	ACETONE
100U	CARBON DISULFIDE
41U	METHYLENE CHLORIDE
41U	TRANS-1,2-DICHLOROETHENE
41U	1,1-DICHLOROETHANE
41U	CIS-1,2-DICHLOROETHENE
41U	2,2-DICHLOROPROPANE
4103	METHYL ETHYL KETONE
41U	BROMOCHLOROMETHANE
41U	CHLOROFORM
41U	1,1,1-TRICHLOROETHANE
41U	1,1-DICHLOROPROPENE
41U	CARBON TETRACHLORIDE
41U	1,2-DICHLOROETHANE
41U	BENZENE
41U	TRICHLOROETHENE(TRICHLOROETHYLENE)
41u	1,2-DICHLOROPROPANE
41U	DIBROMOMETHANE
41U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
41U	CIS-1,3-DICHLOROFROPCNC
100U	METHYL ISOBUTYL KETONE
41U	TOLUENE
41U	TRANS-1,3-DICHLOROPROPENE
41U	1,1,2-TRICHLOROETHANE
41U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
41U	1,3-DICHLOROPROPANE
100U	METHYL BUTYL KETONE
41U	DIBROMOCHLOROMETHANE
41U	CHLOROBENZENE
41U	1,1,1,2-TETRACHLOROETHANE
41U	ETHYL BENZENE
41U	(M- AND/OR P-)XYLENE
41U	O-XYLENE
41U	STYRENE
41U	BROMOFORM
41U	BROMOBENZENE
41U	1,1,2,2-TETRACHLOROETHANE
41U	1,2,3-TRICHLOROPROPANE
41U	O-CHLOROTOLUENE
41U	P-CHLOROTOLUENE
41U	1,3-DICHLOROBENZENE
41U	1,4-DICHLOROBENZENE
41U	1,2-DICHLOROBENZENE
5.3	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYS... MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 **SAMPLE NO.** 84222 SAMPLE TYPE: **SOIL**
SOURCE: NAS PENSACOLA
STATION ID: 32S54

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0955 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1300U	(3-AND/OR 4-)METHYLPHENOL
1300U	1,2,4-TRICHLOROBENZENE
1300U	2,2'-CHLOROISOPROPYLETHER
1300U	2,3,4,6-TETRACHLOROPHENOL
1300U	2,4,5-TRICHLOROPHENOL
1300U	2,4,6-TRICHLOROPHENOL
1300U	2,4-DICHLOROPHENOL
1300U	2,4-DIMETHYLPHENOL
2700U	2,4-DINITROPHENOL
1300U	2,4-DINITROTOLUENE
1300U	2,6-DINITROTOLUENE
1300U	2-CHLORONAPHTHALENE
1300U	2-CHLOROPHENOL
2700U	2-METHYL-4,6-DINITROPHENOL
1300U	2-METHYLNAPHTHALENE
1300U	2-METHYLPHENOL
1300U	2-NITROANILINE
1300U	2-NITROPHENOL
1300U	3,3'-DICHLOROBENZIDINE
1300U	3-NITROANILINE
1300U	4-BROMOPHENYL PHENYL ETHER
1300U	4-CHLORO-3-METHYLPHENOL
1300U	4-CHLOROANILINE
1300U	4-CHLOROPHENYL PHENYL ETHER
1300U	4-NITROANILINE
2700U	4-NITROPHENOL
1300U	ACENAPHTHENE
1300U	ACENAPHTHYLENE
1300U	ANTHRACENE
1300U	BENZO(A)ANTHRACENE
1300U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1300U	BENZO(GHI)PERYLENE
1300U	BENZO-A-PYRENE
1300U	BENZYL BUTYL PHTHALATE
1300U	BIS(2-CHLOROETHOXY) METHANE
1300U	BIS(2-CHLOROETHYL) ETHER
1300U	BIS(2-ETHYLHEXYL) PHTHALATE
1300U	CARBAZOLE
1300U	CHRYSENE
1300U	DI-N-BUTYLPHTHALATE
1300U	DI-N-OCTYLPHTHALATE
1300U	DIBENZO(A,H)ANTHRACENE
1300U	DIBENZOFURAN
1300U	DIETHYL PHTHALATE
1300U	DIMETHYL PHTHALATE
1300U	FLUORANTHENE
1300U	FLUORENE
1300U	HEXACHLOROBENZENE (HCB)
1300U	HEXACHLOROBUTADIENE
1300U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1300U	HEXACHLOROETHANE
1300U	INDENO (1,2,3-CD) PYRENE
1300U	ISOPHORONE
1300U	N-NITROSODI-N-PROPYLAMINE
1300U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1300U	NAPHTHALENE
1300U	NITROBENZENE
2700U	PENTACHLOROPHENOL
1300U	PHENANTHRENE
1300U	PHENOL
1300U	PYRENE
5.3	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84222 SAMPLE TYPE:, SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S54

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0955 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCO-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
5.3	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES

*A-AVERAGE VALUE. *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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1. WHEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYS.. MANAGEMENT SYSTEM
€PA-REGION IV ESD, ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84222 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S54

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0955 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.21U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYS. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84222 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S54

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 0955 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1 OU	SILVER
3.OU	ARSENIC
NA	BORON
2.3	BARIUM
0.50U	BERYLLIUM
0.50U	CADMIUM
1.OU	COBALT
5.3	CHROMIUM
1.OU	COPPER
1.OU	MOLYBDENUM
2.OU	NICKEL
4.OU	LEAD
3.OU	ANTIMONY
4.OU	SELENIUM
4.OU	TIN
1.OU	STRONTIUM
5.OU	TELLURIUM
27	TITANIUM
10U	THALLIUM
1.OU	VANADIUM
1.OU	YTTRIUM
1.1	ZINC
NA	ZIRCONIUM
0.05U	MERCURY
280	ALUMINUM
1.0	MANGANESE

MG/KG	ANALYTICAL RESULTS
50U	CALCIUM
10U	MAGNESIUM
93	IRON
100U	SODIUM
200U	POTASSIUM
5	PERCENT MOISTURE

@**REMARKS***

REMARKS

FOOTNOTES

@A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANAL...S MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS. GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84223 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S55

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PE SACOLA ST: FL
COLLECT N START: 03/08/94 1020 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
37U	CHLOROMETHANE
37u	VINYL CHLORIDE
37u	BROMOMETHANE
37u	CHLOROETHANE
37u	TRICHLOROFLUOROMETHANE
37u	1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)
370U	ACETONE
92U	CARBON DISULFIDE
37u	METHYLENE CHLORIDE
37u	TRANS-1.2-DICHLOROETHENE
37u	1.1-DICHLOROETHANE
37u	CIS-1.2-DICHLOROETHENE
37u	2.2-DICHLOROPROPANE
370U	METHYL ETHYL KETONE
37u	BROMOCHLOROMETHANE
37u	CHLOROFORM
37u	1.1.1-TRICHLOROETHANE
37u	1.1-DICHLOROPROPENE
37u	CARBON TETRACHLORIDE
37u	1.2-DICHLOROETHANE
37U	BENZENE
37u	TRICHLOROETHENE(TRICHLOROETHYLENE)
37u	1.2-DICHLOROPROPANE
37u	DIBROMOMETHANE
37u	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
37u	CIS-1.3-DICHLOROPROPCNE
92U	METHYL ISOBUTYL KETONE
37u	TOLUENE
37u	TRANS-1.3-DICHLOROPROPENE
37u	1.1.2-TRICHLOROETHANE
37u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
37u	1.3-DICHLOROPROPANE
92U	METHYL BUTYL KETONE
37u	DIBROMOCHLOROMETHANE
37u	CHLOROBENZENE
37u	1.1.1.2-TETRACHLOROETHANE
37u	ETHYL BENZENE
37u	(M- AND/OR P-)XYLENE
370	O-XYLENE
37u	
37u	BROMOFORM
37u	BROMOBENZENE
37U	1.1.2.2-TETRACHLOROETHANE
37U	1.2.3-TRICHLOROPROPANE
37u	O-CHLOROTOLUENE
37u	P-CHLOROTOLUENE
37U	1.3-DICHLOROBENZENE
37u	1.4-DICHLOROBENZENE
37u	1.2-DICHLOROBENZENE
3.0	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEM
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS,, MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84223 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S55

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1020 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1300U	(3-AND/OR 4-)METHYLPHENOL
1300U	1,2,4-TRICHLOROBENZENE
1300U	2,2'-CHLOROISOPROPYLETHER
1300U	2,3,4,6-TETRACHLOROPHENOL
1300U	2,4,5-TRICHLOROPHENOL
1300U	2,4,6-TRICHLOROPHENOL
1300U	2,4-DICHLOROPHENOL
1300U	2,4-DIMETHYLPHENOL
2600U	2,4-DINITROPHENOL
1300U	2,4-DINITROTOLUENE
1300U	2,6-DINITROTOLUENE
1300U	2-CHLORONAPHTHALENE
1300U	2-CHLOROPHENOL
2600U	2-METHYL-4,6-DINITROPHENOL
1300U	2-METHYLNAPHTHALENE
1300U	2-METHYLPHENOL
1300U	2-NITROANILINE
1300U	2-NITROPHENOL
1300U	3,3'-DICHLOROBENZIDINE
1300U	3-NITROANILINE
1300U	4-BROMOPHENYL PHENYL ETHER
1300U	4-CHLORO-3-METHYLPHENOL
1300U	4-CHLOROANILINE
1300U	4-CHLOROPHENYL PHENYL ETHER
1300U	4-NITROANILINE
2600U	4-NITROPHENOL
1300U	ACENAPHTHENE
1300U	ACENAPHTHVLENE
1300U	ANTHRACENE
1300U	BENZO(A)ANTHRACENE
1300U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1300U	BENZO(GHI)PERYLENE
1300U	BENZO-A-PYRENE
1300U	BENZYL BUTYL PHTHALATE
1300U	BIS(2-CHLOROETHOXY) METHANE
1300U	BIS(2-CHLOROETHYL) ETHER
1300U	BIS(2-ETHYLHEXYL) PHTHALATE
1300U	CARBAZOLE
1300U	CHRYSENE
1300U	DI-N-BUTYLPHTHALATE
1300U	DI-N-OCTYLPHTHALATE
1300U	DIBENZO(A,H)ANTHRACENE
1300U	DIBENZOFURAN
1300U	DIETHYL PHTHALATE
1300U	DIMETHYL PHTHALATE
1300U	FLUORANTHENE
1300U	FLUORENE
1300U	HEXACHLOROBENZENE (HCB)
1300U	HEXACHLOROBUTADIENE
1300U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1300U	HEXACHLOROETHANE
1300U	INDENO (1,2,3-CD) PYRENE
1300U	ISOPHORONE
1300U	N-NITROSODI-N-PROPYLAMINE
1300U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1300U	NAPHTHALENE
1300U	NITROBENZENE
2600U	PENTACHLOROPHENOL
1300U	PHENANTHRENE
1300U	PHENOL
1300U	PYRENE
3.0	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84223 SAMPLE TYPE: **SOIL**
SOURCE: NAS PENSACOLA
STATION ID: 32S55

PROC ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1020 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
5DU	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	**4.4'-DDT** (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	**4,4'-DDD** (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN 11 (BETA)
SOU	ENDOSULFAN SULFATE
2wu	CHLORDANE (TECH. MIXTURE) **/1**
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCD-1232 (AROCLOR 1232)
500U	PCB-1218 (AROCLOR 1218)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	**CHLORDEUE** *12*
--	**ALPHA-CHLORDENE /2**
--	**BETA CHLORDENE /2**
--	**GAMMA-CHLORDENE /2**
--	**GAMMA-CHLORDANE /2**
--	**TRANS-NONACHLOR /2**
--	**ALPHA-CHLORDANE /2**
--	**CIS-NONACHLOR /2**
--	**OXYCHLORDANE (OC1ACHLOREPOXIDE) /2**
200U	**METHOXYCHLOR**
50U	**ENDRIN KETONE**
3.0	**PERCENT MOISTURE**

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE **•NA-NOT** ANALYZED **•NAI-INTERFERENCES** •J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE **IS KNOWN TO BE** LESS THAN VALUE GIVEN **•L-ACTUAL** VALUE IS KNOWN **TO** BE GREATER THAN VALUE GIVEN C-CONFIRMED BY GC/MS
•U-MATERIAL WAS ANALYZED **FOR BUT** NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT
1. **WHEN NO** VALUE **IS** REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES **OF** TECHNICAL CHLORDANE.

SAMPLE AND ANALYSI MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84223 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCC: NAS PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: 32S55 COLLECTION START: 03/08/94 1020 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.21U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84223 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S55

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1020 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1.0U	SILVER
3.0U	ARSENIC
NA	BORON
1.0U	BARIUM
0.50U	BERYLLIUM
0.50U	CADMIUM
1.0U	COBALT
3.8	CHROMIUM
1.1	COPPER
1.0U	MOLYBDENUM
2.0U	NICKEL
4.0U	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
4.5U	TIN
1.0U	STRONTIUM
5.0U	TELLURIUM
36	TITANIUM
10U	THALLIUM
1.0U	VANADIUM
1.0U	YTTRIUM
2.9	ZINC
NA	ZIRCONIUM
0.05U	MERCURY
210	ALUMINUM
1.8	MANGANESE

MG/KG	ANALYTICAL RESULTS
80	CALCIUM
11	MAGNESIUM
82	IRON
100U	SODIUM
200U	POTASSIUM
3	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALY_ _ MANAGEMENT SYSTEM
EPA-REGION-IV ESD, ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84224 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32532

PROC ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1040 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
42U	CHLOROMCTHANC
42U	VINYL CHLORIDE
42U	BROMOMETHANE
42U	CHLOROETHANE
42U	TRICHLOROFLUOROMETHANE
42U	1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)
420U	ACETONE
110U	CARBON DISULFIDE
42U	METHYLENE CHLORIDE
42U	TRANS-1.2-DICHLOROETHENE
42U	1.1-DICHLOROETHANE
42U	CIS-1,2-DICHLOROETHENE
42U	2.2-DICHLOROPROPANE
420U	METHYL ETHYL KETONE
42U	BROMOCHLOROMETHANE
42U	CHLOROFORM
42U	1,1,1-TRICHLOROETHANE
42U	1.1-DICHLOROPROPENE
42U	CARBON TETRACHLORIDE
13J	1.2-OICHLOROETHANE
42U	BENZENE
42U	TRICHLOROETHENE(TRICHLOROETHYLENE)
42U	1.2-DICHLOROPROPANE
42U	DIBROMOMETHANE
42U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
42U	CIS-1.3-DICHLOROPROPCNC
110U	METHYL ISOBUTYL KETONE
7.0J	TOLUENE
42U	TRANS-1.3-DICHLOROPROPENE
42U	1.1.2-TRICHLOROETHANE
42U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
42U	1.3-DICHLOROPROPANE
110U	METHYL BUTYL KETONE
42U	DIBROMOCHLOROMETHANE
42U	CHLOROBENZENE
42U	1.1.1.2-TETRACHLOROETHANE
42U	ETHYL BENZENE
7.4J	(M- AND/OR P-)XYLENE
42U	O-XYLENE
42U	STYRENE
42U	BROMOFORM
42U	BROMOBENZENE
42U	1.1.2.2-TETRACHLOROETHANE
42U	1.2.3-TRICHLOROPROPANE
42U	O-CHLOROTOLUENE
42U	P-CHLOROTOLUENE
42U	1.3-DICHLOROBENZENE
42U	1.4-DICHLOROBENZENE
42U	1.2-DICHLOROBENZENE
15.4	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR OUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EP -REGION IV ESD, ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/9

PROJECT NO. 94-0323 SAMPLE NO. 84224 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S32

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITV: PENSACOLA ST: FL
COLLCCTION START: 03/08/94 1040 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1EOOU	(3-AND/OR 4-)METHYLPHENOL
1500U	1,2,4-TRICHLOROBENZENE
1500U	2,2'-CHLOROISOPROPYLETHER
1500U	2,3,4,6-TETRACHLOROPHENOL
1500U	2,4,5-TRICHLOROPHENOL
1500U	2,4,6-TRICHLOROPHENOL
1500U	2,4-DICHLOROPHENOL
1500U	2,4-DIMETHYLPHENOL
3100U	2,4-DINITROPHENOL
1500U	2,4-DINITROTOLUENE
1500U	2,6-DINITROTOLUENE
1500U	2-CHLORONAPHTHALENE
1500U	2-CHLOROPHENOL
3100U	2-METHYL-4,6-DINITROPHENOL
1500U	2-METHYLNAPHTHALENE
1500U	2-METHYLPHENOL
1500U	2-NITROANILINE
1500U	2-NITROPHENOL
1500U	3,3'-DICHLOROBENZIDINE
1500U	3-NITROANILINE
1500U	4-BROMOPHENYL PHENYL ETHER
1500U	4-CHLORO-3-METHYLPHENOL
1500U	4-CHLOROANILINE
1500U	4-CHLOROPHENYL PHENYL ETHER
1500U	4-NITROANILINE
3100U	4-NITROPHENOL
1500U	ACENAPHTHENE
1500U	ACENAPHTHYLENE
1500U	ANTHRACEN
1500U	BENZO(A)ANTHRACENE
1500U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1500U	BENZO(GHI)PERYLENE
1500U	BENZO-A-PYRENE
1500U	BENZYL BUTYL PHTHALATE
1500U	BIS(2-CHLOROETHOXY) METHANE
1500U	BIS(2-CHLOROETHYL) ETHER
1500U	BIS(2-ETHYLHEXYL) PHTHALATE
1500U	CARBAZOLE
1500U	CHRYSENE
1500U	DI-N-BUTYLPHTHALATE
1500U	DI-N-OCTYLPHTHALATE
1500U	DIBENZO(A,H)ANTHRACENE
1500U	DIBENZOFURAN
1500U	DIETHYL PHTHALATE
1500U	DIMETHYL PHTHALATE
1500U	FLUORANTHENE
1500U	FLUORENE
1500U	HEXACHLOROBENZENE (HCB)
1500U	HEXACHLOROBUTADIENE
1500U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1500U	HEXACHLOROETHANE
1500U	INDENO (1,2,3-CD) PYRENE
1500U	ISOPHORONE
1500U	N-NITROSODI-N-PROPYLAMINE
1500U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1500U	NAPHTHALENE
1500U	NITROBENZENE
3100U	PENTACHLOROPHENOL
1500U	PHENANTHRENE
1500U	PHENOL
1500U	PYRENE
15.4	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANAL. ,S MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84224 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S32

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1040 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
SOU	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
SOU	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
5011	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
SOU	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
SOU	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
--	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCO-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
560	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
10U	CHLORDENE /2
19U	BETA-CHLORDENE /2
19U	
19U	GAMMA-CHLORDENE /2
10	GAMMA-CHLORDANE /2
7.0	TRANS-NONACHLOR /2
19U	ALPHA-CHLORDANE /2
19U	CIS-NONACHLOR /2
19U	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
15	PERCENT MOISTURE

REMARKS

REMARKS

+FOOTNOTES*
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT C-CONFIRMED BV GC/MS
1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84224 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32532 COLLECTION START: 03/08/94 1040 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.24U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *3-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS .AGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PKOJECT NO. 94-0323 SAMPLE NO. 84224 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32532

PROC ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1040 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
12	SILVER
3.0U	ARSENIC
NA	BORON
53	BARIUM
0.50U	BERYLLIUM
21	CADMIUM
1.3	COBALT
680	CHROMIUM
57	COPPER
1.0U	MOLYBDENUM
46	NICKEL
57	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
12U	TIN
16	STRONTIUM
5.0U	TELLURIUM
54	TITANIUM
10U	THALLIUM
7.4	VANADIUM
1.9	YTTRIUM
120	ZINC
NA	ZIRCONIUM
0.56	MERCURY
7000	ALUMINUY
62	MANGANESE

MG/KG	ANALYTICAL RESULTS
3600	CALCIUM
590	MAGNESIUM
2800	IRON
100U	SODIUM
200U	POTASSIUM
16	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84225 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32551

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITV: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1050 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
55U	CHLOROMETHANE
55u	VINYL CHLORIDE
55u	BROMOMETHANE
55u	CHLOROETHANE
55u	TRICHLOROFLUOROMETHANE
55u	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE I
550U	ACETONE
140U	CARBON DISULFIDE
55u	METHYLENE CHLORIDE
55u	TRANS-1,2-DICHLOROETHENE
55u	1,1-DICHLOROETHANE
55u	CIS-1,2-DICHLOROETHENE
55u	2,2-DICHLOROPROPANE
550U	METHYL ETHYL KETONE
55u	BROMOCHLOROMETHANE
55u	CHLOROFORM
55u	1,1,1-TRICHLOROETHANE
55u	1.1-DICHLOROPROPENE
55U	CARBON TETRACHLORIDE
9.8J	1,2-DICHLOROETHANE
55u	BENZENE
55u	TRICHLOROETHENE(TRICHLOROETHYLENE)
55u	1.2-DICHLOROPROPANE
55u	DIBROMOMETHANE
55U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
55U	CIS-1,3-DICHLOROPROPENE
140U	METHYL ISOBUTYL KETONE
55U	TOLUENE
55u	TRANS-1,3-DICHLOROPROPENE
55u	1,1,2-TRICHLOROETHANE
55u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
55u	1,3-DICHLOROPROPANE
140U	METHYL BUTYL KETONE
55u	DIBROMOCHLOROMETHANE
55U	CHLOROBENZENE
55u	1,1,1,2-TETRACHLOROETHANE
55u	ETHYL BENZENE
55u	(M- AND/OR P-)XYLENE
55u	O-XYLENE
55u	STYREUE
55u	BROMOFORM
55u	BROMOBENZENE
55u	1.1.2.2-TETRACHLOROETHANE
55u	1,2,3-TRICHLOROPROPANE
55U	O-CHLOROTOLUENE
55u	P-CHLOROTOLUENE
55u	1.3-DICHLOROBENZENE
55u	1.4-DICHLOROBENZENE
55u	1.2-DICHLOROBENZENE
17.0	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS.. MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84225 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32501

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: YENSACOLA ST: FL
COLLECTION START: 03/08/94 1050 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1500U	(3-AND/OR 4-)METHYLPHENOL
1500U	1,2,4-TRICHLOROBENZENE
1500U	2,2'-CHLOROISOPROPYLETHER
1500U	2,3,4,6-TETRACHLOROPHENOL
1500U	2,4,5-TRICHLOROPHENOL
1500U	2,4,6-TRICHLOROPHENOL
1500U	2,4-DICHLOROPHENOL
1500U	2,4-DIMETHYLPHENOL
3100U	2,4-DINITROPHENOL
1500U	2,4-DINITROTOLUENE
1500U	2,6-DINITROTOLUENE
1500U	2-CHLORONAPHTHALENE
1500U	2-CHLOROPHENOL
3100U	2-METHYL-4,6-DINITROPHENOL
1500U	2-METHYLNAPHTHALENE
1500U	2-METHYLPHENOL
1500U	2-NITROANILINE
1500U	2-NITROPHENOL
1500U	3,3'-DICHLOROBENZIDINE
1500U	3-NITROANILINE
1500U	4-BROMOPHENYL PHENYL ETHER
1500U	4-CHLORO-3-METHYLPHENOL
1500U	4-CHLOROANILINE
1500U	4-CHLOROPHENYL PHENYL ETHER
1500U	4-NITROANILINE
3100U	4-NITROPHENOL
1500U	ACENAPHTHENE
1500U	ACENAPHTHYLENE
1500U	ANTHRACENE
1500U	BENZO(A)ANTHRACENE
1500U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1500U	BENZO(GHI)PERYLENE
1500U	BENZO-A-PYRENE
1500U	BENZYL BUTYL PHTHALATE
1500U	BIS(2-CHLOROETHOXY) METHANE
1500U	BIS(2-CHLOROETHYL) ETHER
570J	BIS(2-ETHYLHEXYL) PHTHALATE
1500U	CARBAZOLE
1500U	CHRYSENE
1500U	DI-N-BUTYLPHTHALATE
1500U	DI-N-OCTYLPHTHALATE
1500U	DIBENZO(A,H)ANTHRACENE
1500U	DIBENZOFURAN
1500U	DIETHYL PHTHALATE
1500U	DIMETHYL PHTHALATE
1500U	FLUORANTHENE
1500U	FLUORENE
1500U	HEXACHLOROBENZENE (HCB)
1500U	HEXACHLOROBUTADIENE
1500U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1500U	HEXACHLOROETHANE
1500U	INDENO (1,2,3-CD) PYRENE
1500U	ISOPHORONE
1500U	N-NITROSODI-N-PROPYLAMINE
1500U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1500U	NAPHTHALENE
1500U	NITROBENZENE
3100U	PENTACHLOROPHENOL
1500U	PHENANTHRENE
1500U	PHENOL
1500U	PYRENE
17.0	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE Of MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

04/05/94

PESTICIDES/PCB'S DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84225 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S51

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1050 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4.4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
--	CHLORDANE (TECH. MIXTURE) /1
200u	PCB-1242 (AROCLOR 1242)
200u	PCB-1254 (AROCLOR 1254)
200u	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
200u	PCO-1232 (AROCLOR 1232)
200U	PCB-1248 (AROCLOR 1248)
250	PCB-1260 (AROCLOR 1260)
200u	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
10U	CHLORDENE /2
19U	ALPHA-CHLORDENE /2
19U	BETA CHLORDENE /2
19U	GAMMA-CHLORDENE /2
7.8	GAMMA-CHLORDANE /2
5:3	TRANS-NONACHLOR /2
19U	ALPHA-CHLORDANE /2
19U	CIS-NONACHLOR /2
19U	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
17	PERCENT MOISTURE

***REMARKS@**

REMARKS

FOOTNOTES

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1. WHEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANAL , , MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84225 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCC: NAS PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: 32551 COLLECTION START: 03/08/94 1050 STOP: 00/00/00

RCSULTS	UNITS	PARAMCTCR
0.21U	MG/KG	CYANIDE

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84225 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32S51 COLLECTION START: 03/08/94 1050 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
14	SILVER
3.0U	ARSENIC
NA	BORON
58	BARIUM
0.50U	BERYLLIUM
22	CADMIUM
1.4	COBALT
780	CHROMIUM
66	COPPER
1.0	MOLYBDENUM
54	NICKEL
66	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
13U	TIN
19	STRONTIUM
5.0U	TELLURIUM
42	TITANIUM
10U	THALLIUM
6.7	VANADIUM
1.8	YTTRIUM
140	ZINC
NA	ZIRCONIUM
0.50	MERCURY
6600	ALUMINUM
60	MANGANESE

MG/KG	ANALYTICAL RESULTS
4200	CALCIUM
440	MAGNESIUM
2600	IRON
100U	SODIUM
200U	POTASSIUM
17	PERCENT MOISTURE

REMARKS ***REMARKS***

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84226 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S01

PROG ELEM: SSF COLLECTED BY F SLOAN
CITY: PENSACOLA 5 : FL
COLLECTION START: 03/08/94 1 00 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
53U	CHLOROMETHANE
53U	VINYL CHLORIDE
53u	BROMOMETHANE
53u	CHLOROETHANE
53u	TRICHLOROFLUOROMETHANE
53u	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
530U	ACETONE
130U	CARBON DISULFIDE
53u	METHYLENE CHLORIDE
53u	TRANS-1,2-DICHLOROETHENE
53u	1,1-DICHLOROETHANE
53u	CIS-1,2-DICHLOROETHENE
53u	2,2-DICHLOROPROPANE
530U	METHYL ETHYL KETONE
53u	BROMOCHLOROMETHANE
53u	CHLOROFORM
53u	1,1,1-TRICHLOROETHANE
53u	1,1-DICHLOROPROPENE
53u	CARBON TETRACHLORIDE
20J	1,2-DICHLOROETHANE
53u	BENZENE
53U	TRICHLOROETHENE(TRICHLOROETHYLENE)
53U	1,2-DICHLOROPROPANE
53U	DIBROMOMETHANE
53U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
53u	CIS-1,3-DICHLOROPROPENE
130U	METHYL ISOBUTYL KETONE
53u	TOLUENE
53u	TRANS-1,3-DICHLOROPROPENE
53u	1,1,2-TRICHLOROETHANE
53u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
53u	1,3-DICHLOROPROPANE
130U	METHYL BUTYL KETONE
53u	DIBROMOCHLOROMETHANE
53u	CHLOROBENZENE
53u	1,1,1,2-TETRACHLOROETHANE
53U	ETHYL BENZENE
53U	(M- AND/OR P-)XYLENE
53u	O-XYLENE
53u	STYRENE
53u	BROMOFORM
53u	BROMOBENZENE
53u	1,1,2,2-TETRACHLOROETHANE
53u	1,2,3-TRICHLOROPROPANE
53u	O-CHLOROTOLUENE
53u	P-CHLOROTOLUENE
53u	1,3-DICHLOROBENZENE
53u	1,4-DICHLOROBENZENE
53u	1,2-DICHLOROBENZENE
22.0	PERCENT MOISTURE

***FOOTNOTES—*
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SAMPLE AND ANALYSI_ ,ANAGEMENT SYSTEM
€PA-REGION IV ESD. ATHENS. GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84226 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S01

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLLCTION START: 03/08/94 1100 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1700U	(3-AND/OR 4-)METHYLPHENOL
17wu	1,2,4-TRICHLOROBENZENE
1700U	2.2'-CHLOROISOPROPYLETHER
1700U	2.3.4.6-TETRACHLOROPHENOL
1700U	2,4,5-TRICHLOROPHENOL
1700U	2,4,6-TRICHLOROPHENOL
17wu	2,4-DICHLOROPHENOL
1700U	2.4-DIMETHYLPHENOL
3300U	2.4-DINITROPHENOL
1700U	2.4-DINITROTOLUENE
1700U	2,6-DINITROTOLUENE
1700U	2-CHLORONAPHTHALENE
1700U	2-CHLOROPHENOL
3300U	2-METHYL-4,6-DINITROPHENOL
1700U	2-METHVLNAPHTHALENE
1700U	2-METHYLPHENOL
1700U	2-NITROANILINE
1700U	2-NITROPHENOL
1700U	3,3'-DICHLOROBENZIDINE
1700U	3-NITROANILINE
1700U	4-BROMOPHENYL PHENYL ETHER
1700U	4-CHLORO-3-METHYLPHENOL
1700U	4-CHLOROANILINE
1700U	4-CHLOROPHENYL PHENYL ETHER
1700U	4-NITROANILINE
3300U	4-NITROPHENOL
1700U	ACENAPHTHENE
1700U	ACENAPHTHYLENE
1700U	ANTHRACENE
1700U	BENZO(A)ANTHRACENE
1700U	BENZO(B AND/OR KIFLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1700U	DENZO(GHI)PERYLENE
1700U	DENLO-A-PYRENE
1700U	BENZYL BUTYL PHTHALATE
1700U	BIS(2-CHLOROETHOXY) METHANE
1700U	BIS(2-CHLOROETHYL) ETHER
1700U	BIS(2-ETHYLHEXYL) PHTHALATE
1700U	CARBAZOLE
1700U	CHRVSENE
1700U	DI-N-BUTYLPHTHALATE
1700U	DI-N-OCTYLPHTHALATE
1700U	DIBENZO(A,H)ANTHRACENE
1700U	DIBENZOFURAN
1700U	DIETHYL PHTHALATE
1700U	DIMETHYL PHTHALATE
1700U	FLUORANTHENE
1700U	FLUORENE
1700U	HEXACHLOROBENZENE (HCB)
1700U	HEXACHLOROBUTADIENE
1700U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1700U	HEXACHLOROETHANE
1700U	INDENO (1,2,3-CD) PYRENE
1700U	ISOPHORONE
1700U	N-NITROSODI-N-PROPYLAMINE
1700U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1700U	NAPHTHALENE
1700U	NITROBENZENE
3300U	PENTACHLOROPHENOL
1700U	PHENANTHRENE
1700U	PHENOL
1700U	PYRENE
22.0	PERCENT WISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84226 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S01

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1100 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAWA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50u	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200u	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCB-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
160J	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200u	METHOXVCHLOR
50U	ENDRIN KETONE
22	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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1. WHEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84226 SAMPLE TYPE: SOIL
SOURCE: NAS PCNSACOLA
STATION ID: 32501

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PCNSACOLA ST: FL
COLLECTION START: 03/08/94 1100 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.36	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYS.. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84226 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32501

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1100 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
8.2	SILVER
3.0U	ARSENIC
MA	BORON
25	BARIUM
0.50U	BERYLLIUM
23	CADMIUM
1.4	COBALT
910	CHROMIUM
47	COPPER
1.2	MOLYBDENUM
70	NICKEL
46	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
13U	TIN
14	STRONTIUM
5.0U	TELLURIUM
32	TITANIUM
10U	THALLIUM
5.7	VANADIUM
1.3	YTTRIUM
120	ZINC
NA	ZIRCONIUM
0.20	MERCURV
5700	ALUMINUM
52	MANGANESE

MG/KG	ANALYTICAL RESULTS
2200	CALCIUM
1GO	MAGNESIUM
1800	IRON
100U	SODIUM
200U	POTASSIUM
16	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANAL, .S MANAGEMENT SYSTEM
€PA-REGION I V ESD. ATHENS, GA. .

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84227 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S36

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1115 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS	UG/KG	ANALYTICAL RESULTS
49u	CHLOROMCTHANC	49u	CIS-1.3-DICHLOROPROPENE
49u	VINYL CHLORIDE	120U	METHYL ISOBUTYL KETONE
49u	BROMOMETHANE	7.1J	TOLUENE
49u	CHLOROETHANE	49u	TRANS-1.3-DICHLOROPROPENE
49u	TRICHLOROFLUOROMETHANE	49u	1.1.2-TRICHLOROETHANE
49u	1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)	49u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
490U	ACETONE	49u	1.3-DICHLOROPROPANE
120U	CARBON DISULFIDE	120u	METHYL BUTYL KETONE
49u	METHYLENE CHLORIDE	49u	DIBROMOCHLOROMETHANE
49u	TRANS-1.2-DICHLOROETHENE	49u	CHLOROBENZENE
49u	1.1-DICHLOROETHANE	49u	1.1.1.2-TETRACHLOROETHANE
49u	CIS-1.2-DICHLOROETHENE	49u	ETHYL BENZENE
49u	2.2-DICHLOROPROPANE	17J	(M- AND/OR P-)XYLENE
490U	METHYL ETHYL KETONE	49u	O-XYLENE
49u	BROMOCHLOROMETHANE	49u	STVRENE
49u	CHLOROFORM	49u	BROMOFORM
49u	1.1.1-TRICHLOROETHANE	49u	BROMOBENZENE
49u	1.1-DICHLOROPROPENE	49u	1.1.2.2-TETRACHLOROETHANE
49u	CARBON TETRACHLORIDE	49u	1.2.3-TRICHLOROPROPANE
7.8J	1.2-DICHLOROETHANE	49u	O-CHLOROTOLUENE
4QU	BENZENE	49u	P-CHLOROTOLUENE
49u	TRICHLOROETHENE(TRICHLOROETHYLENE)	49u	1.3-DICHLOROBENZENE
49u	1.2-DICHLOROPROPANE	49u	1.4-DICHLOROBENZENE
49u	DIBROMOMETHANE	49u	1.2-DICHLOROBENZENE
4911	BROMODICHLOROMETHANE	14 7	PERCENT MOISTURE

@**FOOTNOTES***
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS _ MANAGEMENT SYSTEM
€PA-REGION I V ESD, ATHENS, GA.

03/30/94

EXTRACTABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84227 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S3G

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLCCTION START: 03/08/94 1115 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1600U	(3-AND/OR 4-)MCTHYLPHCNOL
1600U	1,2,4-TRICHLOROBENZENE
1600U	2,2'-CHLOROISOPROPYLETHER
1600U	2,3,4,6-TETRACHLOROPHENOL
1600U	2,4,5-TRICHLOROPHENOL
1600U	2,4,6-TRICHLOROPHENOL
1600U	2,4-DICHLOROPHENOL
1600U	2,4-DIMETHYLPHENOL
3100U	2,4-DINITROPHENOL
1600U	2,4-DINITROTOLUENE
1600U	2,6-DINITROTOLUENE
1600U	2-CHLORONAPHTHALENE
1600U	2-CHLOROPHENOL
3100U	2-METHYL-4,6-DINITROPHENOL
1600U	2-METHYLNAPHTHALENE
1600U	2-METHYLPHENOL
1600U	2-NITROANILINE
1600U	2-NITROPHENOL
1600U	3,3'-DICHLOROBENZIDINE
1600U	3-NITROANILINE
1600U	4-BROMOPHENYL PHENYL ETHER
1600U	4-CHLORO-3-METHYLPHENOL
1600U	4-CHLOROANILINE
1600U	4-CHLOROPHENYL PHENYL ETHER
1600U	4-NITROANILINE
3100U	4-NITROPHENOL
1600U	ACENAPHTHENE
1600U	ACENAPHTHYLENE
1600U	ANTHRACENE
1600U	BENZO(A)ANTHRACENE
1600U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1600U	BENZO(GHI)PERYLENE
1600U	BENZO-A-PYRENE
1600U	BENZYL BUTYL PHTHALATE
1600U	BIS(2-CHLOROETHOXY) METHANE
1600U	BIS(2-CHLOROETHYL) ETHER
1600U	BIS(2-ETHYLHEXYL) PHTHALATE
1600U	CARBAZOLE
1600U	CHRYSENE
1600U	DI-N-BUTYLPHTHALATE
1600U	DI-N-OCTYLPHTHALATE
1600U	DIBENZO(A,H)ANTHRACENE
1600U	DIBENZOFURAN
1600U	DIETHYL PHTHALATE
1600U	DIMETHYL PHTHALATE
1600U	FLUORANTHENE
1600U	FLUORENE
1600U	HEXACHLOROBENZENE (HCB)
1600U	HEXACHLOROBUTADIENE
1600U	HEXACHLOROCVCLOPENTADIENE (HCCP)
1600U	HEXACHLOROETHANE
1600U	INDENO (1,2,3-CD) PYRENE
1600U	ISOPHORONE
1600U	N-NITROSODI-N-PROPYLAMINE
1600U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1600U	NAPHTHALENE
1600U	NITROBENZENE
3100U	PENTACHLOROPHENOL
16WU	PHENANTHRENE
1600U	PHENOL
1600U	PYRENE
14.7	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES

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*U-MATERIAL WAS ANALYZED FOR BUT *NOT* DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSI. ANAGEMENT SYSTEM
€PA-REGION IV ESD, ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84227 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32536

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1115 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
SOU	D1ELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
SOU	ENDOSULFAN II [BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCO-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLOROENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONAEHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
15	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER JS THE MINIMUM QUANTITATION LIMIT C-CONFIRMED BY GC/MS
1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYS. MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84227 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32536

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PCNSACOLA ST: FL
COLLECTION START: 03/08/94 1115 STOP: 00/00/00

RESULTS	UNITS	PARAMCTCR
0.25	MG/KG	CYANIDE

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84227 SAMPLE TYPE; SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S36

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1115 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
13	SILVER
3.0U	ARSENIC
MA	BORON
41	BARIUM
0.50U	BERYLLIUM
16	CADMIUM
1.1	COBALT
600	CHROMIUM
59	COPPER
1.0U	MOLYBDENUM
43	NICKEL
56	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
13U	TIN
7.6	STRONTIUM
5.0U	TELLURIUM
53	TITANIUM
0U	THALLIUM
6.0	VANADIUM
1.5	YTTRIUM
120	ZINC
NA	ZIRCONIUM
0.56	MERCURY
6700	ALUMINUM
69	MANGANESE

MG/KG	ANALYTICAL RESULTS
1200	CALCIUM
150	MAGNESIUM
2400	IRON
100U	SODIUM
200u	POTASSIUM
14	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84228 SAMPLE TYPE: **SOIL**
SOURCE: NAS PENSACOLA
STATION ID: 32S39

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1130 STOP: **00/00/00**

UG/KG	ANALYTICAL RESULTS
42U	CHLOROMETHANE
42U	VINYL CHLORIDE
42U	BROMOMETHANE
42U	CHLOROETHANE
42U	TRICHLOROFLUOROMETHANE
42U	**1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)**
420U	ACETONE
100U	CARBON DISULFIDE
42U	METHYLENE CHLORIDE
42U	**TRANS-1.2-DICHLOROETHENE**
42U	1,1-DICHLOROETHANE
42U	CIS-1,2-DICHLOROETHENE
42U	2.2-DICHLOROPROPANE
420U	METHYL ETHYL KETONE
42U	BROMOCHLOROMETHANE
42U	CHLOROFORM
42U	1,1,1-TRICHLOROETHANE
42U	1.1-DICHLOROPROPENE
42U	CARBON TETRACHLORIDE
0.75	1.2-DICHLOROETHANE
42U	BENZENE
42U	**TRICHLOROETHENE(TRICHLOROETHYLENE)**
42U	1.2-DICHLOROPROPANE
42U	DIBROMOMETHANE
42U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
42U	CIS-1.3-DICHLOROPROPCNC
100U	METHYL ISOBUTYL KETONE
4.35	TOLUENE
42U	TRANS-1,3-DICHLOROPROPENE
42U	1,1,2-TRICHLOROETHANE
42U	**TETRACHLOROETHENE(TETRACHLOROETHYLENE)**
42U	1.3-DICHLOROPROPANE
100U	METHYL BUTYL KETONE
42U	DIBROMOCHLOROMETHANE
42U	CHLOROBENZENE
42U	1.1.1.2-TETRACHLOROETHANE
42U	ETHYL BENZENE
42U	**(M- AND/**OR P-)XYLENE
42U	**O-XYLENE**
42U	STYRENE
42U	BROMOFORM
42U	BROMOBENZENE
42U	1.1.2.2-TETRACHLOROETHANE
42U	1.2.3-TRICHLOROPROPANE
42U	O-CHLOROTOLUENE
42U	P-CHLOROTOLUENE
42u	**1.3-DICHLOROBENZENE**
42U	**1.4-DICHLOROBENZENE**
42U	**1.2-DICHLOROBENZENE**
7.8	**PERCENT MOISTURE**

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/30/94

EXTRACTABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84228 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S39

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PLNSACOLA ST: FL
COLLECTION START: 03/08/94 1130 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1400U	(3-AND/OR 4-)METHYLPHENOL
1400U	1,2,4-TRICHLOROBENZENE
1400U	2,2'-CHLOROISOPROPYLETHER
1400U	2,3,4,6-TETRACHLOROPHENOL
1400U	2,4,5-TRICHLOROPHENOL
1400U	2,4,6-TRICHLOROPHENOL
1400U	2,4-DICHLOROPHENOL
1400U	2,4-DIMETHYLPHENOL
2900U	2,4-DINITROPHENOL
1400U	2,4-DINITROTOLUENE
1400U	2,6-DINITROTOLUENE
1400U	2-CHLORONAPHTHALENE
1400U	2-CHLOROPHENOL
2900U	2-METHYL-4,6-DINITROPHENOL
1400U	2-METHYLNAPHTHALENE
1400U	2-METHYLPHENOL
1400U	2-NITROANILINE
1400U	2-NITROPHENOL
1400U	3,3'-DICHLOROBENZIDINE
1400U	3-NIIROANILINE
1400U	4-BROMOPHENYL PHENYL ETHER
1400U	4-CHLORO-3-METHYLPHENOL
1400U	4-CHLOROANILINE
1400U	4-CHLOROPHENYL PHENYL ETHER
1400U	4-NITROANILINE
2900U	4-NITROPHENOL
1400U	ACENAPHTHENE
1400U	ACENAPHTHYLENE
1400U	ANTHRACENE
1400U	BENZO(A)ANTHRACENE
1400U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1400U	BENZO(GHI)PERYLENE
1400U	BENZO-A-PVREME
1400U	BENZYL BUTYL PHTHALATE
1400U	BIS(2-CHLOROETHOXY) METHANE
1400U	BIS(2-CHLOROETHYL) ETHER
1400U	BIS(2-ETHYLHEXYL) PHTHALATE
1400U	CARBAZOLE
1400U	CHRVSENE
1400U	DI-N-BUTYLPHTHALATE
1400U	DI-N-OCTYLPHTHALATE
1400U	DIBENZO(A,H)ANTHRACENE
1400U	DIBENZOFURAN
1400U	DIETHVL PHTHALATE
1400U	DIMETHYL PHTHALATE
1400U	FLUORANTHENE
1400U	FLUORENE
1400U	HEXACHLOROBENZENE (HCB)
1400U	HEXACHLOROBUTADIENE
1400U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1400U	HEXACHLOROETHANE
1400U	INDENO (1,2,3-CD) PYRENE
1400U	ISOPHORONE
1400U	N-NITROSODI-N-PROPYLAMINE
1400U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1400U	NAPHTHALENE
1400U	NITROBENZENE
2900U	PENTACHLOROPHENOL
1400U	PHENANTHRENE
1400U	PHENOL
1400U	PVRENE
7.9	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

04/05/94

PESTICIDES/PCB'S DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84228 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32539

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1130 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50UJ	ALDRIN
50UJ	HEPTACHLOR
50UJ	HEPTACHLOR EPOXIDE
50UJ	ALPHA-BHC
50UJ	BETA-BHC
50UJ	GAMMA-BHC (LINDANE)
50UJ	DELTA-BHC
50UJ	ENDOSULFAN I (ALPHA)
50UJ	DIELDRIN
50UJ	4,4'-DDT (P,P'-DDT)
50UJ	4,4'-DDE (P,P'-DDE)
50UJ	4,4'-DDD (P,P'-DDD)
50UJ	ENDRIN
50UJ	ENDOSULFAN II (BETA)
50UJ	ENDOSULFAN SULFATE
200UJ	CHLORDANE (TECH. MIXTURE) /1
500UJ	PCB-1242 (AROCLOR 1242)
500UJ	PCB-1254 (AROCLOR 1254)
500UJ	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500UJ	PCB-1232 (AROCLOR 12321
500UJ	PCB-1248 (AROCLOR 1248)
500UJ	PCB-1260 (AROCLOR 1260)
500UJ	PCB-1016 (AROCLOR 1016)
3000UJ	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200UJ	METHOXYCHLOR
50UJ	ENDRIN KETONE
7.8	PERCENT MOISTURE

REMARKS

QUANT. IS SUSPECT BASED ON QC DATA

REMARKS

FOOTNOTES

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SAMPLE AND ANALYSI_ ..ANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS. GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84228 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: 32539 COLLECTION START: 03/08/94 1130 STOP: 00/00/00

RCSULTS	UNITS	PARAMCTCR
0.21U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84228 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32539

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1130 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
4.3	SILVER
3.0U	ARSENIC
NA	BORON
18	BARIUM
0.50U	BERYLLIUM
6.2	CADMIUM
1.0U	COBALT
240	CHROMIUM
20	COPPER
1.0U	MOLYBDENUM
12	NICKEL
34	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
7.5U	TIN
10	STRONTIUM
5.0U	TELLURIUM
26	TITANIUM
10U	THALLIUM
2.3	VANADIUM
1.0U	YTTRIUM
47	ZINC
NA	ZIRCONIUM
0.16	MERCURY
2300	ALUMINUM
38	MANGANESE

MG/KG	ANALYTICAL RESULTS
1100	CALCIUM
83	MAGNESIUM
1000	IRON
100U	SODIUM
200U	POTASSIUM
6	PERCENT MOISTURE

REMARKS

**REMARKS*

FOOTNOTES
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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84229 SAMPLE TYPE; SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S06

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1220 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
39u	CHLOROMETHANE
39u	VINYL CHLORIDE
39u	BROMOMETHANE
39u	CHLOROETHANE
39u	TRICHLOROFLUOROMETHANE
39u	1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)
390U	ACETONE
98U	CARBON DISULFIDE
39u	METHYLENE CHLORIDE
39u	TRANS-1.2-DICHLOROETHENE
39u	1.1-DICHLOROETHANE
39u	CIS-1.2-DICHLOROETHENE
39u	2.2-DICHLOROPROPANE
390U	METHYL ETHYL KETONE
39u	BROMOCHLOROMETHANE
39u	CHLOROFORM
39u	1.1.1-TRICHLOROETHANE
39u	1.1-DICHLOROPROPENE
39u	CARBON TETRACHLORIDE
39u	1.2-DICHLOROETHANE
39u	BENZENE
39u	TRICHLOROETHENE(TRICHLOROETHYLENE)
39u	1.2-DICHLOROPROPANE
39u	DIBROMOMETHANE
39U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
39u	CIS-1.3-DICHLOROPROPENE
98U	METHYL ISOBUTYL KETONE
39u	TOLUENE
39u	TRANS-1.3-DICHLOROPROPENE
39u	1.1.2-TRICHLOROETHANE
39u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
39u	1.3-DICHLOROPROPANE
98U	METHYL BUTYL KETONE
39u	DIBROMOCHLOROMETHANE
39u	CHLOROBENZENE
39u	1.1.1.2-TETRACHLOROETHANE
39u	ETHYL BENZENE
39u	(M- AND/OR P-)XYLENE
39u	O-XYLENE
39u	STYRENE
39u	BROMOFORM
39u	BROMOBENZENE
39U	1.1.2.2-TETRACHLOROETHANE
39u	1.2.3-TRICHLOROPROPANE
39u	O-CHLOROTOLUENE
39u	P-CHLOROTOLUENE
39u	1.3-DICHLOROBENZENE
39U	1.4-DICHLOROBENZENE
39u	1.2-DICHLOROBENZENE
1 9	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSI_ ,ANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84229 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32S06 COLLCCTION START: 03/08/94 1220 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1300U	(3-AND/OR 4-)METHYLPHENOL
1300U	1,2,4-TRICHLOROBENZENE
1300U	2,2'-CHLOROISOPROPYLETHER
1300U	2,3,4,6-TETRACHLOROPHENOL
1300U	2,4,5-TRICHLOROPHENOL
1300U	2,4,6-TRICHLOROPHENOL
1300U	2,4-DICHLOROPHENOL
1300U	2,4-DIMETHYLPHENOL
2600U	2.4-DINITROPHENOL
1300U	2.4-DINITROTOLUENE
1300U	2,6-DINITROTOLUENE
1300U	2-CHLORONAPHTHALENE
1300U	2-CHLOROPHENOL
2600U	2-METHYL-4,6-DINITROPHENOL
1300U	2-METHYLNAPHTHALENE
1300U	2-METHYLPHENOL
1300U	2-NITROANILINE
1300U	2-NITROPHENOL
1300U	3,3'-DICHLOROBENZIDINE
1300U	3-NITROANILINE
1300U	4-BROMOPHENYL PHENYL ETHER
1300U	4-CHLORO-3-METHYLPHENOL
1300U	4-CHLOROANILINE
1300U	4-CHLOROPHENYL PHENYL ETHER
1300U	4-NITROANILINE
2600U	4-NITROPHENOL
1300U	ACENAPHTHENE
1300U	ACENAPHTHVLENE
1300U	ANTHRACENE
1300U	BENZO(A)ANTHRACENE
1300U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1300U	BENZO(GHI)PERYLENE
1300U	BENZO-A-PYRENE
1300U	BENZYL BUTYL PHTHALATE
1300U	BIS(2-CHLOROETHOXY) METHANE
1300U	BIS(2-CHLOROETHYL) ETHER
1300U	BIS(2-ETHYLHEXYL) PHTHALATE
1300U	CARBAZOLE
1300U	CHRYSENE
1300U	DI-N-BUTYLPHTHALATE
1300U	DI-N-OCTYLPHTHALATE
1300U	DIBENZO(A,H)ANTHRACENE
1300U	DIBENZOFURAN
1300U	DIETHYL PHTHALATE
1300U	DIMETHYL PHTHALATE
1300U	FLUORANTHENE
1300U	FLUORENE
1300U	HEXACHLOROBENZENE (HCB)
1300U	HEXACHLOROBUTADIENE
1300U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1300U	HEXACHLOROETHANE
1300U	INDENO (1,2,3-CD) PYRENE
1300U	ISOPHORONE
1300U	N-NITROSODI-N-PROPYLAMINE
1300U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1300U	NAPHTHOLENE
1300U	NITROBENZENE
2600U	PENTACHLOROPHENOL
1300U	PHENANTHRENE
1300U	PHENOL
1300U	PYRENE
1.9	PERCENT MOISTURE

REMARKS

REMARKS

*****FOOTNOTES*****
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR OUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALY, MANAGEMENT SVSTEM
EPA-REGION IV ESD, ATHENS; GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84229 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S06

PROG ELEM: SSF COLLECTED BV: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1220 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
SOU	HEPTACHLOR
SOU	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
SOU	DELTA-BHC
SOU	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4.4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500u	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCO-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLOROENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLOROANE (OCTACHLOREPOXIDE) /2
200U	METHOXVCHLOR
50U	ENDRIN KETONE
1.9	PERCENT MOISTURE

FOOTNOTES
•A-AVERAGE VALUE •NA-NOT ANALYZED •NAI-INTERFERENCES •J-ESTIMATED VALUE •N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
•K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN •L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
•U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT. C-CONFIRMED BY GC/MS
1. WEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84229 SAMPLE TYPE: SOIL
SOURCE: NAS PCNSACOLA
STATION ID: 32S06

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1220 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.21U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOW TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84229 SAMPLE TYPE: SOIL PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32506 COLLECTION START: 03/08/94 1220 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1 OU	SILVER
3.OU	ARSENIC
NA	BORON
1.OU	BARIUM
0.50U	BERYLLIUM
0.50U	CADMIUM
1.OU	COBALT
1.OU	CHROMIUM
1.OU	COPPER
1.OU	MOLYBDENUM
2.OU	NICKEL
4.OU	LEAD
3.OU	ANTIMONY
4.OU	SELENIUM
4.OU	TIN
23	STRONTIUM
5.OU	TELLURIUM
2.9	TITANIUM
10U	THALLIUM
1.OU	VANADIUM
1.OU	YTTRIUM
1.2	ZINC
NA	ZIRCONIUM
0.05U	MERCURY
100	ALUMINUM
3.5	MANGANESE

MG/KG	ANALYTICAL RESULTS
4700	CALCIUM
40	MAGNESIUM
150	IRON
100U	SODIUM
200U	POTASSIUM
2	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSI_ ,ANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84230 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S67

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1420 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
43u	CHLOROMETHANE
43u	VINYL CHLORIDE
43u	BROMOMETHANE
43u	CHLOROETHANE
43u	TRICHLOROFLUOROMETHANE
43u	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
430U	ACETONE
110U	CARBON DISULFIDE
43u	METHYLENE CHLORIDE
43u	TRANS-1.2-DICHLOROETHENE
43u	1.1-DICHLOROETHANE
43u	CIS-1.2-DICHLOROETHENE
43u	2.2-DICHLOROPROPANE
430U	METHYL ETHYL KETONE
43u	BROMOCHLOROMETHANE
43u	CHLOROFORM
43u	1,1,1-TRICHLOROETHANE
43u	1.1-DICHLOROPROPENE
43u	CARBON TETRACHLORIDE
43u	1.2-DICHLOROETHANE
43u	BENZENE
43u	TRICHLOROETHENE(TRICHLOROETHYLENE)
43U	1.2-DICHLOROPROPANE
43u	DIBROMOMETHANE
4311	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
43u	CIS-1.3-DICHLOROPROPCNC
110U	METHYL ISOBUTYL KETONE
43u	TOLUENE
43u	TRANS-1.3-DICHLOROPROPENE
43u	1,1,2-TRICHLOROETHANE
43u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
43u	1.3-DICHLOROPROPANE
110U	METHYL BUTYL KETONE
43u	DIBROMOCHLOROMETHANE
43u	CHLOROBENZENE
43u	1.1.1.2-TETRACHLOROETHANE
43u	ETHYL BENZENE
43u	(M- AND/OR P-)XYLENE
43u	O-XYLENE
43u	STYRENE
43u	BROMOFORM
43u	BROMOBENZENE
43u	1.1.2.2-TETRACHLOROETHANE
43u	1.2.3-TRICHLOROPROPANE
43u	O-CHLOROTOLUENE
43u	P-CHLOROTOLUENE
43u	1.3-DICHLOROBENZENE
43u	1.4-DICHLOROBENZENE
43u	1.2-DICHLOROBENZENE
2.3	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84230 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S07

PROG ELEM: SSF COLLECTLD BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1420 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1400U	(3-AND/OR 4-)METHYLPHENOL
1400U	1,2,4-TRICHLOROBENZENE
1400U	2,2'-CHLOROISOPROPYLETHER
1400U	2,3,4,6-TETRACHLOROPHENOL
1400U	2,4,5-TRICHLOROPHENOL
1400U	2,4,6-TRICHLOROPHENOL
1400U	2,4-DICHLOROPHENOL
1400U	2,4-DIMETHYLPHENOL
2700U	2,4-DINITROPHENOL
1400U	2,4-DINITROTOLUENE
1400U	2,6-DINITROTOLUENE
1400U	2-CHLORONAPHTHALENE
1400U	2-CHLOROPHENOL
2700U	2-METHYL-4,6-DINITROPHENOL
1400U	2-METHYLNAPHTHALENE
1400U	2-METHYLPHENOL
1400U	2-NITROANILINE
1400U	2-NITROPHENOL
1400U	3,3'-DICHLOROBENZIDINE
1400U	3-NITROANILINE
1400U	4-BROMOPHENYL PHENYL ETHER
1400U	4-CHLORO-3-METHYLPHENOL
1400U	4-CHLOROANILINE
1400U	4-CHLOROPHENYL PHENYL ETHER
1400U	4-NITROANILINE
2700U	4-NITROPHENOL
1400U	ACENAPHTHENE
1400U	ACENAPHTHYLENE
1400U	ANTHRACENE
1400U	BENZO(A)ANTHRACENE
1400U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1400U	BENZO(GHI)PERYLENE
1400U	BENZO-A-PYRENE
1400U	BENZYL BUTYL PHTHALATE
1400U	BIS(2-CHLOROETHOXY) METHANE
1400U	BIS(2-CHLOROETHYL) ETHER
1400U	BIS(2-ETHYLHEXYL) PHTHALATE
1400U	CARBAZOLE
1400U	CHRYSENE
1400U	DI-N-BUTYLPHTHALATE
1400U	DI-N-OCTYL PHTHALATE
1400U	DIBENZO(A,H)ANTHRACENE
1400U	DIBENZOFURAN
1400U	DIETHYL PHTHALATE
1400U	DIMETHYL PHTHALATE
1400U	FLUORANTHENE
1400U	FLUORENE
1400U	HEXACHLOROBENZENE (HCB)
1400U	HEXACHLOROBUTADIENE
1400U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1400U	HEXACHLOROETHANE
1400U	INDENO (1,2,3-CD) PYRENE
1400U	ISOPHORONE
1400U	N-NITROSODI-N-PROPYLAMINE
1400U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1400U	NAPHTHALENE
1400U	NITROBENZENE
2700U	PENTACHLOROPHENOL
1400U	PHENANTHRENE
1400U	PHENOL
1400U	PYRENE
2.3	PERCENT MOISTURE

+REMARKS*

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

04/05/94

PESTICIDES/PCB'S DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84230 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32567

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1420 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
60U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCB-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-MACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	YETHOXVCHLOR
50U	ENDRIN KETONE
2.3	PERCENT MOISTURE

FOOTNOTES

*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN C-CONFIRMED BY GC/MS
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT
1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84230 SAMPLE TVPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32567

PROG ELEM: SSF COLLECTED BV: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1420 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.21U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84230 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32567

PROC ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1420 STOP: 00/00/00

ANALYTICAL RESULTS

MG/KG	
1.0U	SILVER
3.0U	ARSENIC
NA	BORON-
1.GU	BARIUM
0.50U	BERYLLIUM
0.50U	CADMIUM
1.0U	COBALT
1.0	CHROMIUM
1.0U	COPPER
1.0U	MOLYBDENUM
2.0U	NICKEL
4.0U	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
4.0U	TIN
1.0U	STRONTIUM
5.0U	TELLURIUM
5.1	TITAN1UM
10U	THALLIUM
1.0U	VANAOIUM
1.0U	YTTRIUM
1.0U	ZINC-
NA	ZIRCONIUM
0.05U	MERCURY
96	ALUMINUM
1.0U	MANGANESE

ANALYTICAL RESULTS

MG/KG	
150U	CALCIUM
10U	MAGNCSIUM
81	1RON
100U	SODIUM
200u	POTASSIUM
3	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE. *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS.. MANAGEMENT SYS EM
EPA-REGION IV ESD, ATHENS, GA.

PURGEABLE ORGANICS DA A REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 8423 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32568

PROG ELEM: S F COLLECTED BY: F SLOAN
CITY: PENSAC LA TIME STOP: 00/00/00
COLLECTION S ART: 03/08/94

UG/KG	ANALYTICAL RESULTS
3CU	CHLOROMETHANE
36U	VINYL CHLORIDE
36U	BROMOMETHANE
36U	CHLOROETHANE
36U	TRICHLOROFLUORCMETHANE
36U	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
360U	ACETONE
91U	CARBON DISULFIDE
36U	METHYLENE CHLORIDE
36U	TRANS-1,2-DICHLOROETHENE
36U	1.1-DICHLOROETHANE
36U	CIS-1.2-DICHLOROETHENE
36U	2.2-DICHLOROPROPANE
360U	METHYL ETHYL KETONE
36U	BROMOCHLOROMETHANE
36U	CHLOROFORM
36U	1,1,1-TRICHLOROETHANE
36U	1.1-DICHLOROPROPENE
36U	CARBON TETRACHLORIDE
36U	1.2-DICHLOROETHANC
36U	BENZENE
36U	TRICHLOROETHENE(TRICHLOROETHYLENE)
36U	1.2-DICHLOROPROPANE
36U	DIBROMOMETHANE
36U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
3GU	CIS-1.3-DICHLOROPROPCNE
91U	METHYL ISOBUTYL KETONE
36U	TOLUENE
36U	TRANS-1.3-DICHLOROPROPENE
36U	1,1,2-TRICHLOROETHANE
36U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
36U	1.3-DICHLOROPROPANE
91U	METHYL BUTYL KETONE
36U	DIBROMOCHLOROMETHANE
36U	CHLOROBENZENE
36U	1.1.1.2-TETRACHLOROETHANE
36U	ETHYL BENZENE
36U	(M- AND/OR P-)XYLENE
36U	O-XYLENE
36U	STYRENE
36U	BROMOFORM
36U	BROMOBENZENE
36U	1.1.2.2-TETRACHLOROETHANE
36U	1,2,3-TRICHLOROPROPANE
36U	O-CHLOROTOLUENE
36U	P-CHLOROTOLUENE
36U	1.3-DICHLOROBENZENE
36U	1.4-DICHLOROBENZENE
36U	1.2-DICHLOROBENZENE
1.7	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84231 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S08

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1440 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1300U	(3-AND/OR 4-)METHYLPHENOL
1300U	1,2,4-TRICHLOROBENZENE
1300U	2,2'-CHLOROISOPROPYLETHER
1300U	2,3,4,6-TETRACHLOROPHENOL
1300U	2,4,5-TRICHLOROPHENOL
1300U	2,4,6-TRICHLOROPHENOL
1300U	2,4-DICHLOROPHENOL
1300U	2,4-DIMETHYLPHENOL
2700U	2,4-DINITROPHENOL
1300U	2,4-DINITROTOLUENE
1300U	2,6-DINITROTOLUENE
1300U	2-CHLORONAPHTHALENE
1300U	2-CHLOROPHENOL
2700U	2-METHYL-4,6-DINITROPHENOL
1300U	2-METHYLNAPHTHALENE
1300U	2-METHYLPHENOL
1300U	2-NITROANILINE
1300U	2-NITROPHENOL
1300U	3,3'-DICHLOROBENZIDINE
1300U	3-NITROANILINE
1300U	4-BROMOPHENYL PHENYL ETHER
1300U	4-CHLORO-3-METHYLPHENOL
1300U	4-CHLOROANILINE
1300U	4-CHLOROPHENYL PHENYL ETHER
1300U	4-NITROANILINE
2700U	4-NITROPHENOL
1300U	ACENAPHTHENE
1300U	ACENAPHTHYLENE
1300U	ANTHRACENE
1300U	BENZO(A)ANTHRACENE
1300U	BENZO(B AND/OR IE

UG/KG	ANALYTICAL RESULTS
1300U	BENZO(GHI)PERYLENE
1300U	BENZO-A-PYRENE
1300U	BENZYL BUTYL PHTHALATE
1300U	BIS(2-CHLOROETHOXY) METHANE
1300U	BIS(2-CHLOROETHYL) ETHER
1300U	BIS(2-ETHYLHEXYL) PHTHALATE
1300U	CARBAZOLE
1300U	CHRYSENE
1300U	DI-N-BUTYLPHTHALATE
1300U	DI-N-OCTYLPHTHALATE
1300U	DIBENZO(A,H)ANTHRACENE
1300U	DIBENZOFURAN
1300U	DIETHYL PHTHALATE
1300U	DIMETHYL PHTHALATE
1300U	FLUORANTHENE
1300U	FLUORENE
1300U	HEXACHLOROBENZENE (HCB)
1300U	HEXACHLOROBUTADIENE
1300U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1300U	HEXACHLOROETHANE
1300U	INDENO (1,2,3-CD) PYRENE
1300U	ISOPHORONE
1300U	N-NITROSODI-N-PROPYLAMINE
13wu	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1300U	NAPHTHALENE
1300U	NITROBENZENE
2700U	PENTACHLOROPHENOL
1300U	PHENANTHRENE
1300U	PHENOL
1300U	PYRENE
1.7	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALVZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

04/05/94

PESTICIDES/PCB'S DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO, 84231 SAMPLE TYPE; SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S68

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION STAR : 03/08/94 1440 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
SOU	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200u	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCD-1232 AROCLOR 1232)
500U	PCB-1248 AROCLOR 1248)
500U	PCB-1260 AROCLOR 1260)
500U	PCB-1016 AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
1.7	PERCENT MOISTURE

FOOTNOTES

*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN *C-CONFIRMED BY GC/MS
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT
1. WHEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS. GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84231 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32568

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1440 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.21U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84231 SAMPLE TYPE: SOIL
SOURCE: NAS PENSACOLA
STATION ID: 32S68

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1440 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1.0U	SILVER
3.0U	ARSENIC
NA	BORON
1.0U	BARIUM
0.50U	BERYLLIUM
0.50U	CADMIUM
1.0U	COBALT
1.8	CHROMIUM
1.0U	COPPER
1.0U	MOLYBDENUM
2.0U	NICKEL
4.0U	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
4.5U	TIN
1.0U	STRONTIUM
5.0U	TELLURIUM
5.3	TITANIUM
10U	THALLIUM
1.0U	VANADIUM
1.0U	VTTRIUM
1.0U	ZINC
NA	ZIRCONIUM
0.05U	MERCURV
77	ALUMINUM
1.0	MANGANESE

MG/KG	ANALYTICAL RESULTS
50U	CALCIUM
10U	MAGNESIUM
100	IRON
100U	SODIUM
200U	POTASSIUM
2	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
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SAMPLE AND ANALYS. .ANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84232 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: SLUDGE 2

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1440 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
89U	CHLOROMETHANE
89U	VINYL CHLORIDE
89U	BROMOMETHANE
89U	CHLOROETHANE
89U	TRICHLOROFLUOROMETHANE
89U	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
890U	ACETONE
220u	CARBON DISULFIDE
89U	METHYLENE CHLORIDE
89U	TRANS-1.2-DICHLOROETHENE
89U	1.1-DICHLOROETHANE
89U	CIS-1,2-DICHLOROETHENE
89U	2,2-DICWLOROPROPANE
890U	METHYL ETHYL KETONE
89U	BROMOCHLOROMETHANE
89U	CHLOROFORM
89U	1,1,1-TRICHLOROETHANE
89U	1,1-DICHLOROPROPENE
89U	CARBON TETRACHLORIDE
89U	1.2-DICHLOROETHANE
89U	BENZENE
89U	TRICHLOROETHENE(TRICHLOROETHYLENE)
89U	1.2-DICHLOROPROPANE
89U	DIBROMOMETHANE
89U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
89U	CIS-1.3-DICHLOROPROPCNC
220u	METHYL ISOBUTYL KETONE
89U	TOLUENE
89U	TRANS-1.3-DICHLOROPROPENE
89U	1,1,2-TRICHLOROETHANE
89U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
89U	1.3-DICHLOROPROPANE
220u	METHYL BUTYL KETONE
89U	DIBROMOCHLOROMETHANE
89U	CHLOROBENZENE
89U	1.1.1.2-TETRACHLOROETHANE
89U	ETHYL BENZENE
89U	(M- AND/OR P-)XYLENE
89U	O-XYLENE
89U	STYRENE
89U	BROMOFORM
89U	BROMOBENZENE
89U	1.1.2.2-TETRACHLOROETHANE
89U	1.2.3-TRICHLOROPROPANE
89U	O-CHLOROTOLUENE
89U	P-CHLOROTOLUENE
89U	1.3-DICHLOROBENZENE
89U	1.4-DICHLOROBENZENE
89U	1,2-DICHLOROBENZENE
53.4	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84232 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: SLUDGE 2

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1440 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
2600U	(3-AND/OR 4-)METHYLPHENOL
2600U	1,2,4-TRICHLOROBENZENE
2600U	2,2'-CHLOROISOPROPYLETHER
2600U	2,3,4,6-TETRACHLOROPHENOL
2600U	2,4,5-TRICHLOROPHENOL
2600U	2,4,6-TRICHLOROPHENOL
2600U	2,4-DICHLOROPHENOL
2600U	2,4-DIMETHYLPHENOL
5300U	2,4-DINITROPHENOL
2600U	2,4-DINITROTOLUENE
2600U	2,6-DINITROTOLUENE
2600U	2-CHLORONAPHTHALENE
2600U	2-CHLOROPHENOL
5300U	2-METHYL-4,6-DINITROPHENOL
2600U	2-METHYLNAPHTHALENE
2600U	2-METHYLPHENOL
2600U	2-NITROANILINE
2600U	2-NITROPHENOL
2600U	3,3'-DICHLOROBENZIDINE
2600U	3-NITROANILINE
2600U	4-BROMOPHENYL PHENYL ETHER
2600U	4-CHLORO-3-METHYLPHENOL
2600U	4-CHLOROANILINE
2600U	4-CHLOROPHENYL PHENYL ETHER
2600U	4-NITROANILINE
5300U	4-NITROPHENOL
2600U	ACENAPHTHENE
2600U	ACENAPHTHYLENE
2600U	ANTHRACENE
2600U	BENZO(A)ANTHRACENE
2600U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
2600U	BENZO(GHI)PERYLENE
2600U	BENZO-A-PYRENE
2600U	BENZYL BUTYL PHTHALATE
2600U	BIS(2-CHLOROETHOXY) METHANE
2600U	BIS(2-CHLOROETHYL) ETHER
2600U	BIS(2-ETHYLHEXYL) PHTHALATE
2600U	CARBAZOLE
2600U	CHRYSENE
2600U	DI-N-BUTYLPHTHALATE
2600U	DI-N-OCTYLPHTHALATE
2600U	DIBENZO(A,H)ANTHRACENE
2600U	DIBENZOFURAN
2600U	DIETHYL PHTHALATE
2600U	DIMETHYL PHTHALATE
2600U	FLUORANTHENE
2600U	FLUORENE
2600U	HEXACHLOROBENZENE (HCB)
2600U	HEXACHLOROBUTADIENE
2600U	HEXACHLOROCYCLOPENTADIENE (HCCP)
2600U	HEXACHLOROETHANE
2600U	INDENO (1,2,3-CD) PYRENE
2600U	ISOPHORONE
2600U	N-NITROSODI-N-PROPYLAMINE
2600U	N-NITROSODIPHENYLAMINE/DIPHEN LAMINE
2600U	NAPHTHALENE
2600U	NITROBENZENE
5300U	PENTACHLOROPHENOL
2600U	PHENANTHRENE
2600U	PHENOL
2600U	PYRENE
53.4	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS.. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS. GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84232 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: SLUDGE 2

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1440 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BWC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500u	PCB-1254 (AROCLOR 1254)
500u	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCD-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
53	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT C-CONFIRMED BY GC/MS
1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSI, MANAGEMENT SYSTEM
€PA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84232 SAMPLE TYPE. SLUDGE PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCC: NAS PENSACOLA CITY: PENSACOLA ST; FL
STATION ID: SLUDGE 2 COLLECTION START: 03/08/94 1440 STOP: 00/00/00

RCSULTS	UNITS	PARAMCTCR
0.31U	MG/KG	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-ESD,REG IV
ATHENS GEORGIA

03/31/94

DATA CLP
RTIN; SHEET
SOIL/SLUDGE

SAMPLE NO.: 94C84232 SAMPLE YPE: SLUDGE

PROJECT NO.: 94-0323 PROGRAM ELEMENT: SST
SWRCE : NAS PENSACOLA
CITY: PENSACOLA STATE: FL

STATION I.D · SLUDGE 2
STORET STATION NO:

SAMPLE COLLECTION: START DATE/TIME 03/08/94 1440
SAMPLE COLLECTION: STOP DATE/TIME 00/00/00

COLLECTED BY: F SLOAN RECEIVED FROM: F SLOAN
SAMPLE REC'D: DATE/TIME 03/11/94 0940 REC'D BY: D COLOUITT
SEALED: YES
CHEMIST: FVC

REMARK:
REMARK:

SAMPLE LOG VERIFIED BY: VAX SAMPLE DATA VERIFIED BY: MAW

REMARKS

*****ANALYTICAL RESULTS*****

RESULTS	UNITS	PARAMETER
0.05U	MG/L	SILVER
0.15U	MG/L	ARSENIC
0.2011	MG/L	BARIUM
0.025U	MG/L	CADMIUM
0.05U	MG/L	CHROMIUM
0.20U	MG/L	LEAD
0.20U	MG/L	SELENIUM
NA	MG/L	MERCURY
0.10U	MG/L	NICKEL
0.20U	MG/L	ANTIMONY
0.025U	MG/L	BERYLLIUM
0.50U	MG/L	THALLIUM

****

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES
*J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN
*L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84232 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: SLUDGE 2

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITV: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1440 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1.0U	SILVER
5.0U	ARSCNIC
MA	BORON
5.8	BARIUM
0.50U	BERYLLIUM
0.65	CADMIUM
1.2	COBALT
28	CHROMIUM
9.2	COPPER
1.1	MOLYBDENUM
4.2	NICKEL
17	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
3.5U	TIN
16	STRONTIUM
5.0U	TELLURIUM
33	TITANIUM
10U	THALLIUM
12	VANADIUM
2.9	YTTRIUM
29	ZINC
NA	ZIRCONIUM
0.12	YERCURV
4100	ALUMINUM
22	MANGANESE

MG/KG	ANALYTICAL RESULTS
2600	CALCIUM
1000	MAGNCSIUM
4300	IRON
100U	SODIUM
200U	POTASSIUM
36	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR OUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84233 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: 32561

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1520 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
37u	CHLOROMETHANE
37u	VINYL CHLORIDE
37u	BROMOMETHANE
37u	CHLOROETHANE
37u	TRICHLOROFLUOROMETHANE
37u	1.1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
370U	ACETONE
91U	CARBON DISULFIDE
37u	METHYLENE CHLORIDE
37u	TRANS-1.2-DICHLOROETHENE
37u	1.1-DICHLOROETHANE
37u	CIS-1,2-DICHLOROETHENE
37u	2,2-DICHLOROPROPANE
370U	METHYL ETHYL KETONE
37u	BROMOCHLOROMETHANE
37U	CHLOROFORM
37u	1,1,1-TRICHLOROETHANE
37u	1,1-DICHLOROPROPENE
37u	CARBON TETRACHLORIDE
37u	1,2-DICHLOROETHANE
37u	BENZENE
37u	TRICHLOROETHENE(TRICHLOROETHYLENE)
37u	1.2-DICHLOROPROPANE
37u	DIBROMOMETHANE
37U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
37u	CIS-1.3-DICHLOROPROPCNE
91U	METHYL ISOBUTYL KETONE
37u	TOLUENE
37u	TRANS-1,3-DICHLOROPROPENE
37U	1,1,2-TRICHLOROETHANE
37u	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
37U	1.3-DICHLOROPROPANE
91U	METHYL BUTYL KETONE
37u	DIBROMOCHLOROMETHANE
37u	CHLOROBENZENE
37u	1.1.1.2-TETRACHLOROETHANE
37u	ETHYL BENZENE
37u	(M- AND/OR P-)XYLENE
37u	O-XYLENE
37u	STYRENE-
37u	BROMOFORM
37u	BROMOBENZENE
37u	1.1.2.2-TETRACHLOROETHANE
37u	1,2,3-TRICHLOROPROPANE
37U	O-CHLOROTOLUENE
37u	P-CHLOROTOLUENE
37u	1.3-DICHLOROBENZENE
37u	1.4-DICHLOROBENZENE
37u	1.2-DICHLOROBENZENE
2.4	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/30/94

PROJECT NO. 94-0323 SAMPLE NO. 84233 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: 32SG1

FROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLCCTION START: 03/08/94 1520 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1300U	(3-AND/OR 4-)METHYLPHENOL
1300U	1.2.4-TRICHLOROBENZENL
1300U	2.2'-CHLOROISOPROPYLETHER
1300U	2.3.4.6-TETRACHLOROPHENOL
1300U	2,4,5-TRICHLOROPHENOL
1300U	2,4,6-TRICHLOROPHENOL
1300U	2,4-DICHLOROPHENOL
1300U	2,4-DIMETHYLPHENOL
2600U	2.4-DINITROPHENOL
1300U	2,4-DINITROTOLUENE
1300U	2.6-DINITROTOLUENE
1300U	2-CHLORONAPHTHALENE
1300U	2-CHLOROPHENOL
2600U	2-METHVL-4.6-DINITROPHENOL
1300U	2-ME THVLNAPHTHALENE
1300U	2-METHVLPHENOL
1300U	2-NITROANILINE
1300U	2-NITROPHENOL
1300U	3.3'-DICHLOROBENZIDINE
1300U	3-NITROANILINE
1300U	4-BROMOPHENYL PHENYL ETHER
1300U	4-CHLORO-3-METHYLPHENOL
1300U	4-CHLOROANILINE
1300U	4-CHLOROPHENVL PHENYL ETHER
1300U	4-NITROANILINE
2600U	4-NITROPHENOL
1300U	ACENAPHTHENE
1300U	ACENAPHTHYLENE
1300U	ANTHRACENE
1300U	BENZO(A)ANTHRACENE
1300U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1300U	BENZO(CHI)PERYLENE
1300U	BENZO-A-PYRENE
1300U	BENZYL BUTYL PHTHALATE
1300U	BIS(2-CHLOROETHOXY) METHANE
1300U	BIS(2-CHLOROETHYL) ETHER
1300U	BIS(2-ETHYLHEXYL) PHTHALATE
1300U	CARBAZOLE
1300U	CHRYSENE
1300U	DI-N-BUTYLPHTHALATE
1300U	DI-N-OCTVLPHTHALATE
1300U	DIBENZO(A,H)ANTHRACENE
1300U	DIBENZOFURAN
1300U	DIETHYL PHTHALATE
1300U	DIMETHYL PHTHALATE
1300U	FLUORANTHENE
1300U	FLUORENE
1300U	HEXACHLOROBENZENE (HCB)
1300U	HEXACHLOROBUTADIENE
1300U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1300U	HEXACHLOROETHANE
1300U	INDENO (1,2,3-CD) PYRENE
1300U	ISOPHORONE
1300U	N-NITROSODI-N-PROPYLAMINE
1300U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1300U	NAPHTHALENE
1300U	NITROBENZENE
2600U	PENTACHLOROPHENOL
1300U	PHENANTHRENE
1300U	PHENOL
1300U	PYRENE
2.4	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYS . MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84233 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: 32S61

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1520 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
50U	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
50U	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
50U	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200U	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 1254)
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCB-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
2.4	PERCENT MOISTURE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
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1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS OATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84233 SAMPLE TYPE: SLUDGE PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: 32561 COLLECTION START: 03/08/94 1520 STOP: 00/00/00

RESULTS UNITS FARAMCTCR
0.21U MG/KG CYANIDE

+*+FOOTNOTES***
+A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84233 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: 32S61

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START; 03/08/94 1520 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1.011	SILVER
3.0U	ARSCNIC
NA	BORON
1.0U	BARIUM
0.50U	BERYLLIUM
0.50U	CADMIUM
1.0U	COBALT
1.2	CHROMIUM
1.0U	COPPER
1.0U	MOLYBDENUM
2.0U	NICKEL
4.0U	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
4.5U	TIN
1.5U	STRONTIUM
5.0U	TELLURIUM
4.9	TITANIUM
10U	THALLIUM
1.0U	VANADIUM
1.0U	YTTRIUM
1.0U	ZINC
NA	ZIRCONIUM
0.05U	MERCURY
74	ALUMINUM
1 0U	MANGANESE

MG/KG	ANALYTICAL RESULTS
100	CALCIUM
12	MAGNESIUM
69	IRON
10U	SODIUM
200U	POTASSIUM
3	PERCENT MOISTURE

REMARKS

REMARKS

•••FOOTNOTES•••
•A-AVERAGE VALUE •NA-NOT ANALYZED •NAI-INTERFERENCES •J-ESTIMATED VALUE ON PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
•K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN, •L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
•U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

04/05/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84234 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: SLUDGE 1

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1535 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
G3U	CHLOROMCTHANC
63U	VINYL CHLORIDE
63U	BROMOMETHANE
63U	CHLOROETHANE
63U	TRICHLOROFLUOROMETHANE
63U	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
630U	ACETONE
160U	CARBON DISULFIDE
63U	METHYLENE CHLORIDE
63U	TRANS-1,2-DICHLOROETHENE
63U	1,1-DICHLOROETHANE
63U	CIS-1,2-DICHLOROETHENE
63U	2.2-DICHLOROPROPANE
630U	METHYL ETHYL KETONE
63U	BROMOCHLOROMETHANE
63U	CHLOROFORM
63U	1.1.1-TRICHLOROETHANE
63U	1.1-DICHLOROPROPENE
63U	CARBON TETRACHLORIDE
63U	1,2-DICHLOROETHANE
63U	BENZENE
63U	TRICHLOROETHENE(TRICHLOROETHYLENE)
63U	1,2-DICHLOROPROPANE
63U	DIBROMOMETHANE
63U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
G3U	CIS-1.3-DICHLOROPROPCNC
160U	METHYL ISOBUTYL KETONE
63U	TOLUENE
63U	TRANS-1.3-DICHLOROPROPENE
63U	1,1,2-TRICHLOROETHANE
63U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
63U	1,3-DICHLOROPROPANE
160U	METHYL BUTYL KETONE
63U	DIBROMOCHLOROMETHANE
63U	CHLOROBENZENE
63U	1.1.1.2-TETRACHLOROETHANE
63U	ETHYL BENZENE
63U	(M- AND/OR P-)XYLENE
63U	O-XYLENE
63U	STYRENE
63U	BROMOFORM
63U	BROMOBENZENE
63U	1.1.2.2-TETRACHLOROETHANE
63U	1,2,3-TRICHLOROPROPANE
63U	O-CHLOROTOLUENE
63U	P-CHLOROTOLUENE
63U	1.3-DICHLOROBENZENE
63U	1.4-DICHLOROBENZENE
63U	1,2-DICHLOROBENZENE
20.8	PERCENT MOISTURE

FOOTNOTES
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SAMPLE AND ANALYSI, ,ANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/30/94

EXTRACTABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO, 84234 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: SLUDGE 1

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1535 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
1500U	(3-AND/OR 4-)METHYLPHENOL
1500U	1,2,4-TRICHLOROBENZENE
1500U	2,2'-CHLOROISOPROPYLETHER
1500U	2,3,4,6-TETRACHLOROPHENOL
1500U	2,4,5-TRICHLOROPHENOL
1500U	2,4,6-TRICHLOROPHENOL
1500U	2,4-DICHLOROPHENOL
1500U	2,4-DIMETHYLPHENOL
3000U	2,4-DINITROPHENOL
1500U	2,4-DINITROTOLUENE
1500U	2,6-DINITROTOLUENE
1500U	2-CHLORONAPHTHALENE
1500U	2-CHLOROPHENOL
3000U	2-METHYL-4,6-DINITROPHENOL
1500U	2-METHYLNAPHTHALENE
1500U	2-METHYLPHENOL
1500U	2-NITROANILINE
1500U	2-NITROPHENOL
1500U	3,3'-DICHLOROBENZIDINE
1500U	3-NITROANILINE
1500U	4-BROMOPHENYL PHENYL ETHER
1500U	4-CHLORO-3-METHYLPHENOL
1500U	4-CHLOROANILINE
1500U	4-CHLOROPHENYL PHENYL ETHER
1500U	4-NITROANILINE
3000U	4-NITROPHENOL
1500U	ACENAPHTHENE
1500U	ACENAPHTHYLENE
1500U	ANTHRACENE
1500U	BENZO(A)ANTHRACENE
1500U	BENZO(B AND/OR K)FLUORANTHENE

UG/KG	ANALYTICAL RESULTS
1500U	BENZO(GHI)PERYLENE
1500U	BENZO-A-PYRENE
1500U	BENZYL BUTYL PHTHALATE
1500U	BIS(2-CHLOROETHOXY) METHANE
1500U	BIS(2-CHLOROETHYL) ETHER
1500U	BIS(2-ETHYLHEXYL) PHTHALATE
1500U	CARBAZOLE
1500U	CHRYSENE
1500U	DI-N-BUTYLPHTHALATE
1500U	DI-N-OCTYLPHTHALATE
1500U	DIBENZO(A,H)ANTHRACENE
1500U	DIBENZOFURAN
1500U	DIETHYL PHTHALATE
1500U	DIMETHYL PHTHALATE
1500U	FLUORANTHENE
1500U	FLUORENE
1500U	HEXACHLOROBENZENE (HCB)
1500U	HEXACHLOROBUTADIENE
1500U	HEXACHLOROCYCLOPENTADIENE (HCCP)
1500U	HEXACHLOROETHANE
1500U	INDENO (1,2,3-CD) PYRENE
1500U	ISOPHORONE
1500U	N-NITROSODI-N-PROPYLAMINE
1500U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1500U	NAPHTHALENE
1500U	NITROBENZENE
3000U	PENTACHLOROPHENOL
1500U	PHENANTHRENE
1500U	PHENOL
1500U	PYRENE
20.8	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84234 SAMPLE TYPE: SLUDGE
SOURCE: NAS PENSACOLA
STATION ID: SLUDGE 1

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/08/94 1535 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
50U	ALDRIN
50U	HEPTACHLOR
50U	HEPTACHLOR EPOXIDE
SOU	ALPHA-BHC
50U	BETA-BHC
50U	GAMMA-BHC (LINDANE)
50U	DELTA-BHC
50U	ENDOSULFAN I (ALPHA)
50U	DIELDRIN
SOU	4,4'-DDT (P,P'-DDT)
50U	4,4'-DDE (P,P'-DDE)
50U	4,4'-DDD (P,P'-DDD)
50U	ENDRIN
SOU	ENDOSULFAN II (BETA)
50U	ENDOSULFAN SULFATE
200u	CHLORDANE (TECH. MIXTURE) /1
500U	PCB-1242 (AROCLOR 1242)
500U	PCB-1254 (AROCLOR 12541
500U	PCB-1221 (AROCLOR 1221)

UG/KG	ANALYTICAL RESULTS
500U	PCO-1232 (AROCLOR 1232)
500U	PCB-1248 (AROCLOR 1248)
500U	PCB-1260 (AROCLOR 1260)
500U	PCB-1016 (AROCLOR 1016)
3000U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLOROENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOR:EPOXIDE) /2
200U	METHOXYCHLOR
50U	ENDRIN KETONE
21	PERCENT MOISTURE

FOOTNOTES
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SAMPLE AND ANALYS. ANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. **94-0323** SAMPLE **NO. 84234** SAMPLE TYPE: SLUDGE PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCC: **NAS** PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: SLUDGE 1 COLLECTION START: 03/08/94 1535 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
0.25U	MG/KG	CYANIDE

*****FOOTNOTES*****
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-ESD,REG IV
ATHENS GEORGIA

03/31/94

TCLP
DATA REPORTING SHEET
SEDIMENT/SOIL/SLUDGE

SAMPLE NO.: 94C84234 SAMPLE TYPE: SLUDGE

PROJCCT NO.: 94-0020 PROGRAM ELEMENT: SSF
SOUPCE: NAS PENSACOLA
CITY: PENSACOLA STATE: FL

STATION I.O.: SLUDGE 1
STORET STATION NO:

SAMPLE COLLECTION: START DATE/TIME 03/08/94 1535
SAMPLE COLLECTION: STOP DATE/TIME 00/00/00

COLLECTED BY: F SLOAN RECEIVED FROM: F SLOAN
SAMPLE REC'D: DATE/TIME 03/11/94 0940 REC'D BY: D COLQUITT
SEALED: VES
CHEMIST: FVC

REMARK :
RCMARK :

SAMPLE LOG VERIFIED BY: VAX SAMPLE DATA VERIFIED BY: MAW

REMARKS

*****ANALYTICAL RESULTS*****

RESULTS	UNITS	PARAMETER
0.05U	MG/L	SILVER
0.15U	MG/L	ARSENIC
O 2011	MG/L	BARIUM
0.025U	MG/L	CADMIUM
0.05U	MG/L	CHROMIUM
0.20U	MG/L	LEAD
0.20U	MG/L	SELENIUM
NA	MG/L	MERCURY
0.10U	MG/L	NICKEL
0.20U	MG/L	ANTIMONY
0.025U	MG/L	BERVLLIUM
0.50U	MG/L	THALLIUM

****

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES
*J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN
*L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR OUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84234 SAMPLE TYPE: SLUDGE PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: SLUDGE 1 COLLECTION START: 03/08/94 1535 STOP: 00/00/00

MG/KG	ANALYTICAL RESULTS
1.0U	SILVER
3.0U	ARSENIC
NA	BORON
1.0U	BARIUM
0.50U	BERYLLIUM
0.50U	CADMIUM
1.0U	COBALT
3.6	CHROMIUM
1.2	COPPER
1.0U	MOLYBDENUM
2.0U	NICKEL
4.0U	LEAD
3.0U	ANTIMONY
4.0U	SELENIUM
4.5U	TIN
1.2	STRONTIUM
5.0U	TELLURIUM
20	TITANIUM
10U	THALLIUM
1.8	VANADIUM
1.0U	YTTRIUM
4.4	ZINC
NA	ZIRCONIUM
0.05U	MERCURY
540	ALUMINUM
29	MANGANESE

MG/KG	ANALYTICAL RESULTS
130	CALCIUM
110	MAGNESIUM
510	IRON
100U	SODIUM
200U	POTASSIUM
21	PERCENT MOISTURE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE YXNIMUM QUANTITATION LIMIT.

SAMPLE AND ANAL..S MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS. GA.

03/15/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84235 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32007

PROC ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1115 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
5.0U	CHLOROMETHANE
5.0U	VINYL CHLORIDE
5.0U	BROMOMETHANE
5.0U	CHLOROETHANE
5.0U	TRICHLOROFLUOROMETHANE
5.0U	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
50U	ACETONE
12u	CARBON DISULFIDE
5.0U	METHYLENE CHLORIDE
5.0U	TRANS-1,2-DICHLOROETHENE
5.0U	1,1-DICHLOROETHANE
5.0U	CIS-1,2-DICHLOROETHENE
5.0U	2,2-DICHLOROPROPANE
50U	UETHYL ETHYL KETONE
5.0U	BROMOCHLOROMETHANE
5.0U	CHLOROFORM
5.0U	1,1,1-TRICHLOROETHANE
5.0U	1,1-DICHLOROPROPENE
5.0U	CARBON TETRACHLORIDE
5.0U	1,2-DICHLOROETHANE
5.0U	BENZENE
5.0U	TRICHLOROETHENE(TRICHLOROETHYLENE)
5.0U	1,2-DICHLOROPROPANE
5.0U	DIBROMOMETHANE
5.0U	BROMODICHLOROMETHANE

UG/L	ANALYTICAL RESULTS
5.0U	CIS-1,3-DICHLOROPROPENE
12u	METHYL ISOBUTYL KETONE
5.0U	TOLUENE
5.0U	TRANS-1,3-DICHLOROPROPENE
5.0U	1,1,2-TRICHLOROETHANE
5.0U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
5.0U	1,3-DICHLOROPROPANE
12u	METHYL BUTYL KETONE
5.0U	DIBROMOCHLOROMETHANE
5.0U	CHLOROBENZENE
5.0U	1,1,1,2-TETRACHLOROETHANE
5.[illegible]	ETHYL BENZENE
5.[illegible]	(M- AND/OR P-)XYLENE
5.0	O-XYLENE
5.0U	STYRENE
5.0U	BROMOFORM
5.0U	BROMOBENZENE
5.0U	1,1,2,2-TETRACHLOROETHANE
5.0U	1,2,3-TRICHLOROPROPANE
5.0U	O-CHLOROTOLUENE
5.0U	P-CHLOROTOLUENE
5.0U	1,3-DICHLOROBENZENE
5.0U	1,4-DICHLOROBENZENE
5.0U	1,2-DICHLOROBENZENE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSI, ..ANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/28/94

EXTRACTABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84235 SAMPLE TYPE; GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32607

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1115 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10U	(3-AND/OR 4-)METHYLPHENOL
10U	1,2,4-TRICHLOROBENZENE
10U	2,2'-CHLOROISOPROPYLETHER
10U	2,3,4,6-TETRACHLOROPHENOL
10U	2,4,5-TRICHLOROPHENOL
10U	2,4,6-TRICHLOROPHENOL
10U	2.4-DICHLOROPHENOL
10U	2.4-DIMETHYLPHENOL
20u	2.4-DINITROPHENOL
10U	2,4-DINITROTOLUENE
10U	2,6-DINITROTOLUENE
10U	2-CHLORONAPHTHALENE
10U	2-CHLOROPHENOL
20u	2-METHYL-4.6-DINITROPHENOL
10U	2-METHYLNAPHTHALENE
10U	2-METHYLPHENOL
10U	2-NITROANILINE
10U	2-NITROPHENOL
10U	3.3'-DICHLOROBENZIDINE
10U	3-NITROANILINE
10U	4-BROMOPHENYL PHENYL ETHER
10U	4-CHLORO-3-METHYLPHENOL
10U	4-CHLOROANILINE
10U	4-CHLOROPHENYL PHENYL ETHER
10U	4-NITROANILINE
20u	4-NITROPHENOL
10U	ACENAPHTHENE
10U	ACENAPHTHVLENE
10U	ANTHRACENE
10U	BENZO(A)ANTHRACENE
10U	BENZO(B AND/OR K)FLUORANTHENE

UG/L	ANALYTICAL RESULTS
10U	BENZO(GHI)PERYLENE
10U	BENZO-A-PYRENE
10U	BENZYL BUTYL PHTHALATE
10U	BIS(2-CHLOROETHOXY) METHANE
10U	BIS(2-CHLOROETHYL) ETHER
10U	BIS(2-ETHYLHEXYL) PHTHALATE
10U	CARBAZOLE
10U	CHRYSENE
10U	DI-N-BUTYLPHTHALATE
1011	DI-N-OCTVIPHTHALATE
10U	DIBENZO(A,H)ANTHRACENE
10U	DIBENZOFURAN
10U	DIETHYL PHTHALATE
10U	DIMETHYL PHTHALATE
10U	FLUORANTHENE
10U	FLUORENE
10U	HEXACHLOROBENZENE (HCB)
10U	HEXACHLOROBUTADIENE
10U	HEXACHLOROCYCLOPENTADIENE (HCCP)
10U	HEXACHLOROETHANE
10U	INDENO (1,2,3-CD) PYRENE
10U	ISOPHORONE
10U	N-NITROSODI-N-PROPYLAMINE
10U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
10U	NAPHTHALENE
10U	NITROBENZENE
20U	PENTACHLOROPHENOL
10U	PHENANTHRENE
10U	PHENOL
10U	PYRENE

REMARKS

REMARKS

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SAMPLE AND ANALYSIS mANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/31/94

PESTICIDES/PCB'S DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84235 SAMPLE TYPE: CROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G07

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITV: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1115 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
0.50U	ALDRIN
0.50U	HEPTACHLOR
0.50U	HEPTACHLOR EPOXIDE
0.50U	ALPHA-BHC
0.50U	BETA-BHC
0.50U	GAMMA-BHC (LINDANE)
0.50U	DELTA-BHC
0.50U	ENDOSULFAN I ALPHA)
0.50U	DIELDRIN
0.50U	4,4'-DDT (P,P -DDT)
0.50U	4,4'-DDE (P,P -DDE)
0.50U	4,4'-DDD (P,P -DDD)
0.50U	ENDRIN
0.SOU	ENDOSULFAN II (BETA)
0.50U	ENDOSULFAN SU FATE
1.0U	CHLORDANE (TE H. MIX1URE) /1
2.0U	PCB-1242 (ARO LOR 1242)
2.0U	PCB-1254 (ARO LOR 1254)
2.0U	PCB-1221 (ARO LOR 1221)

UG/L	ANALYTICAL RESULTS
2.0U	PCD-1232 (AROCLOR 1232)
2.0U	PCB-1248 (AROCLOR 1248)
2.0U	PCB-1260 (AROCLOR 1260)
2.0U	PCB-1016 (AROCLOR 1016)
20U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
1.0U	METHOXYCHLOR
0.50U	ENDRIN KETONE

REMARKS

REMARKS

FOOTNOTES

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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT C-CONFIRMED BY GC/MS
1. WEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYS. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS. GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84235 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: 32G07 COLLECTION START: 03/09/94 1115 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
4U	UG/L	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84235 SAMPLE TYPE: GROUNDWA PROC ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32G07 COLLECTION START: 03/09/94 1115 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10UU	SILVER
30U	ARSENIC
NA	BORON
10U	BARIUM
5.0U	BERYLLIUM
5.0U	CADMIUM
10U	COBALT
10U	CHROMIUM
10U	COPPER
10U	MOLYBDENUM
20U	NICKEL
45	LEAD
30U	ANTIMONY
40U	SELENIUM
25U	TIN
26	STRONTIUM
50U	TELLURIUM
10U	TITANIUM
100U	THALLIUM
10U	VANADIUM
10U	YTTRIUM
10U	ZINC
NA	ZIRCONIUM
0.2U	MERCURY
430	ALUMINIUM
10U	MANGANESE

MG/L	ANALYTICAL RESULTS
7.2	CALCIUM
0.16	MAGNESJUM
0.35	IRON
3.0	SODIUM
2.0U	POTASSIUM

REMARKS* ***REMARKS***

FOOTNOTES
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*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANAL . S MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 03/15/94

PROJECT NO. 94-0323 SAMPLE NO. 84236 SAMPLE TYPE: CROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32G15 COLLECTION START: 03/09/94 1340 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS	UG/L	ANALYTICAL RESULTS
5.OU	CHLOROMETHANE	5.OU	CIS-1.3-DICHLOROPROPENE
5.OU	VINYL CHLORIDE	12u	METHYL ISOBUTYL KETONE
5.OU	BROMOMETHANE	5.OU	TOLUENE
5.OU	CHLOROETHANE	5.OU	TRANS-1.3-DICHLOROPROPENE
5.OU	TRICHLOROFLUOROMETHANE	5.OU	1.1.2-TRICHLOROETHANE
5.OU	1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)	5.OU	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
50U	ACETONE	5.OU	1.3-DICHLOROPROPANE
12u	CARBON DISULF IDE	12u	METHYL BUTYL KETONE
5.OU	METHYLENE CHLORIDE	5.OU	DIBROMOCHLOROMETHANE
5.OU	TRANS-1.2-DICHLOROETHENE	7.0	CHLOROBENZENE
1.4J	1.1-DICHLOROETHANE	5.OU	1.1.1.2-TETRACHLOROETHANE
5.OU	CIS-1.2-DICHLOROETHENE	5.OU	ETHYL BENZENE
5.OU	2.2-DICHLOROPROPANE	5.OU	(M- AND/OR P-)XYLENE
50U	METHYL ETHYL KETONE	5.OU	O-XYLENE
5.OU	BROMOCHLOROMETHANE	5.OU	STYRENE
5.OU	CHLOROFORM	5.OU	BROMOFORM
5.OU	1.1.1-TRICHLOROETHANE	5.OU	BROMOBENZENE
5.OU	1.1-DICHLOROPROPENE	5.OU	1.1.2.2-TETRACHLOROETHANE
5.OU	CARBON TETRACHLORIDE	5.OU	1.2.3-TRICHLOROPROPANE
5.OU	1.2-DICHLOROETHANE	5.OU	O-CHLOROTOLUENE
5.OU	BENZENE	5.OU	P-CHLOROTOLUENE
5.OU	TRICHLOROETHENE(TRICHLOROETHYLENE)	0.55J	1.3-DICHLOROBENZENE
5.OU	1.2-DICHLOROPROPANE	1.2J	1.4-DICHLOROBENZENE
5.OU	DIBROMOMETHANE	1.5J	1.2-DICHLOROBENZENE
5.OU	BROMODICHLOROMETHANE		

REMARKS **REMARKS**

FOOTNOTES
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/28/94

EXTRACTABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84236 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G15

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1340 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10U	(3-AND/OR 4-)METHYLPHENOL
10U	1,2,4-TRICHLOROBENZENE
10U	2,2'-CHLOROISOPROPYLETHER
10U	2,3,4,6-TETRACHLOROPHENOL
10U	2,4,5-TRICHLOROPHENOL
10U	2,4,6-TRICHLOROPHENOL
10U	2,4-DICHLOROPHENOL
10U	2,4-DIMETHYLPHENOL
20u	2,4-DINITROPHENOL
10U	2,4-DINITROTOLUENE
10U	2,6-DINITROTOLUENE
10U	2-CHLORONAPHTHALENE
10U	2-CHLOROPHENOL
20u	2-METHYL-4,6-DINITROPHENOL
10U	2-METHYLNAPHTHALENE
10U	2-METHYLPHENOL
10U	2-NITROANILINE
10U	2-NITROPHENOL
10U	3,3'-DICHLOROBENZIDINE
10U	3-NITROANILINE
10U	4-BROMOPHENYL PHENYL ETHER
10U	4-CHLORO-3-METHYLPHENOL
10U	4-CHLOROANILINE
10U	4-CHLOROPHENYL PHENYL ETHER
1011	4-NITROANILINE
20u	4-NITROPHENOL
10U	ACENAPHTHENE
10U	ACENAPHTHYLENE
10U	ANTHRACENE
10U	BENZO(A)ANTHRACENE
10U	BENZO(B AND/OR K)FLUORANTHENE

UG/L	ANALYTICAL RESULTS
10U	BENZO(GHI)PERYLENE
10U	BENZO-A-PYRENE
10U	BENZYL BUTYL PHTHALATE
10U	BIS(2-CHLOROETHOXY) METHANE
10U	BIS(2-CHLOROETHYL) ETHER
10U	BIS(2-ETHYLHEXYL) PHTHALATE
10U	CARBAZOLE
10U	CHRYSENE
10U	DI-N-BUTYLPHTHALATE
10U	DI-N-OCTYLPHTHALATE
10U	DIBENZO(A,H)ANTHRACENE
10U	DIBENZOFURAN
10U	DIETHVL PHTHALATE
10U	DIMETHYL PHTHALATE
10U	FLUORANTHENE
10U	FLUORENE
10U	HEXACHLOROBENZENE (HCB)
10U	HEXACHLOROBUTADIENE
10U	HEXACHLOROCYCLOPENTADIENE (HCCP)
10U	HEXALHLOROETHANE
10U	INDENO (1,2,3-CD) PYRENE
10U	ISOPHORONE
10U	N-NITROSODI-N-PROPYLAMINE
10U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
1011	NAPHTHALENE
10U	NITROBENZENE
20u	PENTACHLOROPHENOL
10U	PHENANTHRENE
10U	PHENOL
10U	PYRENE

REMARKS

REMARKS

FOOTNOTES
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SAMPLE AND ANALYSI ANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/31/94

PESTICIDES/PCB'S DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84236 SAMPLE TYPE; GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G15

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITV: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1340 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
0.50U	ALDRIN
0.50U	HEPTACHLOR
0.50U	HEPTACHLOR EPOXIDE
0.50U	ALPHA-BHC
0.50U	BETA-BHC
0.50U	GAMMA-BHC (LINDANE)
0.50U	DELTA-BHC
0.50U	ENDOSULFAN I (ALPHA)
0.50U	DIELDRIN
0.50U	4,4'-DDT (P,P'-DDT)
0.50U	4,4'-DDE (P,P'-DDE)
0.50U	4,4'-DDD (P,P'-DDD)
0.50U	ENDRIN
0.50U	ENDOSULFAN II (BETA)
0.50U	ENDOSULFAN SULFATE
1.0U	CHLORDANE (TECH. MIXTURE) /1
2.0U	PCB-1242 (AROCLOR 1242)
2.0U	PCB-1254 (AROCLOR 1254)
2.0U	PCB-1221 (AROCLOR 1221)

UG/L	ANALYTICAL RESULTS
2.0U	PCB-1232 (AROCLOR 1232)
2.0U	PCB-1248 (AROCLOR 1248)
2.0U	PCB-1260 (AROCLOR 1260)
2.0U	PCB-1016 (AROCLOR 1016)
20U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
1.0U	METHOXYCHLOR
0.50U	ENDRIN KETONE

@**REMARKS***

REMARKS

@FOOTNOTES*
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT C-CONFIRMED BV GC/MS
1. WHEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84236 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: 32615 COLLECTION START: 03/09/94 1340 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
10	UG/L	CYANIDE

***FOOTNOTES—*
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84236 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G15

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1340 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10U	SILVER
30U	ARSENIC
NA	BORON
10U	BARIUM
5.0U	BERYLLIUM
5.0U	CADMIUM
10U	COBALT
10U	CHROMIUM
10U	COPPER
10U	MOLYBDENUM
20U	NICKEL
5.0U	LEAD
30U	ANTIMONY
40U	SELENIUM
25U	TIN
70	STRONTIUM
50U	TELLURIUM
10U	TITANIUM
100U	THALLIUM
11	VANADIUM
10U	YTTRIUM
10U	ZINC
NA	ZIRCONIUM
0.2U	MERCURY
100U	ALUMINUM
10U	MANGANESE

MG/L	ANALYTICAL RESULTS
5.9	CALCIUM
9.8	MAGNESIUM
0.37	IRON
59	SODIUM
6.8	POTASSIUM

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALY..S MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/15/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84237 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G03

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1430 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
5.0U	CHLOROMETHANE
5.0U	VINYL CHLORIDE
5.0U	BROMOMETHANE
5.0U	CHLOROETHANE
5.0U	TRICHLOROFLUOROMETHANE
5.0U	**1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)**
50U	ACETONE
12u	CARBON DISULFIDE
5.0U	METHYLENE CHLORIDE
5.0U	**TRANS-1.2-DICHLOROETHENE**
5.0U	1.1-DICHLOROETHANE
5.0U	CIS-1.2-DICHLOROETHENE
5.0U	2.2-DICHLOROPROPANE
50U	METHYL ETHYL KETONE
5.0U	BROMOCHLOROMETHANE
5.0U	CHLOROFORM
5.0U	1.1.1-TRICHLOROETHANE
5.0U	**1.1-DICHLOROPROPENE**
5.0U	CARBON TETRACHLORIDE
5.0U	1.2-DICHLOROETHANE
5.0U	BENZENE
5.0U	**TRICHLOROETHENE(TRICHLOROETHYLENE)**
5.0U	1.2-DICHLOROPROPANE
5.0U	DIBROMOMETHANE
5.0U	BROMODICHLOROMETHANE

UG/L	ANALYTICAL RESULTS
5.0U	CIS-1.3-DICHLOROPROPENC
12u	METHYL ISOBUTYL KETONE
5.0U	TOLUENE
5.0U	**TRANS-1.3-DICHLOROPROPENE**
5.0U	1.1.2-TRICHLOROETHANE
5.0U	**TETRACHLOROETHENE(TETRACHLOROE**THYLLNE**)**
5.0U	1.3-DICHLOROPROPANE
12u	METHYL BUTYL KETONE
5.0U	DIBROMOCHLOROMETHANE
5.0U	CHLOROBENZENE
5.0U	**1.1.1.2-TETRACHLOROETHANE**
5.0U	ETHYL BENZENE
5.0U	**(M- AND/OR** P-)XYLENE
5.0U	**O-XYLENE**
5.0U	STYRENE
5.0U	BROMOFORM
5.0U	BROMOBENZENE
5.0U	1.1.2.2-TETRACHLOROETHANE
5.0U	1.2.3-TRICHLOROPROPANE
5.0U	0-CHLOROTOLUENE
5.0U	P-CHLOROTOLUENE
5.0U	1.3-DICHLOROBENZENE
5.0U	1.4-DICHLOROBENZENE
5.0U	1.2-DICHLOROBENZENE

••*REMARKS*••

REMARKS

•••**FOOTNOTES***••
•A-AVERAGE VALUE •NA-NOT ANALYZED •NAI-INTERFERENCES •J-ESTIMATED VALUE •N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/28/94

PROJECT NO. 94-0323 SAMPLE NO. 84237 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G03

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1430 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10U	(3-AND/OR 4-)METHYLPHENOL
10U	1,2,4-TRICHLOROBENZENE
10U	2,2'-CHLOROISOPROPYLETHER
10U	2,3,4,6-TETRACHLOROPHENOL
10U	2,4,5-TRICHLOROPHENOL
10U	2,4,6-TRICHLOROPHENOL
10U	2,4-DICHLOROPHENOL
10U	2,4-DIMETHYLPHENOL
20u	2,4-DINITROPHENOL
1011	2,4-DINITROTOLUENE
10U	2,6-DINITROTOLUENE
10U	2-CHLORONAPHTHALENE
10U	2-CHLOROPHENOL
20u	2-METHYL-4,6-DINITROPHENOL
10U	2-METHYLNAPHTHALENE
10U	2-METHYLPHENOL
10U	2-NITROANILINE
10U	2-NITROPHENOL
10U	3,3'-DICHLOROBENZIDINE
10U	3-NITROANILINE
10U	4-BROMOPHENYL PHENYL ETHER
10U	4-CHLORO-3-METHYLPHENOL
10U	4-CHLOROANILINE
10U	4-CHLOROPHENYL PHENYL ETHER
10U	4-NITROANILINE
20u	4-NITROPHENOL
10U	ACENAPHTHENE
10U	ACENAPHTHVLENE
10U	ANTHRACENE
10U	BENZO(A)ANTHRACENE
10U	BENZO(B AND/OR K)FLUORANTHENE

UG/L	ANALYTICAL RESULTS
10U	BENZO(GHI)PERYLENE
10U	BENZO-A-PYRENE
10U	BENZYL BUTYL PHTHALATE
10U	BIS(2-CHLOROETHOXY) METHANE
10U	BIS(2-CHLOROETHYL) ETHER
10U	BIS(2-ETHYLHEXYL) PHTHALATE
10U	CARBAZOLE
10U	CHRYSENE
10U	DI-N-BUTYLPHTHALATE
10U	DI-N-OCTYLPHTHALATE
10U	DIBENZO(A,H)ANTHRACENE
10U	DIBENZOFURAN
10U	DIETHYL PHTHALATE
10U	DIMETHYL PHTHALATE
10U	FLUORANTHENE
10U	FLUORENE
10U	HEXACHLOROBENZENE (HCB)
10U	HEXACHLOROBUTADIENE
10U	HEXACHLOROCYCLOPENTADIENE (HCCP)
10U	HEXACHLOROETHANE
10U	INDENO (1,2,3-CD) PYRENE
10U	ISOPHORONE
10U	N-NITROSODI-N-PROPYLAMINE
10U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
10U	NAPHTHALENE
10U	NITROBENZENE
20U	PENTACHLOROPHENOL
10U	PHENANTHRENE
10U	PHENOL
10U	PYRENE

REMARKS

REMARKS

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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 03/31/94

PROJECT NO. 94-0323 SAMPLE NO. 84237 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G03

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1430 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
0.50U	ALDRIN
0.50U	HEPTACHLOR
0.50U	HEPTACHLOR EPOXIDE
0.50U	ALPHA-BHC
0.50U	BETA-BHC
0.50U	GAMMA-BHC (LINDANE)
0.50U	DELTA-BHC
0.50U	ENDOSULFAN I (ALPHA)
0.50U	DIELDRIN
0.50U	4,4'-DDT (P,P'-DDT)
0.50U	4,4'-DDE (P,P'-DDE)
0.50U	4,4'-DDD (P,P'-DDD)
0.50U	ENDRIN
0.50U	ENDOSULFAN II (BETA)
0.50U	ENDOSULFAN SULFATE
1.0U	CHLORDANE (TECH. MIXTURE) /1
2.0U	PCB-1242 (AROCLOR 1242)
2.0U	PCB-1254 (AROCLOR 1254)
2.0U	PCB-1221 (AROCLOR 1221)

UG/L	ANALYTICAL RESULTS
2.0U	PCB-1232 (AROCLOR 1232)
2.0U	PCB-1248 (AROCLOR 1248)
2.0U	PCB-1260 (AROCLOR 1260)
2.0U	PCB-1016 (AROCLOR 1016)
20U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
1.0U	METHOXYCHLOR
0.50U	ENDRIN KETONE

REMARKS

REMARKS

FOOTNOTES
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1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYS.. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. **94-0323** SAMPLE NO. **84237** SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PCNSACOLA ST: FL
STATION ID: **32G03** COLLECTION START: **03/09/94** 1430 STOP: **00/00/00**

RESULTS	UNITS	PARAMETER
4U	UG/L	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSI_ ..ANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84237 SAMPLE TYPE: GROUNOWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32G03 COLLECTION START: 03/09/94 1430 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10U	SILVER
30U	ARSENIC
NA	BORON
10U	BARIUM
5.0U	BERYLLIUM
5.0U	CADMIUM
10U	COBALT
10U	CHROMIUM
10U	COPPER
10U	MOLYBDENUM
20U	NICKEL
5.3	LEAD
30U	ANTIMONY
40U	SELENIUM
25U	TIN
17	STRONTIUM
50U	TELLURIUM
10U	TITANIUM
100U	THALLIUM
10U	VANADIUM
10U	YTTRIUM
13	ZINC
NA	ZIRCONIUM
0.2U	MERCURY
610	ALUMINUM
10	MANGANESE

MG/L	ANALYTICAL RESULTS
3.3	CALCIUM
1.2	MAGNESIUM
0.84	IRON
5.3	SODIUM
2.0U	POTASSIUM

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/15/94

PURGEABLE ORGANICS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84238 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32GM66

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1525 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
5.0U	CHLOROMETHANE
130J	VINYL CHLORIDE
5.0U	BROMOMETHANE
5.0U	CHLOROETHANE
5.0U	TRICHLOROFLUOROMETHANE
210J	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
50U	ACETONE
12U	CARBON DISULFIDE
5.0U	METHYLENE CHLORIDE
4.9J	TRANS-1,2-DICHLOROETHENE
54	1,1-DICHLOROETHANE
780	CIS-1,2-DICHLOROETHENE
5.0U	2,2-DICHLOROPROPANE
50U	METHYL ETHYL KETONE
5.0U	BROMOCHLOROMETHANE
5.0U	CHLOROFORM
5.0U	1,1,1-TRICHLOROETHANE
5.0U	1,1-DICHLOROPROPENE
5.0U	CARBON TETRACHLORIDE
1.2J	1,2-DICHLOROETHANE
2.0J	BENZENE
4800	TRICHLOROETHENE(TRICHLOROETHYLENE)
5.0U	1,2-DICHLOROPROPANE
5.0U	DIBROMOMETHANE
5.0U	BROMODICHLOROMETHANE

UG/L	ANALYTICAL RESULTS
5.0U	CIS-1,3-DICHLOROPROPENE
3.0J	METHYL ISOBUTYL KETONE
9.8	TOLUENE
5.0U	TRANS-1,3-DICHLOROPROPENE
5.0U	1,1,2-TRICHLOROETHANE
2.6J	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
5.0U	1,3-DICHLOROPROPANE
12U	METHYL BUTYL KETONE
5.0U	DIBROMOCHLOROMETHANE
120J	CHLOROBENZENE
5.0U	1,1,1,2-TETRACHLOROETHANE
5.0U	ETHYL BENZENE
4.9J	(M- AND/OR P-)XYLENE
4.3J	O-XYLENE
5.0U	STYRENE
5.0U	BROMOFORM
5.0U	BROMOBENZENE
5.0U	1,1,2,2-TETRACHLOROETHANE
5.0U	1,2,3-TRICHLOROPROPANE
2.4J	O-CHLOROTOLUENE
5.0U	P-CHLOROTOLUENE
27	1,3-DICHLOROBENZENE
54	1,4-DICHLOROBENZENE
74	1,2-DICHLOROBENZENE

REMARKS

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FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYS.. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/15/94

MISCELLANEOUS PURGEABLE ORGANICS - DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84238 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32GM66 COLLECTION START: 03/09/94 1525 STOP: 00/00/00

ANALYTICAL RESULTS UG/L

1.2J	1.3.5-TRIMETHYLBENZENE
6.1	1.2.4-TRIMETHYLBENZENE
20JN	DICHLOROTRIFLUOROETHANE

***FOOTNOTES*
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT
*R-QC INDICATES THAT DATA UNUSABLE. COMPOUND MAY OR MAY NOT BE PRESENT. RESAMPLING AND REANALYSIS IS NECESSARY FOR VERIFICATION.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/28/94

PROJECT NO. 94-0323 SAMPLE NO. 84238 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA SI: FL
STATION ID: 32GM66 COLLECTION START: 03/09/94 1525 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
1.2J	(3-AND/OR 4-)METHYLPHENOL
10U	1,2,4-TRICHLOROBENZENE
10U	2,2'-CHLOROISOPROPYLETHER
10U	2,3,4,6-TETRACHLOROPHENOL
10U	2,4,5-TRICHLOROPHENOL
10U	2,4,6-TRICHLOROPHENOL
10U	2,4-DICHLOROPHENOL
75	2,4-DIMETHYLPHENOL
20u	2,4-DINITROPHENOL
1011	2,4-DINITROTOLUENE
10U	2,6-DINITROTOLUENE
10U	2-CHLORONAPHTHALENE
10U	2-CHLOROPHENOL
20u	2-METHYL-4,6-DINITROPHENOL
5.0J	2-METHVLNAPHTHALENE
5.6J	2-METHYLPHENOL
10U	2-NITROANILINE
10U	2-NITROPHENOL
10U	3,3'-DICHLOROBENZIDINE
10U	3-NIIROANILINE
10U	4-BROMOPHENYL PHENVL ETHER
10U	4-CHLORO-3-METHYLPHENOL
10U	4-CHLOROANILINE
10U	4-CHLOROPHENYL PHENYL ETHER
10U	4-NITROANILINE
20u	4-NITROPHENOL
2.3J	ACENAPHTHENE
10U	ACENAPHTHVLENE
10U	ANTHRACENE
10U	BENZO(A)ANTHRACENE
10U	BENZO(B AND/OR K)FLUORANTHENE

UG/L	ANALYTICAL RESULTS
10U	BENZO(GHI)PERYLENE
10U	BENZO-A-PYRENE
10U	BENZYL BUTYL PHTHALATE
10U	BIS(2-CHLOROETHOXY) METHANE
10U	BIS 2-CHLOROETHYL) ETHER
10U	BIS(2-ETHYLHEXYL) PHTHALATE
1.4J	CARBAZOLE
10U	CHRYSENE
10U	DI N-BUTVLPHTHALATE
1011	DI-N-OCTYLPHTHALATE
10U	DIBENZO(A,H)ANTHRACENE
1.4J	DIBENZOFURAN
10U	DIETHYL PHTHALATE
10U	DIMETHVL PHTHALATE
10U	FLUORANTHENE
1.7J	FLUORENE
10U	HEXACHLOROBENZENE (HCB)
10U	HEXACHLOROBUTADIENE
10U	HEXACHLOROCYCLOPENTADIENE (HCCP)
10U	HEXACHLOROETHANE
10U	INDENO (1,2,3-CD) PYRENE
10U	ISOPHORONE
10U	N-NITROSODI-N-PROPYLAMINE
10U	N-NITROSODIPHENYLAMINE/DIPHENYLAMINE
53	NAPHTHALENE
10U	NITROBENZENE
20u	PENTACHLORDPHENOL
10U	PHENANTHRENE
10U	PHENOL
10U	PVRENE

REMARKS

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/28/94

MISCELLANEOUS EXTRACTABLE COMPOUNDS - DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84238 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PLNSACOLA ST: FL
STATION ID: 32GM66 COLLECTION START: 03/09/94 1525 STOP: 00/00/00

ANALVTICAL RCSULTS UG/L

70JN	DIMETHYLPHENOL (NOT 2.4) (2 ISOMERS)
10JN	ETHVLDIMETHYLBENZLNE
10JN	TRIMETHYLPHENOL
2DJN	(DIMETHYLETHYL)PHENOL
10J	1 UNIDENTIFIED COMPOUND

FOOTNOTES
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*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.
*R-QC INDICATES THAT DATA UNUSABLE. COMPOUND MAY OR MAY NOT BE PRESENT. RESAMPLING AND REANALYSIS IS NECESSARY FOR VERIFICATION

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 03/31/94

PROJECT NO. 94-0323 SAMPLE NO. 84238 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32GM66

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PLNSACOLA ST: FL
COLLECTION START: 03/09/94 1525 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
0.50U	ALDRIN
0.50U	HEPTACHLOR
0.50U	HEPTACHLOR EPOXIDE
0.50U	ALPHA-BHC
0.50U	BETA-BHC
0.50U	GAMMA-BHC (LINDANE)
0.50U	DELTA-BHC
0.50U	ENDOSULFAN I (ALPHA)
0.50U	DIELDRIN
0.50U	4,4'-DDT (P,P'-DDT)
0.50U	4,4'-DDE (P,P'-DDE)
0.50U	4,4'-DDD (P,P'-DDD)
0.50U	ENDRIN
0.50U	ENDOSULFAN II (BETA)
0.50U	ENDOSULFAN SULFATE
1.0U	CHLORDANE (TECH. MIXTURE) /1
2.0U	PCB-1242 (AROCLOR 1242)
2.0U	PCB-1254 (AROCLOR 1254)
2.0U	PCB-1221 (AROCLOR 1221)

UG/L	ANALYTICAL RESULTS
2.0U	PCO-1232 (AROCLOR 1232)
2.0U	PCB-1248 (AROCLOR 1248)
2.0U	PCB-1260 (AROCLOR 1260)
2.0U	PCB-1016 (AROCLOR 1016)
20U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
1.0U	METHOXYCHLOR
0.50U	ENDRIN KETONE

*+*REMARKS***

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE OIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT. C-CONFIRMED BY GC/MS
1. WHEN NO VALUE IS REPORTED. SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84238 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32GM66 COLLECTION START: 03/09/94 1525 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
10U	UG/L	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALY. MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84238 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32GW66 COLLECTION START: 03/09/94 1525 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
100U	SILVER
300U	ARSENIC
NA	BORON
100U	BARIUM
84	BERYLLIUM
50U	CADMIUM
100U	COBALT
370	CHROMIUM
100U	COPPER
100U	MOLYBDENUM
200U	NICKEL
5.0U	LEAD
300u	ANTIMONY
4wu	SELENIUM
250U	TIN
2200	STRONTIUM
500U	TELLURIUM
100U	TITANIUM
1000U	THALLIUM
550	VANADIUM
1300	YTTRIUM
530	ZINC
NA	ZIRCONIUM
0.2U	MERCURY
1000000	ALUMINUM
2000	MANGANESE

MG/L	ANALYTICAL RESULTS
240	CALCIUM
260	MAGNESIUM
420	IRON
670	SODIUM
110	POTASSIUM

REMARKS

REMARKS

FOOTNOTES
•A-AVERAGE VALUE •NA-NOT ANALYZED •NAI-INTERFERENCES •J-ESTIMATED VALUE •N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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•U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

EXTRACTABLE ORGANICS DATA REPORT 03/28/94

PROJECT NO. 94-0323 SAMPLE NO. 84239 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32G02 COLLECTION START: 03/09/94 1530 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10U	(3-AND/OR 4-)METHYLPHENOL
10U	1,2,4-TRICHLOROBENZENE
10U	2,2'-CHLOROISOPROPYLETHER
10U	2,3,4,6-TETRACHLOROPHENOL
10U	2,4,5-TRICHLOROPHENOL
10U	2,4,6-TRICHLOROPHENOL
10U	2,4-DICHLOROPHENOL
10U	2,4-DIMETHYLPHENOL
20U	2,4-DINITROPHENOL
10U	2,4-DINITROTOLUENE
10U	2,6-DINITROTOLUENE
10U	2-CHLORONAPHTHALENE
10U	2-CHLOROPHENOL
20U	2-METHYL-4,6-DINITROPHENOL
10U	2-METHYLNAPHTHALENE
10U	2-METHYLPHENOL
10U	2-NITROANILINE
10U	2-NITROPHENOL
10U	3,3'-DICHLOROBENZIDINE
10U	3-NITROANILINE
10U	4-BROMOPHENYL PHENYL ETHER
10U	4-CHLORO-3-METHYLPHENOL
10U	4-CHLOROANILINE
10U	4-CHLOROPHENYL PHENYL ETHER
10U	4-NITROANILINE
20U	4-NITROPHENOL
10U	ACENAPHTHENE
10U	ACENAPHTHVLENE
10U	ANTHRACENE
10U	BENZO(A)ANTHRACENE
10U	BENZO(B AND/OR KIFLUORANTHENE

REMARKS

UG/L	ANALYTICAL RESULTS
10U	BENZO(GHI)PERYLENE
10U	BENZO-A-PYRENE
10U	BENZYL BUTYL PHTHALATE
10U	BIS(2-CHLOROETHOXY) METHANE
10U	BIS(2-CHLOROETHYL) ETHER
10U	BIS(2-ETHYLHEXYL) PHTHALATE
10U	CARBAZOLE
10U	CHRYSENE
10U	DI-N-BUTYLPHTHALATE
1011	DI-N-OCTYLPHTHALATE
10U	DIBENZO(A,H)ANTHRACENE
10U	DIBENZOFURAN
10U	DIETHYL PHTHALATE
10U	DIMETHYL PHTHALATE
10U	FLUORANTHENE
10U	FLUORENE
10U	HEXACHLOROBENZENE (HCB)
10U	HEXACHLOROBUTADIENE
10U	HEXACHLOROCYCLOPENTADIENE (HCCP'I
10U	HEXACHLOROE1HANE
10U	INDENO (1,2,3-CD) PYRENE
10U	ISOPHORONE
10U	N-NITROSODI-N-PROPYLAMINE
10U	N-NITROSODIPHENYLAMINE/DIPHENYLAM'INE
10U	NAPHTHALENE
10U	NITROBENZENE
20U	PENTACHLOROPHENOL
10U	PHENANTHRENE
10U	PHENOL
10U	PVRENE

REMARKS

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PESTICIDES/PCB'S DATA REPORT 03/31/94

PROJECT NO. 94-0323 SAMPLE NO. 84239 SAMPLE TYPE: GROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G02

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1530 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
0.50U	ALDRIN
0.50U	HEPTACHLOR
0.50U	HEPTACHLOR EPOXIDE
0.50U	ALPHA-BHC
0.50U	BETA-BHC
0.50U	GAMMA-BHC (LINDANE)
0.50U	DELTA-BHC
0.50U	ENDOSULFAN I (ALPHA)
0.50U	DIELDRIN
0.50U	4,4'-DDT (P,P'-DDT)
0.50U	4,4'-DDE (P,P'-DDE)
0.50U	4,4'-DDD (P,P'-DDD)
0.50U	ENDRIN
0.50U	ENDOSULFAN II (BETA)
0.50U	ENDOSULFAN SULFATE
1.0U	CHLORDANE (TECH. MIXTURE) /1
2.0U	PCB-1242 (AROCLOR 1242)
2.0U	PCB-1254 (AROCLOR 1254)
2.0U	PCB-1221 (AROCLOR 1221)

UG/L	ANALYTICAL RESULTS
2.0U	PCB-1232 (AROCLOR 1232)
2.0U	PCB-1248 (AROCLOR 1248)
2.0U	PCB-1260 (AROCLOR 1260)
2.0U	PCB-1016 (AROCLOR 1016)
20U	TOXAPHENE
--	CHLORDENE /2
--	ALPHA-CHLORDENE /2
--	BETA CHLORDENE /2
--	GAMMA-CHLORDENE /2
--	GAMMA-CHLORDANE /2
--	TRANS-NONACHLOR /2
--	ALPHA-CHLORDANE /2
--	CIS-NONACHLOR /2
--	OXYCHLORDANE (OCTACHLOREPOXIDE) /2
1.0U	METHOXYCHLOR
0.50U	ENDRIN KETONE

.**REMARKS*** ***REMARKS***

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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1. WHEN NO VALUE IS REPORTED, SEE CHLORDANE CONSTITUENTS. 2. CONSTITUENTS OR METABOLITES OF TECHNICAL CHLORDANE.

SAMPLE AND ANALYSIS MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

03/30/94

SPECIFIED ANALYSIS DATA REPORT

PROJECT NO. 94-0323 SAMPLE NO. 84239 SAMPLE TYPE: GROUNDWA PROG ELEM: SSF COLLECTED BY: F SLOAN
SOURCE: NAS PENSACOLA CITY: PENSACOLA ST: FL
STATION ID: 32G02 COLLECTION START: 03/09/94 1530 STOP: 00/00/00

RESULTS	UNITS	PARAMETER
40	UG/L	CYANIDE

FOOTNOTES
*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
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SAMPLE AND ANALY. MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

METALS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84239 SAMPLE TYPE; CROUNDWA
SOURCE: NAS PENSACOLA
STATION ID: 32G02

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/09/94 1530 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
10U	SILVER
	ARSENIC
	BORON
30	BARIUM
5.0U	BERYLLIUM
5.0U	CADMIUM
10U	COBALT
10U	CHROMIUM
10U	COPPER
10U	MOLYBDENUM
20U	NICKEL
5.0U	LEAD
30U	ANTIMONY
40U	SELENIUM
25U	TIN
190	STRONTIUM
50U	TELLURIUM
10U	TITANIUM
100U	THALLIUM
10U	VANADIUM
10U	YTTRIUM
10U	ZINC
NA	ZIRCONIUM
0.2U	MERCURY
290	ALUMINUM
69	MANGANESE

MG/L	ANALYTICAL RESULTS
54	CALCIUM
1.7	MAGNESIUM
1.6	IRON
8.6	SODIUM
2.0U	POTASSIUM

REMARKS

REMARKS

FOOTNOTES
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OK-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
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SAMPLE AND ANAL.,.S MANAGEMENT SYSTEM
EPA-REGION IV ESD. ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 03/15/94

PROJECT NO. 94-0323 SAMPLE NO. 84240 SAMPLE TYPE: BLKWA
SOURCE: NAS PENSACOLA
STATION ID: TB

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/10/94 0700 STOP: 00/00/00

UG/L	ANALYTICAL RESULTS
5.OU	CHLOROMETHANE
5.OU	VINYL CHLORIDE
5.OU	BROMOMETHANE
5.OU	CHLOROETHANE
5.OU	TRICHLOROFLUOROMETHANE
5.OU	1,1-DICHLOROETHENE(1,1-DICHLOROETHYLENE)
50U	ACETONE
12u	CARBON DISULFIDE
5.OU	METHYLENE CHLORIDE
5.OU	TRANS-1.2-DICHLOROETHENE
5.OU	1,1-DICHLOROETHANE
5.OU	CIS-1,2-DICHLOROETHENE
5.OU	2,2-DICHLOROPROPANE
50U	METHYL ETHYL KETONE
5.OU	BROMOCHLOROMETHANE
5.OU	CHLOROFORM
5.OU	1,1,1-TRICHLOROETHANE
5.OU	1,1-DICHLOROPROPENE
5.OU	CARBON TETRACHLORIDE
5 OU	1,2-DICHLOROETHANE
5.OU	BENZENE
5.OU	TRICHLOROETHENE(TRICHLOROETHYLENE)
5.OU	1,2-DICHLOROPROPANE
5.OU	DIBROMOMETHANE
5.OU	BROMODICHLOROMETHANE

UG/L	ANALYTICAL RESULTS
5.OU	CIS-1.3-DICHLOROPROPENE
12u	METHYL ISOBUTYL KETONE
5.OU	TOLUENE
5.OU	TRANS-1.3-DICHLOROPROPENE
5.OU	1,1,2-TRICHLOROETHANE
5.OU	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
5.OU	1,3-DICHLOROPROPANE
12u	METHYL BUTYL KETONE
5.OU	DIBROMOCHLOROMETHANE
5.OU	CHLOROBENZENE
5.OU	1,1,1,2-TETRACHLOROETHANE
5.OU	ETHYL BENZENE
5.OU	(M- AND/OR P-)XYLENE
5.OU	O-XYLENE
5.OU	STYRENE
5.OU	BROMOFORM
5.OU	BROMOBENZENE
5.OU	1.1.2.2-TETRACHLOROETHANE
5.OU	1.2.3-TRICHLOROPROPANE
5.OU	O-CHLOROTOLUENE
5.OU	P-CHLOROTOLUENE
5.OU	1.3-DICHLOROBENZENE
5.OU	1.4-DICHLOROBENZENE
5.OU	1.2-DICHLOROBENZENE

REMARKS

REMARKS

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*A-AVERAGE VALUE *NA-NOT ANALYZED *NAI-INTERFERENCES *J-ESTIMATED VALUE *N-PRESUMPTIVE EVIDENCE OF PRESENCE OF MATERIAL
*K-ACTUAL VALUE IS KNOWN TO BE LESS THAN VALUE GIVEN *L-ACTUAL VALUE IS KNOWN TO BE GREATER THAN VALUE GIVEN
*U-MATERIAL WAS ANALYZED FOR BUT NOT DETECTED. THE NUMBER IS THE MINIMUM QUANTITATION LIMIT.

SAMPLE AND ANALY. . MANAGEMENT SYSTEM
EPA-REGION IV ESD, ATHENS, GA.

PURGEABLE ORGANICS DATA REPORT 04/05/94

PROJECT NO. 94-0323 SAMPLE NO. 84241 SAMPLE TYPE: SEDIMBLK
SOURCE: NAS PENSACOLA
STATION ID: TB2

PROG ELEM: SSF COLLECTED BY: F SLOAN
CITY: PENSACOLA ST: FL
COLLECTION START: 03/10/94 0710 STOP: 00/00/00

UG/KG	ANALYTICAL RESULTS
31U	CHLOROMCTHANC
31U	VINYL CHLORIDE
31U	BROMOMETHANE
31U	CHLOROETHANE
31U	TRICHLOROFLUOROMETHANE
31U	1.1-DICHLOROETHENE(1.1-DICHLOROETHYLENE)
310U	ACETONE
77u	CARBON DISULFIDE
31U	METHYLENE CHLORIDE
31U	TRANS-1.2-DICHLOROETHENE
31U	1.1-DICHLOROETHANE
31U	CIS-1.2-DICHLOROETHENE
31U	2.2-DICHLOROPROPANE
310U	METHYL ETHYL KETONE
31U	BROMOCHLOROMETHANE
31U	CHLOROFORM
31U	1.1.1-TRICHLOROETHANE
31U	1.1-DICHLOROPROPENE
31U	CARBON TETRACHLORIDE
31U	1.2-DICHLOROETHANE
31U	BENZENE
31U	TRICHLOROETHENE(TRICHLOROETHYLENE)
31U	1.2-DICHLOROPROPANE
31U	DIBROMOMETHANE
31U	BROMODICHLOROMETHANE

UG/KG	ANALYTICAL RESULTS
31U	CIS-1.3-DICHLOROPROPENC
77u	METHYL ISOBUTYL KETONE
31U	TOLUENE
31U	TRANS-1.3-DICHLOROPROPENE
31U	1.1.2-TRICHLOROETHANE
31U	TETRACHLOROETHENE(TETRACHLOROETHYLENE)
31U	1.3-DICHLOROPROPANE
77u	METHYL BUTYL KETONE
31U	DIBROMOCHLOROMETHANE
31U	CHLOROBENZENE
31U	1.1.1.2-TETRACHLOROETHANE
31U	ETHYL BENZENE
31U	(M- AND/OR P-)XYLENE
31U	O-XYLENE
31U	STYRENE
31U	BROMOFORM
31U	BROMOBENZENE
31U	1.1.2.2-TETRACHLOROETHANE
31U	1.2.3-TRICHLOROPROPANE
31U	O-CHLOROTOLUENE
31U	P-CHLOROTOLUENE
31u	1.3-DICHLOROBENZENE
31U	1.4-DICHLOROBENZENE
31U	1.2-DICHLOROBENZENE
10.0	PERCENT MOISTURE

FOOTNOTES
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SAMPLE-NUMBERING SYSTEM

The following sample-numbering system was developed for the identification of sample locations and matrices. Sample number 33G05F001 is used as an example and is explained below.

The first two characters of the sample number indicate the site number. Therefore, the 33 in 33G05001F indicates that the sample came from Site 33, which is a part of OU 10. Sample 08G05F001 would come from Site 8.

The third character indicates the matrix where the sample was collected. At OU 10, samples were collected from the soil (S), sediment (M), surface water (W), or groundwater (G). For example, sample 33G05F001 was collected from the groundwater.

The fourth and fifth characters indicate the sample station number, which may be shown on an accompanying map or figure. For example, sample 33G05F001 was collected from sample station 05. The station number is unique at a particular site.

The sixth character is used for QNQC samples only and indicates the type of QNQC sample that was collected in the field. This character will not be found for non-QNQC samples. If this character is not present, it will simply be replaced by the character which would normally follow. A character is never left blank in a sample number. Each type of QNQC sample is listed below and is explained in the next section.

D	Duplicate
E	Equipment Rinsate
M	Matrix Spike
S	Matrix Spike Duplicate
T	Trip Blank
F	Filtered
El	Equipment rinsate

For example, sample 33G05F001 is a filtered sample from a groundwater monitoring well.

The seventh and eighth characters (or sixth and seventh characters for non-QNQC samples) are always numbers. Their meaning depends on the matrix sampled. For soil, sediment and surface water it indicates the depth at which the sample was collected in feet. For groundwater, it indicates the number of the sampling round. For example, sample 33G05F002 was collected from groundwater during sampling round 2. Sample 33S05003 was collected from the soil at a depth of three feet.

Some samples will have up to three additional characters used for QA/QC samples only. These indicate a QA/QC sample created and assigned in the laboratory instead of in the field. The QNQC abbreviations are listed below and are explained in the next section.

MS	Matrix spike
MSD	Matrix Spike Duplicate
F	Filtered
DL	Dilution
RE	Sample rerun
REMS	Re-extracted matrix spike/matrix spike duplicate

Specific exceptions to this sample-numbering system are listed below:

33GGM71
33GGM78
33GGM69
33GGM13
33GGM66
33GGM69
33GGM72
33GGM76
33GGM77
33GGM78
33GR3

All of the above samples were collected from groundwater. The last four digits indicate the well number when: the sample was collected. For example, sample 33GGM71 was collected from monitoring well GM-71. These samples were collected during the first sampling round.

Other specific exceptions are described below:

33GROU	OU 10 Grout QNQC sample
33GUN	OU 10 QNQC sample from decontamination spray gun
33MUD	OU 10 drilling mud QNQC sample
33SAND	OU 10 sand filter pack QMQC sample
33BENT	OU 10 bentonite QMQC sample
NASPAF	Field blank QNQC series
TB	Trip Blank
RB	Rinsate Blank

DATA QUALIFIER AND SAMPLE IDENTIFICATION QUALIFIER DEFINITIONS

Data qualifiers are listed adjacent to the data and indicate any special circumstances involved after analysis of the sample, usually assigned during the data validation process. They are explained below:

DATA QUALIFIER DEFINITIONS

U The compound was analyzed for, but was not detected above the reported sample quantitation limit.

J The compound was positively detected, however, the reported concentration is considered to approximate the concentration within the sample.

UJ The compound was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual

limit of quantitation necessary to accurately and precisely measure the compound in the sample.

R The sample results are rejected due to serious deficiencies in the ability to analyze the Sample and meet QC criteria. The presence or absence of the compound cannot be verified.

SAMPLE IDENTIFICATION QUALIFIER DEFINITIONS

These are always listed next to the actual sample number. They refer to special procedures that were performed before the sample was analyzed.

MS Matrix Spike: The laboratory injects a known amount of a certain compound into the sample to demonstrate acceptable compound recovery by the laboratory at the time of sample analysis. It is used to help determine the accuracy of the entire analytical process.

MSD Matrix Spike Duplicate: The laboratory injects a known amount of a certain compound into the sample with the matrix spike as a means of comparing the accuracy with the matrix spike.

DL Dilution: The sample contained too high a concentration of a compound to be detected on the instrument and had to be diluted.

RE Sample rerun: If a problem was suspected during the original analysis, the sample was analyzed a second time.

REMS Re-extracted matrix spike/matrix spike duplicate: The original sample was re-extracted and had a matrix spike or a matrix spike duplicate analyzed.

APPENDIX S

SPLIT SAMPLING ANALYTICAL RESULTS — MARCH 1994

PENSACOLA, SITE 32

CT5213	SAMPLE ID ---->	032-S-0006-00	032-G-0007-00	032-S-0039-00	032-S-0067-00		
	ORIGINAL ID ----->	32S06	32G07	32S39	32S67		
	LAB SAMPLE ID --->	39535-004	39546-002	39535-005	39535-006		
	SAMPLE DATE ----->	03/08/94	03/09/94	03/08/94	03/08/94		
	UNITS ---------->	MG/KG	UG/L	MG/KG	MG/KG		
Method	Parameter						
CYANIDE	Cyanide	0 J	1. U	0 14 J	0.08 W		

CT5213	SAMPLE ID ------->	032-S-0006-00		032-G-0007-00		032-S-0039-00		032-S-0067-00			
	ORIGINAL ID ----->	32506		32G07		32S39		32S67			
	LAB SAMPLE ID --->	39535-004		39546-002		39535-005		39535-006			
	SAMPLE DATE ----->	03/08/94		03/09/94		03/08/94		03/08/94			
	UNITS ---------->	MG/KG		UG/L		MG/KG		MG/KG			
Method	Parameter										
METAL	Aluminum	1210.	J	285.		2410.	J	56.9	J		
METAL	Antimony	2.6	U	15.6	U	3.5	U	3.1	U		
METAL	Arsenic	0.40	U	2.9	U	0.76	U	0.57	U		
METAL	Barium	7.	U	4.2	U	25.2	U	0.96	U		
METAL	Beryllium	0.07	U	0.4	U	0.09	U	0.08	U		
METAL	Cadmium	1.5	J	2.1	U	7.1	J	0.41	UJ		
METAL	Calcium	5320.		7310.		928.		29.4	U		
METAL	Chromium	60.3	J	3.1		235.	J	1.8	U		
METAL	Cobalt	1.5	U	3.2	U	1.6	U	1.2	U		
METAL	Copper	6.7	J	2.6	U	23.6	J	2.8	J		
METAL	Iron	730.		327.		1030.		72.1			
METAL	Lead	30.	J	2.5		23.5	J	0.84	J		
METAL	Magnesium	72.0		482.		62.		11.2			
METAL	Manganese	30.1	J	2.2		54.7	J	0.42	U		
METAL	Mercury	0.03	U	0.1	U	0.3		0.03	U		
METAL	Nickel	2.	U	10.4	U	16.9		1.8	U		
METAL	Potassium	120.		709.		144.	U	127.	U		
METAL	Selenium	0.28	UJ	1.7	U	0.38	UJ	0.33	UJ		
METAL	Silver	0.63	U	3.6	U	6.5	J	0.71	UJ		
METAL	Sodium	44.5	U	2930.		8.9	U	4.4	U		
METAL	Thallium	0.46	U	2.8	U	0.62	U	0.55	U		
METAL	Vanadium	2.		3.3	U	2.3		0.65	U		
METAL	Zinc	19.7	J	5.3		51.	J	1.8	J		

+** Validation Complete ***

PENSACOLA, SITE 32

	SAMPLE ID ------->	032-S-0006-00		032-K-0006-00MS		032-X-0006-00MSD		032-6-0007-00		032-S-0039-00		032-S-0067-00	
	ORIGINAL ID ----->	32S06		32S06MS		32S06MSD		32607		32S39		32S67	
	LAB SAMPLE ID --->	39535-1		39535-ins		39535-1MSD		39546-1		39535-2		39535-3	
	SAMPLE DATE ----->	03/08/94						03/08/94		03/08/94		03/08/94	
	DATE EXTRACTED -->	03/10/94		03/10/94		03/10/94		03/14/94		03/10/94		03/10/94	
	DATE ANALYZED --->	03/14/94		03/15/94		03/15/94		03/15/94		03/15/94		03/15/94	
	UNITS ---------->	UG/KG		UG/KO		UG/KG		UG/L		UG/KG		UG/KG	
Method	Parameter												
PEST	alpha-BHC	1.8	U	1.7	U	1.7	U	0.06	U	1.9	U	1.7	U
PEST	beta-BHC	1.8	U	1.7	U	1.7	U	0.06	U	1.9	U	1.7	U
PEST	delta-BHC	1.8	U	1.7	U	1.7	U	0.06	U	1.9	U	1.7	U
PEST	gamma-BHC (Lindane)	1.8	U	17.		16.		0.06	U	1.9	U	1.7	U
PEST	Heptachlor	1.8	U	16.		15.		0.06	U	1.9	U	1.7	U
PEST	Aldrin	1.8	U	14.		14.		0.06	U	1.9	U	1.7	U
PEST	Heptachlor epoxide	1.8	U	1.7	U	1.7	U	0.06	U	1.9	U	1.7	U
PEST	Endosulfan I	1.8	U	1.7	U	1.7	U	0.06	U	1.9	U	1.7	U
PEST	Dicldrin	3.4	U	32.		31.		0.12	U	3.7	U	3.3	U
PEST	4,4'-DDE	3.4	U	3.4	U	3.4	U	0.12	U	3.7	U	3.3	U
PEST	Endrin	3.4	U	34.		33.		0.12	U	3.7	U	3.3	U
PEST	Endosulfan II	3.4	U	3.4	U	3.4	U	0.12	U	3.7	U	3.3	U
PEST	4,4'-DDD	3.4	U	3.4	U	3.4	U	0.12	U	3.7	U	3.3	U
PEST	Endosulfan sulfate	3.4	U	3.4	U	3.4	U	0.12	U	3.7	U	3.3	U
PEST	4,4'-DDT	3.4	U	35.		33.		0.12	U	3.7	U	3.3	U
PEST	Methoxychlor	18.	U	17.	U	17.	U	0.6	U	19.	U	17.	U
PEST	Endrin ketone	3.4	U	3.4	U	3.4	U	0.12	U	3.7	U	3.3	U
PEST	Endrin aldehyde	3.4	U	3.4	U	3.4	U	0.12	U	3.7	U	3.3	U
PEST	alpha-Chlordane	1.8	U	1.7	U	1.7	U	0.06	U	2.1		1.7	U
PEST	gamma-Chlordane	1.8	U	1.7	U	1.7	U	0.06	U	23		1.7	U
	Toxaphene	180.	U	170.	U	170.	U	6.	U	190.	U	170.	U
	Aroclor-1016	34.	U	34.	U	34.	U	1.2	U	37.	U	33.	U
	Aroclor-1221	69.	U	69.	U	68.	U	2.4	U	75.	U	68.	U
	Aroclor-1232	34.	U	34.	U	34.	U	1.2	U	37.	U	33.	U
	Aroclor-1242	34.	U	34.	U	34.	U	1.2	U	37.	U	33.	U
	Aroclor-1248	34.	U	34.	U	34.	U	1.2	U	37.	U	33.	U
	Aroclor-1254	34.	U	34.	U	34.	U	1.2	U	37.	U	33.	U
	Aroclor-1260	34.	U	34.	U	34.	U	1.2	U	35.	J	33.	U

+++ Validation Complete ***

PKNSACOLA, SITE 32

CT521	SAMPLE ID ------->	BLK-W-PBLK-67		BLK-W-PBLK-68		BLK-W-PIBK-I1		BLK-W-PIBK-I2		BLK-W-PIBK-IA		BLK-W-PIBK-IB	
	ORIGINAL ID ----->	PBLK67		PBLK68		PIBLKI1		PIBLKI2		PIBLKIA		PIBLKIB	
	LAB SAMPLE ID --->	BP2811		BP2814		PIBLKI1		PIBLKI2		PIBLKIA		PIBLKIB	
	SAMPLE DATE ----->												
	DATE EXTRACTED -->	03/10/94		03/14/94									
	DATE ANALYZED --->	03/14/94		03/15/94		03/08/94		03/09/94		13/14/94		03/14/94	
	UNITS ---------->	UG/KG		UG/L		UG/L		UG/L		JG/L		UG/L	
Method	Parameter												
PEST	alpha-BHC	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	beta-BHC	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	delta-BHC	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	gamma-BHC (Lindane)	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	Heptachlor	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	Aldrin	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	Heptachlor epoxide	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	Endosulfan I	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	Dieldrin	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	4,4'-DDE	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	Endrin	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	Endosulfan II	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	4,4'-DDD	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	Endosulfan sulfate	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	4,4'-DDT	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	Methoxychlor	17.	U	0.5	U	0.5	U	0.5	U	0.5	U	0.5	U
PEST	Endrin ketone	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	Endrin aldehyde	3.3	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
PEST	alpha-Chlordane	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	gamma-Chlordane	1.7	U	0.05	U	0.05	U	0.05	U	0.05	U	0.05	U
PEST	Toxaphene	170.	U	5.	U	5.	U	5.	U	5.	U	5.	U
PEST	Aroclor-1016	33.	U	1.	U	1.	U	1.	U	1.	U	1.	U
PEST	Aroclor-1221	67.	U	2.	U	2.	U	2.	U	2.	U	2.	U
PEST	Aroclor-1232	33.	U	1.	U	1.	U	1.	U	1.	U	1.	U
PEST	Aroclor-1242	33.	U	1.	U	1.	U	1.	U	1.	U	1.	U
PEST	Aroclor-1248	33.	U	1.	U	1.	U	1.	U	1.	U	1.	U
PEST	Aroclor-1254	33.	U	1.	U	1.	U	1.	U	1.	U	1.	U
PEST	Aroclor-1260	33.	U	1.	U	1.	U	1.	U	1.	U	1.	U

*** Validation Complete ***

CT521	SAMPLE ID ------->	BLK-W-PIBK-IC		BLK-W-PIBK-ID	
	ORIGINAL ID ----->	PIBLKIC		PIBLKID	
	LAB SAMPLE ID --->	PIBLKIC		PIBLKID	
	SAMPLE DATE ----->				
	DATE ANALYZED --->	03/15/94		03/15/94	
	UNITS ---------->	UG/L		UG/L	
Method	Parameter				
PEST	alphe-BHC	0.05	U	0.05	U
PEST	beta-BHC	0.05	U	0.05	U
PEST	delta-BHC	0.05	U	0.05	U
PEST	gamma-BHC (Lindane)	0.05	U	0.05	U
PEST	Heptachlor	0.05	U	0.05	U
PEST	Aldrin	0.05	U	0.05	U
PEST	Heptachlor epoxide	0.05	U	0.05	U
PEST	Endosulfan I	0.05	U	0.05	U
PEST	Dieldrin	0.1	U	0.1	U
PEST	4,4'-DDE	0.1	U	0.1	U
PEST	Endrin	0.1	U	0.1	U
PEST	Endosulfan II	0.1	U	0.1	U
PEST	4,4'-DDD	0.1	U	0.1	U
PEST	Endosulfan sulfate	0.1	U	0.1	U
PEST	4,4'-DDT	0.1	U	0.1	U
PEST	Methoxychlor	0.5	U	0.5	U
PEST	Endrin ketone	0.1	U	0.1	U
PEST	Endrin aldehyde	0.1	U	0.1	U
PEST	alpha-Chlordane	0.05	U	0.05	U
PEST	gamma-Chlordane	0.05	U	0.05	U
PEST	Toxaphene	5.	U	5.	U
PEST	Aroclor-1016	1.	U	1.	U
PEST	Aroclor-1221	2.	U	2.	U
PEST	Aroclor-1232	1.	U	1.	U
PEST	Aroclor-1242	1.	U	1.	U
PEST	Aralor-1248	1.	U	1.	U
PEST	Aroclor-1254	1.	U	1.	U
PEST	Aralor-1260	1.	U	1.	U

*** Validation Complete ***

CT521	SAMPLE ID ------->	032-S-0006-00		032-G-0007-00		032-S-0039-00		032-S-0067-00		032-S-0067-00MS		032-S-0067-00MSD	
	ORIGINAL ID ----->	32S06		32G07		32S39		32S67		H30903MS		H30903MSD	
	LAB SAMPLE ID --->	39535-1		59546-1		39535-2		39535-3		39535-3MS		39535-3MSD	
	SAMPLE DATE ----->	03/08/94		03/08/94		03/08/94		03/08/94		03/08/94		03/08/94	
	DATE EXTRACTED -->	03/10/94		03/14/94		03/10/94		03/10/94		03/10/94		03/10/94	
	DATE ANALYZED --->	03/17/94		03/17/94		03/17/94		03/17/94		03/17/94		03/17/94	
	UNITS ---------->	UG/KG		UG/L		UG/KG		UG/KG		UG/KG		UG/KG	
Method	Parameter												
SVOA	Phenol	340.	U	11.	U	370.	U	340.	U	1300.		1300.	
SVOA	bis(2-Chloroethyl)ether	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2-Chlorophenol	340.	U	11.	U	370.	U	340.	U	1400.		1300.	
SVOA	1,3-Dichlorobenzene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	1,4-Dichlorobenzene	340.	U	11.	U	370.	U	340.	U	830.		800.	
SVOA	1,2-Dichlorobenzene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2-Methylphenol	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2,2'-oxybis(1-Chloropropane)	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	4-Methylphenol	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	N-Nitroso-di-n-propylamine	340.	U	11.	U	370.	U	340.	U	1000.		990.	
SVOA	Hexachloroethane	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Nitrobenzene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Isophorone	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2-Nitrophenol	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2,4-Dimethylphenol	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	bis(2-Chloroethoxy)methane	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2,4-Dichlorophenol	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	1,2,4-Trichlorobenzene	340.	U	11.	U	370.	U	340.	U	1000.		1000.	
SVOA	Naphthalene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
WON	4-Chloroaniline	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Hexachlorobutadiene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	4-Chloro-3-methylphenol	340.	U	11.	U	370.	U	340.	U	1600.		1500.	
SVOA	2-Methylnaphthalene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Hexachlorocyclopentadiene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2,4,6-Trichlorophenol	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2,4,5-Trichlorophenol	820.	U	20.	U	900.	U	820.	U	810.	U	810.	U
SVOA	2-Chloronaphthalene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2-Nitroaniline	020.	U	28.	U	900.	U	820.	U	810.	U	010.	U
SVOA	Dimethylphthalate	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Acenaphthylene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2,6-Dinitrotoluene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	3-Nitroaniline	020.	U	28.	U	900.	U	820.	U	810.	U	810.	U
SVOA	Acenaphthene	340.	U	11.	U	370.	U	340.	U	1100.		1000.	
SVOA	2,4-Dinitrophenol	820.	U	28.	U	900.	U	020.	U	810.	U	810.	U
SVOA	4-Nitrophenol	820.	U	28.	U	900.	U	820.	U	2100.		1700.	
SVOA	Dibenzofuran	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	2,4-Dinitrotoluene	340.	U	11.	U	370.	U	340.	U	1200.		1000.	
SVOA	Diethylphthalate	340.	U	5.	U	370.	U	340.	U	340.	U	330.	U

*** Validation Complete ***

CT521	SAMPLE ID -------:	032-S-0006-00		032-G-0007-00		032-S-0039-00		032-S-0067-00		032-S-0067-00MS		032-S-0067-00MSD	
	ORIGINAL ID -----:	32806		32G07		32839		32S67		H30903MS		H30903MSD	
	LAB SAMPLE ID ---:	39535-1		39546-1		39535-2		3953s-3		39535-3MS		39535-3MSD	
	SAMPLE DATE -----:	03/08/94		03/08/94		03/08/94		03/08/94		03/08/94		03/08/94	
	DATE EXTRACTED --:	03/10/94		03/14/94		03/10/94		03/10/94		03/10f 94		03/10/94	
	DATE ANALYZED ---:	03/17/94		03/17/94		03/17/94		03/17/94		03/17/94		03/17/94	
	UNITS ----------:	UG/KG		UG/L		UG/KG		UG/KG		UG/KG		UG/KG	
Method	Parameter												
SVOA	4-Chlorophenyl-phenylether	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Fluorene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	4-Nitroaniline	820.	U	28.	U	900.	U	820.	U	810.	U	810.	U
SVOA	4,6-Dinitro-2-mmethylphenol	820.	U	28.	U	900.	U	820.	U	810.	U	810.	U
SVOA	N-Nitrosodiphenylamine	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	4-Bromophenyl-phenylether	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Hexachlorobenzene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Pentachlorophenol	820.	U	28.	U	900.	U	820.	U	2100.		1900.	
SVOA	Phenanthrene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Anthracene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Carbazole	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Di-n-butylphthalate	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Fluoranthene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Pyrene	340.	U	11.	U	370.	U	340.	U	1300.		1200.	
SVOA	Butylbenzylphthalatc	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	3,3'-Dichlorobenzidine	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Benzo(a)anthracene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Chrysene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	bis(2-Ethylhexyl)phthalate (BEHP)	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Di-n-octylphthalate	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Benzo(b)fluoranthene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Benzo(k)fluoranthene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Benzo(a)pyrene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Indeno(1,2,3-cd)pyrene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Dibenzo(a,h)anthracene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U
SVOA	Benzo(g,h,i)perylene	340.	U	11.	U	370.	U	340.	U	340.	U	330.	U

+++ Validation Complete +++

CT521	SAMPLE ID ------->	BLK-W-SBLK-FI		BLK-W-SBLK-FJ					
	ORIGINAL ID ----->	SBLKFI		SBLKFJ					
	LAB SAMPLE ID --->	B-A1380		B-A1377					
	SAMPLE DATE ----->								
	DATE EXTRACTED -->	03/14/94		03/10/94					
	DATE ANALYZED --->	03/17/94		03/17/94					
	UNITS ---------->	UG/L		UG/KG					
Method	Parameter								
SVOA	Phenol	10.	U	330.	U				
SVOA	bis(2-Chloroethyl)ether	10.	U	330.	U				
SVOA	2-Chlorophenol	10.	U	330.	U				
SVOA	1,3-Dichlorobenzene	10.	U	330.	U				
SVOA	1,4-Dichlorobenzene	10.	U	330.	U				
SVOA	1,2-Dichlorobenzene	10.	U	330.	U				
SVOA	2-Methylphenol	10.	U	330.	U				
SVOA	2,2'-oxybis(1-Chloropropane)	10.	U	330.	U				
SVOA	4-Methylphenol	10.	U	330.	U				
SVOA	N-Nitroso-di-n-propylamine	10.	U	330.	U				
SVOA	Hexachloroethane	10.	U	330.	U				
SVOA	Nitrobenzene	10.	U	330.	U				
SVOA	Isophorone	10.	U	330.	U				
SVOA	2-Nitrophenol	10.	U	330.	U				
SVOA	2,4-Dimethylphenol	10.	U	330.	U				
SVOA	bis(2-Chloroethoxy)methane	10.	U	330.	U				
SVOA	2,4-Dichlorophenol	10.	U	330.	U				
SVOA	1,2,4-Trichlorobenzene	10.	U	330.	U				
SVOA	Naphthalene	10.	U	330.	U				
SVOA	4-Chloroaniline	10.	U	330.	U				
SVOA	Hexachlorobutadiene	10.	U	330.	U				
SVOA	4-Chloro-3-methylphenol	10.	U	330.	U				
SVOA	2-Methylnaphthalene	10.	U	330.	U				
SVOA	Hexachlorocyclopentadiene	10.	U	330.	U				
SVOA	2,4,6-Trichlorophenol	10.	U	330.	U				
SVOA	2,4,5-Trichlorophenol	25.	U	800.	U				
SVOA	2-Chloronaphthalene	10.	U	330.	U				
SVOA	2-Nitroaniline	25.	U	800.	U				
SVOA	Dimethylphthalate	10.	U	330.	U				
SVOA	Acenaphthylene	10.	U	330.	U				
SVOA	2,6-Dinitrotoluene	10.	U	330.	U				
SVOA	3-Nitroaniline	25.	U	000.	U				
SVOA	Acenaphthene	10.	U	330.	U				
SVOA	2,4-Dinitrophenol	25.	U	800.	U				
SVOA	4-Nitrophenol	25.	U	800.	U				
SVOA	Dibenzofuran	10.	U	330.	U				
SVOA	2,4-Dinitrotoluene	10.	U	330.	U				
SVOA	Diethylphthalate	5.	U	330.	U				

*** Validation Complete ***

CT521	SAMPLE ID -------:	BLK-W-SBLK-FI		BLK-W-SBLK-FJ					
	ORIGINAL ID -----:	SBLKFI		SBLKFJ					
	LAB SAMPLE ID ---:	B-A1380		E-A1377					
	SAMPLE DATE -----:								
	DATE EXTRACTED --:	03/14/94		03/10/94					
	DATE ANALYZED ---:	03/17/94		03/17/94					
	UNITS ----------:	UG/L		UG/KG					
Method	Parameter								
SVOA	4-Chlorophenyl-phenylether	10.	U	330.	U				
SVOA	Fluorene	10.	U	330.	U				
SVOA	4-Nitroaniline	25.	U	800.	U				
SVOA	4,6-Dinitro-2-mmethylphenol	25.	U	800.	U				
SVOA	N-Nitrosodiphenylamine	10.	U	330.	U				
SVOA	4-Bromophenyl-phenylether	10.	U	330.	U				
SVOA	Hexachlorobenzene	10.	U	330.	U				
SVOA	Pentachlorophenol	2s.	U	800.	U				
SVOA	Phenanthrene	10.	U	330.	U				
SVOA	Anthracene	10.	U	330.	U				
SVOA	Carbazole	10.	U	330.	U				
SVOA	Di-n-butylphthalate	10.	U	330.	U				
SVOA	Fluoranthm	10.	U	330.	U				
SVOA	Pyrm	10.	U	330.	U				
SVOA	Butylbenzylphthalate	10.	U	330.	U				
SVOA	3,3'-Dichlorobenzidine	10.	U	330.	U				
SVOA	Benzo(a)anthracene	10.	U	330.	U				
SVOA	Chrysene	10.	U	330.	U				
SVOA	bis(2-Ethylhexyl)phthalate (BEHP)	10.	U	330.	U				
SVOA	Di-n-octylphthalate	10.	U	330.	U				
SVOA	Benzo(b)fluoranthene	10.	U	330.	U				
SVOA	Benzo(k)fluoranthene	10.	U	330.	U				
SVOA	Benzo(a)pyrene	10.	U	330.	U				
SVOA	Indeno(1,2,3-cd)pyrene	10.	U	330.	U				
SVOA	Dibenzo(a,h)anthracene	10.	U	330.	U				
SVOA	Benzo(g,h,i)perylene	10.	U	330.	U				

++ Validation Complete ++

PENSACOLA, SITE 32

CT521	SAMPLE ID ------->	032-T-0001-00		032-S-0006-00		032-S-0006-00MS		032-S-0006-00MSD		032-G-0007-00		032-S-0039-00	
	ORIGINAL ID ----->	326101		32506		32S06MS		32S06MSD		32607		32539	
	LAB SAMPLE ID --->	39546-4		3953s-7		39535-7MS		39535-7MSD		39546-3		39535-8	
	SAMPLE DATE ----->			03/08/94								03/08/94	
	DATE ANALYZED --->	03/15/94		03/15/94		03/14/94		03/14/94		03/15/94		03/15/94	
	UNITS ---------->	UG/L		UG/KG		UG/KG		UG/KG		UG/L		UG/KG	
Method	Parameter												
VOA	Chloromethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Bromomethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Vinyl chloride	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Chloroethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Methylene chloride	7.	J	20.	U	12.	U	12.	U	6.	U	21.	U
VOA	Acetone	10.	U	40.	U	40.	U	51.	U	10.	U	170.	U
VOA	Carbon disulfide	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	1,1-Dichloroethylene	10.	U	10.	U	64.		62.		10.	U	11.	U
VOA	1,1-Dichloroethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	1,2-Dichloroethene	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Chloroform	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	1,2-Dichloroethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	2-Butanone (MEK)	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	1,1,1-Trichloroethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Carbon tctrrchloride	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Bromodichloromethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	1,2-Dichloropropane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	cis-1,3-Dichloropropene	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Trichloroethylene	10.	U	10.	U	56.		50.		10.	U	11.	U
VOA	Dibromochloromethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	1,1,2-Trichloroethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Benzene	10.	U	10.	U	55.		57.		10.	U	11.	U
VOA	trans-1,3-Dichloropropene	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Bromoform	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	4-Methyl-2-Pentanone (MIBK)	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	2-Hexanone	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Tetrachloroethylene	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	1,1,2,2-Tetrachloroethane	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Toluene	10.	U	10.	U	58.		58.		10.	U	11.	U
VOA	Chlorobenzene	10.	U	10.	U	57.		56.		10.	U	11.	U
VOA	Ethylbenzene	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Styrene	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U
VOA	Xylene (total)	10.	U	10.	U	11.	U	11.	U	10.	U	11.	U

*** Validation Complete ***

PENSACOLA, SITE 32

CT521	SAMPLE ID ------->	032-5-0067-00		032-T-ST01-00		VBL-T-T521-01		VBL-T-T521-02		VBL-T-T521-03	
	ORIGINAL ID ----->	32S67		32ST01		VBLKCX		VBLKDM		VBLKDN	
	LAB SAMPLE ID --->	39535-9		39535-10		BC031594B		BD031494C		BD031594A	
	SAMPLE DATE ----->	03/08/94		03/08/94							
	DATE ANALYZED --->	03/14/94		03/15/94		03/15/94		03/14/94		03/15/94	
	UNITS ---------->	UG/KG		UG/L		UG/L		UG/KG		UG/KG	
Method	Parameter										
VOA	Chloromethane	10.	U	10.	U	10.	U	10.	U	10.	U
MIA	Bromomethane	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Vinyl chloride	10.	U	10.	U	10.	U	10.	U	10.	U
MIA	Chloroethane	10.	U	10.	U	10.	U	10.	U	10.	U
MA	Methylene chloride	9.	U	7.	U	6.	J	2.	J	4.	J
MA	Acetone	13.	U	10.	U	10.	U	7.	J	10.	J
VOA	Carbon disulfide	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	1,1-Dichloroethylene	10.	U	10.	U	10.	U	10.	U	10.	U
MIA	1,1-Dichloroethane	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	1,2-Dichloroethene	10.	U	10.	U	10.	U	10.	U	10.	U
MIA	Chloroform	10.	U	10.	U	10.	U	10.	U	10.	U
MA	1,2-Dichloroethane	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	2-Butanone (MEK)	10.	U	10.	U	10.	U	10.	U	10.	U
VIM	1,1,1-Trichloroethane	10.	U	10.	U	10.	U	10.	U	10.	U
MA	Carbon tetrachloride	10.	U	10.	U	10.	U	10.	U	10.	U
VIM	Bromodichloromethane	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	1,2-Dichloropropane	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	cis-1,3-Dichloropropene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Trichloroethylene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Dibromochloromethane	10.	U	10.	U	10.	U	10.	U	10.	U
MA	1,1,2-Trichloroethane	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Benzene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	trans-1,3-Dichloropropene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Bromoform	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	4-Methyl-2-Pentanone (MIBK)	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	2-Hexanone	10.	U	10.	U	10.	U	10.	U	10.	U
MIA	Tetrachloroethylene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	1,1,2,2-Tetrachloroethane	10.	U	10.	U	10.	U	10.	U	10.	U
MIA	Toluene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Chlorobenzene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Ethylbenzene	10.	U	10.	U	10.	U	10.	U	10.	U
VOA	Styrene	10.	U	10.	U	10.	U	10.	U	10.	U
MA	*Xylene* (total)	10.	U	10.	U	10.	U	10.	U	10.	U

+++ Validation Complete ***

APPENDIX T

DETECTED CONCENTRATIONS — QA/QC SAMPLES

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	PBLK07	PBLK08	PBLK08	PBLK10	PBLK18	PBLK22
	LAB ID:	532529		532531	532809	533393	
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
P 4,4'-DDD	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P 4,4'-DDE	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P 4,4'-DDT	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Aldrin	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Aroclor 1016	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1221	UG/KG	67.0000 U	67.0000 U	67.0000 U	67.0000 U	67.0000 U	67.0000 U
P Aroclor 1232	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1242	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1248	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1254	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1260	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Dieldrin	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endosulfan I	UG/KG	1.7000 U	1.7000 U	1.7000 U	0.0630 J	1.7000 U	1.7000 U
P Endosulfan II	UG/KG	3.3000 U	3.3000 U	3.3000 U	0.1500 J	3.3000 U	3.3000 U
P Endosulfan sulfate	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin aldehyde	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin ketone	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Heptachior	UG/KG	1.7000 U	0.0730 J	0.1500 J	0.1100 J	1.7000 U	0.0900 J
P Heptachlor epoxide	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Methoxychlor	UG/KG	17.0000 U	17.0000 U	17.0000 U	7.3000 J	17.0000 U	17.0000 U
P Toxaphene	UG/KG	170.0000 U	170.0000 U	170.0000 U	170.0000 U	170.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.7000 U	0.0710 J	1.7000 U	1.7000 U	0.0350 J	1.7000 U
P alpha-Chlordane	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P beta-BHC	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P delta-BHC	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.7000 U	0.1000 J	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P gamma-Chlordane	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	0.2000 J	1.7000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	PBLK27 529994	PBLK27 533532	PBLK29 533694	PBLK30 533691	PBLK32 533891	PBLK39 530274
P 4,4'-DDD	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P 4,4'-DDE	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.2000 U	0.0510 J	3.3000 U
P 4,4'-DDT	UG/KG	0.1300 J	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P Aldrin	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Aroclor 1016	UG/KG	33.0000 U	33.0000 U	33.0000 U	32.0000 U	33.0000 U	33.0000 U
P Aroclor 1221	UG/KG	67.0000 U	67.0000 U	67.0000 U	66.0000 U	67.0000 U	67.0000 U
P Aroclor 1232	UG/KG	33.0000 U	33.0000 U	33.0000 U	32.0000 U	33.0000 U	33.0000 U
P Aroclor 1242	UG/KG	33.0000 U	33.0000 U	33.0000 U	32.0000 U	33.0000 U	33.0000 U
P Aroclor 1248	UG/KG	33.0000 U	33.0000 U	33.0000 U	32.0000 U	33.0000 U	33.0000 U
P Aroclor 1254	UG/KG	33.0000 U	33.0000 U	33.0000 U	32.0000 U	33.0000 U	33.0000 U
P Aroclor 1260	UG/KG	33.0000 U	33.0000 U	33.0000 U	32.0000 U	33.0000 U	33.0000 U
P Dieldrin	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P Endosulfan I	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Endosulfan II	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P Endosulfan sulfate	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P Endrin	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P Endrin aldehyde	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P Endrin ketone	UG/KG	0.0910 J	3.3000 U	3.3000 U	3.2000 U	3.3000 U	3.3000 U
P Heptachlor	UG/KG	0.1800 J	0.0690 J	0.0900 J	1.7000 U	0.1700 J	0.0690 J
P Heptachlor epoxide	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Methoxychlor	UG/KG	17.0000 U	17.0000 U	17.0000 U	17.0000 U	17.0000 U	17.0000 U
P Toxaphene	UG/KG	170.0000 U	170.0000 U	170.0000 U	170.0000 U	170.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P alpha-Chlordane	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P beta-BHC	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P delta-BHC	UG/KG	1.7000 U	1.7000 U	0.0700 J	0.0410 J	1.7000 U	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P gama-Chlordane	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	0.0720 J	1.7000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	PBLK40 530287 .	PBLK43 530372 .	PBLK51 530426 .	PBLK52 530432 .	PBLK61 534658 .	PBLK61 547300 .
P 4,4'-DDD	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P 4,4'-DDE	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P 4,4'-DDT	UG/KG	0.2600 J	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P Aldrin	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U
P Aroclor 1016	UG/KG	33.0000 U	33.0000 U	3.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1221	UG/KG	67.0000 U	67.0000 U	6.0000 U	67.0000 U	67.0000 U	67.0000 U
P Aroclor 1232	UG/KG	33.0000 U	33.0000 U	3.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1242	UG/KG	33.0000 U	33.0000 U	3.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1248	UG/KG	33.0000 U	33.0000 U	3.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1254	UG/KG	33.0000 U	33.0000 U	3.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1260	UG/KG	33.0000 U	33.0000 U	3.0000 U	33.0000 U	33.0000 U	33.0000 U
P Dieldrin	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P Bndoeulfan I	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U
P Endosulfan II	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P Endoeulfan sulfate	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin aldehyde	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin ketone	UG/KG	3.3000 U	3.3000 U	3.0000 U	3.3000 U	3.3000 U	3.3000 U
P Heptachlor	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	0.1500 J	0.1200 J
P Heptachlor epoxide	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U
P Methoxychlor	UG/KG	17.0000 U	17.0000 U	1.0000 U	17.0000 U	17.0000 U	17.0000 U
P Toxaphene	UG/KG	170.0000 U	170.0000 U	17.0000 U	170.0000 U	170.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U
P alpha-Chlordane	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U
P beta-BHC	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U
P delta-BHC	UG/KG	1.7000 U	1.7000 U	1.0000 U	0.0400 J	0.0350 J	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U
P gamma-Chlordane	UG/KG	1.7000 U	1.7000 U	1.0000 U	1.7000 U	1.7000 U	1.7000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	PBLK62 547190	PBLK65 547285	PBLK73 547349	PBLK80 550103	PBLK85 531365	PBLK86 550260
P 4,4'-DDD	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P 4,4'-DDE	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	0.1700 J	3.3000 U
P 4,4'-DDT	UG/KG	3.3000 U	3.3000 U	0.0790 J	0.1000 J	3.3000 U	3.3000 U
P Aldrin	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Aroclor 1016	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1221	UG/KG	67.0000 U	67.0000 U	67.0000 U	67.0000 U	67.0000 U	67.0000 U
P Aroclor 1232	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1242	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1248	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1254	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Aroclor 1260	UG/KG	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U	33.0000 U
P Dieldrin	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endosulfan I	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Endosulfan II	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endosulfan sulfate	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin aldehyde	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Endrin ketone	UG/KG	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U	3.3000 U
P Heptachlor	UG/KG	1.7000 U	1.7000 U	1.7000 U	0.1200 J	1.7000 U	1.7000 U
P Heptachlor epoxide	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P Methoxychlor	UG/KG	17.0000 U	17.0000 U	17.0000 U	17.0000 U	17.0000 U	17.0000 U
P Toxaphene	UG/KG	170.0000 U	170.0000 U	170.0000 U	170.0000 U	170.0000 U	170.0000 U
P alpha-BHC	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P alpha-Chlordane	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P beta-BHC	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P delta-BHC	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P gamma-BHC (Lindane)	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U
P gamma-Chlordane	UG/KG	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U	1.7000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in eignificant digite.

	ENSAFE ID:	SBLK02	SBLK03	SBLK06	SBLK09	SBLK09	SBLK14
	LAB ID:	SBLK02	SBLK03	SBLK06	SBLK 09	SBLK09	SBLK14
	DATE RCVD:						
NAME	WITS						
S 1,2,4-Trichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,2-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,3-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,4-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4,5-Trichlorophenol	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U
S 2,4,6-Trichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dimethylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dinitrophenol	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U
S 2,4-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,6-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Chloronaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Chlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Methylnaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Nitroaniline	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U
S 2-Nitrophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 3-Nitroaniline	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U
S 4-Bromophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chloro-3-methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chloroaniline	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Nitroaniline	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U
S 4-Nitrophenol	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U
S Acenaphthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK02	SBLK03	SBLK06	SBLK09	SBLK09	SBLK14
	LAB ID:	SBLK02	SBLK03	SBLK06	SBLK 09	SBLK09	SBLK14
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Acenaphthylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(a)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(a)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(b,k)fluoranthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(g,h,i)perylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Butylbenzylphthalate	UG/KG	95.0000 J	330.0000 U	330.0000 U	330.0000 U	330.0000 J	190.0000 J
S Carbazole	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Chrysene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Di-n-butylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Di-n-octylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dibenz(a,h)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dibenzofuran	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Diethylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 J	330.0000 U
S Dimethylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Fluoranthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Fluorene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorobutadiene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorocyclopentadiene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachloroethane	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Ieophorone	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S **N-Nitroso-di-n-propylamine**	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S N-Nitrosodiphenylamine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Naphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Nitrobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Pentachlorophenol	UG/KG	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U	800.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK02	SBLK03	SBLK06	SBLK09	SBLK09	SBLK14
	LAB ID:	SBLK02	SBLK03	SBLK06	SBLK 09	SBLK09	SBLK14
	DATE RCVD:						
NAME	WITS						
S Phenanthrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Phenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 J	330.0000 U
S bis(2-chloroethyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330,0000 U	330.0000 U	330.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK21	SBLK22	SBLK22	SBLK23	SBLK26	SBLK34
	LAB ID:	SBLK21	SBLK 22	SBLK22	SBLK23	SBLK26	SBLK34
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S 1,2,4-Trichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,2-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,3-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,4-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4,5-Trichlorophenol	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S 2,4,6-Trichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dimethylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dinitrophenol	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S 2,4-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,6-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Chloronaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Chlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Methylnaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Nitroaniline	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S 2-Nitrophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 3-Nitroaniline	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S 4-Bromophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chloro-3-methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chloroaniline	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Nitroaniline	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S 4-Nitrophenol	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S Acenaphthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK21	SBLK22	SBLK22	SBLK23	SBLK26	SBLK34
	LAB ID:	SBLK21	SBLIC 22	SBLK22	SBLK23	SBLK26	SBLK34
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Acenaphthylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(a)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(a)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(b,k)fluoranthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(g,h,i)perylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Butylbenzylphthalate	UG/KG	330.0000 U	330.0000 U	78.0000 J	330.0000 U	87.0000 J	330.0000 U
S Carbazole	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Chrysene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Di-n-butylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Di-n-octylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dibenz(a,h)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dibenzofuran	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Diethylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dimethylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Fluoranthene	UG/KG	330.0000 U	330,0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Fluorene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorobutadiene	UG/KG	330.0000 U	330,0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorocyclopentadiene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachloroethane	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Isophorone	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S N-Nitrosodiphenylamine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Naphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Nitrobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Pentachlorophenol	UG/KG	790.0000 U	790.0000 U	790.0000 U	800.0000 U	790.0000 U	790.0000 U
S Phenanthrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK21	SBLK22	SBLK22	SBLK23	SBLK26	SBLK34
	LAB ID:	SBLK21	SBLK 22	SBLK22	SBLK23	SBLK26	SBLK34
	DATE RCVD:						
NAME	UNITS						
S Phenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	330.0000 U	62.0000 J	330.0000 U	61.0000 J	330.0000 U	330.0000 U
S bis(2-chloroethyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK45	SBLK51	SBLK66	SBLK71	SBLK74	SBLK80
	LAB ID:	SBLK45			SBLK71	SBLK74	SBLK80
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S 1,2,4-Trichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,2-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,3-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 1,4-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4,5-Trichlorophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S 2,4,6-Trichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dimethylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,4-Dinitrophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S 2,4-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2,6-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Chloronaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Chlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Methylnaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 2-Nitroaniline	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S 2-Nitrophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 3-Nitroaniline	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S 4-Bromophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chloro-3-methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chloroaniline	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S 4-Nitroaniline	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S 4-Nitrophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S Acenaphthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presen ɜd at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK45	SBLK51	SBLK66	SBLK71	SBLK74	SBLK80
	LAB ID:	SBLK45			SBLK71	SBLK74	SBLK80
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Acenaphthylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(a)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(a)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(b,k)fluoranthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Benzo(g,h,i)perylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Butylbenzylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	220.0000 J
S Carbazole	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Chrysene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Di-n-butylphthalate	UG/KG	330.0000 U	330.0000 U	69.0000 J	330.0000 U	330.0000 U	330.0000 U
S Di-n-octylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dibenz(a,h)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dibenzofuran	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Diethylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Dimethylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Fluoranthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Fluorene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorobutadiene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachlorocyclopentadiene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Hexachloroethane	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Isophorone	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S N-Nitrosodiphenylamine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Naphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Nitrobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Pentachlorophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	800.0000 U	800.0000 U
S Phenanthrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK45	SBLK51	SBLK66	SBLK71	SBLK74	SBLK80
	LAB ID:	SBLK45			SBLK71	SBLK74	SBLK80
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Phenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S Pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	330.0000 U	330.0000 U	330,0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U'	330.0000 U	330.0000 U	330.0000 U
S bis(2-chloroethyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U	330.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK84	SBLK90	SBLKS7	SBLKS7	VBLK03	VBLKA5
	LAB ID:	SBLK84	SBLK90	SBLK 97	SBLK97	VBLK03	VBLKA5
	DATE RCVD:						
NAME	UNITS						
S 1,2,4-Trichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 1,2-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 1,3-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 1,4-Dichlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2,4,5-Trichlorophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U		
S 2,4,6-Trichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2,4-Dichlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2,4-Dimethylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2,4-Dinitrophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U		
S 2,4-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2,6-Dinitrotoluene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2-Chloronaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2-Chlorophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2-Methylnaphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 2-Nitroaniline	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U		
S 2-Nitrophenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 3,3'-Dichlorobenzidine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 3-Nitroaniline	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U		
S 4,6-Dinitro-2-methylphenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U		
S 4-Bromophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 4-Chloro-3-methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 4-Chloroaniline	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 4-Chlorophenyl-phenylether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 4-Methylphenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
S 4-Nitroaniline	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U		
S 4-Nitrophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U		
S Acenaphthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK84	SBLK90	SBLK97	SBLK97	VBLK03	VBLKA5
	LAB ID:	SBLK84	SBLK90	SBLK97	SBLK97	VBLK03	VBLKA5
	DATE RCM:	.	.	.	.	.	.
NAME	UNITS						
S Acenaphthylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Benzo(a)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Benzo(a)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Benzo(b,k)fluoranthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Benzo(g,h,i)perylene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Butylbenzylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Carbazole	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Chrysene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Di-n-butylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Di-n-octylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Dibenz(a,h)anthracene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Dibenzofuran	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Diethylphthalate	UG/KG	34.0000 J	330.0000 U	51.0000 J	330.0000 U	.	.
S Dimethylphthalate	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Fluoranthene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Fluorene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Hexachlorobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Hexachlorobutadiene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Hexachlorocyclopentadiene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Hexachloroethane	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Indeno(1,2,3-cd)pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Isophorone	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S N-Nitroso-di-n-propylamine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S N-Nitrosodiphenylamine	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Naphthalene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Nitrobenzene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S Pentachlorophenol	UG/KG	800.0000 U	800.0000 U	790.0000 U	790.0000 U	.	.
S Phenanthrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK84	SBLKS0	SBLKS7	SBLKS7	VBLK03	VBLKA5
	LAB ID:	SBLK84	SBLKS0	SBLK 97	SBLKS7	VBLK03	VBLKA5
	DATE RCVD:						
NAME	UNITS						
S Phenol	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	
S Pyrene	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	
S bis(2-Chloroethoxy)methane	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	.
S bis(2-Ethlyhexy)phthalate	UG/KG	330.0000 U	330.0000 U	41.0000 J	33.0000 J	.	.
S bis (2-chloroethyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U	.	
S bis(2-chloroisopropyl)ether	UG/KG	330.0000 U	330.0000 U	330.0000 U	330.0000 U		
V 1,1,1-Trichloroethane	UG/KG	.	.		.	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Acetone	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Benzene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Bromoform	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Bromomethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Chloroethane	UG/KG	.	.		.	10.0000 U	10.0000 U
V Chloroform	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Chloromethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	.	.		.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK84	SBLKS0	SBLKS7	SBLK97	VBLK03	VBLKA5
	LAB ID:	SBLK84	SBLKS0	SBLK 97	SBLK97	VBLK03	VBLKA5
	DATE RCVD:						
NAME	UNITS						
V Ethylbenzene	UG/KG		.	.	.	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	.	.	.	.	12.0000 J	13.0000 J
V Styrene	UG/KG		.	.	.	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Toluene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Trichloroethene	UG/KG		.	.	.	10.0000 U	10.0000 U
V Vinyl Chloride	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG		.	.	.	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of ppendix. Data not shown in significant d gits.

	ENSAFE ID:	VBLKA7	VBLKA7	VBLKA8	VBLKA8	VBLKB1	VBLKB2
	LAB ID:		VBLKA7	VBLK A8	VBLKA8	VBLKB1	VBLKB2
	DATE RCVD:						
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/KG	23.0000 J	23.0000 J	10.0000 U	13.0000 J	10.0000 U	14.0000 J
V Benzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/KG	1.0000 J	1.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	18.0000 J	18.0000 J	7.0000 J	5.0000 J	4.0000 J	17.0000 J
V Styrene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKA7	VBLKA7	VBLKA8	VBLKA8	VBLKB1	VBLKB2
	LAB ID:		VBLKA7	VBLK A8	VBLKA8	VBLKB1	VBLKB2
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	VBLKC3	VBLKC3 / VBLKC3	VBLKC8 / VBLKC8	VBLKCZ	VBLKD4	VBLKD6 / VBLKD6
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Acetone	UG/KG	22.0000 J	22.0000 J	10.0000 U	1200.0000 U	2100.0000 J	10.0000 U
V Benzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Chloroform	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Methylene Chloride	UG/KG	21.0000 J	21.0000 J	26.0000 J	140.0000 J	610.0000 J	28.0000 J
V Styrene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKC3	VBLKC3	VBLKC8	VBLKCZ	VBLKD4	VBLKD6
	LAB ID:		VBLKC3	VBLKC8			VBLKD6
	DATE RCVD:						
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	1200.0000 U	1200.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKE2	VBLKE6	VBLKF1	VBLKF2	VBLKG4	VBLKG9
	LAB ID:		VBLKE6	VBLKF1	VBLKF2	VBLKG4	VBLKG9
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10,0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/KG	8.0000 J	11.0000 J	19.0000 J	38.0000 J	14.0000 J	16.0000 J
V Benzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	16.0000 J	12.0000 J	16.0000 J	22.0000 J	17.0000 J	23.0000 J
V Styrene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKE2	VBLKE6	VBLKF1	VBLKF2	VBLKG4	VBLKG9
	LAB ID:		VBLKE6	VBLKF1	VBLKF2	VBLKG4	VBLKG9
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10,0000 U	10.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKH2		VBLKH3		VBLKH9		VBLKI3		VBLKI3		VBLKI7	
	LAB ID:	VBLKH2		VBLKH3		VBLKH9		VBLK I3		VBLKI3		VBLKI7	
	DATE RCVD:												
NAME	UNITS												
V 1,1,1-Trichloroethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 1,1,2-Trichloroethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 1,1-Dichloroethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 1,1-Dichloroethene	UG/KG	10.0000	3	10.0000	5	1200.0000	3	10.0000	5	10.0000	3	10.0000	3
V 1,2-Dichloroethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 1,2-Dichloroethene (total)	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 1,2-Dichloropropane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 2-Butanone	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 2-Hexanone	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V 4-Methyl-2-Pentanone	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Acetone	UG/KG	10.0000	3	10.0000	3	1200.0000	3	11.0000	3	11.0000	3	10.0000	3
V Benzene	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	1	10.0000	1	10.0000	3
V Bromodichloromethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Bromoform	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Bromomethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Carbon Disulfide	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Carbon Tetrachloride	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Chlorobenzene	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Chloroethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Chloroform	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Chloromethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Dibromochloromethane	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	3	10.0000	3
V Ethylbenzene	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	3	10.0000	1	10.0000	3
V Methylene Chloride	UG/KG	3.0000	3	2.0000	3	540.0000	3	5.0000	3	5.0000	3	8.0000	3
V Styrene	UG/KG	10.0000	1	10.0000	1	1200.0000	1	10.0000	3	10.0000	3	10.0000	1
V Tetrachloroethene	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	1	10.0000	3	10.0000	3
V Toluene	UG/KG	10.0000	3	10.0000	3	1200.0000	3	10.0000	1	10.0000	1	10.0000	3
			3		3		3		3		3		3
									3		3		
									3		3		

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKH2	VBLKH3	VBLKH9	VBLKI3	VBLKI3	VBLKI7
	LAB ID:	VBLKH2	VBLKH3	VBLKH9	VBLK I3	VBLKI3	VBLKI7
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKJ4	VBLKJ9	VBLKL3	VBLKL3	VBLKL5	VBLKP4
	LAB ID:	VBLKJ4	VBLKJ9	VBLK L3	VBLKL3	VBLKL5	VBLKP4
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	4.0000 J
V Benzene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	27.0000 J	290.0000 J	30.0000 J	30.0000 J	2.0000 J	3.0000 J
V Styrene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKJ4	VBLKJ9	VBLKL3	VBLKL3	VBLKL5	VBLKP4
	LAB ID:	VBLKJ4	VBLKJ9	VBLK L3	VBLKL3	VBLKL5	VBLKP4
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	1200.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	VBLKP5 / VBLKP5	VBLKR6 / VBLKRC	VBLKR9 / VBLKR9	VBLKW5 / VBLKW5	VBLKW6 / VBLKW6	VBLKW9 / VBLKW9
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/KG	11.0000 J	10.0000 J	15.0000 J	10.0000 U	10.0000 U	12.0000 J
V Benzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	9.0000 J	3.0000 J	4.0000 J	5.0000 J	6.0000 J	9.0000 J
V Styrene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10,0000 U	10.0000 U	10.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKP5	VBLKR6	VBLKR9	VBLKW5	VBLKW6	VBLKW9
	LAB ID:	VBLKP5	VBLKR6	VBLKR9	VBLKW5	VBLKW6	VBLKW9
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1.3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKX1	VBLKX4	VBLKX8	VBLKY3	VBLKY9	VBLKY9
	LAB ID:	VBLKX1	VBLKX4		VBLKY3		VBLKY9
	DATE RCVD:						
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 2-Butanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 2-Hexanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Acetone	UG/KG	21.0000 J	7.0000 J	21.0000 J	10.0000 U	18.0000 J	1200.0000 U
V Benzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Bromoform	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Bromomethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Chlorobenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Chloroethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Chloroform	UG/KG	10.0000 U	10.0000 U	2.0000 J	10.0000 U	2.0000 J	1200.0000 U
V Chloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Ethylbenzene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Methylene Chloride	UG/KG	4.0000 J	2.0000 J	13.0000 J	6.0000 J	36.0000 J	1200.0000 U
V Styrene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Toluene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKX1	VBLKX4	VBLKX8	VBLKY3	VBLKY9	VBLKY9
	LAB ID:	VBLKX1	VBLKX4		VBLKY3		VBLKY9
	DATE RCVD:						
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Vinyl Chloride	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V Xylene (total)	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1200.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	ENSAFE ID: VBLKZ2 LAB ID: VBLKZ2 DATE RCVD:						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U						
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U						
V 1,1,2-Trichloroethane	UG/KG	10.0000 U						
V 1,1-Dichloroethane	UG/KG	10.0000 U						
V 1,1-Dichloroethene	UG/KG	10.0000 U						
V 1,2-Dichloroethane	UG/KG	10.0000 U						
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U						
V 1,2-Dichloropropane	UG/KG	10.0000 U						
V 2-Butanone	UG/KG	10.0000 U						
V 2-Hexanone	UG/KG	10.0000 U						
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U						
V Acetone	UG/KG	25.0000 J						
V Benzene	UG/KG	10.0000 U						
V Bromodichloromethane	UG/KG	10.0000 U						
V Bromoform	UG/KG	10.0000 U						
V Bromomethane	UG/KG	10.0000 U						
V Carbon Disulfide	UG/KG	10.0000 U						
V Carbon Tetrachloride	UG/KG	10.0000 U						
V Chlorobenzene	UG/KG	10.0000 U						
V Chloroethane	UG/KG	10.0000 U						
V Chloroform	UG/KG	10.0000 U						
V Chloromethane	UG/KG	10.0000 U						
V Dibrmochloromethane	UG/KG	10.0000 U						
V Ethylbenzene	UG/KG	10.0000 U						
V Methylene Chloride	UG/KG	11.0000 J						
V Styrene	UG/KG	10.0000 U						
V Tetrachloroethene	UG/KG	10.0000 U						
V Toluene	UG/KG	10.0000 U						

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Dsta not shown in significant digits.

	ENSAFE ID:	VBLKZ2					
	LAB ID:	VBLKZ2					
	DATE RCVD:		.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/KG	10.0000 U	.	.	.	.	
V Vinyl Chloride	UG/KG	10.0000 U	.		.	.	
V Xylene (total)	UG/KG	10.0000 U		.	.	.	
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	.	.	.	.	
V trans-1,3-Dichloropropene	UG/KG	10.0000 U		.	.	.	

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	BS 520459	BS 520463	BS 530846	BS 534055	BS 546691	BSD 520460
ENSAFE ID:							
LAB ID:							
DATE RCVD:							
S 1,2,4-Trichlorobenzene	UG/L	.	45.0000	38.0000	25.0000	44.0000	.
S 1,2-Dichlorobenzene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 1,3-Dichlorobenzene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 1,4-Dichlorobenzene	UG/L	.	35.0000	31.0000	18.0000	41.0000	.
S 2,4,5-Trichlorophenol	UG/L	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.
S 2,4,6-Trichlorophenol	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 2,4-Dichlorophenol	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 2,4-Dimethylphenol	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 2,4-Dinitrophenol	UG/L	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.
S 2,4-Dinitrotoluene	UG/L	.	34.0000	37.0000	29.0000	42.0000	.
S 2,6-Dinitrotoluene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 2-Chloronaphthalene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 2-Chlorophenol	UG/L	.	55.0000	61.0000	33.0000	55.0000	.
S 2-Methylnaphthalene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 2-Methylphenol	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 2-Nitroaniline	UG/L	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.
S 2-Nitrophenol	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 3-Nitroaniline	UG/L	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	.	68.0000	76.0000	59.0000	68.0000	.
S 4-Chloroaniline	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 4-Methylphenol	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S 4-Nitroaniline	UG/L	.	25.0000 U	1.0000 J	25.0000 U	25.0000 U	.
S 4-Nitrophenol	UG/L	.	54.0000	67.0000	53.0000	66.0000	.
S Acenaphthene	UG/L	.	41.0000	40.0000	30.0000	46.0000	.

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	BS	BS	BS	BS	BS	BSD
	LAB ID:	520459	520463	530846	534055	546691	520460
	DATE RCVD:	.	.	.	.	.	
NAME	UNITS						
S Acenaphthylene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Anthracene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Benzo(a)anthracene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Benzo(a)pyrene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Benzo(b,k)fluoranthene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
S Benzo(g,h,i)perylene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Butylbenzylphthalate	UG/L	.	10.0000 U	10.0000 U	72.0000 U	10.0000 U	.
S Carbazole	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Chrysene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Di-n-butylphthalate	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Di-n-octylphthalate	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Dibenzofuran	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Diethylphthalate	UG/L	.	10.0000 U	10.0000 U	3.0000 U	10.0000 U	.
S Dimethylphthalate	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Fluoranthene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Fluorene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Hexachlorobenzene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Hexachlorobutadiene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Hexachloroethane	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Isophorone	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	.	35.0000	34.0000	29.0000	51.0000	.
S N-Nitrosodiphenylamine	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Naphthalene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Nitrobenzene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
S Pentachlorophenol	UG/L	.	78.0000	51.0000	48.0000	60.0000	.
S Phenanthrene	UG/L	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	BS	BS	BS	BS	BS	BSD
	LAB ID:	520459	520463	530846	534055	546691	520460
	DATE RCVD:						
NAME	UNITS						
S Phenol	UG/L		57.0000	63.0000	36.0000	55.0000	
S Pyrene	UG/L		46.0000	51.0000	41.0000	41.0000	
S bis(2-Chloroethoxy)methane	UG/L		10.0000 U	10.0000 U	10.0000 U	10.0000 U	
S bis(2-Ethlyhexy)phthalate	UG/L		10.0000 U	1.0000 U	3.0000 U	1.0000 U	
S bis(2-chloroethyl)ether	UG/L		10.0000 U	10.0000 U	10.0000 U	10.0000 U	
3 bis(2-chloroisopropyl)ether	UG/L		10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V 1,1,1-Trichloroethane	UG/L	10.0000 U					10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U					10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U					10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U					10.0000 U
V 1,1-Dichloroethene	UG/L	51.0000					54.0000
V 1,2-Dichloroethane	UG/L	10.0000 U					10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U					10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U					10.0000 U
V 2-Butanone	UG/L	10.0000 U					10.0000 U
V 2-Hexanone	UG/L	10.0000 U					10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U					10.0000 U
V Acetone	UG/L	10.0000 U					10.0000 U
V Benzene	UG/L	48.0000					53.0000
V Bromodichloromethane	UG/L	10.0000 U					10.0000 U
V Bromofo m	UG/L	10.0000 U					10.0000 U
V Bromomethane	UG/L	10.0000 U					10.0000 U
V Carbon Disulfide	UG/L	10.0000 U					10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U					10.0000 U
V Chlorobenzene	UG/L	50.0000					51.0000
V Chloroethane	UG/L	10.0000 U					10.0000 U
V Chlorofo m	UG/L	10.0000 U					10.0000 U
V Chloromethane	UG/L	10.0000 U					10.0000 U
V Dibromochloromethane	UG/L	10.0000 U					10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	BS	BS	BS	BS	BS	BSD
	LAB ID:	520459	520463	530846	534055	546691	520460
	DATE RCVD:						
NAME	UNITS						
V Ethylbenzene	UG/L	10.0000 U	.	.	.	.	10.0000 U
V Methylene Chloride	UG/L	1.0000 U	.	.	.	.	1.0000 U
V Styrene	UG/L	10.0000 U	.	.	.	.	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	.	.	.	.	10.0000 U
V Toluene	UG/L	51.0000	.	.	.	.	50.0000
V Trichloroethene	UG/L	49.0000	.	.	.	.	50.0000
V Vinyl Chloride	UG/L	10.0000 U	.	.	.	.	10.0000 U
V Xylene (total)	UG/L	10.0000 U	.	.	.	.	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	.	.	.	.	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	.	.	.	.	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	BSD 520465	BSD 530848	BSD 534056	BSD 546692	HAZWRAPBLK 529590	HAZWRAPBLK HAZWRAPBLK
ENSAFE ID:		BSD	BSD	BSD	BSD	HAZWRAPBLK	HAZWRAPBLK
LAB ID:		520465	530848	534056	546692	529590	HAZWRAPBLK
DATE RCVD:							
S 1,2,4-Trichlorobenzene	UG/L	48.0000	36.0000	23.0000	41.0000		
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 1,4-Dichlorobenzene	UG/L	38.0000	29.0000	21.0000	37.0000		
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		
S 2,4-Dinitrotoluene	UG/L	40.0000	36.0000	19.0000	45.0000		
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2-Chlorophenol	UG/L	60.0000	58.0000	34.0000	53.0000		
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 4-Chloro-3-methylphenol	UG/L	77.0000	71.0000	42.0000	69.0000		
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		
S 4-Nitrophenol	UG/L	67.0000	54.0000	31.0000	71.0000		
S Acenaphthene	UG/L	47.0000	39.0000	25.0000	48.0000		

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID. LAB ID: DATE RCVD UNITS	BSD 520465	BSD 530848	BSD 534056	BSD 546692	HAZWRAPBLK 529590	HAZWRAPBLK HAZWRAPBLK
S Acenaphthylene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω		
S Anthracene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω		
S Benzo(a)anthracene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω		
S Benzo(a)pyrene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω		
S Benzo(b,k)fluoranthene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Benzo(g,h,i)perylene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Butylbenzylphthalate	UG/L	10.0000 ω	10.0000 ω	1.0000 ω	10.0000 U		
S Carbazole	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Chrysene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Di-n-butylphthalate	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Di-n-octylphthalate	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Dibenz(a,h)anthracene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Dibenzofuran	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Diethylphthalate	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	1.0000 J		
S Dimethylphthalate	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Fluoranthene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Fluorene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Hexachlorobenzene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Hexachlorobutadiene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Hexachlorocyclopentadiene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Hexachloroethane	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Isophorone	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S N-Nitroso-di-n-propylamine	UG/L	43.0000	31.0000	27.0000	51.0000		
S N-Nitrosodiphenylamine	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 U		
S Naphthalene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω		
S Nitrobenzene	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω		
S Pentachlorophenol	UG/L	93.0000	47.0000	30.0000	63.0000		
S Phenanthrene	UG/L	10.0000 ω	10.0000 ω	10.0000	10.0000 ω		

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	BSD	BSD	BSD	BSD	HAZWRAP BLK	HAZWRAPBLK
	LAB ID:	520465	530848	534056	546692	52890	HAZWRAPBLK
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Phenol	UG/L	67.0000	64.0000	37.0000	55.0000	.	.
S Pyrene	UG/L	60.0000	48.0000	27.0000	41.0000	.	.
S bis(2-Chloroethoxy)methane	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω	.	.
S bis(2-Ethlyhexy)phthalate	UG/L	2.0000 ω	2.0000 ω	10.0000 ω	4.0000 ω	.	.
S bis(2-chloroethyl)ether	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω	.	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 ω	10.0000 ω	10.0000 ω	10.0000 ω	.	.
V 1,1,1-Trichloroethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 1,1,2,2-Tetrachloroethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 1,1,2-Trichloroethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 1,1-Dichloroethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 1,1-Dichloroethene	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 1,2-Dichloroethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 1,2-Dichloroethene (total)	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 1,2-Dichloropropane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 2-Butanone	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 2-Hexanone	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V 4-Methyl-2-Pentanone	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Acetone	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Benzene	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Bromodichloromethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Bromoform	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Bromomethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Carbon Disulfide	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Carbon Tetrachloride	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Chlorobenzene	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Chloroethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Chloroform	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Chloromethane	UG/L	.	.	.	.	10.0000 U	10.0000 ω
V Dibromochloromethane	UG/L					10.0000 U	10.0000 ω

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	BSD	BSD	BSD	BSD	HAZWRAPBLK	HAZWRAPBLK
	LAB ID:	520465	530848	534056	546692	529590	HAZWRAPBLK
	DATE RCVD:						
NAME	UNITS						
V Ethylbenzene	UG/L	.	.	.	.	10.0000 u	10.0000 U
V Methylene Chloride	UG/L	.	.	.	.	1.0000 J	1.0000 J
V Styrene	UG/L	.	.		.	10.0000 u	10.0000 U
V Tetrachloroethene	UG/L	.	.	.	.	10.0000 u	10.0000 U
V Toluene	UG/L	.	.	.	.	10.0000 u	10.0000 u
V Trichloroethene	UG/L	.	.	.	.	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	.	.	.	.	10.0000 U	10.0000 U
V Xylene (total)	UG/L	.	.	.	.	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	.	.	.	.	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	.	.	.	.	10.0000 u	10.0000 u

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD:	UNITS	LABPURE 524179	LABPURE 525448	LABPURE 529765	LABPURE 530116	OG5WL 524190	OG5WLMS 524168
V 1,1,1-Trichloroethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	48.0000
V 1,2-Dichloroethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	5.0000 U	10.0000 U
V Benzene		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	48.0000
V Bromodichloromethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	47.0000
V Chloroethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride		UG/L	3.0000 J	7.0000 J	3.0000 J	3.0000 'J	2.0000 U	7.0000 U
V Styrene		UG/L	10.0000 U	10.0000 U	10,0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	48.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	LABPURE	LABPURE	LABPURE	LABPURE	OG5WL	OG5WLMS
	LAB ID:	524179	525448	529765	530116	524190	524168
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	50.0000
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD:	UNITS	OG5WLMSD 524169	PBLK01 531980	PBLK02 529105	PBLK05 532163	PBLK14 533200	PBLK17 533416
P 4,4'-DDD		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.0031 J
P 4,4'-DDE		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDT		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Aldrin		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Aroclor 1016		UG/L	.	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1221		UG/L	.	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
P Aroclor 1232		UG/L	.	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1242		UG/L	.	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1248		UG/L	.	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1254		UG/L	.	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1260		UG/L	.	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Dieldrin		UG/L	.	0.1000 U	0.1000 U	0.0017 J	0.1000 U	0.1000 U
P Endosulfan I		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Endoeulfan II		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan sulfate		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin aldehyde		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin ketone		UG/L	.	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Heptachlor		UG/L	.	0.0500 U		0.0500 U	0.0070 J	0.0500 U
P Heptachlor epoxide		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Methoxychlor		UG/L	.	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U
P Toxaphene		UG/L	.	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
P alpha-BHC		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P alpha-Chlordane		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0020 J	0.0500 U
P beta-BHC		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P delta-BHC		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-BHC (Lindane)		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-Chlordane		UG/L	.	0.0500 U	0.0500 U	0.0500 U	0.0016 J	0.0500 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	OG5WLMSD	PBLK01	PBLK02	PBLK05	PBLK14	PBLK17
	LAB ID:	524169	531980	529105	532163	533200	533416
	DATE RCVD:						
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	.	.	.	.	.
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	.	.	.	.	
V 1,1,2-Trichloroethane	UG/L	10.0000 U	.	.	.	.	.
V 1,1-Dichloroethane	UG/L	10.0000 U	.	.	.	.	.
V 1,1-Dichloroethene	UG/L	48.0000	.	.	.	.	.
V 1,2-Dichloroethane	UG/L	10.0000 U	.	.	.	.	
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	.	.	.	.	.
V 1,2-Dichloropropane	UG/L	10.0000 U	.	.	.	.	.
V 2-Butanone	UG/L	10.0000 U	.	.	.	.	.
V 2-Hexanone	UG/L	10.0000 U	.		.	.	.
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	.	.	.	.	.
V Acetone	UG/L	9.0000 U	.	.	.	.	.
V Benzene	UG/L	56.0000	.	.	.	.	.
V Bromodichloromethane	UG/L	10.0000 U	.	.	.	.	.
V Bromoform	UG/L	10.0000 U	.	.	.	.	.
V Bromomethane	UG/L	10.0000 U	.	.	.	.	.
V Carbon Disulfide	UG/L	10.0000 U	.	.	.	.	.
V Carbon Tetrachloride	UG/L	10.0000 U	.	.	.	.	.
V Chlorobenzene	UG/L	51.0000	.	.	.	.	.
V Chloroethane	UG/L	10.0000 U	.	.	.	.	.
V Chloroform	UG/L	10.0000 U	.	.	.	.	.
V Chloromethane	UG/L	10.0000 U	.	.	.	.	.
V Dibromochloromethane	UG/L	10.0000 U	.	.	.	.	.
V Ethylbenzene	UG/L	10.0000 U	.	.	.	.	.
V Methylene Chloride	UG/L	6.0000 U	.	.	.	.	.
V Styrene	UG/L	10.0000 U	.	.	.	.	.
V Tetrachloroethene	UG/L	10.0000 U	.	.	.	.	.
V Toluene	UG/L	53.0000	.	.	.	.	.
V Trichloroethene	UG/L	52.0000	.	.	.	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	OG5WLMSD	PBLK01	PBLK02	PBLK05	PBLK14	PBLK17
	LAB ID:	524169	531980	529105	532163	533200	533416
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U		.	.	.	.
V Xylene (total)	UG/L	10.0000 U		.	.	.	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U		.	.	.	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U		.	.	.	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	PBLK18 529463	PBLK21 536141	PBLK22 534418	PBLK28 530174	PBLK31 530178	PBLK35 520735
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Aldrin	UG/L	*0.0500* U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Dieldrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Endosulfan II	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Heptachlor	UG/L	0.0500 U	0.0056 J	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Methoxychlor	UG/L	*0.5000* U	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U
P Toxaphene	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
P alpha-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	PBLK36 530365	PBLK39 534031	PBLK46 530434	PBLK47 534123	PBLK50 546924	PBLK52 534817
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Aldrin	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Dieldrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Endosulfan II	UG/L	0.0035 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.0240 U
P Heptachlor	UG/L	0.0021 U	0.0035 U	0.0500 U	0.0500 U	0.0500 U	0.0021 U
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Methoxychlor	UG/L	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U
P Toxaphene	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
P alpha-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0032 U
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID. / LAB ID: / DATE RCVD / UNITS	PBLK56 534566	PBLK59 534648	PBLK60 534657	PBLK60 534657 R1	PBLK73 530889	PBLK74 534920
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	0.0023 J	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Aldrin	UG/L	0.0500 U	0.0500 U	0.0023 J	0.0034 J	0.0300 U	0.0300 U
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Dieldrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U
P Endosulfan II	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.0100 J	0.0098 J	0.1000 U	0.1000 U
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Heptachlor	UG/L	0.0500 U	0.0010 J	0.0080 J	0.0068 J	0.0300 U	0.0300 U
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U
P Methoxychlor	UG/L	0.5000 U	0.5000 U	0.0400 J	0.0450 J	0.5000 U	0.5000 U
P Toxaphene	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
P alpha-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U
P gamma-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0300 U	0.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	PBLK75	PBLK78	PBLK82	PBLK83	PBLK87	PBLK89
ENSAFE ID:		PBLK75	PBLK78	PBLK82	PBLK83	PBLK87	PBLK89
LAB ID:			534924	535143	522844	531536	531912
DATE RCVD:		.	.	.	.	.	.
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Aldrin	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Dieldrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Endosulfan II	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Heptachlor	UG/L	0.0500 U	0.0022 J	0.0018 J	0.0500 U	0.0500 U	0.0500 U
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Methoxychlor	UG/L	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U
P Toxaphene	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
P alpha-BHC	UG/L	0.0500 U	0.0500 U	*0.0500* U	0.0500 U	0.0500 U	0.0500 U
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0500 U	*0.0500* U	0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	*0.0500* U	0.0500 U	0.0500 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	PBLK90 / 528393 / .	PBLK98 / .	PIBLKIG / PIBLK IG / .	PIBLKPU / PIBLK PU / .	SBLK06 / SBLK06 / .	SBLK09 / SBLK09 / .
						.	.
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P Aldrin	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	0.5000 U	0.5000 U	.	.
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	1.0000 U	1.0000 U	.	.
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	0.5000 U	0.5000 U	.	.
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	0.5000 U	0.5000 U	.	.
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	0.5000 U	0.5000 U	.	.
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	0.5000 U	0.5000 U	.	.
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	0.5000 U	0.5000 U	.	.
P Dieldrin	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P Endosulfan II	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P Endrin	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.0500 U	0.0500 U	.	.
P Heptachlor	UG/L	0.0014 J	0.0500 U	0.0200 U	0.0200 U	.	.
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P Methoxychlor	UG/L	0.5000 U	0.5000 U	0.2000 U	0.2000 U	.	.
P Toxaphene	UG/L	5.0000 U	5.0000 U	2.0000 U	2.0000 U	.	.
P alpha-BHC	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.
P gamma-Chlordane	UG/L	0.0500 U	0.0500 U	0.0200 U	0.0200 U	.	.

NAS I'ENSACOLA OU-10 AND SITE 13

QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digite.

	ENSAFE ID:	PBLK90	PBLK98	PIBLKIG	PIBLKPU	SBLKO6	SBLK09
	LAB ID:	528393		PIBLK IG	PIBLK PU	SBLKO6	SBLK09
	DATE RCVD:						
NAME	UNITS						
S 1,2,4-Trichlorobenzene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	.	.	.	.	25.0000 U	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	.	.	.	.	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	.	.	.	.	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	.	.	.	.	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	.	.	.	.	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	.	.	.	.	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	.	.	.	.	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	.	.	.	.	25.0000 U	25.0000 U
S Acenaphthene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Acenaphthylene	UG/L	.	.	.	.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	PBLK90	PBLK98	PIBLKIG	PIBLKPU	SBLKO6	SBLK09
	LAB ID:	528393		PIBLK IG	PIBLK PU	SBLKO6	SBLK09
	DATE RCVD:			.	.	.	.
NAME	UNITS						
S Anthracene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Butylbenzylphthalate	UG/L	.	.	.	.	10.0000 U	*6.0000* J
S Carbazole	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Chrysene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Dibenz (a,h)anthracene	UG/L		.	.	.	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	.	.	.	.	10.0000 U	1.0000 J
S Dimethylphthalate	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Fluoranthene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Fluorene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Ieophorone	UG/L	.	.	.	.	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	.	.	.	.	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Naphthalene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	.	.	.	.	25.0000 U	25.0000 U
S Phenanthrene	UG/L	.	.	.	.	10.0000 U	10.0000 U
S Phenol	UG/L	.	.	.	.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	PBLK90	PBLK98	PIBLKIG	PIBLKPU	SBLK06	SBLK09
	LAB ID:	528393		PIBLK IG	PIBLK PU	SBLK06	SBLK09
	DATE RCVD:						
NAME	UNITS						
S Pyrene	UG/L			.	.	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L		.	.	.	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L		.	.	.	2.0000 J	2.0000 J
S bis(2-chloroethyl)ether	UG/L			.	.	10.0000 U	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L				.	10.0000 u	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK17	SBLK23	SBLK25	SBLK31	SBLK32	SBLK35
	LAB ID:	SBLK17	SBLK23	SBLK25	SBLK31	SBLK32	SBLK35
	DATE RCVD:						
NAME	UNITS						
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK17	SBLK23	SBLK25	SBLK31	SBLK32	SBLK35
	LAB ID:	SBLK17	SBLK23	SBLK25	SBLK31	SBLK32	SBLK35
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Butylbenzylphthalate	UG/L	10.0000 U	5.0000 J	10.0000 U	10.0000 U	10.0000 U	3.0000 J
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1.0000 J
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	10.0000 U	2.0000 J	1.0000 J	10.0000 U	10.0000 U	1.0000 J
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Isophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK17	SBLK23	SBLK25	SBLK31	SBLK32	SBLK35
	LAB ID:	SBLK17	SBLK23	SBLK25	SBLK31	SBLK32	SBLK35
	DATE RCVD:	-	-	-	-	-	-
NAME	UNITS						
S Phenol	UG/L	10.0000 ω	10.0000 ω	3.0000 J	10.0000 ω	10.0000 U	10.0000 ω
S Pyrene	UG/L	10.0000 ω	10.0000 ω	10.0000 U	10.0000 ω	10.0000 U	10.0000 ω
S bis(2-Chloroethoxy)methane	UG/L	10.0000 ω	10.0000 ω	10.0000 U	10.0000 ω	10.0000 U	10.0000 ω
S bis(2-Ethlyhexy)phthalate	UG/L	10.0000 ω	10.0000 ω	3.0000 J	9.0000 5	4.0000 J	2.0000 5
S bis(2-chloroethyl)ether	UG/L	10.0000 ω	10.0000 ω	10.0000 U	10.0000 ω	10.0000 U	10.0000 ω
S bis(2-chloroisopropyl)ether	UG/L	10.0000 ω	10.0000 ω	10.0000 U	10.0000 ω	10.0000 U	10.0000 ω

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK36	SBLK37	SBLK41	SBLK43	SBLK47	SBLK53
	LAB ID:	SBLK36	SBLK37	SBLK41	SBLK43	SBLK47	SBLK53
	DATE RCVD:						
NAME	UNITS						
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	SBLK36 / SBLK36	SBLK37 / SBLK37	SBLK41 / SBLK41	SBLK43 / SBLK43	SBLK47 / SBLK47	SBLK53 / SBLK53
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Butylbenzylphthalate	UG/L	8.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1.0000 J
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chryeene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Ieophoronc	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK36	SBLK37	SBLK41	SBLK43	SBLK47	SBLK53
	LAB ID:	SBLK36	SBLK37	SBLK41	SBLK43	SBLK47	SBLK53
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Phenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	2.0000 J	8.0000 J	3.0000 J	2.0000 J	7.0000 J	43.0000 J
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK54	SBLK54	SBLK57	SBLK60	SBLK62	SBLK64
	LAB ID:	SBLK 54	SBLK54	SBLK57	SBLK60	SBLK62	SBLK64
	DATE RCVD:						
NAME	WITS						
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	'10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	SBLK54	SBLK54	SBLK57	SBLKCO	SBLK62	SBLK64
ENSAFE ID:		SBLK54	SBLK54	SBLK57	SBLKCO	SBLK62	SBLK64
LAB ID:		SBLK 54	SBLK54	SBLK57	SBLKCO	SBLK62	SBLK64
DATE RCVD:							
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Butylbenzylphthalate	UG/L	1.0000 J	1.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	10.0000 U	10.0000 U	2.0000 J	10.0000 U	10.0000 U	10.0000 U
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10,0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Ieophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK54	SBLK54	SBLK57	SBLK60	SBLK62	SBLK64
	LAB ID:	SBLK 54	SBLK54	SBLK57	SBLK60	SBLK62	SBLK64
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Phenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	10.0000 U	10.0000 U	2.0000 J	2.0000 J	7.0000 J	3.0000 J
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK68	SBLK69	SBLK69	SBLK73	SBLK81	SBLK83
	LAB ID:	SBLK68	SBLK 69	SBLK69	SBLK73	SBLK81	SBLK83
	DATE RCVD:						
NAME	UNITS						
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK68	SBLK69	SBLK69	SBLK73	SBLK81	SBLK83
	LAB ID:	SBLK68	SBLK 69	SBLK69	SBLK73	SBLK81	SBLK83
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Acenaphthylene	UG/L	10.0000 u	10.0000 u	10.0000 u	10.0000 U	10.0000 U	10.0000 U
s Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 u
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 u	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U
S Butylbenzylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	3.0000 J
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1.0000 J
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.00'00 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Isophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 u	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
s Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK68	SBLK69	SBLK69	SBLK73	SBLK81	SBLK83
	LAB ID:	SBLK68	SBLK 69	SBLK69	SBLK73	SBLK81	SBLK83
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Phenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	1.0000 J	1.0000 J	1.0000 J	45.0000 J	7.0000 J	5.0000 J
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	SBLK96 / SBLK96 / .	SBLK97 / SBLK 97 / .	SBLK97 / SBLK97 / .	SBLK99 / SBLK99 / .	VBLKAA / VBLK AA / .	VBLKAA / VBLKAA / .
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	SBLK96	SBLK97	SBLK97	SBLK99	VBLKAA	VBLKAA
	LAB ID:	SBLK96	SBLK 97	SBLK97	SBLK99	VBLK AA	VBLKAA
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Butylbenzylphthalate	UG/L	10.0000 U	10.0000 U	2.0000 J	10.0000 U		.
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	1.0000 J	10.0000 U		.
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Dibenz (a,h)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Diethylphthalate	UG/L	1.0000 J	1.0000 J	2.0000 J	10.0000 U		.
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Isophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD:	UNITS	SBLK96 / SBLK96	SBLK97 / SBLK 97	SBLK97 / SBLK97	SBLK99 / SBLK99	VBLKAA / VBLK M	VBLKAA / VBLKAA
S Phenol		UG/L	1.0000 J	10.0000 U	10.0000 U	10.0000 U	.	
S Pyrene		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	
S bis (2-Chloroethoxy)methane		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	
S bis(2-Ethlyhexy)phthalate		UG/L	10.0000 U	6.0000 J	33.0000	1.0000 J		
S bis (2-chloroethyl)ether		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S bis(2-chloroisopropyl)ether		UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
V 1,1,1-Trichloroethane		UG/L	.				10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane		UG/L	.		.	.	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane		UG/L			.	.	10.0000 U	10.0000 U
V 1,1-Dichloroethane		UG/L	.	.	.	.	10.0000 U	10.0000 U
V 1,1-Dichloroethene		UG/L	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloroethane		UG/L	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)		UG/L	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloropropane		UG/L	.	.	.	.	10.0000 U	10.0000 U
V 2-Butanone		UG/L	.	.	.	.	10.0000 U	10.0000 U
V 2-Hexanone		UG/L	.	.		.	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone		UG/L	.	.	.	.	10.0000 J	2.0000 J
V Acetone		UG/L	.	.		.	10.0000 J	8.0000 J
V Benzene		UG/L	.		.	.	10.0000 J	0.1000 J
V Bromodichloromethane		UG/L	.	.	.	.	10.0000 U	10.0000 U
V Bromoform		UG/L	.	.		.	10.0000 U	10.0000 U
V Bromomethane		UG/L	.	.	.	.	10.0000 U	10.0000 U
V Carbon Disulfide		UG/L	.	.	.	.	10.0000 J	0.0700 J
V Carbon Tetrachloride		UG/L	.	.	.	.	10.0000 U	10.0000 U
V Chlorobenzene		UG/L	.	.	.	.	10.0000 J	0.1000 J
V Chloroethane		UG/L	.	.	.	.	10.0000 U	10.0000 U
V Chloroform		UG/L	.	.	.	.	10.0000 J	0.1000 J
V Chloromethane		UG/L	.	.	.	.	10.0000 U	10.0000 U
V Dibromochloromethane		UG/L	.	.	.	.	10.0000 U	10.0000 U

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	SBLK96	SBLK97	SBLK97	SBLK99	VBLKAA	VBLKAA
	LAB ID:	SBLK96	SBLK 97	SBLK97	SBLK99	VBLK AA	VBLKAA
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Ethylbenzene	UG/L	.	.	.	.	10.0000 J	0.2000 J
V Methylene Chloride	UG/L	.	.	.	.	10.0000 J	0.6000 J
V Styrene	UG/L	.	.	.	.	10.0000 J	0.2000 J
V Tetrachloroethene	UG/L	.	.	.	.	10.0000 U	10.0000 U
V Toluene	UG/L	.	.	.	.	10.0000 J	0.1000 J
V Trichloroethene	UG/L	.	.	.	.	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	.	.	.	.	10.0000 U	10.0000 U
V Xylene (total)	UG/L	.	.	.	.	10.0000 J	0.4000 J
V cis-1,3-Dichloropropene	UG/L	.	.	.	.	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	.	.	.	.	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKAC	VBLKAE	VBLKAO	VBLKBF	VBLKBJ	VBLKC5
	LAB ID:	VBLKAC	VBLKAE	VBLKAO	VBLKBF	VBLKBJ	VBLK C5
	DATE RCVD:						
NAME	WITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	18.0000 J
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	7.0000 J	7.0000 J	4.0000 J	2.0000 J	2.0000 J	1.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKAC	VBLKAE	VBLKAO	VBLKBF	VBLKBJ	VBLKC5
	LAB ID:	VBLKAC	VBLKAE	VBLKAO	VBLKBF	VBLKBJ	VBLK C5
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	VBLKC5 / VBLKC5	VBLKCC / VBLKC6	VBLKCB / VBLKCB	VBLKCC / VBLKCC	VBLKD3 / VBLKD3	VBLKDC / VBLKDC
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10,0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	18.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	1.0000 J	10.0000 U	8.0000 J	5.0000 J	10.0000 U	1.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKC5	VBLKC6	VBLKCB	VBLKCC	VBLKD3	VBLKDC
	LAB ID:	VBLKC5	VBLKC6	VBLKCB	VBLKCC	VBLKD3	VBLKDC
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	VBLKDG	VBLKDI	VBLKDW	VBLKG3	VBLKG3	VBLKG9
ENSAFE ID:		VBLKDG	VBLKDI	VBLKDW	VBLK G3	VBLKG3	VBLKG9
LAB ID:		VBLKDG	VBLKDI	VBLKDW	VBLK G3	VBLKG3	VBLKG9
DATE RCVD:		.	.	.	.	.	.
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	10.0000 U	1.0000 J	10.0000 U	3.0000 J	2.0000 J	10.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKDG	VBLKDI	VBLKDW	VBLKG3	VBLKG3	VBLKG9
	LAB ID:	VBLKDG	VBLKDI	VBLKDW	VBLK G3	VBLKG3	VBLKG9
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKI2	VBLKK7	VBLKK8	VBLKK9	VBLKKP	VBLKKS
	LAB ID:	VBLKI2	VBLKK7	VBLKK8	VBLKK9	VBLKKP	VBLKKS
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	10.0000 U	1.0000 J	1.0000 J	10.0000 U	3.0000 J	5.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKI2	VBLKK7	VBLKK8	VBLKK9	VBLKKP	VBLKKS
	LAB ID:	VBLKI2	VBLKK7	VBLKK8	VBLKK9	VBLKKP	VBLKKS
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKL1	VBLKL2	VBLKL3	VBLKO5	VBLKSB	VBLKSG
	LAB ID:	VBLKL1	VBLKL2	VBLKL3	VBLKO5	VBLKSB	VBLKSG
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	2.0000 J	2.0000 J	2.0000 J	2.0000 J	5.0000 J	4.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKL1	VBLKL2	VBLKL3	VBLKO5	VBLKSB	VBLKSG
	LAB ID:	VBLKL1	VBLKL2	VBLKL3	VBLKO5	VBLKSB	VBLKSG
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10,0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKVR	VBLKW9	VBLKWN	VBLKWX	VBLKX1	VBLKXY
	LAB ID:	VBLKVR	VBLKW9	VBLKWN	VBLKWX	VBLKX1	VBLKXY
	DATE RCVD:	.	.	.	.	.	.
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 J	10.0000 U	10.0000 U	10.0000 J	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	1.0000 J	10.0000 J	4.0000 J	6.0000 J	1.0000 J	2.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKVR	VBLKW9	VBLKWN	VBLKWX	VBLKX1	VBLKXY
	LAB ID:	VBLKVR	VBLKW9	VBLKWN	VBLKWX	VBLKX1	VBLKXY
	DATE RCVD:						
NAME	WITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKY8	VBLKZ1
	LAB ID:	VBLKY8	VBLKZ1
	DATE RCVD:		
NAME	UNITS		
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10,0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	1.0000 J	1.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	VBLKY8	VBLKZ1					
	LAB ID:	VBLKY8	VBLKZ1					
	DATE RCVD:							
NAME	UNITS							
V Trichloroethene	UG/L	10.0000 U	10.0000 U					
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U					
V Xylene (total)	UG/L	10.0000 U	10.0000 U					
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U					
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U					

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001MS	13M61001MSD	13M63M01	13M63S01	13M64001MS	13M64001MSD
	LAB ID:	532105	532106	532108	532109	532102	532103
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
P 4,4'-DDD	UG/KG	.	.	0.5600 J	0.6800 J		.
P 4,4'-DDE	UG/KG	.	.	6.6000	4.4000 J		.
P 4,4'-DDT	UG/KG	.	.	30.0000	30.0000		.
P Aldrin	UG/KG	.	.	11.0000	12.0000		.
P Aroclor 1016	UG/KG	.	.	33.0000 U	33.0000 U		.
P Aroclor 1221	UG/KG	.	.	67.0000 U	67.0000 U		.
P Aroclor 1232	UG/KG	.	.	33.0000 U	*33.0000* U		.
P Aroclor 1242	UG/KG	.	.	33.0000 U	*33.0000* U		.
P Aroclor 1248	UG/KG	.		33.0000 U	33.0000 U		.
P Aroclor 1254	UG/KG	.		33.0000 U	33.0000 U		.
P Aroclot 1260	UG/KG	.		*33.0000* U	33.0000 U		.
P Dieldrin	UG/KG	.		27.0000	27.0000		.
P Endosulfan I	UG/KG	.		1.7000 U	1.7000 U		.
P Endosulfan II	UG/KG	.		3.3000 U	*3.3000* U		.
P Endosulfan sulfate	UG/KG	.		3.3000 U	3.3000 U		.
P Endrin	UG/KG	.		29.0000	30.0000		.
P Endrin aldehyde	UG/KG	.		*3.3000* U	*3.3000* U		.
P Endrin ketone	UG/KG	.		0.2300 J	0.2900 J		.
P Heptachlor	UG/KG	.		12.0000 J	13.0000 J		.
P Heptachlor epoxide	UG/KG	.		1.7000 U	1.7000 U		.
P Methoxychlor	UG/KG	.		17.0000 U	17.0000 U		.
P Toxaphene	UG/KG	.		170.0000 U	170.0000 U		.
P alpha-BHC	UG/KG	.		1.7000 U	1.7000 U		.
P alpha-Chlordane	UG/KG	.		1.7000 U	1.7000 U		.
P beta-BHC	UG/KG	.		1.7000 U	1.7000 U		.
P delta-BHC	UG/KG	.		1.7000 U	1.7000 U		.
P gamma-BHC (Lindane)	UG/KG	.		12.0000	13.0000		.
P gamma-Chlordane	UG/KG			1.7000 U	1.7000 U		.
S 1,2,4-Trichlorobenzene	UG/KG	2200.0000	1900.0000				.

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001MS	13M61001MSD	13M63M01	13M63S01	13M64001MS	13M64001MSD
	LAB ID:	532105	532106	532108	532109	532102	532103
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
S 1,2-Dichlorobenzene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S 1,3-Dichlorobenzene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S 1,4-Dichlorobenzene	UG/KG	1800.0000	1600.0000		.	.	
S 2,4,5-Trichlorophenol	UG/KG	1100.0000 U	1100.0000 U		.	.	
S 2,4,6-Trichlorophenol	UG/KG	440.0000 U	440.0000 U		.	.	
S 2,4-Dichlorophenol	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 2,4-Dimethylphenol	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 2,4-Dinitrophenol	UG/KG	1100.0000 U	1100.0000 U	.	.	.	
S 2,4-Dinitrotoluene	UG/KG	1800.0000	1600.0000	.	.	.	
S 2,6-Dinitrotoluene	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 2-Chloronaphthalene	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 2-Chlorophenol	UG/KG	2500.0000	2300.0000	.	.	.	
S 2-Methylnaphthalene	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 2-Methylphenol	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 2-Nitroaniline	UG/KG	1100.0000 U	1100.0000 U	.	.	.	
S 2-Nitrophenol	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 3,3'-Dichlorobenzidine	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 3-Nitroaniline	UG/KG	1100.0000 U	1100.0000 U	.	.	.	
S 4,6-Dinitro-2-methylphenol	UG/KG	1100.0000 U	1100.0000 U	.	.	.	
S 4-Bromophenyl-phenylether	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 4-Chloro-3-methylphenol	UG/KG	3300.0000	3100.0000	.	.	.	
S 4-Chloroaniline	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 4-Chlorophenyl-phenylether	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 4-Methylphenol	UG/KG	440.0000 U	440.0000 U	.	.	.	
S 4-Nitroaniline	UG/KG	1100.0000 U	1100.0000 U	.	.	.	
S 4-Nitrophenol	UG/KG	4000.0000 J	3500.0000	.	.	.	
S Acenaphthene	UG/KG	1900.0000	1800.0000	.	.	.	
S Acenaphthylene	UG/KG	440.0000 U	440.0000 U	.	.	.	
S Anthracene	UG/KG	440.0000 U	440.0000 U	.	.	.	

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001MS	13M61001MSD	13M63M01	13M63S01	13M64001MS	13M64001MSD
	LAB ID:	532105	532106	532108	532109	532102	532103
	DATE RCVD:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	WITS						
S Benzo(a)anthracene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Benzo(a)pyrene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Benzo(b,k)fluoranthene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Benzo(g,h,i)perylene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Butylbenzylphthalate	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Carbazole	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Chrysene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Di-n-butylphthalate	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Di-n-octylphthalate	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Dibenz(a,h)anthracene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Dibenzofuran	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Diethylphthalate	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Dimethylphthalate	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Fluoranthene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Fluorene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Hexachlorobenzene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Hexachlorobutadiene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Hexachlorocyclopentadiene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Hexachloroethane	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Indeno(1,2,3-cd)pyrene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Ieophorone	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S N-Nitroso-di-n-propylamine	UG/KG	1900.0000	1700.0000	.	.	.	.
S N-Nitrosodiphenylamine	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Naphthalene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Nitrobenzene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Pentachlorophenol	UG/KG	3500.0000	3200.0000	.	.	.	.
S Phenanthrene	UG/KG	440.0000 U	440.0000 U	.	.	.	.
S Phenol	UG/KG	2600.0000	2400.0000	.	.	.	.
S Pyrene	UG/KG	1900.0000	1800.0000	.	.	.	.

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCW: / WITS	13M61001MS	13M61001MSD	13M63M01	13M63S01	13M64001MS	13M64001MSD
	LAB ID:	532105	532106	532108	532109	532102	532103
	DATE RCW:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
S bin(2-Chloroethoxy)methane	UG/KG	440.0000 U	440.0000 U	.	.		
S bis(2-Ethlyhexy)phthalate	UG/KG	440.0000 U	440.0000 U	.	.		
S bis(2-chloroethyl)ether	UG/KG	440.0000 U	440.0000 U	.	.		
S bis(2-chloroisopropyl)ether	UG/KG	440.0000 U	440.0000 U	.	.		
V 1,1,1-Trichloroethane	UG/KG		.	.	.	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG		.	.	.	59.0000	57.0000
V 1,2-Dichloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 2-Butanone	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 2-Hexanone	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Acetone	UG/KG	.	.	.	.	17.0000 U	18.0000 U
V Benzene	UG/KG	.	.	.	.	68.0000	66.0000
V Bromodichloromethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Bromoform	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Bromomethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Chlorobenzene	UG/KG	.	.	.	.	59.0000	58.0000
V Chloroethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Chloroform	UG/KG	.	.	.	.	1.0000 U	2.0000 U
V Chloromethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Ethylbenzene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Methylene Chloride	UG/KG	.	.	.	.	27.0000 U	17.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M61001MS	13M61001MSD	13M63M01	13M63S01	13M64001MS	13M64001MSD
	LAB ID:	532105	532106	532108	532109	532102	532103
	DATE RCW:	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93	01/15/93
NAME	UNITS						
V Styrene	UG/KG	.	.	.	.	10.0000 U	10.0000 U
V Tetrachloroethene	UG/KG			.	.	10.0000 U	10.0000 U
V Toluene	UG/KG	.		.	.	58.0000	56.0000
V Trichloroethene	UG/KG	.		.	.	66.0000	*63.0000*
V Vinyl Chloride	UG/KG	.		.	.	10.0000 U	10.0000 U
V Xylene (total)	UG/KG	.		.	.	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	.		.	.	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	.		.	.	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	13M65D01 / 533375 / 01/15/93	13S02M03 / 12/09/92	13S02M06 / 12/09/92	13S02S03 / 12/09/92	13S02S06 / 12/09/92	13S04D06 / 533000 / 01/20/93
I Aluminum	MG/KG	9170.0000	.	.	.	.	76.0000
I Antimony	MG/KG	14.4000 U	.		.	.	12.0000 U
I Arsenic	MG/KG	6.2000	.	.	.	.	3.2000 U
I Barium	MG/KG	12.3000 J	.	.	.	.	1.3000 U
I Beryllium	MG/KG	0.7600 U	.	.	.	.	0.6300 U
I Cadmium	MG/KG	2.1000	.	.	.	.	1.1000 U
I Calcium	MG/KG	6170.0000	.	.	.	.	27.3000 U
I Chromium	MG/KG	46.6000		.	.	.	2.1000 U
I Cobalt	MG/KG	3.0000 U	.	.	.	.	2.5000 U
I Copper	MG/KG	17.7000	.	.	.	.	5.3000 U
I Cyanide	MG/KG	0.6300 U	.	.	.	.	0.5300 U
I Iron	MG/KG	10300.0000	.	.	.	.	46.9000
I Lead	MG/KG	48.3000	.	.	.	.	0.9700
I Magnesium	MG/KG	1910.0000	.	.	.	.	116.0000 U
I Manganese	MG/KG	57.0000	.	.	.	.	1.3000 U
I Mercury	MG/KG	0.1300 U			.	.	0.1100 U
I Nickel	MG/KG	9.6000 U	.	.	.	.	8.0000 U
I Potassium	MG/KG	996.0000 U	.		.	.	831.0000 U
I Selenium	MG/KG	0.7600 U	.	.	.	.	4.2000 U
I Silver	MG/KG	2.5000 U	.	.	.	.	2.1000 U
I sodium	MG/KG	413.0000 J	.	.	.	.	102.0000 J
I Thallium	MG/KG	0.5000 U	.	.	.	.	0.6300 U
I Vanadium	MG/KG	23.6000	.	.	.	.	3.2000 U
I Zinc	MG/KG	55.5000	.		.		0.6300 U
P 4,4'-DDD	UG/KG	3.5000 U	.	3.4000 U	.	3.3000 U	3.4000 U
P 4,4'-DDE	UG/KG	0.3300 J	.	1.1000 J	.	1.2000 J	3.4000 U
P 4,4'-DDT	UG/KG	0.2000 J	.	32.0000	.	33.0000	3.4000 U
P Aldrin	UG/KG	1.8000 U	.	13.0000	.	13.0000	1.8000 U

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M65D01	13S02M03	13S02M06	13S02S03	13S02S06	13S04D06
	LAB ID:	533375					533008
	DATE RCVD:	01/15/93	12/09/92	12/09/92	12/09/92	12/09/92	01/20/93
NAME	UNITS						
P Aroclor 1016	UG/KG	35.0000 U		34.0000 U		33.0000 U	34.0000 U
P Aroclor 1221	UG/KG	72.0000 U		70.0000 U		67.0000 U	69.0000 U
P Aroclor 1232	UG/KG	35.0000 U		34.0000 U		33.0000 U	34.0000 U
P Aroclor 1242	UG/KG	35.0000 U		34.0000 U		33.0000 U	34.0000 U
P Aroclor 1248	UG/KG	35.0000 U		34.0000 U		33.0000 U	34.0000 U
P Aroclor 1254	UG/KG	35.0000 U		34.0000 U		33.0000 U	34.0000 U
P Aroclor 1260	UG/KG	4.4000 J		34.0000 U		33.0000 U	34.0000 U
P Dieldrin	UG/KG	3.5000 U		29.0000		31.0000	3.4000 U
P Endosulfan I	UG/KG	1.8000 U		1.8000 U		1.7000 U	1.8000 U
P Endosulfan II	UG/KG	3.5000 U		3.4000 U		0.0880 J	3.4000 U
P Endosulfan sulfate	UG/KG	3.5000 U		3.4000 U		3.3000 U	3.4000 U
P Endrin	UG/KG	3.5000 U		31.0000		33.0000	3.4000 U
P Endrin aldehyde	UG/KG	3.5000 U		3.4000 U		3.3000 U	3.4000 U
P Endrin ketone	UG/KG	3.5000 U		0.7500 J		0.9000 J	3.4000 U
P Heptachlor	UG/KG	0.2000 J		13.0000 J		14.0000 J	0.1300 J
P Heptachlor epoxide	UG/KG	1.8000 U		0.1200 J		1.7000 U	1.8000 U
P Methoxychlor	UG/KG	18.0000 U		18.0000 U		17.0000 U	18.0000 U
P Toxaphene	UG/KG	180.0000 U		180.0000 U		170.0000 U	180.0000 U
P alpha-BHC	UG/KG	1.8000 U		1.8000 U		1.7000 U	1.8000 U
P alpha-Chlordane	UG/KG	1.8000 U		1.8000 U		0.1400 J	1.8000 U
P beta-BHC	UG/KG	1.8000 U		1.8000 U		1.7000 U	1.8000 U
P delta-BHC	UG/KG	1.8000 U		1.8000 U		1.7000 U	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.8000 U		14.0000 J		15.0000 J	1.8000 U
P gamma-Chlordane	UG/KG	1.8000 U		1.8000 U		1.7000 U	1.8000 U
S 1,2,4-Trichlorobenzene	UG/KG	320.0000 U	1600.0000		1700.0000		340.0000 U
S 1,2-Dichlorobenzene	UG/KG	320.0000 U	340.0000 U		340.0000 U		340.0000 U
S 1.3-Dichlorobenzene	UG/KG	320.0000 U	340.0000 U		340.0000 U		340.0000 U
S 1.4-Dichlorobenzene	UG/KG	320.0000 U	1400.0000		1400.0000		340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	790.0000 U	820.0000 U		820.0000 U		830.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13M65D01	13S02M03	13S02M06	13S02S03	13S02S06	13S04D06
ENSAFE ID:		13M65D01	13S02M03	13S02M06	13S02S03	13S02S06	13S04D06
LAB ID:		533375					533008
DATE RCVD:		01/15/93	12/09/92	12/09/92	12/09/92	12/09/92	01/20/93
S 2,4,6-Trichlorophenol	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 2,4-Dichlorophenol	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 2,4-Dimethylphenol	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 2,4-Dinitrophenol	UG/KG	790.0000 U	820.0000 U		820.0000 U	.	830.0000 U
S 2,4-Dinitrotoluene	UG/KG	320.0000 U	1500.0000		1500.0000	.	340.0000 U
S 2,6-Dinitrotoluene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 2-Chloronaphthalene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 2-Chlorophenol	UG/KG	320.0000 U	2000.0000		2100.0000	.	340.0000 U
S 2-Methylnaphthalene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 2-Methylphenol	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 2-Nitroaniline	UG/KG	790.0000 U	820.0000 U		820.0000 U	.	830.0000 U
S 2-Nitrophenol	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 3-Nitroaniline	UG/KG	790.0000 U	820.0000 U		820.0000 U	.	830.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	790.0000 U	820.0000 U		820.0000 U	.	830.0000 U
S 4-Bromophenyl-phenylether	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 4-Ctiloro-3-methylphenol	UG/KG	320.0000 U	1400.0000		1400.0000	.	340.0000 U
S 4-Chloroaniline	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 4-Methylphenol	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S 4-Nitroaniline	UG/KG	790.0000 U	820.0000 U		820.0000 U	.	830.0000 U
S 4-Nitrophenol	UG/KG	790.0000 U	1300.0000		1400.0000	.	830.0000 U
S Acenaphthene	UG/KG	320.0000 U	1500.0000		1500.0000	.	340.0000 U
S Acenaphthylene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Anthracene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Benzo(a)anthracene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Benzo(a)pyrene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG	320.0000 U	62.0000 J		34.0000 J	.	340.0000 U
S Benzo(g,h,i)perylene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13M65D01	13S02M03	13SO2M06	13S02S03	13502506	13S04D06
ENSAFE ID:		13M65D01	13S02M03	13SO2M06	13S02S03	13502506	13S04D06
LAB ID:		533375					533008
DATE RCVD:		01/15/93	12/09/92	12/09/92	12/09/92	12/09/92	01/20/93
S Butylbenzylphthalate	UG/KG	240.0000 U	340.0000 U		340.0000 U	.	230.0000 U
S Carbazole	UG/KG	320.0000 U	340.0000 U		340.0000 U		340.0000 U
S Chrysene	UG/KG	320.0000 U	340.0000 U		340.0000 U		340.0000 U
S Di-n-butylphthalate	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Di-n-octylphthalate	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Dibenz(a,h)anthracene	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Dibenzofuran	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S Diethylphthalate	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Dimethylphthalate	UG/KG	320.0000 U	340.0000 U		340.0000 U	.	340.0000 U
S Fluoranthene	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S Fluorene	UG/KG	320.0000 U	340,0000 U		340.0000 U	.	340.0000 U
S Hexachlorobenzene	UG/KG	320.0000 U	340.0000 U		340.0000 U		340.0000 U
S Hexachlorobutadiene	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	320.0000 U	340.0000 U		340.0000 U		340.0000 U
S Hexachloroethane	UG/KG	320.0000 U	340.0000 U	.	340.0000 U		340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	320.0000 U	340.0000 U	.	340.0000 U		340.0000 U
S Ieophorone	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	320.0000 U	1600.0000	.	1700.0000	.	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S Naphthalene	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S Nitrobenzene	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S Pentachlorophenol	UG/KG	790.0000 U	2500.0000	.	2500.0000	.	830.0000 U
S Phenanthrene	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S Phenol	UG/KG	320.0000 U	1900.0000	.	2000.0000	.	340.0000 U
S Pyrene	UG/KG	320.0000 U	1200.0000	.	1400.0000	.	340.0000 U
S bis(2-Chloroethoxy)methane	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	45.0000 U	340.0000 U	.	40.0000 U	.	47.0000 U
S bis(2-chloroethyl)ether	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	320.0000 U	340.0000 U	.	340.0000 U	.	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13M65D01	13S02M03	13S02M06	13S02S03	13S02S06	13S04D06
	LAB ID:	533375					533008
	DATE RCVD:	01/15/93	12/09/92	12/09/92	12/09/92	12/09/92	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 1,2-Dichloroethene (total	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 2-Butanone	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 2-Hexanone	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Acetone	UG/KG	7.0000 U	.	.	.	.	96.0.000U
V Benzene	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Bromoform	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Bromomethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Chlorobenzene	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Chloroethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Chloroform	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Chloromethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Ethylbenzene	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Methylene Chloride	UG/KG	8.0000 U	.	.	.	.	20.0000 U
V Styrene	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Toluene	UG/KG	10.0000 U	.	.	.	.	11.0000 U
V Trichloroethene	UG/KG	10.0000 U	.	.	.	.	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in eignificant digits.

NAME	UNITS						
	ENSAFE ID:	13M65D01	13S02M03	13S02M06	13S02S03	13S02S06	13S04D06
	LAB ID:	533375					533008
	DATE RCVD:	01/15/93	12/09/92	12/09/92	12/09/92	12/09/92	01/20/93
V Vinyl Chloride	UG/KG	10.0000 U					11.0000 U
v Xylene (total)	UG/KG	10.0000 U					11.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U					11.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U					11.0000 u

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not ehown in significant digits.

NAME	UNITS	13S06D02	13S64D08	13S65D06	13S66D08	13S68M04	13S68S04
ENSAFE ID:		13S06D02	13S64D08	13S65D06	13S66D08	13S68M04	13S68S04
LAB ID:		529876	532097	532101	532117	533211	533212
DATE RCVD:		12/18/92	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93
I Aluminum	MG/KG	180.0000	12000.0000	285.0000	209.0000		
I Antimony	MG/KG	12.5000 U	12.2000 U	11.8000 U	11.8000 U		
I Arsenic	MG/KG	0.6300 U	2.8000	0.7800 J	0.6200 U		
I Barium	MG/KG	2.3000 J	7.3000 J	1.2000 U	1.2000 U		
I Beryllium	MG/KG	0.2100 U	0.6400 U	0.6200 U	0.6200 U		
I Cadmium	MG/KG	1.0000 U	1.1000 U	1.0000 U	1.0000 U		
I Calcium	MG/KG	9110.0000	532.0000 J	45.1000 J	86.6000 J		
I Chromium	MG/KG	1.3000 U	12.6000	2.1000 U	2.5000		
I Cobalt	MG/KG	1.9000 U	2.6000 U	2.5000 U	2.5000 U		
I Copper	MG/KG	1.0000 U	5.4000 U	5.2000 U	5.2000 U		
I Cyanide	MG/KG	0.5200 U	1,8000	0.5200 U	0.5200 U		
I Iron	MG/KG	236.0000	7800.0000	190.0000	165.0000		
I Lead	MG/KG	3.6000	11.5000	4.2000	1.2000		
I Magnesium	MG/KG	39.6000 J	119.0000 U	114.0000 U	115.0000 U		
I Manganese	MG/KG	3.4000	30.3000	1.2000 U	1.6000 J		
I Mercury	MG/KG	0.1000 U	0.1100 U	0.1000 U	0.1000 U		
I Nickel	MG/KG	2.3000 U	8.2000 U	7.8000 U	7.9000 U		
I Potaseium	MG/KG	285.0000 U	849.0000 U	815.0000 U	821.0000 U		
I Selenium	MG/KG	0.6300 U	0.8600 U	0.8200 U	0.6200 U		
I Silver	MG/KG	1.3000 U	2.1000 U	2.1000 U	2.1000 U		
I Sodium	MG/KG	188.0000 J	176.0000 J	88.8000 J	115.0000 J		
I Thallium	MG/KG	0.4200 U	0.6400 U	0.6200 U	0.6200 U		
I Vanadium	MG/KG	1.0000 U	19.5000	3.1000 U	3.1000 U		
I Zinc	MG/KG	6.1000	5.7000	2.3000 J	2.9000 J		
P 4,4'-DDD	UG/KG	3.3000 U	0.9700 J	3.4000 U	3.4000 U		
P 4,4'-DDE	UG/KG	3.3000 U	0.1400 J	3.4000 U	3.4000 U		
P 4,4'-DDT	UG/KG	0.1000 J	0.1300 J	3.4000 U	3.4000 U		
P Aldrin	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key pteeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	13S06D02	13S64D08	13S65D06	13S66D08	13868104	13S68S04
ENSAFE ID:		13S06D02	13S64D08	13S65D06	13S66D08	13868104	13S68S04
LAB ID:		529076	532097	532101	532117	533211	533212
DATE RCVD:		12/18/92	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93
	UG/KG	33.0000 U	35.0000 U	34.0000 U	34.0000 U		.
	UG/KG	68.0000 U	71.0000 U	69.0000 U	70.0000 U		
	UG/KG	33.0000 U	35.0000 U	34.0000 U	34.0000 U		.
	UG/KG	33.0000 U	35.0000 U	34.0000 U	34.0000 U		.
	UG/KG	33.0000 U	35.0000 U	34.0000 U	34.0000 U		.
	UG/KG	33.0000 U	35.0000 U	34.0000 U	34.0000 U		.
	UG/KG	33.0000 U	35.0000 U	34.0000 U	14.0000 J		.
	UG/KG	3.3000 U	0.1200 J	3.4000 U	3.4000 U		.
P Endosulfan I	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		.
P Endosulfan II	UG/KG	3.3000 U	3.5000 U	3.4000 U	3.4000 U		.
P Endosulfan sulfate	UG/KG	3.3000 U	3.5000 U	3.4000 U	3.4000 U		.
P Endrin	UG/KG	3.3000 U	3.5000 U	3.4000 U	3.4000 U		.
P Endrin aldehyde	UG/KG	3.3000 U	3.5000 U	3.4000 U	3.4000 U		.
P Endrin ketone	UG/KG	3.3000 U	3.5000 U	3.4000 U	3.4000 U		.
P Heptachlor	UG/KG	1.7000 U	1.8000 U	0.0560 J	0.0940 J		.
P Heptachlor epoxide	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		.
P Methoxychlor	UG/KG	17.0000 U	18.0000 U	18.0000 U	18.0000 U		.
P Toxaphene	UG/KG	170.0000 U	180.0000 U	180.0000 U	180.0000 U		.
P alpha-BHC	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		.
P alpha-Chlordane	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		.
P beta-BHC	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		.
P delta-BHC	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		.
P gamma-BHC (Lindane)	UG/KG	1.7000 U	1.8000 U	1.8000 U	1.8000 U		.
P gamma-Chlordane	UG/KG	1.7000 U	1.8000 U	1.8000 U	0.0680 J		.
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U		.
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U		.
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U		.
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U		.
S 2,4,5-Trichlorophenol	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U		.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S06D02	13S64D08	13S65D06	13S66D08	13S68M04	13S68S04
	LAB ID:	529876	532097	532101	532117	533211	533212
	DATE RCVD:	12/18/92	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2,4-Dichlorophenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2,4-Dimethylphenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2,4-Dinitrophenol	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U	.	.
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2-Chloronaphthalene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2-Chlorophenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2-Methylnaphthalene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2-Methylphenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 2-Nitroaniline	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U	.	.
S 2-Nitrophenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 3-Nitroaniline	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U	.	.
S 4,6-Dinitro-2-methylphenol	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U	.	.
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 4-Chloroaniline	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 4-Methylphenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S 4-Nitroaniline	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U	.	.
S 4-Nitrophenol	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U	.	.
S Acenaphthene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Acenaphthylene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Anthracene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Benzo(a)anthracene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Benzo(a)pyrene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Benzo(b,k)fluoranthene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Benzo(g,h,i)perylene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. **Data** not shown in significant digits.

	ENSAFE ID:	13S06D02	13S64D08	13S65D06	13S66D08	13S68M04	13S68S04
	LAB ID:	529876	532097	532101	532117	533211	533212
	DATE RCVD:	12/18/92	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	340.0000 U	180.0000 U	330.0000 U	340.0000 U	.	.
S Carbazole	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Chrysene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Di-n-butylphthalate	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Di-n-octylphthalate	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Dibenz(a,h)anthracene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Dibenzofuran	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Diethylphthalate	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Dimethylphthalate	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Fluoranthene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Fluorene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Hexachlorobenzene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Hexachlorobutadiene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Hexachloroethane	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Indeno(1,2,3-cd)pyrene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Ieophorone	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Naphthalene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Nitrobenzene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Pentachlorophenol	UG/KG	820.0000 U	860.0000 U	800.0000 U	830.0000 U	.	.
S Phenanthrene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Phenol	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S Pyrene	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S bis(2-Ethlyhexy)phthalate	UG/KG	340.0000 U	60.0000 U	330.0000 U	340.0000 U	.	.
S bis(2-chloroethyl)ether	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.
S bis(2-chloroisopropyl)ether	UG/KG	340.0000 U	350.0000 U	330.0000 U	340.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S06D02	13S64D08	13S65D06	13S66D08	13S68M04	13S68S04
	LAB ID:	529876	532097	532101	532117	533211	533212
	DATE RCVD:	12/18/92	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	45.0000	37.0000
V 1,2-Dichloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	66.0000 U	51.0000 U	54.0000 U	75.0000 U	42.0000 U	40.0000 U
V Benzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	57.0000	49.0'000
V Bromodichloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	55.0000	55.0000
V Chloroethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chloroform	UG/KG	10.0000 U	11.0000 U	2.0000 U	10.0000 U	11.0000 U	11.0000 U
V Chloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	22.0000 U	23.0000 U	16.0000 U	20.0000 U	21.0000 U	9.0000 U
V Styrene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	53.0000	51.0000
V Trichloroethene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	55.0000	52.0000

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S06D02	13S64D08	13S65D06	13S66D08	13S68M04	13S68S04
	LAB ID:	529876	532097	532101	532117	533211	533212
	DATE RCVD:	12/18/92	01/15/93	01/15/93	01/15/93	01/20/93	01/20/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	10.0000 U	10.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID:	13S69004MS	13S69004MSD	13S70002MS	13S70002MSD	33BENT01	33GROU01
	LAB ID:	532988	532989	532951	532952	529929	529932
	DATE RCVD: WITS	01/20/93	01/20/93	01/20/93	01/21/93	12/18/92	12/18/92
I Aluminum	MG/KG					10400.0000	19800.0000
I Antimony	MG/KG					12.4000 U	11.5000 U
I Arsenic	MG/KG					7.3000	8.6000
I Barium	MG/KG					69.1000	81.9000
I Beryllium	MG/KG					0.9300 J	0.8000 J
I Cadmium	MG/KG					1.1000 U	1.0000 U
I Calcium	MG/KG					3210.0000	404000.000
I Chromium	MG/KG					2.2000 U	49.3000
I Cobalt	MG/KG					2.4000 U	3.5000 J
I Copper	MG/KG					8.7000	10.2000
I Cyanide	MG/KG					0.5400 U	0.5000 U
I Iron	MG/KG					4400.0000	19200.0000
I Lead	MG/KG					41.4000	2.0000
I Magnesium	MG/KG					1860.0000	16000.0000
I Manganese	MG/KG					46.1000	97.8000
I Mercury	MG/KG					0.1100 U	0.1000 U
I Nickel	MG/KG					8.7000 U	11.8000
I Potassium	MG/KG					1050.0000 U	2950.0000
I Selenium	MG/KG					0.6500 U	3.0000 U
I Silver	MG/KG					2.2000 U	2.0000 U
I Sodium	MG/KG					15900.0000	988.0000 J
I Thallium	MG/KG					0.6500 U	0.6000 U
I Vanadium	MG/KG					4.1000 U	59.3000
I Zinc	MG/KG					63.1000	32.9000
P 4,4'-DDD	UG/KG	4.3000 U	4.3000 U	0.5600 J	0.5600 J	4.4000 U	4.0000 U
P 4,4'-DDE	UG/KG	1.5000 J	1.4000 J	0.6400 J	2.0000 J	4.4000 U	4.0000 U
P 4,4'-DDT	UG/KG	34.0000 J	34.0000 J	37.0000	39.0000	0.2200 J	0.1100 J
P Aldrin	UG/KG	13.0000 J	13.0000 J	15.0000	16.0000	2.2000 U	2.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	13S69004MS / 532995 / 01/20/93	13S69004MSD / 532996 / 01/20/93	13S70002MS / 532951 / 01/20/93	13S70002MSD / 532952 / 01/21/93	33BENT01 / 529929 / 12/18/92	33GROU01 / 529932 / 12/18/92
P Aroclor 1016	UG/KG	43.0000 U	43.0000 U	34.0000 U	34.0000 U	44.0000 U	40.0000 U
P Aroclor 1221	UG/KG	87.0000 U	88.0000 U	70.0000 U	70.0000 U	88.0000 U	81.0000 U
P Aroclor 1232	UG/KG	43.0000 U	43.0000 U	34.0000 U	34.0000 U	44.0000 U	40.0000 U
P Aroclor 1242	UG/KG	43.0000 U	43.0000 U	34.0000 U	34.0000 U	44.0000 U	40.0000 U
P Aroclor 1248	UG/KG	43.0000 U	43.0000 U	34.0000 U	34.0000 U	44.0000 U	40.0000 U
P Aroclor 1254	UG/KG	43.0000 U	43.0000 U	34.0000 U	34.0000 U	44.0000 U	40.0000 U
P Aroclor 1260	UG/KG	5.1000 J	4.2000 J	34.0000 U	34.0000 U	44.0000 U	40.0000 U
P Dieldrin	UG/KG	32.0000 J	32.0000 J	32.0000	36.0000	0.4300 J	0.1700 J
P Endosulfan I	UG/KG	2.2000 U	2.2000 U	1.8000 U	1.8000 U	2.2000 U	2.0000 U
P Endosulfan II	UG/KG	4.3000 U	4.3000 U	3.4000 U	3.4000 U	0.1900 J	4.0000 U
P Endosulfan sulfate	UG/KG	4.3000 U	4.3000 U	3.4000 U	3.4000 U	4.4000 U	4.0000 U
P Endrin	UG/KG	33.0000 J	32.0000 J	34.0000	36.0000	0.2900 J	4.0000 U
P Endrin aldehyde	UG/KG	4.3000 U	4.3000 U	3.4000 U	0.2300 J	4.4000 U	4.0000 U
P Endrin ketone	UG/KG	4.3000 U	4.3000 U	3.4000 U	3.4000 U	4.4000 U	4.0000 U
P Heptachlor	UG/KG	14.0000 J	14.0000 J	17.0000 J	17.0000 J	2.2000 U	0.0690 J
P Heptachlor epoxide	UG/KG	2.2000 U	2.2000 U	1.8000 U	1.8000 U	2.2000 U	2.0000 U
P Methoxychlor	UG/KG	22.0000 U	22.0000 U	18.0000 U	18.0000 U	22.0000 U	20.0000 U
P Toxaphene	UG/KG	220.0000 U	220.0000 U	180.0000 U	180.0000 U	220.0000 U	200.0000 U
P alpha-BHC	UG/KG	2.2000 U	2.2000 U	1.8000 U	1.8000 U	2.2000 U	2.0000 U
P alpha-Chlordane	UG/KG	2.2000 U	2.2000 U	1.8000 U	1.8000 U	2.2000 U	2.0000 U
P beta-BHC	UG/KG	2.2000 U	2.2000 U	1.8000 U	1.8000 U	2.2000 U	2.0000 U
P delta-BHC	UG/KG	2.2000 U	2.2000 U	1.8000 U	1.8000 U	2.2000 U	2.0000 U
P gamma-BHC (Lindane)	UG/KG	14.0000 J	15.0000 J	16.0000	17.0000	0.0550 J	2.0000 U
P gamma-Chlordane	UG/KG	2.2000 U	'2.2000 U	1.8000 U	1.8000 U	0.0880 J	2.0000 U
S 1,2,4-Trichlorobenzene	UG/KG	1500.0000	1400.0000			960.0000 U	330.0000 U
S 1,2-Dichlorobenzene	UG/KG	350.0000 U	350.0000 U			360.0000 U	330.0000 U
S 1,3-Dichlorobenzene	UG/KG	350.0000 U	350.0000 U			360.0000 U	330.0000 U
S 1,4-Dichlorobenzene	UG/KG	1200.0000	1200.0000			360.0000 U	330.0000 U
S 2,4,5-Trichlorophenol	UG/KG	840.0000 U	840.0000 U			860.0000 U	800.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificsnt digits.

	ENSAFE ID:	13S69004MS	13S69004MSD	13S70002MS	13S70002MSD	33BENT01	33GROU01
	LAB ID:	532988	532989	532951	532952	529929	529932
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/21/93	12/18/92	12/18/92
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 2,4-Dichlorophenol	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 2,4-Dimethylphenol	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 2,4-Dinitrophenol	UG/KG	840.0000 U	840.0000 U	.	.	860.0000 U	800.0000 U
S 2,4-Dinitrotoluene	UG/KG	1500.0000	1500.0000	.	.	360.0000 U	330.0000 U
S 2,6-Dinitrotoluene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 2-Chloronaphthalene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 2-Chlorophenol	UG/KG	1900.0000	1700.0000	.	.	360.0000 U	330.0000 U
S 2-Methylnaphthalene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 2-Methylphenol	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 2-Nitroaniline	UG/KG	840.0000 U	840.0000 U	.	.	860.0000 U	800.0000 U
S 2-Nitrophenol	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 3-Nitroaniline	UG/KG	840.0000 U	840.0000 U	.	.	860.0000 U	800.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	840.0000 U	840.0000 U	.	.	860.0000 U	800.0000 U
S 4-Bromophenyl-phenylether	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 4-Chloro-3-methylphenol	UG/KG	2000.0000	1800.0000	.	.	360.0000 U	330.0000 U
S 4-Chloroaniline	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 4-Methylphenol	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S 4-Nitroaniline	UG/KG	840.0000 U	840.0000 U	.	.	860.0000 U	800.0000 U
S 4-Nitrophenol	UG/KG	1700.0000	1400.0000	.	.	860.0000 U	800.0000 U
S Acenaphthene	UG/KG	1200.0000	1200.0000	.	.	360.0000 U	330.0000 U
S Acenaphthylene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Anthracene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Benzo(a)anthracene	UG/KG	350.0000 U	350.0000 U	.	.	3130.0000 U	330.0000 U
S Benzo(a)pyrene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Benzo(b,k)fluoranthene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Benzo(g,h,i)perylene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S69004MS	13S69004MSD	13S70002MS	13S70002MSD	33BENT01	33GROU01
	LAB ID:	532988	532989	532951	532952	529929	529932
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/21/93	12/18/92	12/18/92
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	240.0000 U	170.0000 U	.	.	360.0000 U	330.0000 U
S Carbazole	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Chrysene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Di-n-butylphthalate	UG/KG	350.0000 U	350.0000 U	.	.	54.0000 U	39.0000 U
S Di-n-octylphthalate	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Dibenz (a,h) anthracene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Dibenzofuran	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Diethylphthalate	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Dimethylphthalate	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Fluoranthene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Fluorene	UG/KG	350,0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Hexachlorobenzene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Hexachlorobutadiene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Hexachlorocyclopentadiene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Hexachloroethane	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Ieophorone	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S N-Nitroeo-di-n-propylamine	UG/KG	1300.0000	1100.0000	.	.	360.0000 U	330.0000 U
S N-Nitrosodiphenylamine	UG/KG	350.0000 U	350.0000 U	.	.	56.0000 J	330.0000 U
S Naphthalene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Nitrobenzene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Pentachlorophenol	UG/KG	2600.0000	2000.0000	.	.	860.0000 U	800.0000 U
S Phenanthrene	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S Phenol	UG/KG	2000.0000	1900.0000	.	.	360.0000 U	330.0000 U
S Pyrene	UG/KG	1600.0000	1600.0000	.	.	360.0000 U	330.0000 U
S bie(2-Chloroethoxy)methane	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	46.0000 U	40.0000 U	.	.	66.0000 U	330.0000 U
S bis(2-chloroethyl)ether	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	350.0000 U	350.0000 U	.	.	360.0000 U	330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	WITS	13S69004MS	13S69004MSD	13S70002MS	13S70002MSD	33BENT01	33GROU01
ENSAFE ID:		13S69004MS	13S69004MSD	13S70002MS	13S70002MSD	33BENT01	33GROU01
LAB ID:		533214	533215	532951	532952	529929	529932
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/21/93	12/18/92	12/18/92
V 1,1,1-Trichloroethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 1.1-Dichloroethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/KG	53.0000	50.0000	.	.	11.0000 U	10.0000 U
V 1.2-Dichloroethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 1.2-Dichloroethene (total)	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 1.2-Dichloropropane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 2-Butanone	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 2-Hexanone	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Acetone	UG/KG	130.0000 U	68.0000 U	.	.	11.0000 U	59.0.000 U
V Benzene	UG/KG	67.0000	59.0000	.	.	11.0000 U	10.0000 U
V Bromodichloromethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Bromoform	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Bromomethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Carbon Disulfide	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Carbon Tetrachloride	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Chlorobenzene	UG/KG	61.0000	57.0000	.	.	11.0000 U	10.0000 U
V Chloroethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Chloroform	UG/KG	1.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Chloromethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Dibromochloromethane	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Ethylbenzene	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Methylene Chloride	UG/KG	100.0000 U	19.0000 U	.	.	6.0000 U	16.0000 U
V Styrene	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Tetrachloroethene	UG/KG	11.0000 U	11.0000 U	.	.	11.0000 U	10.0000 U
V Toluene	UG/KG	58.0000	56.0000	.	.	11.0000 U	10.0000 U
V Trichloroethene	UG/KG	65.0000	56.0000	.	.	11.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S69004MS	13S69004MSD	13S70002MS	13S70002MSD	33BENT01	33GROU01
	LAB ID:	533214	533215	532951	532952	529929	529932
	DATE RCM:	01/20/93	01/20/93	01/20/93	01/21/93	12/18/92	12/18/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	11.0000 U	11.0000 U		.	11.0000 U	10.0000 U
V Xylene (total)	UG/KG	11.0000 U	11.0000 U		.	11.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U		.	11.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	11.0000 U	11.0000 U		.	11.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33M03D01	33M03D01DL	33MUD001	33MUDM01	33MUDS01	33S01M02
ENSAFE ID:		33M03D01	33M03D01DL	33MUD001	33MUDM01	33MUDS01	33S01M02
LAB ID:		533352	533352 R1D50	531055	531056	531057	
DATE RCM:		01/22/93	01/22/93	01/06/93	01/06/93	01/06/93	12/10/92
I Aluminum	MG/KG	756.0000	.	9270.0000		.	.
I Antimony	MG/KG	14.7000 U		274.0000 U		.	.
I Arsenic	MG/KG	0.7700 U	.	14.7000 U			.
I Barium	MG/KG	1.5000 U	.	223.0000 J		.	.
I Beryllium	MG/KG	0.7700 U		14.4000 U			.
I Cadmium	MG/KG	1.3000 U	.	24.0000 U		.	.
I Calcium	MG/KG	121.0000 J		9530.0000 J		.	.
I Chromium	MG/KG	5.3000		48.1000 U		.	.
I Cobalt	MG/KG	3.1000 U		57.7000 U		.	.
I Copper	MG/KG	6.4000 U		120.0000 U		.	.
I Cyanide	MG/KG	0.6400 U		12.5000 U		.	.
I Iron	MG/KG	453.0000		5340.0000		.	.
I Lead	MG/KG	2.8000		33.8000		.	.
I Magnesium	MG/KG	142.0000 U		2650.0000 U		.	.
I Manganese	MG/KG	4.0000		134.0000		.	.
I Mercury	MG/KG	0.1300 U		2.5000 U		.	.
I Nickel	MG/KG	9.8000 U		183.0000 U		.	.
I Potassium	MG/KG	1020.0000 U		19000.0000 U		.	.
I Selenium	MG/KG	0.7700'U		14.7000 U		.	.
I Silver	MG/KG	2.6000 U		48.0000 U		.	.
I Sodium	MG/KG	145.0000 J		13200.0000 J		.	.
I Thallium	MG/KG	0.7700 U		14.7000 U		.	.
I Vanadium	MG/KG	3.9000 U		72.1000 U			.
I Zinc	MG/KG	5.6000		52.3000 J			.
P 4,4'-DDD	UG/KG	140.0000 J	150.0000 J	81.0000 U	82.0000 U	3.0000 J	.
P 4,4'-DDE	UG/KG	120.0000 J	130.0000 J	2.3000 J	82.0000 U	81.0000 U	.
P 4,4'-DDT	UG/KG	69.0000 J	72.0000 J	81.0000 U	690.0000 J	840.0000 J	.
P Aldrin	UG/KG	0.1300 J	11.0000 U	42.0000 U	280.0000	340.0000	.

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	33M03D01	33M03D01DL	33MUD001	33MUDM01	33MUDS01	33S01M02
	LAB ID:	533352 R1	533352 R1D50	531055	531062	531063	
	DATE RCVD:	01/22/93	01/22/93	01/06/93	01/06/93	01/06/93	12/10/92
NAME	UNITS						
P Aroclor 1016	UG/KG	42.0000 U	210.0000 U	810.0000 U	820.0000 U	810.0000 U	.
P Aroclor 1221	UG/KG	86.0000 U	430.0000 U	1700.0000 U	1700.0000 U	1600.0000 U	.
P Aroclor 1232	UG/KG	42.0000 U	210.0000 U	810.0000 U	820.0000 U	810.0000 U	.
P Aroclor 1242	UG/KG	42.0000 U	210.0000 U	810.0000 U	820.0000 U	810.0000 U	.
P Aroclor 1248	UG/KG	42.0000 U	210.0000 U	810.0000 U	820.0000 U	810.0000 U	.
P Aroclor 1254	UG/KG	42.0000 U	210.0000 U	810.0000 U	820.0000 U	810.0000 U	.
P Aroclor 1260	UG/KG	42.0000 U	210.0000 U	810.0000 U	820.0000 U	810.0000 U	.
P Dieldrin	UG/KG	0.6100 J	21.0000 U	81.0000 U	610.0000	720.0000	.
P Endosulfan I	UG/KG	2.2000 U	11.0000 U	42.0000 U	42.0000 U	42.0000 U	
P Endosulfan II	UG/KG	4.2000 U	21.0000 U	81.0000 U	82.0000 U	81.0000 U	
P Endosulfan sulfate	UG/KG	4.2000 U	21.0000 U	81.0000 U	82.0000 U	81.0000 U	
P Endrin	UG/KG	4.2000 U	21.0000 U	81.0000 U	610.0000	720.0000 J	
P Endrin aldehyde	UG/KG	0.1500 J	21.0000 U	81.0000 U	16.0000 J	18.0000 J	
P Endrin ketone	UG/KG	4.2000 U	21.0000 U	81.0000 U	13.0000 J	17.0000 J	
P Heptachlor	UG/KG	0.1800 J	11.0000 U	1.8000 J	300.0000	350.0000	
P Heptachlor epoxide	UG/KG	0.0920 J	0.2800 J	42.0000 U	42.0000 U	42.0000 U	.
P Methoxychlor	UG/KG	22.0000 U	110.0000 U	420.0000 U	420.0000 U	7.3000 J	
P Toxaphene	UG/KG	220.0000 U	1100.0000 U	4200.0000 U	4200.0000 U	4200.0000 U	
P alpha-BHC	UG/KG	2.2000 U	11.0000 U	42.0000 U	42.0000 U	42.0000 U	
P alpha-Chlordane	UG/KG	0.9200 J	0.8400 J	42.0000 U	42.0000 U	42.0000 U	.
P beta-BHC	UG/KG	0.3500 J	11.0000 U	42.0000 U	42.0000 U	42.0000 U	
P delta-BHC	UG/KG	2.2000 U	11.0000 U	42.0000 U	42.0000 U	42.0000 U	
P gamma-BHC (Lindane)	UG/KG	2.2000 U	11.0000 U	42.0000 U	280.0000	340.0000	
P gamma-Chlordane	UG/KG	0.2200 J	11.0000 U	42.0000 U	42.0000 U	42.0000 U	.
S 1,2,4-Trichlorobenzene	UG/KG	420.0000 U		8200.0000 U	35000.0000	36000.0000	
S 1,2-Dichlorobenzene	UG/KG	420.0000 U		8200.0000 U	8200.0000 U	8200.0000 U	.
S 1,3-Dichlorobenzene	UG/KG	420.0000 U		8200.0000 U	8200.0000 U	8200.0000 U	.
S 1,4-Dichlorobenzene	UG/KG	420.0000 U		8200.0000 U	30000.0000	30000.0000	.
S 2,4,5-Trichlorophenol	UG/KG	1000.0000 U		20000.0000 U	20000.0000 U	20000.0000 U	.

NAY PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33M03D01	33M03D01DL	33MUD001	33MUDM01	33MUDS01	33S01M02
	LAB ID:	533352	533352 R1D50	531055	531059	531060	
	DATE RCVD:	01/22/93	01/22/93	01/06/93	01/06/93	01/06/93	12/10/92
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 2,4-Dichlorophenol	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 2,4-Dimethylphenol	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 2,4-Dinitrophenol	UG/KG	1000.0000 U	.	20000.0000 U	20000.0000 U	20000.0000 U	.
S 2,4-Dinitrotoluene	UG/KG	420.0000 U	.	8200.0000 U	36000.0000	35000.0000	.
S 2,6-Dinitrotoluene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 2-Chloronaphthalene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 2-Chlorophenol	UG/KG	420.0000 U	.	8200.0000 U	*50000.0000*	48000.0000	.
S 2-Methylnaphthalene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 2-Methylphenol	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 2-Nitroaniline	UG/KG	1000.0000 U	.	20000.0000 U	20000.0000 U	20000.0000 U	.
S 2-Nitrophenol	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 3,3'-Dichlorobenzidine	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 3-Nitroaniline	UG/KG	1000.0000 U	.	20000.0000 U	20000.0000 U	20000.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/KG	1000.0000 U	.	20000.0000 U	20000.0000 U	20000.0000 U	.
S 4-Bromophenyl-phenylether	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 4-Chloro-3-methylphenol	UG/KG	420.0000 U	.	8200.0000 U	58000.0000	54000.0000	.
S 4-Chloroaniline	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 4-Chlorophenyl-phenylether	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 4-Methylphenol	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S 4-Nitroaniline	UG/KG	1000.0000 U	.	20000.0000 U	20000.0000 U	20000.0000 U	.
S 4-Nitrophenol	UG/KG	1000.0000 U	.	20000.0000 U	55000.0000	60000.0000	.
S Acenaphthene	UG/KG	420.0000 U	.	8200.0000 U	35000.0000	34000.0000	.
S Acenaphthylene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000,U	8200.0000 U	.
S Anthracene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Benzo(a)anthracene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Benzo(a)pyrene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Benzo(b,k)fluoranthene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Benzo(g,h,i)perylene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33M03D01	33M03D01DL	33MUD001	33MUDM01	33MUDS01	33S01M02
	LAB ID:	533352	533352 R1D50	531055	531059	531060	
	DATE RCVD:	01/22/93	01/22/93	01/06/93	01/06/93	01/06/93	12/10/92
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	110.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Carbazole	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Chrysene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Di-n-butylphthalate	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Di-n-octylphthalate	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Dibenz(a,h)anthracene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	
S Dibenzofuran	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	
S Diethylphthalate	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	
S Dimethylphthalate	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Fluoranthene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	
S Fluorene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Hexachlorobenzene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Hexachlorobutadiene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Hexachlorocyclopentadiene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Hexachloroethane	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Ieophorone	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S N-Nitroso-di-n-propylamine	UG/KG	420.0000 U	.	8200.0000 U	34000.0000	32000.0000	.
S N-Nitrosodiphenylamine	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Naphthalene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Nitrobenzene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Pentachlorophenol	UG/KG	1000.0000 U	.	20000.0000 U	58000.0000	57000.0000	.
S Phenanthrene	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S Phenol	UG/KG	420.0000 U	.	8200.0000 U	53000.0000	51000.0000	.
S Pyrene	UG/KG	420.0000 U	.	8200.0000 U	37000.0000	33000.0000	.
S bis(2-Chloroethoxy)methane	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S bis(2-Ethylhexy)phthalate	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S bis(2-chloroethyl)ether	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.
S bis(2-chloroisopropyl)ether	UG/KG	420.0000 U	.	8200.0000 U	8200.0000 U	8200.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33M03D01	33M03D01DL	33MUD001	33MUDM01	33MUDS01	33S01M02
	LAB ID:	533352	533352 R1D50	531055	531056	531057	
	DATE RCVD:	01/22/93	01/22/93	01/06/93	01/06/93	01/06/93	12/10/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 1,1,2-Trichloroethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 1,1-Dichloroethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 1,1-Dichloroethene	UG/KG	13.0000 U	.	250.0000 U	1200.0000	1700.0000	120.0000
V 1,2-Dichloroethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 1,2-Dichloroethene (total)	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 1,2-Dichloropropane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 2-Butanone	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 2-Hexanone	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V 4-Methyl-2-Pentanone	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Acetone	UG/KG	37.0000 U	.	250.0000 U	250.0000 U	640.0000	36.0000 U
V Benzene	UG/KG	13.0000 U	.	250.0000 U	1500.0000	1500.0000	110.0000
V Bromodichloromethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Bromoform	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Bromomethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Carbon Disulfide	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Carbon Tetrachloride	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Chlorobenzene	UG/KG	13.0000 U	.	250.0000 U	1500.0000	1400.0000	240.0000
V Chloroethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Chloroform	UG/KG	13.0000 U	.	35.0000 U	27.0000 U	41.0000 U	3.0000 U
V Chloromethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Dibromochloromethane	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Ethylbenzene	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Methylene Chloride	UG/KG	12.0000 U	.	480.0000 J	530.0000 U	1100.0000 J	57.0000 U
V Styrene	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Tetrachloroethene	UG/KG	13.0000 U	.	250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Toluene	UG/KG	13.0000 U	.	66.0000 J	1500.0000	1500.0000	100.0000
V Trichloroethene	UG/KG	13.0000 U	.	250.0000 U	1600.0000	1700.0000	93.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33M03D01	33M03D01DL	33MUD001	33MUDM01	33MUDS01	33S01M02
	LAB ID:	533352	533352 R1D50	531055	531056	531057	
	DATE RCVD:	01/22/93	01/22/93	01/06/93	01/06/93	01/06/93	12/10/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	13.0000 U		250.0000 U	250.0000 U	250.0000 U	19.0000 U
V Xylene (total)	UG/KG	13.0000 U		250.0000 U	250.0000 U	250.0000 U	19,0000 U
V cis-1,3-Dichloropropene	UG/KG	13.0000 U		250.0000 U	250.0000 U	250.0000 U	19.0000 U
V trans-1,3-Dichloropropene	UG/KG	13.0000 U		250.0000 U	250.0000 U	250.0000 U	19.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S01S02	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
	LAB ID:				529101	528694	528695
	DATE RCVD:	12/10/92	12/10/92	12/14/92	12/10/92	12/12/92	12/12/92
NAME	UNITS						
I Aluminum	MG/KG	.	.	.	10300.0000	.	
I Antimony	MG/KG	.	.	.	18.3000 U		
I Arsenic	MG/KG	.	.	.	1.5000 J		
I Barium	MG/KG	.	.	.	118.0000		
I Beryllium	MG/KG	.	.	.	0.3100 U		
I Cadmium	MG/KG	.	.	.	33.7000		
I Calcium	MG/KG	.	.	.	4610.0000		
I Chromium	MG/KG	.	.	.	915.0000		
I Cobalt	MG/KG	.	.	.	2.8000 U		
I Copper	MG/KG	.	.	.	137.0000		
I Cyanide	MG/KG	.	.	.	2.0000		
I Iron	MG/KG	.	.	.	3700.0000		
I Lead	MG/KG	.	.	.	140.0000		
I Magnesium	MG/KG	.	.	.	245.0000 J		
I Manganese	MG/KG	.	.	.	77.4000		
I Mercury	MG/KG	.	.	.	3.1000		
I Nickel	MG/KG	.	.	.	17.8000		
I Potassium	MG/KG	.	.	.	417.0000 U		
I Selenium	MG/KG	.	.	.	0.9200 U		
I Silver	MG/KG	.	.	.	33.4000 U		
I Sodium	MG/KG	.	.	.	150.0000 J		
I Thallium	MG/KG	.	.	.	0.6100 U		
I Vanadium	MG/KG	.	.	.	7.6000 J		
I Zinc	MG/KG	.	.	.	257.0000		
P 4,4'-DDD	UG/KG	.	.	.	25.0000 U	3.600.0 U	3.6000 U
P 4,4'-DDE	UG/KG	.	.	.	5.1000 J	3.6000 U	0.1600 J
P 4,4'-DDT	UG/KG	.	.	.	25.0000 U	26.0000 J	23.0000 J
P Aldrin	UG/KG	.	.	.	13.0000 U	11.0000	10.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S01S02	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
ENSAFE ID:		33S01S02	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
LAB ID:					529101 D50	528700 R1	528701 R1
DATE RCVD:		12/10/92	12/10/92	12/10/92	12/10/92	12/12/92	12/12/92
P Aroclor 1016	UG/KG		.	.	250.0000 U	36.0000 U	36.0000 U
P Aroclor 1221	UG/KG		.	.	510.0000 U	73.0000 U	73.0000 U
P Aroclor 1232	UG/KG			.	250.0000 U	36.0000 U	36.0000 U
P Aroclor 1242	UG/KG			.	250.0000 U	36.0000 U	36.0000 U
P Aroclor 1248	UG/KG			.	250.0000 U	36.0000 U	36.0000 U
P Aroclor 1254	UG/KG			.	250.0000 U	36.0000 U	36.0000 U
P Aroclor 1260	UG/KG	.	.	.	200.0000 J	36.0000 U	36.0000 U
P Dieldrin	UG/KG	.	.	.	110.0000	31.0000	29.0000
P Endosulfan I	UG/KG			.	13.0000 U	1.8000 U	1.9000 U
P Endosulfan II	UG/KG			.	25.0000 U	0.3400 J	0.3400 J
P Endosulfan sulfate	UG/KG		.	.	25.0000 U	0.1200 J	0.1000 J
P Endrin	UG/KG		.	.	1.4000 J	24.0000	22.0000
P Endrin aldehyde	UG/KG		.	.	25.0000 U	0.4500 J	0.7000 J
P Endrin ketone	UG/KG	.	.	.	25.0000 U	3.6000 U	3.6000 U
P Heptachlor	UG/KG	.	.	.	0.4400 J	11.0000	10.0000
P Heptachlor epoxide	UG/KG		.	.	13.0000 U	0.3400 J	0.2600 J
P Methoxychlor	UG/KG		.	.	130.0000 U	18.0000 U	0.5300 J
P Toxaphene	UG/KG		.	.	1300.0000 U	180.0000 U	190.0000 U
P alpha-BHC	UG/KG	.	.	.	13.0000 U	0.2100 J	0.2500 J
P alpha-Chlordane	UG/KG	.	.	.	56.0000 J	0.6300 J	0.6500 J
P beta-BHC	UG/KG	.	.	.	13.0000 U	1.8000 U	1.9000 U
P delta-BHC	UG/KG		.	.	2.9000 J	1.8000 U	1.9000 U
P gama-BHC (Lindane)	UG/KG	.	.	.	13.0000 U	12.0000	11.0000
P gama-Chlordane	UG/KG	.	.	.	73.0000 J	1.0000 J	1.0000 J
S 1,2,4-Trichlorobenzene	UG/KG	.	.	.	10000.0000 U	1400.0000	1300.0000
S 1,2-Dichlorobenzene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 1,3-Dichlorobenzene	UG/KG		.	.	10000.0000 U	360.0000 U	360.0000 U
S 1,4-Dichlorobenzene	UG/KG		.	.	10000.0000 U	1300.0000	1200.0000
S 2,4,5-Trichlorophenol	UG/KG	.	.	.	24000.0000 U	880.0000 U	880.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33S01S02	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
	LAB ID:				529101	528697	528698
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92	12/12/92
S 2,4,6-Trichlorophenol	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 2,4-Dichlorophenol	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 2,4-Dimethylphenol	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 2,4-Dinitrophenol	UG/KG	.	.	.	24000.0000 U	880.0000 U	880.0000 U
S 2,4-Dinitrotoluene	UG/KG	.	.	.	10000.0000 U	1400.0000	1300.0000
S 2,6-Dinitrotoluene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 2-Chloronaphthalene	UG/KG	.		.	10000.0000 U	360.0000 U	360.0000 U
S 2-Chlorophenol	UG/KG	.		.	10000.0000 U	1700.0000	1700.0000
S 2-Methylnaphthalene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 2-Methylphenol	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 2-Nitroaniline	UG/KG	.	.	.	24000.0000 U	880.0000 U	880.0000 U
S 2-Nitrophenol	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	.	.	.	10000.0000 U	360.0060 U	360.0000 U
S 3-Nitroaniline	UG/KG	.	.	.	24000.0000 U	880.0000 U	880.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	.	.	.	24000.0000 U	880.0000 U	880.0000 U
S 4-Bromophenyl-phenylether	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 4-Chloro-3-methylphenol	UG/KG	.	.	.	10000.0000 U	2100.0000	2100.0000
S 4-Chloroaniline	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 4-Chlordphenyl-phenylether	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 4-Methylphenol	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S 4-Nitroaniline	UG/KG	.	.	.	24000.0000 U	880.0000 U	880.0000 U
S 4-Nitrophenol	UG/KG	.	.	.	24000.0000 U	2100.0000	2000.0000
S Acenaphthene	UG/KG	.	.	.	10000.0000 U	1300.0000	1300.0000
S Acenaphthylene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Anthracene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Benzo(a)anthracene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Benzo(a)pyrene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Benzo(b,k)fluoranthene	UG/KG	.	.	.	10000.0000 U	72.0000 J	57.0000 J
S Benzo(g,h,i)perylene	UG/KG	.	.	.	10000.0000 U	120.0000 J	100.0000 J

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S01S02	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
	LAB ID:				529101	528697	528698
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92	12/12/92
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	.	.	.	1300.0000 U	360.0000 U	360.0000 U
S Carbazole	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Chryeene	UG/KG		.	.	10000.0000 U	360.0000 U	360.0000 U
S Di-n-butylphthalate	UG/KG	.	.		10000.0000 U	360.0000 U	360.0000 U
S Di-n-octylphthalate	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Dibenz (a,h) anthracene	UG/KG	.	.		10000.0000 U	360.0000 U	360.0000 U
S Dibenzofuran	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Diethylphthalate	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Dimethylphthalate	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Fluoranthene	UG/KG	.	.	.	10000.0000 U	43.0000 J	360.0000 U
S Fluorene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Hexachlorobenzene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Hexachlorobutadiene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Hexachlorocyclopentadiene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Hexachloroethane	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	.	.	.	10000.0000 U	85.0000 J	75.0000 J
S Ieophorone	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	.	.	.	10000.0000 U	1200.0000	1200.0000
S N-Nitrosodiphenylamine	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Naphthalene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Nitrobenzene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Pentachlorophenol	UG/KG	.	.	.	24000.0000 U	2000.0000	2000.0000
S Phenanthrene	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S Phenol	UG/KG	.	.	.	10000.0000 U	1600.0000	1500.0000
S Pyrene	UG/KG	.	.	.	10000.0000 U	1700.0000	1700.0000
S bis(2-Chloroethoxy)methane	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	.	.	.	3200.0000 U	250.0000 U	230.0000 U
S bis(2-chloroethyl)ether	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	.	.	.	10000.0000 U	360.0000 U	360.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33501502	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
	LAB ID:				529101	528694	528695
	DATE RCVD:	12/10/92	12/10/92	12/10/92	12/10/92	12/12/92	12/12/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 1,1-Dichloroethene	UG/KG	120.0000	11000.0000	11000.0000	15.0000 U	62.0000	59.0000
V 1,2-Dichloroethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 1,2-Dichloropropane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 2-Butanone	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 2-Hexanone	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Acetone	UG/KG	54.0000 U	9500.0000 U	9100.0000 U	27.0000 U	990.0000	1000.0000 U
V Benzene	UG/KG	120.0000	7400.0000	7600.0000	15.0000 U	64.0000	62.0000
V Bromodichloromethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Bromoform	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Bromomethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Carbon Disulfide	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Carbon Tetrachloride	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Chlorobenzene	UG/KG	140.0000	6500.0000	6400.0000	15.0000 U	57.0000	55.0000
V Chloroethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Chloroform	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	2.0000 U	1.0000 U	1.0000 U
V Chloromethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Dibromochloromethane	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Ethylbenzene	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Methylene Chloride	UG/KG	39.0000 U	640.0000 U	780.0000 U	32.0000 U	39.0000 U	67.0000 U
V Styrene	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Tetrachloroethene	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Toluene	UG/KG	110.0000	6900.0000	6700.0000	15.0000 U	57.0000	55.0000
V Trichloroethene	UG/KG	100.0000	6800.0000	7000.0000	15.0000 U	58.0000	54.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S01S02	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
ENSAFE ID:		33S01S02	33S02001MSD	33S02M01	33S03D02	33S10M02	33S10S02
LAB ID:					529101	528694	528695
DATE RCVD:		12/10/92	12/10/92	12/10/92	12/10/92	12/12/92	12/12/92
V Vinyl Chloride	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V Xylene (total)	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V cis-1,3-Dichloropropene	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U
V trans-1,3-Dichloropropene	UG/KG	19.0000 U	1300.0000 U	1300.0000 U	15.0000 U	11.0000 U	11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
	LAB ID:	530138	530145	530146	530664	529346	529536
	DATE RCVD:	12/19/92	12/19/92	12/19/92	12/18/92	12/16/92	12/17/92
NAME	UNITS						
I Aluminum	MG/KG	2240.0000			.	.	8750.0000
I Antimony	MG/KG	11.9000 U	.		.	.	13.0000 U
I Arsenic	MG/KG	0.8000 J	.		.	.	1.4000 J
I Barium	MG/KG	18.4000 J	.		.	.	15.0000 J
I Beryllium	MG/KG	0.4200 U	.		.	.	0.2200 U
I Cadmium	MG/KG	1.0000 U			.	.	1.1000 U
I Calcium	MG/KG	250.0000 J	.		.	.	670.0000 J
I Chromium	MG/KG	12.7000			.	.	9.7000
I Cobalt	MG/KG	2.3000 U	.		.	.	1.9000 U
I Copper	MG/KG	5.2000 U	.		.	.	5.1000 J
I Cyanide	MG/KG	0.5200 U	.		.	.	0.5400 U
I Iron	MG/KG	1800.0000	.		.	.	4330.0000
I Lead	MG/KG	41.3000	.		.	.	40.7000
I Magnesium	MG/KG	98.2000 J			.	.	172.0000 J
I Manganese	MG/KG	6.3000	.		.	.	194.0000
I Mercury	MG/KG	0.1000 U	.		.	.	0.1100 U
I Nickel	MG/KG	8.3000 U	.		.	.	2.4000 U
I Potassium	MG/KG	1010.0000 U	.		.	.	295.0000 U
I Selenium	MG/KG	0.6200 U	.		.	.	0.6500 U
I Silver	MG/KG	2.1000 U	.		.	.	1.3000 U
I sodium	MG/KG	96.4000 J	.		.	.	85.7000 J
I Thallium	MG/KG	0.6200 U	.		.	.	0.4300 U
I Vanadium	MG/KG	4.0000 U	.		.	.	11.4000
I Zinc	MG/KG	16.4000	.		.	.	408.0000
P 4,4'-DDD	UG/KG	3.4000 U	.		.	.	2.0000 J
P 4,4'-DDE	UG/KG	0.6600 J	.		.	.	190.0000 J
P 4,4'-DDT	UG/KG	1.8000 J	.		.	.	120.0000 J
P Aldrin	UG/KG	1.7000 U	.		.	.	1.8000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
ENSAFE ID:		33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
LAB ID:		530138	530145	530146	530664	529346	529536
DATE RCVD:		12/19/92	12/19/92	12/19/92	12/18/92	12/16/92	12/17/92
P Aroclor 1016	UG/KG	34.0000 U	.		.	.	36.0000 U
P Aroclor 1221	UG/KG	69.0000 U		.	.	.	73.0000 U
P Aroclor 1232	UG/KG	34.0000 U	.		.	.	36.0000 U
P Aroclor 1242	UG/KG	34.0000 U			.	.	36.0000 U
P Aroclor 1248	UG/KG	34.0000 U			.	.	36.0000 U
P Aroclor 1254	UG/KG	34.0000 U	.	.	.	.	36.0000 U
P Aroclor 1260	UG/KG	34.0000 U	.	.	.	.	36.0000 U
P Dieldrin	UG/KG	1.9000 J	.	.	.	.	3.6000 U
P Endosulfan I	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P Endosulfan II	UG/KG	0.2100 J	.	.	.	.	3.6000 U
P Endosulfan sulfate	UG/KG	3.4000 U	.	.	.	.	5.4000 J
P Endrin	UG/KG	3.4000 U	.	.	.	.	0.0610 J
P Endrin aldehyde	UG/KG	3.4000 U	.	.	.	.	1.1000 J
P Endrin ketone	UG/KG	0.6200 J	.	.	.	.	0.1800 J
P Heptachlor	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P Heptachlor epoxide	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P Methoxychlor	UG/KG	7.1000 J	.	.	.	.	3.1000 J
P Toxaphene	UG/KG	170.0000 U	.	.	.	.	180.0000 U
P alpha-BHC	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P alpha-Chlordane	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P beta-BHC	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P delta-BHC	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P gamma-BHC (Lindane)	UG/KG	1.7000 U	.	.	.	.	1.8000 U
P gamma-Chlordane	UG/KG	0.2900 J			.		1.8000 U
S 1,2,4-Trichlorobenzene	UG/KG	340.0000 U	1900.0000	1700.0000	.	2500.0000 U	350.0000 U
S 1,2-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 1,3-Dichlorobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 1,4-Dichlorobenzene	UG/KG	340.0000 U	1500.0000	1400.0000	.	2500.0000 U	350.0000 U
S 2,4,5-Trichlorophenol	UG/KG	830.0000 U	840.0000 U	850.0000 U	.	6200.0000 U	860.0000 U

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
ENSAFE ID:		33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
LAB ID:		530138	530148	530149	530664	529346	529536
DATE RCVD:		12/19/92	12/19/92	12/19/92	12/18/92	12/16/92	12/17/92
S 2,4,6-Trichlorophenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 2,4-Dichlorophenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 2,4-Dimethylphenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 2,4-Dinitrophenol	UG/KG	830.0000 U	840.0000 U	850.0000 U	.	6200.0000 U	860.0000 U
S 2,4-Dinitrotoluene	UG/KG	340.0000 U	1500.0000	1400.0000	.	2500.0000 U	350.0000 U
S 2,6-Dinitrotoluene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 2-Chloronaphthalene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 2-Chlorophenol	UG/KG	340.0000 U	2700,0000	2400.0000	.	2500.0000 U	350.0000 U
S 2-Methylnaphthalene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 2-Methylphenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 2-Nitroaniline	UG/KG	830.0000 U	840.0000 U	850.0000 U	.	6200.0000 U	860.0000 U
S 2-Nitrophenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 3-Nitroaniline	UG/KG	830.0000 U	840.0000 U	850.0000 U	.	6200.0000 U	860.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	830.0000 U	840.0000 U	850.0000 U	.	6200.0000 U	860.0000 U
S 4-Bromophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 4-Chloro-3-methylphenol	UG/KG	340.0000 U	3300.0000 J	3100.0000 J	.	2500.0000 U	350.0000 U
S 4-Chloroaniline	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 4-Methylphenol	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S 4-Nitroaniline	UG/KG	830.0000 U	840.0000 U	850.0000 U	.	6200.0000 U	860.0000 U
S 4-Nitrophenol	UG/KG	830.0000 U	3300.0000 J	3200.0000 J	.	6200.0000 U	860.0000 U
S Acenaphthene	UG/KG	340.0000 U	1800.0000	1600.0000	.	2500.0000 U	350.0000 U
S Acenaphthylene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Anthracene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Benzo(a)anthracene	UG/KG	52.0000 J	350.0000 U	350.0000 U	.	3100.0000 J	120.0000 J
S Benzo(a)pyrene	UG/KG	46.0000 J	350.0000 U	350.0000 U	.	3400.0000 J	120.0000 J
S Benzo(b,k)fluoranthene	UG/KG	94.0000 J	350.0000 U	350.0000 U	.	5900.0000 J	300,0000 J
S Benzo(g,h,i)perylene	UG/KG	61.0000 J	350.0000 U	350.0000 U	.	2600.0000 J	220.0000 J

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

NAME	UNITS	33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
ENSAFE ID:		33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
LAB ID:		530138	530148	530149	530664	529346	529536
DATE RCVD:		12/19/92	12/19/92	12/19/92	12/18/92	12/16/92	12/17/92
S Butylbenzylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Carbazole	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Chrysene	UG/KG	42.0000 J	350.0000 U	350.0000 U	.	2800.0000 J	74.0000 J
S Di-n-butylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Di-n-octylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Dibenz (a,h) anthracene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	610.0000 J	47.0000 J
S Dibenzofuran	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Diethylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Dimethylphthalate	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Fluoranthene	UG/KG	56.0000 J	350.0000 U	350.0000 U	.	3400.0000 J	120.0000 J
S Fluorene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Hexachlorobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Hexachlorobutadiene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Hexachlorocyclopentadiene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Hexachloroethane	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Indeno(1,2,3-cd) pyrene	UG/KG	46.0000 J	350.0000 U	350.0000 U	.	3000.0000 J	250.0000 J
S Isophorone	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	340.0000 U	1600.0000	1400.0000	.	2500.0000 U	350.0000 U
S N-Nitrosodiphenylamine	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Naphthalene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Nitrobenzene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S Pentachlorophenol	UG/KG	830.0000 U	3700.0000 J	3500.0000 J	.	6200.0000 U	860.0000 U
S Phenanthrene	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	760.0000 J	350.0000 U
S Phenol	UG/KG	340.0000 U	2900.0000 J	2600.0000	.	2500.0000 U	350.0000 U
S Pyrene	UG/KG	56.0000 J	1900.0000	1800.0000	.	4400.0000 J	130.0000 J
S bis(2-Chloroethoxy)methane	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S bie (2-Ethlyhexy) phthalate	UG/KG	79.0000 U	46.0000 U	56.0000 U	.	2500.0000 U	43.0000 U
S bis(2-chloroethyl) ether	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U
S bis(2-chloroisopropyl) ether	UG/KG	340.0000 U	350.0000 U	350.0000 U	.	2500.0000 U	350.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
	LAB ID:	530138	530145	530146	530664	529346	529536
	DATE RCVD:	12/19/92	12/19/92	12/19/92	12/18/92	12/16/92	12/17/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 1,1,2-Trichloroethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300,0000 U	.	11.0000 U
V 1,1-Dichloroethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 1,1-Dichloroethene	UG/KG	10.0000 U	62.0000	67.0000	11000.0000	.	11.0000 U
V 1,2-Dichloroethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 1,2-Dichloroethene (total)	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 1,2-Dichloropropane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 2-Butanone	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 2-Hexanone	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V 4-Methyl-2-Pentanone	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Acetone	UG/KG	21.0000 U	6.0000 U	11.0000 U	1300.0000 U	.	6.0000 U
V Benzene	UG/KG	10.0000 U	58.0000	57.0000	6900.0000	.	11.0000 U
V Bromodichloromethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Bromoform	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Bromomethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Carbon Disulfide	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Carbon Tetrachloride	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Chlorobenzene	UG/KG	10.0000 U	55.0000	56.0000	6600.0000	.	11.0000 U
V Chloroethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Chloroform	UG/KG	10.0000 U	11.0000 U	2.0000 U	1300.0000 U	.	11.0000 U
V Chloromethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Dibromochloromethane	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Ethylbenzene	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Methylene Chloride	UG/KG	5.0000 U	25.0000 U	44.0000 U	480.0000 U	.	45.0000 U
V Styrene	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Tetrachloroethene	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U	.	11.0000 U
V Toluene	UG/KG	10.0000 U	57.0000	59.0000	6500.0000	.	11.0000 U
V Trichloroethene	UG/KG	10.0000 U	51.0000	53.0000	6600.0000	.	11.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S12D02	33S12M04	33S12S04	33S14M02	33S20D02	33S23D02
	LAB ID:	530138	530145	530146	530664	529346	529536
	DATE RCVD:	12/19/92	12/19/92	12/19/92	12/18/92	12/16/92	12/17/92
NAME	UNITS						
V Vinyl Chloride	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U		11.0000 U
V Xylene (total)	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U		11.0000 U
V cis-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U		11.0000 U
V trans-1,3-Dichloropropene	UG/KG	10.0000 U	11.0000 U	11.0000 U	1300.0000 U		11.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCW:	UNITS	33S23D02DL / 529536 D11 / 12/17/92	33S23M02 / 529542 / 12/17/92	33S23M04 / 529545 / 12/17/92	33S23S02 / 529543 / 12/17/92	33S23S04 / 529546 / 12/17/92	33833D02 / 533006 / 01/20/93
I Aluminum		MG/KG	.	.	.	.	.	6320.0000
I Antimony		MG/KG	.	.	.	.	.	13.2000 U
I Arsenic		MG/KG	.	.	.	.	.	1.4000 J
I Barium		MG/KG	.	.	.	.	.	29.5000 J
I Beryllium		MG/KG	.	.	.	.	.	0.7000 U
I Cadmium		MG/KG	.	.	.	.	.	23.4000
I Calcium		MG/KG	.	.	.	.	.	6890.0000
I Chromium		MG/KG	.	.	.	.	.	722.0000
I Cobalt		MG/KG	.	.	.	.	.	2.8000 U
I Copper		MG/KG	.	.	.	.	.	47.8000
I Cyanide		MG/KG	.	.	.	.	.	0.5800 U
I Iron		MG/KG	.	.	.	.	.	2270.0000
I Lead		MG/KG	.	.	.	.	.	58.9000
I Magnesium		MG/KG	.	.	.	.	.	265.0000 J
I Manganese		MG/KG	.	.	.	.	.	36.2000
I Mercury		MG/KG	.	.	.	.	.	0.4100
I Nickel		MG/KG	.	.	.	.	.	42.2000
I Potassium		MG/KG	.	.	.	.	.	916.0000 U
I Selenium		MG/KG	.	.	.	.	.	0.9300 U
I Silver		MG/KG	.	.	.	.	.	9.5000 U
I Sodium		MG/KG	.	.	.	.	.	142.0000 J
I Thallium		MG/KG	.	.	.	.	.	0.7000 U
I Vanadium		MG/KG	.	.	.	.	.	6.0000 J
I Zinc		MG/KG	.	.	.	.	.	123.0000
P 4,4'-DDD		UG/KG	8.1000 J	4.1000 J	.	1.7000 J	.	3.8000 U
P 4,4'-DDE		UG/KG	490.0000 J	42.0000	.	44.0000	.	3.3000 J
P 4,4'-DDT		UG/KG	100.0000 J	45.0000 J	.	47.0000 J	.	3.8000 U
P Aldrin		UG/KG	0.7200 J	12.0000	.	13.0000	.	0.5100 J

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S23D02DL	33S23M02	33S23M04	33S23S02	33S23S04	33833D02
ENSAFE ID:		33S23D02DL	33S23M02	33S23M04	33S23S02	33S23S04	33833D02
LAB ID:		529536 D11	529548	529545	529549	529546	533006 R1
DATE RCVD:		12/17/92	12/17/92	12/17/92.	12/17/92	12/17/92	01/20/93
P Aroclor 1016	UG/KG	360.0000 U	37.0000 U		36.0000 U		38.0000 U
P Aroclor 1221	UG/KG	730.0000 U	74.0000 U		73.0000 U		78.0000 U
P Aroclor 1232	UG/KG	360.0000 U	37.0000 U		36.0000 U		38.0000 U
P Aroclor 1242	UG/KG	360.0000 U	37.0000 U		36.0000 U		38.0000 U
P Aroclor 1248	UG/KG	360.0000 U	37.0000 U		36.0000 U		38.0000 U
P Aroclor 1254	UG/KG	360.0000 U	37.0000 U		36.0000 U		38.0000 U
P Aroclor 1260	UG/KG	360.0000 U	37.0000 U		36.0000 U		77.0000
P Dieldrin	UG/KG	36.0000 U	28.0000		29.0000		15.0000 J
P Endosulfan I	UG/KG	18.0000 U	1.9000 U		1.9000 U		2.0000 U
P Endosulfan II	UG/KG	36.0000 U	3.7000 U		3.6000 U		3.1000 J
P Endosulfan sulfate	UG/KG	36.0000 U	3.7000 U		3.6000 U		0.6700 J
P Endrin	UG/KG	36.0000 U	28.0000		29.0000 J		0.3200 J
P Endrin aldehyde	UG/KG	36.0000 U	0.6800 J		0.6400 J		2.9000 J
P Endrin ketone	UG/KG	36.0000 U	0.5300 J		0.5700 J		3.8000 U
P Heptachlor	UG/KG	18.0000 U	12.0000 J		13.0000 J		0.2100 J
P Heptachlor epoxide	UG/KG	18.0000 U	1.9000 U		0.1600 J		0.3000 J
P Methoxychlor	UG/KG	180.0000 U	19.0000 U		19.0000 U		20.0000 U
P Toxaphene	UG/KG	1800.0000 U	190.0000 U		190.0000 U		200.0000 U
P alpha-BHC	UG/KG	18.0000 U	1.9000 U		1.9000 U		0.0790 J
P alpha-Chlordane	UG/KG	0.9700 J	1.9000 U		1.9000 U		5.0000 J
P beta-BHC	UG/KG	18.0000 U	1.9000 U		1.9000 U		2.0000 U
P delta-BHC	UG/KG	18.0000 U	1.9000 U		0.2700 J		0.4100 J
P gamma-BHC (Lindane)	UG/KG	18.0000 U	13.0000 J		13.0000 J		2.0000 U
P gamma-Chlordane	UG/KG	18.0000 U	1.9000 U		1.9000 U		5.3000 J
S 1,2,4-Trichlorobenzene	UG/KG			1800.0000		2000.0000	380.0000 U
S 1,2-Dichlorobenzene	UG/KG			370.0000 U		370.0000 U	56.0000 J
S 1,3-Dichlorobenzene	UG/KG	.		370.0000 U		370.0000 U	210.0000 J
S 1,4-Dichlorobenzene	UG/KG	.		1700.0000		1800.0000	360.0000 J
S 2,4,5-Trichlorophenol	UG/KG	.		900.0000 U		890.0000 U	920.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	33S23D02DL	33S23M02	33S23M04	33S23S02	33S23S04	33S33D02
	LAB ID:	529536 D11	529542	529545	529543	529546	533006
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/17/92	12/17/92	01/20/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 2,4-Dichlorophenol	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 2,4-Dimethylphenol	UG/KG	.	.	370.0000 U	.	370.0000 U	67.0000 J
S 2,4-Dinitrophenol	UG/KG	.	.	900.0000 U	.	890.0000 U	920.0000 U
S 2,4-Dinitrotoluene	UG/KG	.	.	1900.0000	.	2100.0000	380.0000 U
S 2,6-Dinitrotoluene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 2-Chloronaphthalene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 2-Chlorophenol	UG/KG	.	.	2600.0000	.	2800.0000	380.0000 U
S 2-Methylnaphthalene	UG/KG	.	.	370.0000 U	.	370.0000 U	350.0000 J
S 2-Methylphenol	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 2-Nitroaniline	UG/KG	.	.	900.0000 U	.	890.0000 U	920.0000 U
S 2-Nitrophenol	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 3-Nitroaniline	UG/KG	.	.	900.0000 U	.	890.0000 U	920.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	.	.	900.0000 U	.	890.0000 U	920.0000 U
S 4-Bromophenyl-phenylether	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 4-Chloro-3-methylphenol	UG/KG	.	.	3000.0000	.	3100.0000 J	380.0000 U
S 4-Chloroaniline	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 4-Methylphenol	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S 4-Nitroaniline	UG/KG	.	.	900.0000 U	.	890.0000 U	920.0000 U
S 4-Nitrophenol	UG/KG	.	.	3000.0000	.	3200.0000 J	920.0000 U
S Acenaphthene	UG/KG	.	.	1800.0000	.	1900.0000	380.0000 U
S Acenaphthylene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Anthracene	UG/KG	.	.	370.0000 U	.	370.0000 U	48.0000 J
S Benzo(a)anthracene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Benzo(a)pyrene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Benzo(b,k)fluoranthene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Benzo(g,h,i)perylene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S23D02DL	33S23M02	33S23M04	33S23S02	33S23S04	33S33D02
	LAB ID:	529536 D11	529542	529545	529543	529546	533006
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/17/92	12/17/92	01/20/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	.	.	370.0000 U	.	370.0000 U	340.0000 U
S Carbazole	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Chrysene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Di-n-butylphthalate	UG/KG	.	.	95.0000 U	.	370.0000 U	380.0000 U
S Di-n-octylphthalate	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Dibenz(a,h)anthracene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Dibenzofuran	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Diethylphthalate	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Dimethylphthalate	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Fluoranthene	UG/KG	.	.	370.0000 U	.	370.0000 U	88.0000 J
S Fluorene	UG/KG	.	.	370.0000 U	.	370.0000 U	53.0000 J
S Hexachlorobenzene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Hexachlorobutadiene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Hexachlorocyclopentadiene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Hexachloroethane	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Isophorone	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	.	.	1900.0000	.	2000.0000	380.0000 U
S N-Nitrosodiphenylamine	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Naphthalene	UG/KG	.	.	370.0000 U	.	370.0000 U	110.0000 J
S Nitrobenzene	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S Pentachlorophenol	UG/KG	.	.	2600.0000	.	2700.0000	920.0000 U
S Phenanthrene	UG/KG	.	.	370.0000 U	.	370.0000 U	140.0000 J
S Phenol	UG/KG	.	.	2900.0000	.	2900.0000	160.0000 J
S Pyrene	UG/KG	.	.	1900.0000	.	2100.0000	79.0000 J
S bis(2-Chloroethoxy)methane	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	.	.	51.0000 U	.	370.0000 U	260.0000 U
S bis(2-chloroethyl)ether	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	.	.	370.0000 U	.	370.0000 U	380.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	33S23D02DL	33S23M02	33S23M04	33S23S02	33S23S04	33S33D02
	LAB ID:	529536 D11	529542	529545	529543	529546	533006
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/17/92	12/17/92	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG		11.0000 U		11.0000 U	.	12.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG		11.0000 U	.	11.0000 U	.	12.0000 U
V 1,1,2-Trichloroethane	UG/KG		11.0000 U	.	11.0000 U	.	12.0000 U
V 1,1-Dichloroethane	UG/KG		11.0000 U	.	11.0000 U	.	12.0000 U
V 1,1-Dichloroethene	UG/KG	.	71.0000		65.0000	.	12.0000 U
V 1,2-Dichloroethane	UG/KG		11.0000 U	.	11.0000 U	.	12.0000 U
V 1,2-Dichloroethene (total)	UG/KG		11.0000 U	.	11.0000 U	.	12.0000 U
V 1,2-Dichloropropane	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V 2-Butanone	UG/KG		11.0000 U	.	11.0000 U	.	27.0000 U
V 2-Hexanone	UG/KG		11.0000 U	.	11.0000 U	.	12.0000 U
V 4-Methyl-2-Pentanone	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Acetone	UG/KG	.	43.0000 U	.	38.0000 U	.	150.0000 U
V Benzene	UG/KG	.	63.0000	.	61.0000	.	12.0000 U
V Bromodichloromethane	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Bromoform	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Bromomethane	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Carbon Disulfide	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Carbon Tetrachloride	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Chlorobenzene	UG/KG		56.0000	.	56.0000	.	410.0000 J
V Chloroethane	UG/KG		11.0000 U		11.0000 U	.	12.0000 U
V Chloroform	UG/KG	.	11.0000 U	.	2.0000 U	.	1.0000 U
V Chloromethane	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Dibromochloromethane	UG/KG	.	11.0000 U	.	11.0000 U	.	12.0000 U
V Ethylbeneene	UG/KG	.	11.0000 U	.	11.0000 U	.	1.0000 U
V Methylene Chloride	UG/KG		52.0000 U	.	34.0000 U	.	19.0000 U
V Styrene	UG/KG	.	11.0000 U	.	11.0000 U		12.0000 U
V Tetrachloroethene	UG/KG		11.0000 U	.	11.0000 U	.	12.0000 U
V Toluene	UG/KG		59.0000	.	58.0000		12.0000 U
V Trichloroethene	UG/KG	.	56.0000	.	56.0000		12.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S23D02DL	33S23M02	33S23M04	33S23S02	33S23S04	33S33D02
	LAB ID:	529536 D11	529542	529545	529543	529546	533006
	DATE RCVD:	12/17/92	12/17/92	12/17/92	12/17/92	12/17/92	01/20/93
NAME	WITS						
V Vinyl Chloride	UG/KG		11.0000 U		11.0000 U		12.0000 U
V Xylene (total)	UG/KG		11.0000 U		11.0000 U		9.0000 J
V cis-1,3-Dichloropropene	UG/KG		11.0000 U		11.0000 U		12.0000 U
V trans-1,3-Dichloropropene	UG/KG		11.0000 U		11.0000 U		12.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S33D02RE	33S37D04	33S37D04RE	33S37D04REMS	33S37D04REMS	33S38004MS
	LAB ID:	533006	533000	533000	533547	533540	532944
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
I Aluminum	MG/KG		1100.0000	.	.	.	.
I Antimony	MG/KG		13.0000 U	.	.	.	.
I Arsenic	MG/KG		1.1000 J	.	.	.	.
I Barium	MG/KG		1.6000 J	.	.	.	.
I Beryllium	MG/KG		0.6800 U	.	.	.	.
I Cadmium	MG/KG		1.1000 U	.	.	.	.
I Calcium	MG/KG	.	2830.0000	.	.	.	.
I Chromium	MG/KG		4.4000	.	.	.	.
I Cobalt	MG/KG		2.7000 U	.	.	.	.
I Copper	MG/KG	.	5.7000 U	.	.	.	.
I Cyanide	MG/KG	.	0.6600	.	.	.	.
I Iron	MG/KG	.	1180.0000	.	.	.	.
I Lead	MG/KG	.	3.7000	.	.	.	.
I Magnesium	MG/KG	.	126.0000 U	.	.		.
I Manganese	MG/KG	.	7.1000	.	.	.	.
I Mercury	MG/KG		0.1100 U	.	.	.	
I Nickel	MG/KG	.	8.7000 U	.	.	.	.
I Potaeeium	MG/KG	.	900.0000 U	.	.	.	.
I Selenium	MG/KG	.	0.6800 U	.	.	.	.
I Silver	MG/KG	.	2.3000 U	.	.	.	.
I Sodium	MG/KG	.	140.0000 J	.	.	.	.
I Thallium	MG/KG	.	0.6800 U	.	.	.	.
I Vanadium	MG/KG	.	3.4000 U	.	.	.	.
I Zinc	MG/KG	.	18.0000	.	.	.	.
P 4,4'-DDD	UG/KG	.	5.0000 U	.	.	.	.
P 4,4'-DDE	UG/KG	.	1.2000 J	.	.	.	.
P 4,4'-DDT	UG/KG	.	5.0000 U	.	.	.	.
P Aldrin	UG/KG	.	0.0930 J	.	.	.	.

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S33D02RE	33S37D04	33S37D04RE	33S37D04REMS	33S37D04REMS	33S38004MS
	LAB ID:	533006	533000 R1	533000	533547	533540	532944
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
P Aroclor 1016	UG/KG		50.0000 U	.	.	.	
P Aroclor 1221	UG/KG		100.0000 U		.	.	.
P Aroclor 1232	UG/KG	.	50.0000 U	.	.	.	.
P Aroclor 1242	UG/KG	.	50.0000 U	.	.	.	.
P Aroclor 1248	UG/KG	.	50.0000 U	.	.	.	
P Aroclor 1254	UG/KG	.	50.0000 U	.	.	.	.
P Aroclor 1260	UG/KG		63.0000	.	.	.	.
P Dieldrin	UG/KG	.	17.0000	.	.	.	.
P Endosulfan I	UG/KG		2.6000 U	.	.	.	.
P Endosulfan II	UG/KG	.	5.0000 U	.	.	.	.
P Endosulfan sulfate	UG/KG	.	1.1000 J	.	.	.	.
P Endrin	UG/KG	.	5.0000 U	.	.	.	.
P Endrin aldehyde	UG/KG		1.3000 J	.	.	.	.
P Endrin ketone	UG/KG	.	5.0000 U	.	.	.	.
P Heptachlor	UG/KG		0.6100 J	.	.	.	.
P Heptachlor epoxide	UG/KG	.	3.3000 J	.	.	.	.
P Methoxychlor	UG/KG	.	26.0000 U	.	.	.	.
P Toxaphene	UG/KG	.	260.0000 U	.	.	.	.
P alpha-BHC	UG/KG	.	2.6000 U	.	.	.	.
P alpha-Chlordane	UG/KG	.	4.7000 J	.	.	.	.
P beta-BHC	UG/KG	.	2.6000 U	.	.	.	.
P delta-BHC	UG/KG	.	2.6000 U	.	.	.	.
P gamma-BHC (Lindane)	UG/KG	.	2.6000 U	.	.	.	.
P gamma-Chlordane	UG/KG	.	28.0000 J	.	.	.	
S 1,2,4-Trichlorobenzene	UG/KG	.	370.0000 U	.	.	.	11000.0000 J
S 1,2-Dichlorobenzene	UG/KG	.	370.0000 U	.	.	.	19000.0000 J
S 1,3-Dichlorobenzene	UG/KG	.	370.0000 U	.	.	.	1600.0000
S 1,4-Dichlorobenzene	UG/KG	.	370.0000 U	.	.	.	6700.0000 J
S 2,4,5-Trichlorophenol	UG/KG	.	900.0000 U	.	.	.	960.0000 U

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S33D02RE	33S37D04	33S37D04RE	33S37D04REMS	33S37D04REMS	33S38004MS
ENSAFE ID:		33S33D02RE	33S37D04	33S37D04RE	33S37D04REMS	33S37D04REMS	33S38004MS
LAB ID:		533006	533000	533000	533547	533548	532944
DATE RCVD:		01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
S 2,4,6-Trichlorophenol	UG/KG		370.0000 U	.	.	.	400.0000 U
S 2,4-Dichlorophenol	UG/KG		370.0000 U	.	.	.	400.0000 U
S 2,4-Dimethylphenol	UG/KG		370.0000 U	.	.	.	400.0000 U
S 2,4-Dinitrophenol	UG/KG		900.0000 U	.	.	.	960.0000 U
S 2,4-Dinitrotoluene	UG/KG		370.0000 U	.	.	.	1400.0000
S 2,6-Dinitrotoluene	UG/KG	.	370.0000 U		.	.	400.0000 U
S 2-Chloronaphthalene	UG/KG		370.0000 U	.	.	.	400.0000 U
S 2-Chlorophenol	UG/KG	.	370.0000 U	.	.	.	2100.0000
S 2-Methylnaphthalene	UG/KG		370.0000 U	.	.	.	130.0000 J
S 2-Methylphenol	UG/KG		370.0000 U	.	.	.	400.0000 U
S 2-Nitroaniline	UG/KG	.	900.0000 U	.	.	.	960.0000 U
S 2-Nitrophenol	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S 3-Nitroaniline	UG/KG	.	900.0000 U	.	.	.	960.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	.	900.0000 U	.	.	.	960.0000 U
S 4-Bromophenyl-phenylether	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S 4-Chloro-3-methylphenol	UG/KG	.	370.0000 U	.	.	.	2400.0000
S 4-Chloroaniline	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S 4-Methylphenol	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S 4-Nitroaniline	UG/KG	.	900.0000 U	.	.	.	960.0000 U
S 4-Nitrophenol	UG/KG	.	900.0000 U	.	.	.	2000.0000
S Acenaphthene	UG/KG	.	370.0000 U	.	.	.	1200.0000
S Acenaphthylene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Anthracene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Benzo(a)anthracene	UG/KG	.	370.0000 U	.	.	.	61.0000 U
S Benzo(a)pyrene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Benzo(b,k)fluoranthene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Benzo(g,h,i)perylene	UG/KG	.	370.0000 U	.	.	.	400.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S33D02RE	33S37D04	33S37D04RE	33S37D04REMS	33S37D04REMS	33S38004MS
	LAB ID:	533006	533000	533000	533547	533540	532944
	DATE RCW:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	.	200.0000 U	.	.	.	160.0000 U
S Carbazole	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Chrysene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Di-n-butylphthalate	UG/KG	.	370.0000 U	.	.	.	86.0000 U
S Di-n-octylphthalate	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Dibenz(a,h)anthracene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Dibenzofuran	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Diethylphthalate	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Dimethylphthalate	UG/KG	.	370.0000 U		.	.	400.0000 U
S Fluoranthene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Fluorene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Hexachlorobenzene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Hexachlorobutadiene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Hexachlorocyclopentadiene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Hexachloroethane	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Ieophorone	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	.	370.0000 U	.	.	.	1400.0000
S N-Nitrosodiphenylamine	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Naphthalene	UG/KG	.	370.0000 U	.	.	.	200.0000 J
S Nitrobenzene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Pentachlorophenol	UG/KG	.	900.0000 U	.	.	.	3300.0000 J
S Phenanthrene	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S Phenol	UG/KG	.	370.0000 U	.	.	.	2400.0000
S Pyrene	UG/KG	.	370.0000 U	.	.	.	1000.0000
S bis(2-Chloroethoxy)methane	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	.	50.0000 U	.	.	.	400.0000 U
S bis(2-chloroethyl)ether	UG/KG	.	370.0000 U	.	.	.	400.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	.	370.0000 U	.	.	.	400.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S33D02RE	33S37D04	33S37D04RE	33S37D04REMS	33S37D04REMS	33S38004MS
	LAB ID:	533006	533000	533000	533547	533548	532944
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 1,1,2,2-Tetrachloroethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 1,1,2-Trichloroethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 1,1-Dichloroethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 1,1-Dichloroethene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	6000.0000	5600.0000	.
V 1,2-Dichloroethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 1,2-Dichloroethene (total)	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 1,2-Dichloropropane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 2-Butanone	UG/KG	29.0000 U	11.0000 U	870.0000 U	1400.0000 U	1400.0000 U	.
V 2-Hexanone	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V 4-Methyl-2-Pentanone	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Acetone	UG/KG	100.0000 U	1100.0000	1400.0000 U	4600.0000	6600.0000	.
V Benzene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	6900.0000	6700.0000	.
V Bromodichloromethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Bromoform	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Bromomethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Carbon Disulfide	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Carbon Tetrachloride	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Chlorobenzene	UG/KG	140.0000 J	11.0000 U	1400.0000 U	6400.0000	6500.0000	.
V Chloroethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Chloroform	UG/KG	29.0000 U	1.0000 U	160.0000 U	1400.0000 U	1400.0000 U	.
V Chloromethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Dibromochloromethane	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Ethylbenzene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Methylene Chloride	UG/KG	19.0000 U	11.0000 U	2200.0000 U	1700.0000 U	2300.0000 U	.
V Styrene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Tetrachloroethene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	.
V Toluene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	6600.0000	6700.0000	.
V Trichloroethene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	6600.0000	6600.0000	.

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S33D02RE	33S37D04	33S37D04RE	33S37D04REMS	33S37D04REMS	33S38004MS
	LAB ID:	533006	533000	533000	533547	533548	532944
	DATE RCVD:	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93	01/20/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	
V Xylene (total)	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	
V cis-1,3-Dichloropropene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	
V trans-1,3-Dichloropropene	UG/KG	29.0000 U	11.0000 U	1400.0000 U	1400.0000 U	1400.0000 U	

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S38004MSD	33S38M04	33S38S04	33S45D02	33S45D02DL	33S50004MS
ENSAFE ID:		33S38004MSD	33S38M04	33S38S04	33S45D02	33S45D02DL	33S50004MS
LAB ID:		532945	533544	533545	547124	547124 D11	547484
DATE RCVD:		01/20/93	01/20/93	01/20/93	04/22/93	04/22/93	04/21/93
I Aluminum	MG/KG	.	.	.	2390.0000		.
I Antimony	MG/KG	.	.	.	9.7000 U		.
I Arsenic	MG/KG	.		.	1.3000 J		.
I Barium	MG/KG	.	.	.	3.8000 J		.
I Beryllium	MG/KG	.		.	0.4200 U		.
I Cadmium	MG/KG	.	.	.	1.1000 U		.
I Calcium	MG/KG		.	.	751.0000 J		.
I Chromium	MG/KG	.	.	.	3.8000		.
I Cobalt	MG/KG	.	.	.	1.9000 U		.
I Copper	MG/KG	.	.	.	5.3000 U		.
I Cyanide	MG/KG	.	.	.	0.5200 U		.
I Iron	MG/KG	.	.	.	1520.0000		.
I Lead	MG/KG	.	.	.	14.5000		.
I Magnesium	MG/KG	.	.	.	100.0000 U		.
I Manganese	MG/KG	.	.		16.5000		.
I Mercury	MG/KG	.	.	.	0.1100 U		.
I Nickel	MG/KG	.	.	.	6.5000 U		.
I Potassium	MG/KG	.	.	.	470.0000 U		.
I Selenium	MG/KG	.	.	.	0.6300 U		.
I Silver	MG/KG	.	.	.	2.1000 U		.
I Sodium	MG/KG	.	.	.	58.9000 J		.
I Thallium	MG/KG	.	.	.	0.8400 U		.
I Vanadium	MG/KG	.	.	.	3.6000 J		.
I Zinc	MG/KG		.	.	15.5000		.
P 4,4'-DDD	UG/KG		.	.	14.0000 J	58.0000 J	.
P 4,4'-DDE	UG/KG	.	.	.	36.0000	35.0000 J	.
P 4,4'-DDT	UG/KG	.	.	.	180.0000 J	190.0000 J	.
P Aldrin	UG/KG	.	.	.	1.8000 U	18.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S38004MSD	33S38M04	33S38S04	33S45D02	33S45D02DL	33S50004MS
	LAB ID:	532945	533544	533545	547124	547124 D11	547484
	DATE RCVD:	01/20/93	01/20/93	01/20/93	04/22/93	04/22/93	04/21/93
NAME	UNITS						
P Aroclor 1016	UG/KG		.	.	35.0000 U	350.0000 U	.
P Aroclor 1221	UG/KG		.	.	70.0000 U	700.0000 U	.
P Aroclor 1232	UG/KG		.	.	35.0000 U	350.0000 U	.
P Aroclor 1242	UG/KG	.	.	.	35.0000 U	350.0000 U	.
P Aroclor 1248	UG/KG	.	.	.	35.0000 U	350.0000 U	.
P Aroclor 1254	UG/KG	.	.	.	35.0000 U	350.0000 U	.
P Aroclor 1260	UG/KG	.	.	.	35.0000 U	350.0000 U	.
P Dieldrin	UG/KG		.	.	8.8000	7.1000 J	.
P Endosulfan I	UG/KG		.	.	0.3400 J	18.0000 U	.
P Endosulfan II	UG/KG		.	.	3.5000 U	35.0000 U	.
P Endosulfan sulfate	UG/KG		.	.	3.5000 U	35.0000 U	.
P Endrin	UG/KG		.	.	3.5000 U	35.0000 U	.
P Endrin aldehyde	UG/KG	.	.	.	0.4900 J	35.0000 U	.
P Endrin ketone	UG/KG		.	.	3.5000 U	35.0000 U	.
P Heptachlor	UG/KG	.	.	.	0.2700 J	18.0000 U	.
P Heptachlor epoxide	UG/KG		.	.	1.8000 U	18.0000 U	.
P Methoxychlor	UG/KG		.	.	18.0000 U	180.0000 U	.
P Toxaphene	UG/KG		.	.	180.0000 U	1800.0000 U	.
P alpha-BHC	UG/KG		.	.	1.8000 U	18.0000 U	.
P alpha-Chlordane	UG/KG		.	.	0.5000 J	18.0000 U	.
P beta-BHC	UG/KG		.	.	1.8000 U	18.0000 U	.
P delta-BHC	UG/KG		.	.	1.8000 U	18.0000 U	.
P gamma-BHC (Lindane)	UG/KG		.	.	1.8000 U	18.0000 U	.
P gamma-Chlordane	UG/KG		.	.	0.5100 J	18.0000 U	.
S 1,2,4-Trichlorobenzene	UG/KG	15000.0000 J	.	.	340.0000 U		.
S 1,2-Dichlorobenzene	UG/KG	21000.0000 J	.	.	340.0000 U		.
S 1,3-Dichlorobenzene	UG/KG	2100.0000	.	.	340.0000 U		.
S 1,4-Dichlorobenzene	UG/KG	7900.0000 J	.	.	340.0000 U		.
S 2,4,5-Trichlorophenol	UG/KG	960.0000 U	.	.	830.0000 U		.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S38004MSD	33S38M04	33S38S04	33S45D02	33S45D02DL	33S50004MS
	LAB ID:	532945	533544	533545	547124	547124 D11	547484
	DATE RCVD:	01/20/93	01/20/93	01/20/93	04/22/93	04/22/93	04/21/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 2,4-Dichlorophenol	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 2,4-Dimethylphenol	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 2,4-Dinitrophenol	UG/KG	960.0000 U		.	830.0000 U	.	.
S 2,4-Dinitrotoluene	UG/KG	1200.0000			340.0000 U	.	.
S 2,6-Dinitrotoluene	UG/KG	400.0000 U		.	340.0000 U	.	.
S 2-Chloronaphthalene	UG/KG	400.0000 U			340.0000 U	.	
S 2-Chlorophenol	UG/KG	1800.0000	.		340.0000 U	.	.
S 2-Methylnaphthalene	UG/KG	210.0000 J			340.0000 U	.	.
S 2-Methylphenol	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 2-Nitroaniline	UG/KG	960.0000 U	.	.	830.0000 U	.	.
S 2-Nitrophenol	UG/KG	400.0000 U		.	340.0000 U	.	.
S 3,3'-Dichlorobenzidine	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 3-Nitroaniline	UG/KG	960.0000 U	.	.	830.0000 U	.	.
S 4,6-Dinitro-2-methylphenol	UG/KG	960.0000 U	.	.	830.0000 U	.	.
S 4-Bromophenyl-phenylether	UG/KG	400.0000 U		.	340.0000 U	.	.
S 4-Chloro-3-methylphenol	UG/KG	2000.0000	.	.	340.0000 U	.	.
S 4-Chloroaniline	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 4-Chlorophenyl-phenylether	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 4-Methylphenol	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S 4-Nitroaniline	UG/KG	960.0000 U	.	.	830.0000 U	.	.
S 4-Nitrophenol	UG/KG	1700.0000	.	.	830.0000 U	.	.
S Acenaphthene	UG/KG	1100.0000	.	.	340.0000 U	.	.
S Acenaphthylene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Anthracene	UG/KG	400.0000 U	.		340.0000 U	.	.
S Benzo(a)anthracene	UG/KG	400.0000 U	.	.	96.0000 J	.	.
S Benzo(a)pyrene	UG/KG	400.0000 U	.	.	94.0000 J	.	.
S Benzo(b,k)fluoranthene	UG/KG	400.0000 U	.	.	230.0000 J		.
S Benzo(g,h,i)perylene	UG/KG	400.0000 U	.	.	340.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S38004MSD	33S38M04	33S38S04	33S45D02	33S45D02DL	33S50004MS
	LAB ID:	532945	533544	533545	547124	547124 D11	547484
	DATE RCVD:	01/20/93	01/20/93	01/20/93	04/22/93	04/22/93	04/21/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	110.0000 U	.	.	340.0000 U	.	.
S Carbazole	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Chrysene	UG/KG	400.0000 U	.	.	96.0000 J	.	.
S Di-n-butylphthalate	UG/KG	76.0000 U	.	.	340.0000 U	.	.
S Di-n-octylphthalate	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Dibenz(a,h)anthracene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Dibenzofuran	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Diethylphthalate	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Dimethylphthalate	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Fluoranthene	UG/KG	400.0000 U	.	.	170.0000 J	.	.
S Fluorene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Hexachlorobenzene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Hexachlorobutadiene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Hexachlorocyclopentadiene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Hexachloroethane	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Indeno(1,2,3-cd)pyrene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Ieophorone	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S N-Nitroso-di-n-propylamine	UG/KG	1200.0000	.	.	340.0000 U	.	.
S N-Nitrosodiphenylamine	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Naphthalene	UG/KG	310.0000 J	.	.	340.0000 U	.	.
S Nitrobenzene	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S Pentachlorophenol	UG/KG	3000.0000	.	.	830.0000 U	.	.
S Phenanthrene	UG/KG	400.0000 U	.	.	68.0000 J	.	.
S Phenol	UG/KG	2000.0000	.	.	340.0000 U	.	.
S Pyrene	UG/KG	930.0000	.	.	140.0000 J	.	.
S bis(2-Chloroethoxy)methane	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S bis(2-Ethlyhexy)phthalate	UG/KG	380.0000 U	.	.	89.0000 J	.	.
S bis(2-chloroethyl)ether	UG/KG	400.0000 U	.	.	340.0000 U	.	.
S bis(2-chloroisopropyl)ether	UG/KG	400.0000 U	.	.	340.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S38004MSD	33S38M04	33S38S04	33545D02	33S45D02DL	33S50004MS
	LAB ID:	532945	533544	533545	547124	547124 D11	547484
	DATE RCVD:	01/20/93	01/20/93	01/20/93	04/22/93	04/22/93	04/21/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 1,1,2-Trichloroethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 1,1-Dichloroethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 1,1-Dichloroethene	UG/KG	.	7800.0000	7000.0000	11.0000 U	.	6400.0000
V 1,2-Dichloroethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 1,2-Dichloroethene (total)	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 1,2-Dichloropropane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 2-Butanone	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 2-Hexanone	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V 4-Methyl-2-Pentanone	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Acetone	UG/KG	.	2100.0000 U	2100.0000 U	62.0000 U	.	1400.0000 U
V Benzene	UG/KG	.	7500.0000	7S00.0000	11.0000 U	.	6600.0000
V Bromodichloromethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Bromoform	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Bromomethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Carbon Disulfide	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Carbon Tetrachloride	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Chlorobenzene	UG/KG	.	12000.0000	11000.0000	11.0000 U	.	6500.0000
V Chloroethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Chloroform	UG/KG	.	2100.0000 U	2100.0000 U	1.0000 J	.	1400.0000 U
V Chloromethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Dibromochloromethane	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Ethylbenzene	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Methylene Chloride	UG/KG	.	370.0000 U	4400.0000 U	27.0000 U	.	570.0000 J
V Styrene	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Tetrachloroethene	UG/KG	.	2100.0000 U	2100.0000 U	11.0000 U	.	1400.0000 U
V Toluene	UG/KG	.	7300.0000	7200.0000	11.0000 U	.	6600.0000
V Trichloroethene	UG/KG		7200.0000	7300.0000	11.0000 U	.	5900.0000

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S38004MSD	33S38M04	33S38S04	33S45D02	33S45D02DL	33S50004MS
	LAB ID:	532945	533544	533545	547124	547124 D11	547484
	DATE RCVD:	01/20/93	01/20/93	01/20/93	04/22/93	04/22/93	04/21/93
NAME	UNITS						
V Vinyl Chloride	UG/KG		2100.0000 U	2100.0000 U	11.0000 U		1400.0000 U
V Xylene (total)	UG/KG		2100.0000 U	270.0000 J	11.0000 U		860.0000 J
V cis-1,3-Dichloropropene	UG/KG		2100.0000 U	2100.0000 U	11,0000 U		1400.0000 U
V trans-1,3-Dichloropropene	UG/KG		2100.0000 U	2100.0000 U	11.0000 U		1400.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33S50004MSD 547485 04/21/93	33S51002MS 546866 R1 04/21/93	33S51002MSD 546867 R1 04/21/93	33S52002MS 546863 04/21/93	33S52002MSD 546864 04/21/93	33S52D02 547129 04/22/93
I Aluminum	MG/KG	.	.	.	.	.	1680.0000
I Antimony	MG/KG	.	.	.	.	.	9.6000 U
I Arsenic	MG/KG	.	.	.	.	.	0.7200 J
I Barium	MG/KG	.	.	.	.	.	12.6000 J
I Beryllium	MG/KG	.	.	.	.	.	0.4200 U
I Cadmium	MG/KG	.	.	.	.	.	1.0000 U
I Calcium	MG/KG	.	.	.	.	.	296.0000 J
I Chromium	MG/KG	.	.	.	.	.	2.8000
I Cobalt	MG/KG	.	.	.	.	.	1.9000 U
I Copper	MG/KG	.	.	.	.	.	7.2000
I Cyanide	MG/KG	.	.	.	.	.	0.5200 U
I Iron	MG/KG	.	.	.	.	.	987.0000
I Lead	MG/KG	.	.	.	.	.	36.3000
I Magnesium	MG/KG	.	.	.	.	.	99.0000 U
I Manganese	MG/KG	.	.	.	.	.	29.4000
I Mercury	MG/KG	.	.	.	.	.	0.1000 U
I Nickel	MG/KG	.	.	.	.	.	6.4000 U
I Potassium	MG/KG	.	.	.	.	.	465.0000 U
I Selenium	MG/KG	.	.	.	.	.	0.6200 U
I Silver	MG/KG	.	.	.	.	.	2.1000 U
I Sodium	MG/KG	.	.	.	.	.	67.9000 J
I Thallium	MG/KG	.	.	.	.	.	0.8300 U
I Vanadium	MG/KG	.	.	.	.	.	2.5000 J
I Zinc	MG/KG	.	.	.	.	.	33.8000
P 4,4'-DDD	UG/KG	.	0.5400 J	0.3200 J	.	.	3.4000 U
P 4,4'-DDE	UG/KG	.	3.2000 J	3.4000 J	.	.	0.2100 J
P 4,4'-DDT	UG/KG	.	27.0000 J	28.0000	.	.	0.4800 J
P Aldrin	UG/KG	.	10.0000	11.0000	.	.	1.8000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S50004MSD	33S51002MS	33S51002MSD	33S52002MS	33S52002MSD	33552802
	LAB ID:	547485	546866 R1	546867 R1	546863	546864	547129
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93
NAME	UNITS						
P Aroclor 1016	UG/KG		35.0000 U	35.0000 U	.		34.0000 U
P Aroclor 1221	UG/KG		71.0000 U	71.0000 U	.		70.0000 U
P Aroclor 1232	UG/KG		35.0000 U	35.0000 U	.		34.0000 U
P Aroclor 1242	UG/KG		35.0000 U	35.0000 U	.	.	34.0000 U
P Aroclor 1248	UG/KG		35.0000 U	35.0000 U	.	.	34.0000 U
P Aroclor 1254	UG/KG		35.0000 U	35.0000 U	.	.	34.0000 U
P Aroclor 1260	UG/KG		35.0000 U	35.0000 U	.		34.0000 U
P Dieldrin	UG/KG		25.0000	26.0000	.	.	0.2200 J
P Endosulfan I	UG/KG		1.8000 U	1.8000 U	.	.	1.8000 U
P Endosulfan II	UG/KG	.	3.5000 U	3.5000 U	.	.	3.4000 U
P Endosulfan sulfate	UG/KG	.	0.3400 J	3.5000 U	.	.	3.4000 U
P Endrin	UG/KG	.	25.0000	25.0000	.	.	3.4000 U
P Endrin aldehyde	UG/KG	.	2.2000 J	2.2000 J	.	.	3.4000 U
P Endrin ketone	UG/KG	.	0.0820 J	0.3200 J	.	.	3.4000 U
P Heptachlor	UG/KG	.	12.0000	12.0000	.	.	1.8000 U
P Heptachlor epoxide	UG/KG	.	1.8000 U	1.8000 U	.	.	1.8000 U
P Methoxychlor	UG/KG	.	18.0000 U	18.0000 U	.	.	18.0000 U
P Toxaphene	UG/KG	.	180.0000 U	180.0000 U	.	.	180.0000 U
P alpha-BHC	UG/KG	.	1.8000 U	1.8000 U	.	.	1.8000 U
P alpha-Chlordane	UG/KG	.	0.2100 J	0.1700 J	.	.	0.9000 J
P beta-BHC	UG/KG	.	1.8000 U	1.8000 U	.	.	1.8000 U
P delta-BHC	UG/KG	.	1.8000 U	1.8000 U	.	.	1.8000 U
P gamma-BHC (Lindane)	UG/KG	.	12.0000	13.0000	.	.	1.8000 U
P gamma-Chlordane	UG/KG	.	1.2000 J	1.5000 J			0.8800 J
S 1,2,4-Trichlorobenzene	UG/KG	.			1500.0000	1700.0000	340.0000 U
S 1,2-Dichlorobenzene	UG/KG	.			340.0000 U	340.0000 U	340.0000 U
S 1,3-Dichlorobenzene	UG/KG	.			340.0000 U	340.0000 U	340.0000 U
S 1,4-Dichlorobenzene	UG/KG	.			1400.0000	1700.0000	340.0000 U
S 2,4,5-Trichlorophenol	UG/KG	.			830.0000 U	830.0000 U	820.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S50004MSD	33S51002MS	33S51002MSD	33S52002MS	33S52002MSD	33S52D02
	LAB ID:	547485	546866 R1	546867 R1	546863	546864	547129
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/KG				340.6000 U	340.0000 U	340.0000 U
S 2,4-Dichlorophenol	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 2,4-Dimethylphenol	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 2,4-Dinitrophenol	UG/KG				830.0000 U	830.0000 U	820.0000 U
S 2,4-Dinitrotoluene	UG/KG				1400.0000	1700.0000	340.0000 U
S 2,6-Dinitrotoluene	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 2-Chloronaphthalene	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 2-Chlorophenol	UG/KG				2200.0000	2600.0000	340.0000 U
S 2-Methylnaphthalene	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 2-Methylphenol	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 2-Nitroaniline	UG/KG				830.0000 U	830.0000 U	820.0000 U
S 2-Nitrophenol	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 3,3'-Dichlorobenzidine	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 3-Nitroaniline	UG/KG				830.0000 U	830.0000 U	820.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG				830.0000 U	830.0000 U	820.0000 U
S 4-Bromophenyl-phenylether	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 4-Chloro-3-methylphenol	UG/KG				2800.0000 J	3100.0000 J	340.0000 U
S 4-Chloroaniline	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 4-Chlorophenyl-phenylether	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 4-Methylphenol	UG/KG				340.0000 U	340.0000 U	340.0000 U
S 4-Nitroaniline	UG/KG				830.0000 U	830.0000 U	820.0000 U
S 4-Nitrophenol	UG/KG				2400.0000	3100.0000 J	820.0000 U
S Acenaphthene	UG/KG				1500.0000	1700.0000	340.0000 U
S Acenaphthylene	UG/KG				340.0000 U	340.0000 U	340.0000 U
S Anthracene	UG/KG				340.0000 U	340.0000 U	340.0000 U
S Benzo(a)anthracene	UG/KG				340.0000 U	340.0000 U	38.0000 J
S Benzo(a)pyrene	UG/KG				340.0000 U	340.0000 U	340.0000 U
S Benzo(b,k)fluoranthene	UG/KG				340.0000 U	37.0000 J	100.0000 J
S Benzo(g,h,i)perylene	UG/KG				340.0000 U	340.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S50004MSD	33S51002MS	33S51002MSD	33S52002MS	33S52002MSD	33S52D02
	LAB ID:	547485	546866 R1	546867 R1	546863	546864	547129
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG		.	.	340.0000 U	340.0000 U	340.0000 U
S Carbazole	UG/KG			.	340.0000 U	340.0000 U	340.0000 U
S Chrysene	UG/KG		.	.	340.0000 U	340.0000 U	37.0000 J
S Di-n-butylphthalate	UG/KG			.	340.0000 U	340.0000 U	340.0000 U
S Di-n-octylphthalate	UG/KG			.	340.0000 U	340.0000 U	340.0000 U
S Dibenz (a,h) anthracene	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Dibenzofuran	UG/KG			.	340.0000 U	340.0000 U	340.0000 U
S Diethylphthalate	UG/KG			.	340.0000 U	340.0000 U	340.0000 U
S Dimethylphthalate	UG/KG		.	.	340.0000 U	340.0000 U	340.0000 U
S Fluoranthene	UG/KG	.		.	41.0000 J	76.0000 J	110.0000 J
S Fluorene	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Hexachlorobenzene	UG/KG	.			340.0000 U	340.0000 U	340.0000 U
S Hexachlorobutadiene	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Hexachlorocyclopentadiene	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Hexachloroethane	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Isophorone	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	.	.	.	1300.0000	1400.0000	340.0000 U
S N-Nitrosodiphenylamine	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Naphthalene	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Nitrobenzene	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S Pentachlorophenol	UG/KG	.	.	.	2300.0000	2900.0000 J	820.0000 U
S Phenanthrene	UG/KG	.	.	.	340.0000 U	51.0000 J	82.0000 J
S Phenol	UG/KG	.	.	.	2500.0000	2800.0000	340.0000 U
S Pyrene	UG/KG	.	.	.	1700.0000	2200.0000	69.0000 J
S bis (2-Chloroethoxy) methane	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	.	.	.	340.0000 U	66.0000 U	340.0000 U
S bis(2-chloroethyl)ether	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	.	.	.	340.0000 U	340.0000 U	340.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S50004MSD	33S51002MS	33S51002MSD	33S52002MS	33S52002MSD	33S52D02
	LAB ID:	547485	546866 R1	546867 R1	546863	546864	547129
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V 1,1,2-Trichloroethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V 1,1-Dichloroethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V 1,1-Dichloroethene	UG/KG	6600.0000	.	.	.	.	10.0000 U
V 1,2-Dichloroethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V 1,2-Dichloropropane	UG/KG	1400.0000 U			.	.	10.0000 U
V 2-Butanone	UG/KG	770.0000 U	.	.	.	.	10.0000 U
V 2-Hexanone	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Acetone	UG/KG	1400.0000 U	.	.	.	.	470.0.000 J
V Benzene	UG/KG	7100.0000	.	.	.	.	10.0000 U
V Bromodichloromethane	UG/KG	1400.0000 U		.	.	.	10.0000 U
V Bromoform	UG/KG	1400.0000 U		.	.	.	10.0000 U
V Bromomethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Carbon Disulfide	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Carbon Tetrachloride	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Chlorobenzene	UG/KG	6600.0000	.	.	.	.	10.0000 U
V Chloroethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Chloroform	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Chloromethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Dibromochlormethane	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Ethylbenzene	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Methylene Chloride	UG/KG	450.0000 J		.	.	.	35.0000 U
V Styrene	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Tetrachloroethene	UG/KG	1400.0000 U	.	.	.	.	10.0000 U
V Toluene	UG/KG	6500.0000	.	.	.	.	10.0000 U
V Trichloroethene	UG/KG	6100.0000		.	.	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S50004MSD	33S51002MS	33S51002MSD	33S52002MS	33S52002MSD	33S52D02
	LAB ID:	547485	546866 R1	546867 R1	546863	546864	547129
	DATE RCVD:	04/21/93	04/21/93	04/21/93	04/21/93	04/21/93	04/22/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	1400.0000 U		.	.	.	10.0000 U
V Xylene (total)	UG/KG	840.0000 J		.	.	.	10.0000 U
V cis-1,3-Dichloropropene	UG/KG	1400.0000 U		.	.	.	10.0000 U
V trans-1,3-Dichloropropene	UG/KG	1400.0000 U		.	.	.	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD	33S57D01
ENSAFE ID:		33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD	33S57D01
LAB ID:		547129	546756	546757	546860	546861	547127
DATE RCVD:		04/22/93	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93
I Aluminum	MG/KG	.	.	.	.	.	2040.0000
I Antimony	MG/KG	.	.	.	.	.	15.9000 U
I Arsenic	MG/KG	.	.	.	.	.	1.9000 J
I Barium	MG/KG	.	.	.	.	.	51.4000 J
I Beryllium	MG/KG	.	.	.	.	.	0.6900 U
I Cadmium	MG/KG	.	.	.	.	.	271.0000
I Calcium	MG/KG	.	.	.	.	.	4290.0000
I Chromium	MG/KG	.	.	.	.	.	4760.0000
I Cobalt	MG/KG	.	.	.	.	.	3.1000 U
I Copper	MG/KG	.	.	.	.	.	446.0000
I Cyanide	MG/KG	.	.	.	.	.	94.8000
I Iron	MG/KG	.	.	.	.	.	5860.0000
I Lead	MG/KG	.	.	.	.	.	923.0000
I Magnesium	MG/KG	.	.	.	.	.	250.0000 J
I Manganese	MG/KG	.	.	.	.	.	33.4000
I Mercury	MG/KG	.	.	.	.	.	2.9000
I Nickel	MG/KG	.	.	.	.	.	11.0000 J
I Potassium	MG/KG	.	.	.	.	.	771.0000 U
I Selenium	MG/KG	.	.	.	.	.	1.0000 U
I Silver	MG/KG	.	.	.	.	.	8.7000
I sodium	MG/KG	.	.	.	.	.	194.0000 J
I Thallium	MG/KG	.	.	.	.	.	1.4000 U
I Vanadium	MG/KG	.	.	.	.	.	3.1000 J
I Zinc	MG/KG	.	.	.	.	.	713.0000
P 4,4'-DDD	UG/KG	.	3.7000 U	3.7000 U	.	.	44.0000
P 4,4'-DDE	UG/KG	.	6.4000	6.9000	.	.	240.0000 J
P 4,4'-DDT	UG/KG	.	41.0000 J	36.0000 J	.	.	240.0000 J
P Aldrin	UG/KG	.	15.0000	15.0000	.	.	0.7700 J

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD	33S57D01
	LAB ID:	547129	546762	546763	546860	546861	547127
	DATE RCVD:	04/22/93	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93
NAME	UNITS						
P Aroclor 1016	UG/KG	.	37.0000 U	37.0000 U	.	.	56.0000 U
P Aroclor 1221	UG/KG	.	75.0000 U	75.0000 U	.	.	110.0000 U
P Aroclor 1232	UG/KG	.	37.0000 U	37.0000 U	.	.	56.0000 U
P Aroclor 1242	UG/KG	.	37.0000 U	37.0000 U	.	.	56.0000 U
P Aroclor 1248	UG/KG	.	37.0000 U	37.0000 U	.	.	56.0000 U
P Aroclor 1254	UG/KG	.	37.0000 U	37.0000 U	.	.	56.0000 U
P Aroclor 1260	UG/KG	.	270.0000	290.0000	.	.	56.0000 U
P Dieldrin	UG/KG	.	38.0000 U	38.0000	.	.	5.6000 U
P Endosulfan I	UG/KG	.	1.9000 U	1.9000 U	.	.	1.4000 J
P Endosulfan II	UG/KG	.	3.7000 U	3.7000 U	.	.	5.6000 U
P Endosulfan sulfate	UG/KG	.	3.7000 U	3.7000 U	.	.	5.6000 U
P Endrin	UG/KG	.	46.0000	47.0000	.	.	5.6000 U
P Endrin aldehyde	UG/KG	.	3.7000 U	3.7000 U	.	.	9.7000 J
P Endrin ketone	UG/KG	.	3.7000 U	3.7000 U	.	.	5.6000 U
P Heptachlor	UG/KG	.	16.0000	15.0000	.	.	0.4200 J
P Heptachlor epoxide	UG/KG	.	1.9000 U	0.2300 J	.	.	2.9000 U
P Methoxychlor	UG/KG	.	19.0000 U	19.0000 U	.	.	18.0000 J
P Toxaphene	UG/KG	.	190.0000 U	190.0000 U	.	.	290.0000 U
P alpha-BHC	UG/KG	.	1.9000 U	1.9000 U	.	.	2.9000 U
P alpha-Chlordane	UG/KG	.	2.8000	3.1000	.	.	2.9000 U
P beta-BHC	UG/KG	.	1.9000 U	1.9000 U	.	.	2.9000 U
P delta-BHC	UG/KG	.	1.9000 U	1.9000 U	.	.	2.9000 U
P gamma-BHC (Lindane)	UG/KG	.	16.0000	16.0000	.	.	2.9000 U
P gamma-Chlordane	UG/KG	.	0.5600 J	0.9100 J	.	.	0.4400 J
S 1,2,4-Trichlorobenzene	UG/KG	.	1200.0000	1100.0000	.	.	560.0000 U
S 1,2-Dichlorobenzene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S 1,3-Dichlorobenzene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S 1,4-Dichlorobenzene	UG/KG	.	1200.0000	1000.0000	.	.	560.0000 U
S 2,4,5-Trichlorophenol	UG/KG	.	900.0000 U	900.0000 U	.	.	1400.0000 U

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD	33857001
ENSAFE ID:		33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD	33857001
LAB ID:		547129	546759	546760	546860	546861	547127
DATE RCVD:		04/22/93	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93
S 2,4,6-Trichlorophenol	UG/KG		370.0000 U	370.0000 U	.		560.0000 U
S 2,4-Dichlorophenol	UG/KG		370.0000 U	370.0000 U			560.0000 U
S 2,4-Dimethylphenol	UG/KG		370.0000 U	370.0000 U	.	.	560.0000 U
S 2,4-Dinitrophenol	UG/KG		900.0000 U	900.0000 U			1400.0000 U
S 2,4-Dinitrotoluene	UG/KG		1200.0000	1100.0000			560.0000 U
S 2,6-Dinitrotoluene	UG/KG		370.0000 U	370.0000 U			560.0000 U
S 2-Chloronaphthalene	UG/KG		370.0000 U	370.0000 U	.		560.0000 U
S 2-Chlorophenol	UG/KG		1600.0000	1400.0000	.	.	560.0000 U
S 2-Methylnaphthalene	UG/KG		370.0000 U	370.0000 U	.	.	560.0000 U
S 2-Methylphenol	UG/KG		370.0000 U	370.0000 U	.	.	560.0000 U
S 2-Nitroaniline	UG/KG	.	900.0000 U	900.0000 U	.	.	1400.0000 U
S 2-Nitrophenol	UG/KG		370.0000 U	370.0000 U	.	.	560.0000 U
S 3,3'-Dichlorobenzidine	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S 3-Nitroaniline	UG/KG	.	900.0000 U	900.0000 U	.	.	1400.0000 U
S 4,6-Dinitro-2-methylphenol	UG/KG	.	900.0000 U	900.0000 U	.	.	1400.0000 U
S 4-Bromophenyl-phenylether	UG/KG		370.0000 U	370.0000 U	.	.	560.0000 U
S 4-Chloro-3-methylphenol	UG/KG	.	2000.0000	1800.0000	.	.	560.0000 U
S 4-Chloroaniline	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S 4-Chlorophenyl-phenylether	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S 4-Methylphenol	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S 4-Nitroaniline	UG/KG	.	900.0000 U	900.0000 U	.		1400.0000 U
S 4-Nitrophenol	UG/KG	.	2400.0000	2300.0000	.	.	1400.0000 U
S Acenaphthene	UG/KG	.	1200.0000	1100.0000	.	.	560.0000 U
S Acenaphthylene	UG/KG	.	370.0000 U	370.0000 U	.	.	91.0000 J
S Anthracene	UG/KG	.	370.0000 U	370.0000 U	.	.	110.0000 J
S Benzo(a)anthracene	UG/KG	.	60.0000 J	120.0000 J	.		490.0000 J
S Benzo(a)pyrene	UG/KG	.	120.0000 J	84.0000 J	.	.	500.0000 J
S Benzo(b,k)fluoranthene	UG/KG	.	310.0000 J	2e0.0000 J	.	.	930.0000
S Benzo(g,h,i)perylene	UG/KG	.	80.0000 J	42.0000 J	.	.	740.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD	33S57D01
	LAB ID:	547129	546759	546760	546860	546861	547127
	DATE RCVD:	04/22/93	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93
NAME	UNITS						
S Butylbenzylphthalate	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Carbazole	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Chrysene	UG/KG	.	72.0000 J	81.0000 J	.	.	380.0000 J
S Di-n-butylphthalate	UG/KG	.	38.0000 U	370.0000 U	.	.	330.0000 J
S Di-n-octylphthalate	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Dibenz(a,h)anthracene	UG/KG	.	46.0000 J	44.0000 J	.	.	180.0000 J
S Dibenzofuran	UG/KG	.	370.0000'U	370.0000 U	.	.	560.0000 U
S Diethylphthalate	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Dimethylphthalate	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Fluoranthene	UG/KG	.	46.0000 J	39.0000 J	.	.	310.0000 J
S Fluorene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Hexachlorobenzene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Hexachlorobutadiene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Hexachlorocycloperitadiene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Hexachloroethane	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG	.	150.0000 J	120.0000 J	.	.	520.0000 J
S Ieophorone	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S N-Nitroso-di-n-propylamine	UG/KG	.	1400.0000	1200.0000	.	.	560.0000 U
S N-Nitrosodiphenylamine	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Naphthalene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Nitrobenzene	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S Pentachlorophenol	UG/KG	.	2100.0000	1900.0000	.	.	1400.0000 U
S Phenanthrene	UG/KG	.	370.0000 U	370.0000 U	.	.	90.0000 J
S Phenol	UG/KG	.	1600.0000	1500.0000	.	.	82.0000 J
S Pyrene	UG/KG	.	1100.0000	1100.0000	.	.	300.0000 J
S bis(2-Chloroethoxy)methane	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG	.	60.0000 U	60.0000 U	.	.	98.0000 U
S bis(2-chloroethyl)ether	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U
S bis(2-chloroisopropyl)ether	UG/KG	.	370.0000 U	370.0000 U	.	.	560.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD '	33S57D01
	LAB ID:	547129	546756	546757	546860	546861	547127
	DATE RCVD:	04/22/93	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 1,1,2-Trichloroethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 1,1-Dichloroethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 1,1-Dichloroethene	UG/KG	35.0000 U	52.0000	63.0000	48.0000	52.0000	17.0000 U
V 1,2-Dichloroethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 1.2-Dichloroethene (total)	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 1,2-Dichloropropane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 2-Butanone	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 2-Hexanone	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V 4-Methyl-2-Pentanone	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Acetone	UG/KG	400.0000 J	64.0000 U	30.0000 U	58.0000 U	180.0000 U	13.0000 U
V Benzene	UG/KG	35.0000 U	55.0000	61.0000	51.0000	50.0000	17.0000 U
V Bromodichloromethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Bromoform	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Bromomethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Carbon Disulfide	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Carbon Tetrachloride	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Chlorobenzene	UG/KG	35.0000 U	56.0000	61.0000	47.0000	46.0000	17.0000 U
V Chloroethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Chloroform	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Chloromethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Dibromochloromethane	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Ethylbenzene	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Methylene Chloride	UG/KG	63.0000 U	27.0000 U	13.0000 U	21.0000 U	51.0000 U	52.0000 U
V Styrene	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Tetrachloroethene	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Toluene	UG/KG	35.0000 U	56.0000	63.0000	50.0000	51.0000	17.0000 U
V Trichloroethene	UG/KG	35.0000 U	57.0000	56.0000	44.0000	43.0000	17.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S52D02RE	33S53002MS	33S53002MSD	33S57B04MS	33S57B04MSD	33S57D01
	LAB ID:	547129	546756	546757	546860	546861	547127
	DATE RCVD:	04/22/93	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93
NAME	UNITS						
V Vinyl Chloride	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V Xylene (total)	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V cis-1,3-Dichloropropene	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U
V trans-1,3-Dichloropropene	UG/KG	35.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U	17.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57D01DL	33SAND01
	LAB ID:	547127 D11	529909
	DATE RCVD:	04/22/93	12/18/92
NAME	UNITS		
I Aluminum	MG/KO	.	137.0000
I Antimony	MG/KG	.	11.4000 U
I Arsenic	MG/KG		0.6000 U
I Barium	MG/KG		1.3000 J
I Beryllium	MG/KG	.	0.4000 U
I Cadmium	MG/KG		1.0000 U
I Calcium	MG/KG	.	185.0000 U
I Chromium	MG/KG	.	2.0000 U
I Cobalt	MG/KG	.	2.2000 U
I Copper	MG/KG	.	5.0000 U
I Cyanide	MG/KG		0.5000 U
I Iron	MG/KG		54.9000
I Lead	MG/KG	.	0.5200 J
I Magnesium	MG/KG	.	85.2000 U
I Manganese	MG/KG	.	1.5000 J
I Mercury	MG/KG		0.1000 U
I Nickel	MG/KG	.	8.0000 U
I Potassium	MG/KG		973.0000 U
I Selenium	MG/KG	.	0.6000 U
I Silver	MG/KG	.	2.0000 U
I Sodium	MG/KG	.	90.1000 J
I Thallium	MG/KG	.	0.6000 U
I Vanadium	MG/KG	.	3.8000 U
I Zinc	MG/KG		1.0000 U
P 4,4'-DDD	UG/KG	44.0000 J	3.2000 U
P 4,4'-DDE	UG/KG	280.0000	3.2000 U
P 4,4'-DDT	UG/KG	260.0000	3.2000 U
P Aldrin	UG/KG	29.0000 U	1.6000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57D01DL	33SAND01				
	LAB ID:	547127 D11	529909				
	DATE RCVD:	04/22/93	12/18/92				
NAME	UNITS						
P Aroclor 1016	UG/KG	560.0000 U	32.0000 U	.	.	.	.
P Aroclor 1221	UG/KG	1100.0000 U	64.0000 U	.	.	.	.
P Aroclor 1232	UG/KG	560.0000 U	32.0000 U	.	.	.	.
P Aroclor 1242	UG/KG	560.0000 U	32.0000 U	.	.	.	.
P Aroclor 1248	UG/KG	560.0000 U	32.0000 U	.	.	.	.
P Aroclor 1254	UG/KG	560.0000 U	32.0000 U	.	.	.	.
P Aroclor 1260	UG/KG	560.0000 U	32.0000 U	.	.	.	.
P Dieldrin	UG/KG	56.0000 U	3.2000 U	.	.	.	
P Endosulfan I	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P Endosulfan II	UG/KG	56.0000 U	3.2000 U	.	.	.	.
P Endosulfan sulfate	UG/KG	56.0000 U	3.2000 U	.	.	.	.
P Endrin	UG/KG	56.0000 U	3.2000 U	.	.	.	.
P Endrin aldehyde	UG/KG	12.0000 J	3.2000 U	.	.	.	.
P Endrin ketone	UG/KG	56.0000 U	3.2000 U	.	.	.	.
P Heptachlor	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P Heptachlor epoxide	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P Methoxychlor	UG/KG	290.0000 U	16.0000 U	.	.	.	.
P Toxaphene	UG/KG	2900.0000 U	160.0000 U	.	.	.	.
P alpha-BHC	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P alpha-Chlordane	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P beta-BHC	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P delta-EHC	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P gama-BHC (Lindane)	UG/KG	29.0000 U	1.6000 U	.	.	.	.
P gama-Chlordane	UG/KG	29.0000 U	1.6000 U	.	.	.	.
S 1,2,4-Trichlorobenzene	UG/KG		330.0000 U	.	.	.	.
S 1,2-Dichlorobenzene	UG/KG		330.0000 U	.	.	.	.
S 1,3-Dichlorobenzene	UG/KG		330.0000 U	.	.	.	.
S 1,4-Dichlorobenzene	UG/KG		330.0000 U		.	.	
S 2,4,5-Trichlorophenol	UG/KG		790.0000 U	.	.	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57D01DL	33SAND01				
	LAB ID:	547127 D11	529909				
	DATE RCVD:	04/22/93	12/18/92				
NAME	WITS						
S 2,4,6-Trichlorophenol	UG/KG	.	330.0000 U	.	.	.	.
S 2,4-Dichlorophenol	UG/KG	.	330.0000 U	.	.	.	.
S 2,4-Dimethylphenol	UG/KG	.	330.0000 U		.	.	.
S 2,4-Dinitrophenol	UG/KG	.	790.0000 U		.	.	.
S 2,4-Dinitrotoluene	UG/KG	.	330.0000 U	.	.	.	.
S 2,6-Dinitrotoluene	UG/KG	.	330.0000 U		.	.	.
S 2-Chloronaphthalene	UG/KG	.	330.0000 U		.	.	.
S 2-Chlorophenol	UG/KG	.	330.0000 U		.	.	.
S 2-Methylnaphthalene	UG/KG	.	330.0000 U	.	.	.	.
S 2-Methylphenol	UG/KG	.	330.0000 U		.	.	.
S 2-Nitroaniline	UG/KG	.	790.0000 U	.	.	.	.
S 2-Nitrophenol	UG/KG	.	330.0000 U	.	.	.	.
S 3,3'-Dichlorobenzidine	UG/KG	.	330.0000 U	.	.	.	.
S 3-Nitroaniline	UG/KG	.	790.0000 U	.	.	.	.
S 4,6-Dinitro-2-methylphenol	UG/KG	.	790.0000 U	.	.	.	.
S 4-Bromophenyl-phenylether	UG/KG	.	330.0000 U	.	.	.	.
S 4-Chloro-3-methylphenol	UG/KG	.	330.0000 U	.	.	.	.
S 4-Chloroaniline	UG/KG	.	330.0000 U	.	.	.	.
S 4-Chlorophenyl-phenylether	UG/KG	.	330.0000 U	.	.	.	.
S 4-Methylphenol	UG/KG	.	330.0000 U	.	.	.	.
S 4-Nitroaniline	UG/KG	.	790.0000 U	.	.	.	.
S 4-Nitrophenol	UG/KG	.	790.0000 U	.	.	.	.
S Acenaphthene	UG/KG	.	330.0000 U	.	.	.	.
S Acenaphthylene	UG/KG	.	330.0000 U	.		.	.
S Anthracene	UG/KG	.	330.0000 U	.	.	.	.
S Benzo(a)anthracene	UG/KG	.	330.0000 U	.		.	.
S Benzo(a)pyrene	UG/KG	.	330.0000 U	.	.	.	.
S Benzo(b,k)fluoranthene	UG/KG	.	330.0000 U	.	.	.	.
S Benzo(g,h,i)perylene	UG/KG	.	330.0000 U	.	.	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57D01DL	33SAND01
	LAB ID:	547127 D11	529909
	DATE RCVD:	04/22/93	12/18/92
NAME	UNITS		
S Butylbenzylphthalate	UG/KG		330.0000 U
S Carbazole	UG/KG		330.0000 U
S Chrysene	UG/KG		330.0000 U
S Di-n-butylphthalate	UG/KG		330.0000 U
S Di-n-octylphthalate	UG/KG		330.0000 U
S Dibenz(a,h)anthracene	UG/KG		330.0000 U
S Dibenzofuran	UG/KG		330.0000 U
S Diethylphthalate	UG/KG		330.0000 U
S Dimethylphthalate	UG/KG		330.0000 U
S Fluoranthene	UG/KG		330.0000 U
S Fluorene	UG/KG		330.0000 U
S Hexachlorobenzene	UG/KG		330.0000 U
S Hexachlorobutadiene	UG/KG		330.0000 U
S Hexachlorocyclopentadiene	UG/KG		330.0000 U
S Hexachloroethane	UG/KG		330.0000 U
S Indeno(1,2,3-cd)pyrene	UG/KG		330.0000 U
S Isophorone	UG/KG		330.0000 U
S N-Nitroso-di-n-propylamine	UG/KG		330.0000 U
S N-Nitrosodiphenylamine	UG/KG		330.0000 U
S Naphthalene	UG/KG		330.0000 U
S Nitrobenzene	UG/KG		330.0000 U
S Pentachlorophenol	UG/KG		790.0000 U
S Phenanthrene	UG/KG		330.0000 U
S Phenol	UG/KG		330.0000 U
S Pyrene	UG/KG		330.0000 U
S bis(2-Chloroethoxy)methane	UG/KG		330.0000 U
S bis(2-Ethlyhexy)phthalate	UG/KG		330.0000 U
S bis(2-chloroethyl)ether	UG/KG		330.0000 U
S bis(2-chloroisopropyl)ether	UG/KG		330.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key pteeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S57D01DL	33SAND01
	LAB ID:	547127 D11	529909
	DATE RCVD:	04/22/93	12/18/92
NAME	UNITS		
V 1,1,1-Trichloroethane	UG/KG	.	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/KG		10.0000 U
V 1,1,2-Trichloroethane	UG/KG	.	10.0000 U
V 1,1-Dichloroethane	UG/KG		10.0000 U
V 1,1-Dichloroethene	UG/KG		10.0000 U
V 1,2-Dichloroethane	UG/KG		10.0000 U
V 1,2-Dichloroethene (total)	UG/KG	.	10.0000 U
V 1,2-Dichloropropane	UG/KG		10.0000 U
V 2-Butanone	UG/KG	.	10.0000 U
V 2-Hexanone	UG/KG	.	10.0000 U
V 4-Methyl-2-Pentanone	UG/KG	.	10.0000 U
V Acetone	UG/KG	.	46.0000 U
V Benzene	UG/KG	.	10.0000 U
V Bromodichloromethane	UG/KG		10.0000 U
V Bromoform	UG/KG	.	10.0000 U
V Bromomethane	UG/KG		10.0000 U
V Carbon Disulfide	UG/KG	.	10.0000 U
V Carbon Tetrachloride	UG/KG	.	10.0000 U
V Chlorobenzene	UG/KG	.	10.0000 U
V Chloroethane	UG/KG	.	10.0000 U
V Chloroform	UG/KG	.	10.0000 U
V Chloromethane	UG/KG		10.0000 U
V Dibromochloromethane	UG/KG	.	10.0000 U
V Ethylbenzene	UG/KG	.	10.0000 U
V Methylene Chloride	UG/KG	.	14.0000 U
V Styrene	UG/KG	.	10.0000 U
V Tetrachloroethene	UG/KG	.	10.0000 U
V Toluene	UG/KG	.	10.0000 U
V Trichloroethene	UG/KG	.	10.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	33S57D01DL	33SAND01
	LAB ID:	547127 D11	529909
	DATE RCVD:	04/22/93	12/18/92
NAME	UNITS		
V Vinyl Chloride	UG/KG		10.0000 U
V Xylene (total)	UG/KG		10.0000 U
V cis-1,3-Dichloropropene	UG/KG		10.0000 U
V trans-1,3-Dichloropropene	UG/KG		10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF13	NASPAP14	NASPAF14MS	NASPAF14MSD	RB121192	RB121192
	LAB ID:	534681	546686	546687	546688	528679	528680
	DATE RCVD:	02/04/93	04/21/93	04/21/93	04/21/93	12/12/92	12/12/92
NAME	UNITS						
I Aluminum	UG/L	83.4000 J	142.0000 J	.	.	.	171.0000 J
I Antimony	UG/L	52.0000 U	36.0000 U	.	.	.	60.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	.	.	.	4.0000 U
I Barium	UG/L	2.0000 U	13.9000 J	.	.	.	2.2000 J
I Beryllium	UG/L	1.0000 U	1.0000 U	.	.	.	1.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	.	.	.	5.0000 U
I Calcium	UG/L	49.0000 J	512.0000 J	.	.	.	93.4000 J
I Chromium	UG/L	6.0000 U	8.0000 U	.	.	.	6.0000 U
I Cobalt	UG/L	9.0000 U	8.0000 U	.	.	.	9.0000 U
I Copper	UG/L	4.0000 U	4.0000 U	.	.	.	5.0000 U
I Cyanide	UG/L	10.0000 U	10.0000 U	.	.	.	10.0000 U
I Iron	UG/L	40.1000 J	62.9000 J	.	.	.	34.0000 U
I Lead	UG/L	2.0000 U	2.0000 U	.	.	.	2.5000 J
I Magnesium	UG/L	47.0000 U	349.0000 J	.	.	.	46.0000 U
I Manganeee	UG/L	1.0000 U	3.9000 J	.	.	.	2.0000 U
I Mercury	UG/L	0.2000 U	0.2000 U	.	.	.	0.2000 U
I Nickel	UG/L	13.0000 U	17.0000 U	.	.	.	11.0000 U
I Potaeeium	UG/L	1440.0000 U	1350.0000 U	.	.	.	1360.0000 U
I Selenium	UG/L	4.0000 U	3.0000 U	.	.	.	4.0000 U
I Silver	UG/L	5.0000 U	7.0000 U	.	.	.	6.0000 U
I Sodium	UG/L	371.0000 J	679.0000 J	.	.	.	315.0000 J
I Thallium	UG/L	3.0000 U	4.0000 U	.	.	.	2.0000 U
I Vanadium	UG/L	4.0000 U	6.0000 U	.	.	.	5.0000 U
I Zinc	UG/L	5.0000 U	13.9000 J	.	.		3.9000 J
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	.	.	0.1000 J	
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	.	.	0.1000 J	
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	.	.	0.1000 J	
P Aldrin	UG/L	0.0500 U	0.0500 U	.	.	0.0500 J	

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID. / LAB ID: / DATE RCVD. / UNITS	NASPAF13 534681 02/04/93	NASPAF14 546686 04/21/93	NASPAF14MS 546687 04/21/93	NASPAF14MSD 546688 04/21/93	RB121192 528679 12/12/92	RB121192 528680 12/12/92
P Aroclor 1016	UG/L	1.0000 U	1.0000 ω	.	.	1.0000 J	.
P Aroclor 1221	UG/L	2.0000 U	2.0000 ω	.	.	2.0000 J	.
P Aroclor 1232	UG/L	1.0000 U	1.0000 ω	.	.	1.0000 J	.
P Aroclor 1242	UG/L	1.0000 U	1.0000 ω	.	.	1.0000 J	.
P Aroclor 1248	UG/L	1.0000 U	1.0000 ω	.	.	1.0000 J	.
P Aroclor 1254	UG/L	1.0000 U	1.0000 ω	.	.	1.0000 J	.
P Aroclor 1260	UG/L	1.0000 U	1.0000 ω	.	.	1.0000 J	.
P Dieldrin	UG/L	0.1000 U	0.1000 ω	.	.	0.1000 J	.
P Endosulfan I	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
P Endosulfan II	UG/L	0.1000 U	0.1000 ω	.	.	0.1000 J	.
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 ω	.	.	0.1000 J	.
P Endrin	UG/L	0.1000 U	0.1000 ω	.	.	0.1000 J	.
P Endrin aldehyde	UG/L	0.1000 U	0.1000 ω	.	.	0.1000 J	.
P Endrin ketone	UG/L	0.1000 U	0.1000 ω	.	.	0.1000 J	.
P Heptachlor	UG/L	0.0010 J	0.0500 ω	.	.	0.0010 J	.
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
P Methoxychlor	UG/L	0.5000 U	0.5000 ω	.	.	0.5000 J	.
P Toxaphene	UG/L	5.0000 U	5.0000 ω	.	.	5.0000 J	.
P alpha-BHC	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
P alpha-Chlordane	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
P beta-BHC	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
P delta-BHC	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
P gamma-Chlordane	UG/L	0.0500 U	0.0500 ω	.	.	0.0500 J	.
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 ω	.	.	10.0000 U	.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 ω	.	.	10.0000 U	.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 ω	.	.	10.0000 U	.
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 ω	.	.	10.0000 U	.
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 ω	.	.	25.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	NASPAF13 534681 02/04/93	NASPAF14 546686 04/21/93	NASPAF14MS 546687 04/21/93	NASPAF14MSD 546688 04/21/93	RB121192 528679 12/12/92	RB121192 528680 12/12/92
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	.
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	.
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	.
S Acenaphthene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Anthracene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	NASPAF13	NASPAF14	NASPAF14MS	NASPAF14MSD	RB121192	RB121192
ENSAFE ID:		NASPAF13	NASPAF14	NASPAF14MS	NASPAF14MSD	RB121192	RB121192
LAB ID:		534681	546686	546687	546688	528679	528680
DATE RCVD:		02/04/93	04/21/93	04/21/93	04/21/93	12/12/92	12/12/92
S Butylbenzylphthalate	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Carbazole	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Chrysene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Dibenz(a,h)anthracene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Diethylphthalate	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Fluoranthene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Fluorene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Isophorone	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Naphthalene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	.
S Phenanthrene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Phenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S Pyrene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	1.0000 U	9.0000 U	.	.	5.0000 J	.
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF13	NASPAP14	NASPAF14MS	NASPAF14MSD	RB121192	RB121192
	LAB ID:	534681	546606	546687	546688	528679	528680
	DATE RCVD:	02/04/93	04/21/93	04/21/93	04/21/93	12/12/92	12/12/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	40.0000	41.0000	10.0000 U	.
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 J	
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V Benzene	UG/L	10.0000 U	10.0000 U	43.0000	45.0000	10.0000 U	
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	43.0000	46.0000	10.0000 U	.
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Methylene Chloride	UG/L	9.0000 U	2.0000 U	13.0000 U	13.0000 U	1.0000 J	
V Styrene	UG/L	10.0000 U	10.0000 U	10,0000 U	10.0000 U	10.0000 U	.
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.
V Toluene	UG/L	10.0000 U	10.0000 U	43.0000	47.0000	3.0000 J	.
V Trichloroethene	UG/L	10.0000 U	10.0000 U	41.0000	46.0000	10.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF13	NASPAF14	NASPAF14MS	NASPAF14MSD	RB121192	RB121192
	LAB ID:	534681	546686	546687	546688	528679	528680
	DATE RCVD:	02/04/93	04/21/93	04/21/93	04/21/93	12/12/92	12/12/92
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	TB120892 / 527770 / 12/09/92	TB121092 / 528678 / 12/12/92	TB60 / 546720 / 04/21/93	TEMPBLANK / 524186 / 11/20/92	TEMPBLANK / 527771 / 12/09/92	TEMPBLANK / 528140 / 12/10/92
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 J	130.0000 J	8.0000 J
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	17.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1.0000 J	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	2.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	2.0000 J	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	4.0000 J	2.0000 J	3.0000 J	7.0000 J	6.0000 J	2.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	2.0000 J	10.0000 U	10.0000 U	4.0000 J	1.0000 J

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	TB120892	TB121092	TB60	TEMPBLANK	TEMPBLANK	TEMPBLANK
ENSAFE ID:		TB120892	TB121092	TB60	TEMPBLANK	TEMPBLANK	TEMPBLANK
LAB ID:		527770	528678	546720	524186	527771	528140
DATE RCVD:		12/09/92	12/12/92	04/21/93	11/20/92	12/09/92	12/10/92
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	lo.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	TRIPBLANK	TRIPBLANK				
	LAB ID:	524167	529357				
	DATE RCVD:	11/20/92	12/16/92				
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U				
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U				
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U				
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U				
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U				
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U				
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U				
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U				
V 2-Butanone	UG/L	10.0000 U	10.0000 U				
V 2-Hexanone	UG/L	10.0000 U	10.0000 U				
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U				
V Acetone	UG/L	10.0000 U	10.0000 U				
V Benzene	UG/L	10.0000 U	10.0000 U				
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U				
V Bromoform	UG/L	10.0000 U	10.0000 U				
V Bromomethane	UG/L	10.0000 U	10.0000 U				
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U				
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U				
V Chlorobenzene	UG/L	10.0000 U	10.0000 U				
V Chloroethane	UG/L	10.0000 U	10.0000 U				
V Chloroform	UG/L	10.0000 U	10.0000 U				
V Chloromethane	UG/L	10.0000 U	10.0000 U				
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U				
V Ethylbenzene	UG/L	10.0000 U	10.0000 U				
V Methylene Chloride	UG/L	6.0000 J	1.0000 J				
V Styrene	UG/L	10.0000 U	10.0000 U				
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U				
V Toluene	UG/L	10.0000 U	10.0000 U				

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	TRIPBLANK	TRIPBLANK			
	LAB ID:	524167	529357			
	DATE RCVD:	11/20/92	12/16/92	.	.	.
NAME	UNITS					
V Trichloroethene	UG/L	10.0000 U	10.0000 U	.	.	.
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	.	.	.
V Xylene (total)	UG/L	10.0000 U	10.0000 U	.	.	.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	.	.	.
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	.	.	.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G07001MS	13G07001MSD	13G07E01	13S01E01	13S04E01	13S60E01
	LAB ID:	533767	533768	533779	527759	532929	531928
	DATE RCVD:	01/27/93	01/27/93	01/27/93	12/09/92	01/20/93	01/14/93
NAME	UNITS						
I Aluminum	UG/L	.	.	229.0000	183.0000 J	228.0000	262.0000
I Antimony	UG/L	.	.	57.0000 U	60.0000 U	57.0000 U	57.0000 U
I Arsenic	UG/L	.	.	3.0000 U	4.0000 U	3.0000 U	3.0000 U
I Barium	UG/L	.	.	6.0000 U	2.5000 J	6.0000 U	6.0000 U
I Beryllium	UG/L	.	.	3.0000 U	1.0000 U	3.0000 U	3.0000 U
I Cadmium	UG/L	.	.	5.0000 U	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	.	.	130.0000 U	1190.0000 J	130.0000 U	153.0000 J
I Chromium	UG/L			10.0000 U	6.0000 U	10.0000 U	10.0000 U
I Cobalt	UG/L	.		12.0000 U	9.0000 U	12.0000 U	12.0000 U
I Copper	UG/L	.	.	25.0000 U	5.0000 U	25.0000 U	25.0000 U
I Cyanide	UG/L	.	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
I Iron	UG/L	.	.	513.0000	34.0000 U	248.0000	65.1000 J
I Lead	UG/L	.	.	2.0000 U	2.0000 U	2.0000 U	2.0000 U
I Magnesium	UG/L	.	.	552.0000 U	591.0000 J	552.0000 U	552.0000 U
I Manganese	UG/L	.	.	7.3000 J	2.0000 U	6.0000 J	6.0000 U
I Mercury	UG/L	.	.	0.2000 U	0.2000 U	0.2000 U	0.2000 U
I Nickel	UG/L	.	.	38.0000 U	11.0000 U	38.0000 U	38.0000 U
I Potassium	UG/L	.	.	3950.0000 U	1360.0000 U	3950.0000 U	3950.0000 U
I Selenium	UG/L	.	.	3.0000 U	4.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	.	.	10.0000 U	6.0000 U	10.0000 U	10.0000 U
I Sodium	UG/L	.	.	407.0000 J	903.0000 J	445.0000 J	465.0000 J
I Thallium	UG/L	.	.	3.0000 U	2.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	.	.	15.0000 U	5.0000 U	15.0000 U	15.0000 U
I Zinc	UG/L			95.6000	21.2000	23.3000	47.3000
P 4,4'-DDD	UG/L	0.0090 J	0.0069 J	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	0.6300	0.6100	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Aldrin	UG/L	0.2900	0.2800	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits

	ENSAFE ID.	13G07001MS	13G07001MSD	13G07E01	13S01E01	13S04E01	13S60E01
	LAB ID:	533773	533774	533779	527759	532929	531928
	DATE RCVD.	01/27/93	01/27/93	01/27/93	12/09/92	01/20/93	01/14/93
NAME	UNITS						
P Aroclor 1221	UG/L	2.0000	2.0000	2.0000	2.0000 U	2.0000	2.0000
P Aroclor 1232	UG/L	1.0000	1.0000	1.0000	1.0000 U	1.0000	1.0000
P Aroclor 1242	UG/L	1.0000	1.0000	1.0000	1.0000 U	1.0000	1.0000
P Aroclor 1248	UG/L	1.0000	1.0000	1.0000	1.0000 U	1.0000	1.0000
P Aroclor 1254	UG/L	1.0000	1.0000	1.0000	1.0000 U	1.0000	1.0000
P Aroclor 1260	UG/L	1.0000	1.0000	1.0000	1.0000 U	1.0000	1.0000
P Dieldrin	UG/L	0.6600	0.6500	0.1000	0.1000 U	0.1000	0.1000
P Endosulfan I	UG/L	0.0500	0.0500	0.0500	0.0500 U	0.0500	0.0500
P Endosulfan II	UG/L	0.1000	0.1000	0.1000	0.1000 U	0.1000	0.1000
P Endosulfan sulfate	UG/L	0.1000	0.1000	0.1000	0.1000 U	0.1000	0.1000
P Endrin	UG/L	0.6800	0.6600	0.1000	0.1000 U	0.1000	0.1000
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.1000	0.1000 U	0.1000	0.1000
P Endrin ketone	UG/L	0.0030 J	0.0030 J	0.1000	0.1000 U	0.1000	0.1000
P Heptachlor	UG/L	0.3100	0.3000	0.0500	0.0024 J	0.0024	0.0500
P Heptachlor epoxide	UG/L	0.0500	0.0500	0.0500	0.0500 U	0.0500	0.0500
P Methoxychlor	UG/L	0.5000	0.5000	0.5000	0.5000 U	0.5000	0.5000
P Toxaphene	UG/L	5.0000	5.0000	5.0000	5.0000 U	5.0000	5.0000
P alpha-BHC	UG/L	0.0500	0.0500	0.0500	0.0500 U	0.0500	0.0500
P alpha-Chlordane	UG/L	0.0500	0.0500	0.0500	0.0500 U	0.0500	0.0500
P beta-BHC	UG/L	0.0500	0.0500	0.0500	0.0500 U	0.0500	0.0500
P delta-BHC	UG/L	0.0500	0.0500	0.0500	0.0500 U	0.0500	0.0500
P gamma-BHC (Lindane)	UG/L	0.3500	0.3400	0.0500	0.0500 U	0.0500	0.0500
P gamma-Chlordane	UG/L	0.0500	0.0500	0.0500	0.0500 U	0.0023	0.0500
S 1,2,4-Trichlorobenzene	UG/L	17.0000	19.0000	10.0000	10.0000 U	10.0000	10.0000
S 1,2-Dichlorobenzene	UG/L	10.0000	10.0000	10.0000	10.0000 U	10.0000	10.0000
S 1,3-Dichlorobenzene	UG/L	10.0000	10.0000	10.0000	10.0000 U	10.0000	10.0000
S 1,4-Dichlorobenzene	UG/L	14.0000	16.0000	10.0000	10.0000 U	10.0000	10.0000
S 2,4,5-Trichlorophenol	UG/L	25.0000	25.0000	25.0000	25.0000 U	25.0000	25.0000
S 2,4,6-Trichlorophenol	UG/L	10.0000	10.0000	10.0000	10.0000 U	10.0000	10.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G07001MS	13G07001MSD	13G07E01	13S01E01	13S04E01	13S60E01
	LAB ID:	533770	533771	533779	527759	532929	531928
	DATE RCVD:	01/27/93	01/27/93	01/27/93	12/09/92	01/20/93	01/14/93
NAME	UNITS						
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	19.0000	22.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	24.0000	28.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	28.0000	32.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000	30.0000	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Acenaphthene	UG/L	19.0000	22.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Butylbenzylphthalate	UG/L	7.0000 U	8.0000 U	2.0000 U	10.0000 U	5.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G07001MS	13G07001MSD	13G07E01	13S01E01	13S04E01	13S60E01
	LAB ID:	533770	533771	533779	527759	532929	531928
	DATE RCVD:	01/27/93	01/27/93	01/27/93	12/09/92	01/20/93	01/14/93
NAME	UNITS						
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	2.0000 U	2.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz (a,h) anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	2.0000 U	2.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Ieophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	20.0000	23.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	29.0000	34.0000	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Phenol	UG/L	24.0000	27.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pyrene	UG/L	17.0000	19.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis (2-Ethlyhexy)phthalate	UG/L	5.0000 U	6.0000 U	2.0000 U	16.0000 U	5.0000 U	3.0000 U
S bie (2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bie (2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	10.0.000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G07001MS	13G07001MSD	13G07E01	13S01E01	13S04E01	13S60E01
	LAB ID:	533767	533768	533779	527759	532929	531928
	DATE RCVD:	01/27/93	01/27/93	01/27/93	12/09/92	01/20/93	01/14/93
NAME	UNITS						
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	46.0000	45.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	49.0000 U	10.0000 U	92.0000	72.0000 U
V Benzene	UG/L	51.0000	51.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromofo m	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10,0000 U
V Chlorobenzene	UG/L	50.0000	50.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	9.0000 U	10.0000 U	3.0000 U	10.0000 U	7.0000 U	1.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	50.0000	51.0000	10.0000 U	2.0000 J	10.0000 U	1.0000 J
V Trichloroethene	UG/L	48.0000	48.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13G07001MS	13G07001MSD	13G07E01	13S01E01	13S04E01	13S60E01
	LAB ID:	533767	533768	533779	527759	532929	531928
	DATE RCVD:	01/27/93	01/27/93	01/27/93	12/09/92	01/20/93	01/14/93
NAME	UNITS						
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	13S60E02	13S60E03	13S60E04	33G00E30	33G00T01	33G00T02
	LAB ID:	532080	532928	532919	547108	533872	534238
	DATE RCVD:	01/15/93	01/20/93	01/20/93	04/22/93	01/27/93	01/29/93
I Aluminum	UG/L	257.0000	280.0000	258.0000	99.4000 J	.	.
I Antimony	UG/L	57.0000 U	57.0000 U	57.0000 U	36.0000 U	.	.
I Arsenic	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	.	.
I Barium	UG/L	6.0000 U	6.0000 U	6.0000 U	2.0000 U	.	.
I Beryllium	UG/L	3.0000 U	3.0000 U	3.0000 U	1.0000 U	.	.
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	.	.
I Calcium	UG/L	146.0000 J	130.0000 U	634.0000 J	115.0000 J	.	.
I Chromium	UG/L	10.0000 U	10.0000 U	10.0000 U	8.0000 U	.	.
I Cobalt	UG/L	12.0000 U	12.0000 U	12.0000 U	8.0000 U	.	.
I Copper	UG/L	25.0000 U	25.0000 U	25.0000 U	4.0000 U	.	.
I Cyanide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
I Iron	UG/L	76.1000 J	159.0000	66.6000 J	41.0000 U	.	.
I Lead	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	.	.
I Magnesium	UG/L	552.0000 U	552.0000 U	552.0000 U	59.0000 U	.	.
I Manganese	UG/L	6.0000 U	6.0000 U	6.0000 U	1.0000 U	.	.
I Mercury	UG/L	0.2000 U	0.2000 U	0.2000 U	0.2000 U	.	.
I Nickel	UG/L	38.0000 U	38.0000 U	38.0000 U	17.0000 U	.	.
I Potassium	UG/L	3950.0000 U	3950.0000 U	3950.0000 U	1350.0000 U	.	.
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	3.0000 U	.	.
I Silver	UG/L	10.0000 U	10.0000 U	10.0000 U	7.0000 U	.	.
I sodium	UG/L	511.0000 J	508.0000 J	457.0000 J	279.0000 J	.	.
I Thallium	UG/L	3.0000 U	3.0000 U	3.0000 U	4.0000 U	.	.
I Vanadium	UG/L	15.0000 U	15.0000 U	15.0000 U	6.0000 U	.	.
I Zinc	UG/L	22.1000	41.6000	18.3000 J	4.0000 U	.	.
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.0310 J	0.1000 U	.	.
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	.	.
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	.	.
P Aldrin	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	13S60E02	13S60E03	13S60E04	33G00E30	33G00T01	33G00T02
	LAB ID:	532080	532928	532919	547108	533872	534238
	DATE RCVD:	01/15/93	01/20/93	01/20/93	04/22/93	01/27/93	01/29/93
NAME	UNITS						
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	.	.
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	.	.
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	.	.
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	.	.
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	.	.
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	.	.
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	.	.
P Dieldrin	UG/L	0.0023 J	0.1000 U	0.1000 U	0.1000 U	.	.
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.
P Endosulfan II	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	.	.
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	.	.
P Endrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	.	.
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	.	.
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	.	.
P Heptachlor	UG/L	0.0500 U	0.0022 J	0.0500 U	0.0500 U	.	.
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.
P Methoxychlor	UG/L	0.5000 U	0.5000 U	0.5000 U	0.5000 U	.	.
P Toxaphene	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	.	.
P alpha-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	.	.
P gamma-Chlordane	UG/L	0.0500 U	0.0023 J	0.0025 J	0.0500 U	.	.
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S60E02	13S60E03	13560304	33G00E30	33G00T01	33G00T02
	LAB ID:	532080	532928	532919	547108	533872	534238
	DATE RCVD:	01/15/93	01/20/93	01/20/93	04/22/93	01/27/93	01/29/93
NAME	UNITS						
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Butylbenzylphthalate	UG/L	3.0000 U	3.0000 U	5.0000 U	10.0000 U		.
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Dibenz(a,h)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Diethylphthalate	UG/L	10.0000 U	1.0000 U	10.0000 U	10.0000 U		.
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Isophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U		.
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Phenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S Pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.
S bis(2-Ethlyhexy)phthalate	UG/L	5.0000 U	1.0000 U	2.0000 U	2.0000 U		
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	13S60E02	13S60E03	13S60E04	33G00E30	33G00T01	33G00T02
	LAB ID:	532080	532928	532919	547108	533872	534238
	DATE RCVD:	01/15/93	01/20/93	01/20/93	04/22/93	01/27/93	01/29/93
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID:	33G00T03	33G00T04	33G00T05	33G02D01	33G02D01F	33G04D01
	LAB ID:	534319	534613	534794	534217	534192	534206
	DATE RCVD:	01/30/93	02/03/93	02/04/93	01/29/93	01/29/93	01/29/93
	UNITS						
I Aluminum	UG/L	.	.	.	1880.0000	269.0000	3040.0000
I Antimony	UG/L	.	.	.	57.0000 U	52.0000 U	57.0000 U
I Arsenic	UG/L	.	.	.	3.2000 J	5.1000 J	4.1000 J
I Barium	UG/L	.	.	.	28.2000 J	23.1000 J	31.8000 J
I Beryllium	UG/L	.	.	.	3.0000 U	1.0000 U	3.0000 U
I Cadmium	UG/L	.	.	.	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	.	.	.	32800.0000	31800.0000	31800.0000
I Chromium	UG/L	.	.	.	19.8000	6.0000 U	10.0000 U
I Cobalt	UG/L	.	.	.	12.0000 U	9.0000 U	12.0000 U
I Copper	UG/L	.	.	.	25.0000 U	4.0000 U	25.0000 U
I Cyanide	UG/L	.	.	.	10.0000 U		10.0000 U
I Iron	UG/L	.	.	.	6300.0000	3590.0000	3170.0000
I Lead	UG/L	.	.	.	17.2000	2.0000 U	5.5000
I Magnesium	UG/L	.	.	.	2070.0000 J	1960.0000 J	3340.0000 J
I Manganese	UG/L	.	.	.	92.7000	86.8000	38.3000
I Mercury	UG/L	.	.	.	0.3100	0.2000 U	0.3100
I Nickel	UG/L	.	.	.	38.0000 U	13.0000 U	38.0000 U
I Potaeeium	UG/L	.	.	.	3950.0000 U	2390.0000 J	4440.0000 J
I Selenium	UG/L	.	.	.	3.0000 U	3.0000 U	3.0000 U
I Silver	UG/L	.	.	.	10.0000 U	5.0000 U	10.0000 U
I Sodium	UG/L	.	.	.	2870.0000 J	3160.0000 J	36000.0000
I Thallium	UG/L	.	.	.	3.0000 U	3.0000 U	3.0000 U
I Vanadium	UG/L	.	.	.	15.0000 U	4.0000 U	15.0000 U
I Zinc	UG/L	.	.	.	34.9000	8.6000 J	16.3000 J
P 4,4'-DDD	UG/L	.	.	.	0.1000 U		0.1000 U
P 4,4'-DDE	UG/L	.	.	.	0.1000 U		0.1000 U
P 4,4'-DDT	UG/L	.	.	.	0.1000 U		0.1000 U
P Aldrin	UG/L	.	.	.	0.0500 U		0.0500 U

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33600T03	3360OT04	33G00T05	33G02D01	33G02D01F	33G04D01
ENSAFE ID:		33600T03	3360OT04	33G00T05	33G02D01	33G02D01F	33G04D01
LAB ID:		534319	534613	534794	534231	534192	534226
DATE RCVD:		01/30/93	02/03/93	02/04/93	01/29/93	01/29/93	01/29/93
P Aroclor 1016	UG/L	.	.	.	1.0000 U	.	1.0000 U
P Aroclor 1221	UG/L	.	.	.	2.0000 U	.	2.0000 U
P Aroclor 1232	UG/L	.	.	.	1.0000 U	.	1.0000 U
P Aroclor 1242	UG/L	.	.	.	1.0000 U	.	1.0000 U
P Aroclor 1248	UG/L	.	.	.	1.0000 U	.	1.0000 U
P Aroclor 1254	UG/L	.	.	.	1.0000 U	.	1.0000 U
P Aroclor 1260	UG/L	.	.	.	1.0000 U	.	1.0000 U
P Dieldrin	UG/L	.		.	0.1000 U	.	0.0028 J
P Endosulfan I	UG/L	.	.	.	0.0500 U	.	0.0500 U
P Endosulfan II	UG/L	.	.	.	0.1000 U	.	0.1000 U
P Endosulfan sulfate	UG/L	.	.	.	0.1000 U	.	0.1000 U
P Endrin	UG/L	.	.	.	0.1000 U	.	0.1000 U
P Endrin aldehyde	UG/L	.	.	.	0.1000 U	.	0.1000 U
P Endrin ketone	UG/L	.	.	.	0.1000 U	.	0.1000 U
P Heptachlor	UG/L	.	.	.	0.0029 J	.	0.0022 J
P Heptachlor epoxide	UG/L	.	.	.	0.0500 U	.	0.0500 U
P Methoxychlor	UG/L	.	.	.	0.5000 U	.	0.5000 U
P Toxaphene	UG/L	.	.	.	5.0000 U	.	5.0000 U
P alpha-BHC	UG/L	.	.	.	0.0500 U	.	0.0500 U
P alpha-Chlordane	UG/L	.	.	.	0.0500 U	.	0.0500 U
P beta-BHC	UG/L	.	.	.	0.0500 U		0.0500 U
P delta-BHC	UG/L	.	.	.	0.0500 U	.	0.0500 U
P gamma-BHC (Lindane)	UG/L	.	.		0.0500 U		0.0500 U
P gamma-Chlordane	UG/L	.	.		0.0500 U	.	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 1,2-Dichlorobenzene	UG/L	.	.		10.0000 U	.	10.0000 U
S 1,3-Dichlorobenzene	UG/L	.	.		10.0000 U	.	10.0000 U
S 1,4-Dichlorobenzene	UG/L	.	.	.	10.0000 U		10.0000 U
S 2,4,5-Trichlorophenol	UG/L	.	.	.	25.0000 U	.	25.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G00T03	33G00T04	33G00T05	33G02D01	33G02D01F	33G04D01
	LAB ID:	534319	534613	534794	534231	534192	534226
	DATE RCVD:	01/30/93	02/03/93	02/04/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 2,4-Dichlorophenol	UG/L		.		10.0000 U	.	10.0000 U
S 2,4-Dimethylphenol	UG/L			.	10.0000 U	.	10.0000 U
S 2,4-Dinitrophenol	UG/L	.		.	25.0000 U	.	25.0000 U
S 2,4-Dinitrotoluene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 2,6-Dinitrotoluene	UG/L			.	10.0000 U	.	10.0000 U
S 2-Chloronaphthalene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 2-Chlorophenol	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 2-Methylnaphthalene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 2-Methylphenol	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 2-Nitroaniline	UG/L	.	.	.	25.0000 U	.	25.0000 U
S 2-Nitrophenol	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 3-Nitroaniline	UG/L	.	.	.	25.0000 U	.	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	.	.	.	25.0000 U	.	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 4-Chloroaniline	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 4-Methylphenol	UG/L	.	.	.	10.0000 U	.	10.0000 U
S 4-Nitroaniline	UG/L	.	.	.	25.0000 U	.	25.0000 U
S 4-Nitrophenol	UG/L	.	.	.	25.0000 U	.	25.0000 U
S Acenaphthene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Acenaphthylene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Anthracene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Benzo(a)anthracene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Benzo(a)pyrene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Benzo(g,h,i)perylene	UG/L	.	.	.	10.0000 U	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33G00T03	33G00T04	33G00T05	33G02D01	33G02D01F	33G04D01
ENSAFE ID:		33G00T03	33G00T04	33G00T05	33G02D01	33G02D01F	33G04D01
LAB ID:		534319	534613	534794	534231	534192	534226
DATE RCVD:		01/30/93	02/03/93	02/04/93	01/29/93	01/29/93	01/29/93
S Butylbenzylphthalate	UG/L	.		.	10.0000 U	.	10.0000 U
S Carbazole	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Chrysene	UG/L	.	.		10.0000 U	.	10.0000 U
S Di-n-butylphthalate	UG/L				10.0000 U	.	10.0000 U
S Di-n-octylphthalate	UG/L	.		.	10.0000 U	.	10.0000 U
S Dibenz(a,h)anthracene	UG/L	.	.		10.0000 U	.	10.0000 U
S Dibenzofuran	UG/L	.	.		10.0000 U	.	10.0000 U
S Diethylphthalate	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Dimethylphthalate	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Fluoranthene	UG/L	.		.	10.0000 U	.	10.0000 U
S Fluorene	UG/L	.	.		10.0000 U	.	10.0000 U
S Hexachlorobenzene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Hexachlorobutadiene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Hexachlorocyclopentadiene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Hexachloroethane	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Isophorone	UG/L	.	.	.	10.0000 U	.	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	.	.	.	10.0000 U	.	10.0000 U
S N-Nitrosodiphenylamine	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Naphthalene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Nitrobenzene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Pentachlorophenol	UG/L	.	.	.	25.0000 U	.	25.0000 U
S Phenanthrene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Phenol	UG/L	.	.	.	10.0000 U	.	10.0000 U
S Pyrene	UG/L	.	.	.	10.0000 U	.	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	.	.	.	10.0000 U	.	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	.	.	.	4.0000 U	.	3.0000 U
S bis(2-chloroethyl)ether	UG/L	.	.	.	10.0000 U	.	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L		.	.	10.0000 U	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preaented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33G00T03	33G00T04	33G00T05	33G02D01	33G02D01F	33G04D01
ENSAFE ID:		33G00T03	33G00T04	33G00T05	33G02D01	33G02D01F	33G04D01
LAB ID:		534319	534613	534794	534231	534192	534226
DATE RCVD:		01/30/93	02/03/93	02/04/93	01/29/93	01/29/93	01/29/93
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	81.0000 U		10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	8.0000 J		10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Methylene Chloride	UG/L	2.0000 U	2.0000 U	2.0000 U	10.0000 U		2.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G00T03	33G00T04	33G00IO5	33G02D01	33G02D01F	33G04D01
	LAB ID:	534319	534613	534794	534231	534192	534226
	DATE RCVD:	01/30/93	02/03/93	02/04/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U		10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	BNSAPE ID:	33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
	LAB ID:	534187	534297	534298	534603	534603	534625
	DATE RCVD:	01/29/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
NAME	UNITS						
I Aluminum	UG/L	234.0000			972.0000	.	65.2000 J
I Antimony	UG/L	52.0000 U			29.0000 U	.	30.0000 U
I Arsenic	UG/L	3.9000 J			18.7000	.	15.0000 U
I Barium	UG/L	32.7000 J			86.2000 J	.	78.6000 J
I Beryllium	UG/L	1.0000 U			3.0000 U	.	1.0000 U
I Cadmium	UG/L	5.0000 U			5.0000 U	.	5.0000 U
I Calcium	UG/L	34100.0000			167000.000	.	161000.000
I Chromium	UG/L	6.0000 U			10.0000 U	.	6.0000 U
I Cobalt	UG/L	9.0000 U			12.0000 U	.	9.0000 U
I Copper	UG/L	4.0000 U			25.0000 U	.	4.0000 U
I Cyanide	UG/L				10.0000 U	.	
I Iron	UG/L	964.0000			4470.0000	.	3970.0000
I Lead	UG/L	2.0000 U			20.0000 U	.	2.6000 J
I Magnesium	UG/L	3570.0000 J			318000.000	.	311000.000
I Manganese	UG/L	38.0000			500.0000	.	453.0000
I Mercury	UG/L	0.2000 U			0.2000 U	.	0.2000 U
I Nickel	UG/L	13.0000 U			38.0000 U	.	13.0000 U
I Potassium	UG/L	4860.0000 J			105000.000	.	100000.000
I Selenium	UG/L	3.0000 U			3.0000 U	.	15.0000 U
I Silver	UG/L	5.0000 U			10.0000 U	.	5.0000 U
I Sodium	UG/L	43100.0000			1360000.00	.	2600000.00
I Thallium	UG/L	3.0000 U			30.0000 U	.	3.0000 U
I Vanadium	UG/L	4.0000 U			15.0000 U	.	4.0000 U
I Zinc	UG/L	8.6000 J			29.3000	.	7.8000 J
P 4,4'-DDD	UG/L		0.1000 U	0.1000 U	0.1000 U	.	
P 4,4'-DDE	UG/L		0.1000 U	0.1000 U	0.1000 U	.	
P 4,4'-DDT	UG/L		0.7800	0.7900	0.1000 U	.	
P Aldrin	UG/L		0.2500 J	0.2400 J	0.0500 U	.	

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
ENSAFE ID:		33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
LAB ID:		534187	534303	534304	534603	534603	534625
DATE RCM:		01/29/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
P Aroclor 1016	UG/L		1.0000 U	1.0000 U	1.0000 U		.
P Aroclor 1221	UG/L		2.0000 U	2.0000 U	2.0000 U		.
P Aroclor 1232	UG/L		1.0000 U	1.0000 U	1.0000 U		.
P Aroclor 1242	UG/L		1.0000 U	1.0000 U	1.0000 U		.
P Aroclor 1248	UG/L		1.0000 U	1.0000 U	1.0000 U		.
P Aroclor 1254	UG/L		1.0000 U	1.0000 U	1.0000 U		.
P Aroclor 1260	UG/L		1.0000 U	1.0000 U	1.0000 U		.
P Dieldrin	UG/L		0.7500	0.7500	0.1000 U		.
P Endosulfan I	UG/L		0.0500 U	0.0500 U	0.0500 U		.
P Endosulfan II	UG/L		0.1000 U	0.1000 U	0.1000 U		.
P Endosulfan sulfate	UG/L		0.1000 U	0.1000 U	0.1000 U		.
P Endrin	UG/L		0.8000	0.8000 J	0.1000 U		.
P Endrin aldehyde	UG/L		0.0140 J	0.0180 J	0.0300 J		.
P Endrin ketone	UG/L		0.1000 U	0.1000 U	0.1000 U		.
P Heptachlor	UG/L		0.2500 J	0.2400 J	0.0500 U		.
P Heptachlor epoxide	UG/L		0.0500 U	0.0500 U	0.0500 U		.
P Methoxychlor	UG/L		0.5000 U	0.5000 U	0.5000 U		.
P Toxaphene	UG/L		5.0000 U	5.0000 U	5.0000 U		.
P alpha-BHC	UG/L		0.0500 U	0.0500 U	0.0500 U		.
P alpha-Chlordane	UG/L		0.0500 U	0.0500 U	0.0500 U		.
P beta-BHC	UG/L		0.0500 U	0.0500 U	0.0500 U		.
P delta-BHC	UG/L		0.0500 U	0.0500 U	0.0500 U		.
P gamma-BHC (Lindane)	UG/L		0.2700 J	0.2700 J	0.0500 U		.
P gamma-Chlordane	UG/L		0.0500 U	0.0500 U	0.0500 U		.
S 1,2,4-Trichlorobenzene	UG/L		25.0000	23.0000	50.0000 U	80.0000 U	.
S 1,2-Dichlorobenzene	UG/L		10.0000 U	10.0000 U	410.0000 J	370.0000 J	.
S 1,3-Dichlorobenzene	UG/L		10.0000 U	10.0000 U	230.0000	220.0000 J	.
S 1,4-Dichlorobenzene	UG/L		21.0000	20.0000	370.0000	350.0000 J	.
S 2,4,5-Trichlorophenol	UG/L		25.0000 U	25.0000 U	120.0000 U	200.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS					
ENSAFE ID:		33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
LAB ID:		534187	534300	534301	534603	534603	534625
DATE RCVD:		01/29/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
S 2,4,6-Trichlorophenol	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 2,4-Dichlorophenol	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 2,4-Dimethylphenol	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 2,4-Dinitrophenol	UG/L	.	25.0000 U	25.0000 U	120.0000 U	200.0000 U	
S 2,4-Dinitrotoluene	UG/L	.	32.0000	33.0000	50.0000 U	80.0000 U	
S 2,6-Dinitrotoluene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 2-Chloronaphthalene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 2-Chlorophenol	UG/L	.	49.0000	47.0000	50.0000 U	80.0000 U	
S 2-Methylnaphthalene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 2-Methylphenol	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 2-Nitroaniline	UG/L	.	25.0000 U	25.0000 U	120.0000 U	200.0000 U	
S 2-Nitrophenol	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 3,3'-Dichlorobenzidine	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 3-Nitroaniline	UG/L	.	25.0000 U	25.0000 U	120.0000 U	200.0000 U	
S 4,6-Dinitro-2-methylphenol	UG/L	.	25.0000 U	25.0000 U	120.0000 U	200.0000 U	
S 4-Brmophenyl-phenylether	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 4-Chloro-3-methylphenol	UG/L	.	58.0000	53.0000	50.0000 U	80.0000 U	
S 4-Chloroaniline	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 4-Chlorophenyl-phenylether	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 4-Methylphenol	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S 4-Nitroaniline	UG/L	.	25.0000 U	25.0000 U	120.0000 U	200.0000 U	
S 4-Nitrophenol	UG/L	.	52.0000	48.0000	120.0000 U	200.0000 U	
S Acenaphthene	UG/L	.	31.0000	30.0000	11.0000 J	11.0000 J	
S Acenaphthylene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S Anthracene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S Benzo(a)anthracene	UG/L	.	10.0000 U	10.0000 U	50.0000 'U	80.0000 U	
S Benzo(a)pyrene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S Benzo(b,k)fluoranthene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	
S Benzo(g,h,i)perylene	UG/L	.	10.0000 U	10.0000 U	50.0000 U	80.0000 U	

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
	LAB ID:	534187	534300	534301	534603	534603	534625
	DATE RCVD:	01/29/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
NAME	UNITS						
S Butylbenzylphthalate	UG/L		1.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Carbazole	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Chrysene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Di-n-butylphthalate	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Di-n-octylphthalate	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Dibenz(a,h)anthracene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Dibenzofuran	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Diethylphthalate	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Dimethylphthalate	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Fluoranthene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Fluorene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Hexachlorobenzene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Hexachlorobutadiene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Hexachlorocyclopentadiene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Hexachloroethane	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Isophorone	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L		25.0000	22.0000	50.0000 U	80.0000 U	.
S N-Nitrosodiphenylamine	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Naphthalene	UG/L		10.0000 U	10.0000 U	18.0000 J	19.0000 J	.
S Nitrobenzene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Pentachlorophenol	UG/L		65.0000	68.0000	120.0000 U	200.0000 U	.
S Phenanthrene	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S Phenol	UG/L		51.0000	47.0000	50.0000 U	80.0000 U	.
S Pyrene	UG/L		31.0000	31.0000	50.0000 U	80.0000 U	.
S bis(2-Chloroethoxy)methane	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L		20.0000 U	1.0000 U	32.0000 U	38.0000 U	.
S bis(2-chloroethyl)ether	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L		10.0000 U	10.0000 U	50.0000 U	80.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
	LAB ID:	534187	534297	534298	534603	534603	534625
	DATE RCVD:	01/29/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V 1,1,2,2-Tetrachloroethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V 1,1,2-Trichloroethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V 1,1-Dichloroethane	UG/L	.	10.0000 U	10.0000 U	200.0000		.
V 1,1-Dichloroethene	UG/L	.	58.0000	*53.0000*	14.0000 U		.
V 1,2-Dichloroethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V 1,2-Dichloroethene (total)	UG/L	.	10.0000 U	10.0000 U	10.0000 J		.
V 1,2-Dichloropropane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V 2-Butanone	UG/L	.	10.0000 U	10.0000 U	160.0000 U		.
V 2-Hexanone	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V 4-Methyl-2-Pentanone	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Acetone	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Benzene	UG/L	.	58.0000	55.0000	14.0000 U		.
V Bromodichloromethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Bromoform	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Bromomethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Carbon Disulfide	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Carbon Tetrachloride	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Chlorobenzene	UG/L	.	51.0000	49.0000	130.0000		.
V Chloroethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Chloroform	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Chloromethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Dibromochloromethane	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Ethylbenzene	UG/L	.	10.0000 U	10.0000 U	14.0000 U		.
V Methylene Chloride	UG/L	.	2.0000 U	*3.0000* U	14.0000 U		.
V Styrene	UG/L	.	10.0000 U	10.0000 U	5.0000 J		.
V Tetrachloroethene	UG/L	.	10.0000 U	10.0000 U	6.0000 J		.
V Toluene	UG/L	.	52.0000	50.0000	4.0000 J		.
V Trichloroethene	UG/L	.	55.0000	51.0000	2.0000 J		.

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
ENSAFE ID:		33G04D01F	33G10001MS	33G10001MSD	33G12D01	33G12D01DL	33G12D01F
LAB ID:		534187	534297	534298	534603	534603	534625
DATE RCVD:		01/29/93	02/03/93	02/03/93	02/03/93	02/03/93	02/03/93
V Vinyl Chloride	UG/L		10.0000 U	10.0000 U	16.0000		.
V Xylene (total)	UG/L		10.0000 U	10.0000 U	4.0000 J		.
V cis-1,3-Dichloropropene	UG/L		10.0000 U	10.0000 U	14.0000 U		.
V trans-1,3-Dichloropropene	UG/L		10.0000 U	10.0000 U	14.0000 U		.

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33G18E01 / 534605 / 02/03/93	33G20001MS / 546725 / 04/27/93	33G20001MSD / 546726 / 04/27/93	33020D01 / 547567 / 04/27/93	33G20D01DL / 547567 / 04/27/93	33G20D01F / 547574 / 04/27/93
I Aluminum	UG/L	80.5000 J	.	.	2870.0000 J	.	56.9000 J
I Antimony	UG/L	29.0000 U		.	36.0000 U	.	36.0000 U
I Arsenic	UG/L	3.0000 U	.	.	3.4000 J	.	3.0000 U
I Barium	UG/L	6.0000 U	.	.	19.8000 J	.	17.6000 J
I Beryllium	UG/L	3.0000 u	.	.	1.0000 U	.	1.0000 U
I Cadmium	UG/L	5.0000 U		.	5.0000 U	.	5.0000 U
I Calcium	UG/L	176.0000 J	.	.	60100.0000	.	67200.0000
I Chromium	UG/L	10.0000 u		.	11.6000	.	8.0000 U
I Cobalt	UG/L	12.0000 u	.	.	8.0000 U	.	8.0000 u
I Copper	UG/L	25.0000 U		.	5.2000 J	.	4.0000 u
I Cyanide	UG/L	10.0000 U		.	10.0000 U	.	2.0000
I Iron	UG/L	74.7000 J	.	.	4650.0000	.	1540.0000
I Lead	UG/L	2.0000 u	.	.	2.0000 u	.	2.0000 U
I Magnesium	UG/L	552.0000 U	.	.	33300.0000	.	37300.0000
I Manganese	UG/L	6.0000 U	.	.	44.1000	.	42.2000
I Mercury	UG/L	0.2000 U	.	.	0.2000 U	.	0.2000 U
I Nickel	UG/L	38.0000 U	.	.	17.0000 U	.	17.0000 U
I Potassium	UG/L	3950.0000 U	.	.	22600.0000	.	25000.0000
I Selenium	UG/L	3.0000 U		.	3.0000 U	.	3.0000 U
I Silver	UG/L	10.0000 u	.	.	7.0000 U	.	7.0000 U
I Sodium	UG/L	2730.0000 J	.	.	272000.000	.	307000.000
I Thallium	UG/L	3.0000 U	.	.	40.0000 U	.	40.0000 U
I Vanadium	UG/L	15.0000 U	.	.	17.4000 J		8.9000 J
I Zinc	UG/L	3.0000 U	.	.	12.0000 J	.	4.0000 u
P 4,4'-DDD	UG/L	0.1000 U	.	.		.	
P 4,4'-DDE	UG/L	0.1000 u	.	.		.	
P 4,4'-DDT	UG/L	0.1000 U	.	.		.	
P Aldrin	UG/L	0.0500 U		.		.	

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	33G18E01	33G20001MS	33G20001MSD	33G20D01	33G20D01DL	33G2OD01F
	LAB ID:	534605	546725	546726	547567	547567	547574
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	04/27/93	04/27/93
P Aroclor 1016	UG/L	1.0000 U	.	.	.	.	.
P Aroclor 1221	UG/L	2.0000 U	.	.	.	.	.
P Aroclor 1232	UG/L	1.0000 U			.	.	.
P Aroclor 1242	UG/L	1.0000 U	.	.		.	.
P Aroclor 1248	UG/L	1.0000 U		.	.	.	.
P Aroclor 1254	UG/L	1.0000 U	.	.	.	.	
P Aroclor 1260	UG/L	1.0000 U	.	.	.	.	
P Dieldrin	UG/L	0.1000 U	.	.	.	.	.
P Endosulfan I	UG/L	0.0500 U		.	.	.	.
P Endosulfan II	UG/L	0.1000 U	.	.	.	.	.
P Endosulfan sulfate	UG/L	0.1000 U	.	.	.	.	.
P Endrin	UG/L	0.1000 U	.	.	.	.	.
P Endrin aldehyde	UG/L	0.1000 U	.	.	.	.	
P Endrin ketone	UG/L	0.1000 U	.	.	.	.	.
P Heptachlor	UG/L	0.0500 U	.	.	.	.	.
P Heptachlor epoxide	UG/L	0.0500 U	.	.	.	.	.
P Methoxychlor	UG/L	0.5000 U	.	.	.	.	.
P Toxaphene	UG/L	5.0000 U	.	.	.	.	.
P alpha-BHC	UG/L	0.0500 U	.	.	.	.	.
P alpha-Chlordane	UG/L	0.0500 U	.	.	.	.	.
P beta-BHC	UG/L	0.0500 U	.	.	.	.	.
P delta-BHC	UG/L	0.0500 U	.	.	.	.	.
P gamma-BHC (Lindane)	UG/L	0.0500 U	.	.	.	.	.
P gamma-Chlordane	UG/L	0.0500 U	.	.			.
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S 1,2-Dichlorobenzene	UG/L	10.0000 U	.	.	78.0000 R	1200.0000 J	.
S 1,3-Dichlorobenzene	UG/L	10.0000 U	.	.	48.0000 R	730.0000 J	.
S 1,4-Dichlorobenzene	UG/L	10.0000 U	.	.	44.0000 R	690.0000 J	.
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	.	.	25.0000 U	500.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	33G18E01	33G20001MS	33G20001MSD	33G20D01	33G20D01DL	33G20D01F
	LAB ID:	534605	546725	546726	547567	547567	547574
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	04/27/93	04/27/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2,4-Dichlorophenol	UG/L	10.0000 U		.	1.0000 R	200.0000 U	.
S 2,4-Dimethylphenol	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2,4-Dinitrophenol	UG/L	25.0000 U		.	25.0000 U	500.0000 U	.
S 2,4-Dinitrotoluene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2,6-Dinitrotoluene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2-Chloronaphthalene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2-Chlorophenol	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2-Methylnaphthalene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2-Methylphenol	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 2-Nitroaniline	UG/L	25.0000 U		.	25.0000 U	500.0000 U	.
S 2-Nitrophenol	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 3-Nitroaniline	UG/L	25.0000 U		.	25.0000 U	500.0000 U	.
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U		.	25.0000 U	500.0000 U	.
S 4-Bromophenyl-phenylether	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 4-Chloro-3-methylphenol	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 4-Chloroaniline	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 4-Methylphenol	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S 4-Nitroaniline	UG/L	25.0000 U		.	25.0000 U	500.0000 U	.
S 4-Nitrophenol	UG/L	25.0000 U		.	25.0000 U	500.0000 U	.
S Acenaphthene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S Acenaphthylene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S Anthracene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S Benzo(a)anthracene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S Benzo(a)pyrene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S Beneo(b,k)fluoranthene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.
S Benzo(g,h,i)perylene	UG/L	10.0000 U		.	10.0000 U	200.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G18E01	33G20001MS	33G20001MSD	33G20D01	33G20D01DL	33G20D01F
	LAB ID:	534605	546725	546726	547567	547567	547574
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	04/27/93	04/27/93
NAME	UNITS						
S Butylbenzylphthalate	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Carbazole	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Chrysene	UG/L	10.0000 u	.	.	10.0000 U	200.0000 U	.
S Di-n-butylphthalate	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Di-n-octylphthalate	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Dibenz(a,h)anthracene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Dibenzofuran	UG/L	10.0000 U	.	.	10.0000 U	200.0000 u	.
S Diethylphthalate	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Dimethylphthalate	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Fluoranthene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Fluorene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Hexachlorobenzene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Hexachlorobutadiene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Hexachlorocyclopentadiene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Hexachloroethane	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Isophorone	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S N-Nitrosodiphenylamine	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Naphthalene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Nitrobenzene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Pentachlorophenol	UG/L	25.0000 U	.	.	25.0000 U	500.0000 U	.
S Phenanthrene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Phenol	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S Pyrene	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S bis(2-Ethlyhexy)phthalate	UG/L	2.0000 U	.	.	10.0000 U	200.0000 U	.
S bis(2-chloroethyl)ether	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	.	.	10.0000 U	200.0000 U	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33G18E01	33G20001MS	33G20001MSD	33G20D01	33G20D01DL	33G20D01F
ENSAFE ID:		33G18E01	33G20001MS	33G20001MSD	33G20D01	33G20D01DL	33G20D01F
LAB ID:		534605	546725	546726	547567	547567	547574
DATE RCVD:		02/03/93	04/27/93	04/27/93	04/27/93	04/27/93	04/27/93
V 1,1,1-Trichloroethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V 1,1,2-Trichloroethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V 1,1-Dichloroethane	UG/L	10.0000 U	230.0000	220.0000	190.0000	.	.
V 1,1-Dichloroethene	UG/L	10.0000 U	150.0000	140.0000	33.0000 U	.	.
V 1,2-Dichloroethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	14.0000 J	13.0000 J	12.0000 J	.	.
V 1,2-Dichloropropane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V 2-Butanone	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V 2-Hexanone	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Acetone	UG/L	31.0000	33.0000 U	33.0000 U	33.0000 U	.	.
V Benzene	UG/L	10.0000 U	160.0000	160.0000	33.0000 U	.	.
V Bromodichloromethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Bromoform	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Bromomethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Carbon Disulfide	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Carbon Tetrachloride	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Chlorobenzene	UG/L	10.0000 U	360.0000	360.0000	200.0000	.	.
V Chloroethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Chloroform	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Chloromethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Dibromochloromethane	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Ethylbenzene	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Methylene Chloride	UG/L	4.0000 U	29.0000 U	41.0000 U	8.0000 U	.	.
V Styrene	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U	.	.
V Tetrachloroethene	UG/L	10.0000 U	5.0000 J	5.0000 J	6.0000 J	.	.
V Toluene	UG/L	10.0000 U	160.0000	170.0000	6.0000 J	.	.
V Trichloroethene	UG/L	10.0000 U	160.0000	160.0000	7.0000 J	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G18E01	33G20001MS	33G20001MSD	33G20D01	33G20D01DL	33G20D01F
	LAB ID:	534605	546725	546726	547567	547567	547574
	DATE RCVD:	02/03/93	04/27/93	04/27/93	04/27/93	04/27/93	04/27/93
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	15.0000 J	33.0000 U	33.0000 U		.
V Xylene (total)	UG/L	10.0000 U	5.0000 J	5.0000 J	4.0000 J		.
V cis-1,3-Dichloropropene	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U		
V trans-1,3-Dichloropropene	UG/L	10.0000 U	33.0000 U	33.0000 U	33.0000 U		

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G20D01RE	33G20T01	33G81001MS	33G81001MSD	33GGM71E1	33GGM78E1
	LAB ID:	547567	547583	532920	532921	534220	534223
	DATE RCVD:	04/27/93	04/27/93	01/27/93	01/27/93	01/29/93	01/29/93
NAME	WITS						
I Aluminum	UG/L	.	.	.	.	126.0000 J	119.0000 J
I Antimony	UG/L	.	.	.	.	57.0000 U	57.0000 U
I Arsenic	UG/L	.	.	.	.	.7.0000 J	3.0000 U
I Barium	UG/L	.	.	.	.	6.0000 U	6.0000 U
I Beryllium	UG/L	.	.	.	.	3.0000 U	3.0000 U
I Cadmium	UG/L	.	.	.	.	5.0000 U	5.0000 U
I Calcium	UG/L	.	.	.	.	130.0000 U	130.0000 U
I Chromium	UG/L	.	.	.	.	10.0000 U	10.0000 U
I Cobalt	UG/L	.	.	.	.	12.0000 U	12.0000 U
I Copper	UG/L	.	.	.	.	25.0000 U	25.0000 U
I Cyanide	UG/L	.	.	.	.	10.0000 U	10.0000 U
I Iron	UG/L	.	.	.	.	46.3000 J	54.3000 J
I Lead	UG/L	.	.	.	.	2.0000 U	2.0000 U
I Magnesium	UG/L	.	.	.	.	552.0000 U	552.0000 U
I Manganese	UG/L	.	.	.	.	6.0000 U	6.0000 U
I Mercury	UG/L	.	.	.	.	0.2000 U	0.2000 U
I Nickel	UG/L	.	.	.	.	38.0000 U	38.0000 U
I Potassium	UG/L	.	.	.	.	3950.0000 U	3950.0000 U
I Selenium	UG/L	.	.	.	.	3.0000 U	3.0000 U
I Silver	UG/L	.	.	.	.	10.0000 U	10.0000 U
I Sodium	UG/L	.	.	.	.	352.0000 J	324.0000 J
I Thallium	UG/L	.	.	.	.	3.0000 U	3.0000 U
I Vanadium	UG/L	.	.	.	.	15.0000 U	15.0000 U
I Zinc	UG/L	.	.	.	.	3.0000 U	3.0000 U
P 4,4'-DDD	UG/L	.	.	.	.	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	.	.	.	.	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	.	.		.	0.1000 U	0.1000 U
P Aldrin	UG/L	.	.	.	.	0.0500 U	0.0500 U

NAY PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G20D01RE	33G20T01	33G81001MS	33(381001MSD	33GGM71E1	33GGM78El
	LAB ID:	547567	547583	532920	532921	534234	534237
	DATE RCVD:	04/27/93	04/27/93	01/27/93	01/27/93	01/29/93	01/29/93
NAME	UNITS						
P Aroclor 1016	UG/L	.	.	.		1.0000 U	1.0000 U
P Aroclor 1221	UG/L	.	.	.		2.0000 U	2.0000 U
P Aroclor 1232	UG/L	.	.	.		1.0000 U	1.0000 U
P Aroclor 1242	UG/L	.	.	.		1.0000 U	1.0000 U
P Aroclor 1248	UG/L	.	.	.		1.0000 U	1.0000 U
P Aroclor 1254	UG/L	.	.	.		1.0000 U	1.0000 U
P Aroclor 1260	UG/L	.	.	.		1.0000 U	1.0000 U
P Dieldrin	UG/L	.	.	.		0.1000 U	0.1000 U
P Endosulfan I	UG/L	.	.	.		0.0500 U	0.0500 U
P Endosulfan II	UG/L	.	.	.		0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	.	.	.		0.1000 U	0.1000 U
P Endrin	UG/L	.	.	.		0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	.	.	.		0.1000 U	0.1000 U
P Endrin ketone	UG/L	.	.	.		0.1000 U	0.1000 U
P Heptachlor	UG/L	.	.	.		0.0017 J	0.0025 J
P Heptachlor epoxide	UG/L	.	.	.		0.0500 U	0.0500 U
P Methoxychlor	UG/L	.	.	.		0.5000 U	0.5000 U
P Toxaphene	UG/L	.	.	.		5.0000 U	5.0000 U
P alpha-BHC	UG/L	.	.	.		0.0500 U	0.0500 U
P alpha-Chlordane	UG/L	.	.	.		0.0500 U	0.0500 U
P beta-BHC	UG/L	.	.	.		0.0500 U	0.0500 U
P delta-BHC	UG/L	.	.	.		0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L	.	.	.		0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	.	.	.		0.0500 U	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	1.0000 J		41.0000	38.0000	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	580.0000 J		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	420.0000 J		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	360.0000 J		33.0000	30.0000	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	1.0000 J		25.0000 U	25.0000 U	25.0000 U	25.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G20D01RE	33G20T01	33G81001MS	33G81001MSD	33GGM71El	33GGM78El
	LAB ID:	547567	547583	532920	532921	534234	534237
	DATE RCVD:	04/27/93	04/27/93	01/27/93	01/27/93	01/29/93	01/29/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	11.0000 J	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	10.0000 U	.	37.0000	36.0000	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	2.0000 J	.	55.0000	52.0000	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	25.0000 U	25.0000 U	25,0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	.	61.0000	58.0000	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	.	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000 U	.	60.0000	55.0000	25.0000 U	25.0000 U
S Acenaphthene	UG/L	10.0000 U	.	39.0000	36.0000	10.0000 U	10.0000 U
S Acenaphthylene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME		ENSAFE ID:	33G20D01RE	33G20T01	33G81001MS	33G81001MSD	33GGM71E1	33GGM78El
		LAB ID:	547567	547583	532920	532921	534234	534237
		DATE RCVD:	04/27/93	04/27/93	01/27/93	01/27/93	01/29/93	01/29/93
	UNITS							
S Butylbenzylphthalate	UG/L		10.0000 U	.	2.0000 U	2.0000 U	10.0000 U	10.0000 U
S Carbazole	UG/L		10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L		10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L		10.0000 U		2.0000 U	2.0000 U	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L		10.0000 U		1.0000 U	2.0000 U	10.0000 U	10.0000 U
S Dimethylphthalate	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L		10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L		10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Ieophorone	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L		10.0000 U		37.0000	35.0000	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L		10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L		25.0000 U	.	84.0000	80.0000	25.0000 U	25.0000 U
S Phenanthrene	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Phenol	UG/L		10.0000 U		50.0000	49.0000	10.0000 U	10.0000 U
S Pyrene	UG/L		10.0000 U		29.0000	32.0000	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L		2.0000 U	.	8.0000 U	4.0000 U	1.0000 U	10.0000 U
S bis(2-chloroethyl)ether	UG/L		10.0000 U		10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L		10.0000 U	.	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33G20D01RE	33G20T01	33G81001MS	33G81001MSD	33GGM71E1	33GGM78E1
	LAB ID:	547567	547583	532920	532921	534234	534237
	DATE RCVD:	04/27/93	04/27/93	01/27/93	01/27/93	01/29/93	01/29 /93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 2-Butanone	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Acetone	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Benzene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Bromoform	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Bromomethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Chloroethane .	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Chloroform	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Chloromethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	.	14.0000 U		.	10.0000 U	3.0000 U
V Styrene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Toluene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
V Trichloroethene	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	33G20D01RE	33G20T01	33G81001MS	33G81001MSD	33GGM71E1	33GGM78E1
	LAB ID:	547567	547583	532920	532921	534234	534237
	DATE RCVD:	04/27/93	04/27/93	01/27/93	01/27/93	01/29/93	01/29/93
NAME	UNITS						
V Vinyl Chloride	UG/L		10.0000 U		.	10.0000 U	10.0000 U
V Xylene (total)	UG/L		10.0000 U		.	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L		10.0000 U		.	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L		10.0000 U		.	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GUN001	33M01E01	33S00E01	33S00E02	33S00E03	33S00E04
	LAB ID:	531913	533162	529307	529533	529855	530192
	DATE RCVD:	01/14/93	01/21/93	12/16/92	12/17/92	12/18/92	12/19/92
NAME	UNITS						
I Aluminum	UG/L	312.0000	264.0000	193.0000 J	207.0000	208.0000	205.0000
I Antimony	UG/L	57.0000 U	57.0000 U	57.0000 U	57.0000 U	57.0000 U	57.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	4.0000 U	4.0000 U	4.0000 U	4.0000 U
I Barium	UG/L	28.2000 J	6.0000 U	3.0000 U	3.0000 U	3.0000 U	3.0000 U
I Beryllium	UG/L	3.0000 U	3.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U	5.0000 U
I Calcium	UG/L	2280.0000 J	130.0000 U	924.0000 U	924.0000 U	924.0000 U	924.0000 U
I Chromium	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
I Cobalt	UG/L	12.0000 U	12.0000 U	11.0000 U	11.0000 U	11.0000 U	11.0000 U
I Copper	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
I Cyanide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
I Iron	UG/L	540.0000	231.0000	144.0000	28.0000 U	28.0000 U	183.0000
I Lead	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U	3.0000 J
I Magnesium	UG/L	1540.0000 J	552.0000 U	425.0000 U	425.0000 U	425.0000 U	425.0000 U
I Manganese	UG/L	18.1000	6.0000 U	7.0000 U	7.0000 U	7.0000 U	7.0000 U
I Mercury	UG/L	0.2000 U	0.2000 U	0.2000 U	0.2000 U	0.2000 U	0.2000 U
I Nickel	UG/L	38.0000 U	38.0000 U	40.0000 U	40.0000 U	40.0000 U	40.0000 U
I Potassium	UG/L	3950.0000 U	3950.0000 U	4860.0000 U	4860.0000 U	4860.0000 U	4860.0000 U
I Selenium	UG/L	3.0000 U	3.0000 U	4.0000 U	4.0000 U	4.0000 U	4.0000 U
I Silver	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
I Sodium	UG/L	33300.0000	468.0000 J	409.0000 J	414.0000 J	357.0000 J	378.0000 J
I Thallium	UG/L	3.0000 U	3.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
I Vanadium	UG/L	15.0000 U	15.0000 U	19.0000 U	19.0000 U	19.0000 U	19.0000 U
I Zinc	UG/L	111.0000	40.9000	5.0000 U	5.0000 U	5.0000 U	5.5000 J
P 4,4'-DDD	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	0.1000 U	0.0013 J	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Aldrin	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GUN001	33M01E01	33S00E01	33S00E02	33S00E03	33S00E04
	LAB ID:	531913	533162	529307	529533	529855	530192
	DATE RCVD:	01/14/93	01/21/93	12/16/92	12/17/92	12/18/92	12/19/92
NAME	UNITS						
P Aroclor 1016	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1,0000 U	1.0000 U
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	1,0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U	1.0000 U
P Dieldrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan I	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Endosulfan II	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Endrin ketone	UG/L	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U	0.1000 U
P Heptachlor	UG/L	0.0018 J	0.0018 J	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P Methoxychlor	UG/L	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U	0.5000 U
P Toxaphene	UG/L	5.0000 U	5.0000 U	5.0000 U	5.0000 U	s.0000 U	5.0000 U
P alpha-BHC	UG/L	0.0500 U	0.0014 J	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P beta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P delta-BHC	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
ENSAFE ID:		33GUN001	33M01E01	33S00E01	33S00E02	33S00E03	33S00E04
LAB ID:		531913	533162	529307	529533	529655	530192
DATE RCVD:		01/14/93	01/21/93	12/16/92	12/17/92	12/18/92	12/19/92
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33GUN001 / 531913 / 01/14/93	33M01E01 / 533162 / 01/21/93	33S00E01 / 529307 / 12/16/92	33S00E02 / 529533 / 12/17/92	33S00E03 / 529855 / 12/18/92	33S00E04 / 530192 / 12/19/92
S Butylbenzylphthalate	UG/L	10.0000 U	4.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Carbazole	UG/L	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	10.0000 u	10.0000 u	10.0000 U	10.0000 U	10.0000 u	10.0000 U
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	1.0000 u	1.0000 u
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 u	10.0000 u
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10,0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 u
S Hexachlorobenzene	UG/L	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Ieophorone	UG/L	10.0000 u	10.0000 u	10.0000 u	10.0000 u	10.0000 u	10.0000 u
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 u	10.0000 u	10.0000 U	10.0000 u	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 u	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Phenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 u	10.0000 U
S Pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	3.0000 U	2.0000 U	5.0000 u	9.0000 U	5.0000 U	4.0000 U
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 u	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	33GUN001	33M01E01	33S00E01	33S00E02	33S00E03	33S00E04
ENSAFE ID:		33GUN001	33M01E01	33S00E01	33S00E02	33S00E03	33S00E04
LAB ID:		531913	533162	529307	529533	529855	530192
DATE RCVD:		01/14/93	01/21/93	12/16/92	12/17/92	12/18/92	12/19/92
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	32.0000	10.0000 U	14.0000 U	16.0000 U	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	1.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	2.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	1.0000 U	7.0000 U	2.0000 U	1.0000 U	2.0000 U	4.0000 J
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	2.0000 J	1.0000 J	10.0000 U	2.0000 J
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33GUN001	33M01E01	33S00E01	33S00E02	33S00E03	33S00E04
	LAB ID:	531913	533162	529307	529533	529855	530192
	DATE RCVD:	01/14/93	01/21/93	12/16/92	12/17/92	12/18/92	12/19/92
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID:	33S00E30	33S00E31	33S00E32	33S00E32MS	33S00E32MSD	33TANK01
	LAB ID:	546724	546745	547109	546728	546729	531930
	DATE RCVD:	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93	01/14/93
	UNITS						
I Aluminum	UG/L	94.6000 J	96.3000 J	102.0000 J	.	.	506.0000
I Antimony	UG/L	36.0000 U	36.0000 U	36.0000 U	.	.	57.0000 U
I Arsenic	UG/L	3.0000 U	3.0000 U	3.0000 U	.	.	3.0000 U
I Barium	UG/L	3.2000 J	2.3000 J	2.3000 J	.	.	6.0000 U
I Beryllium	UG/L	1.0000 U	1.0000 U	1.0000 U	.	.	3.0000 U
I Cadmium	UG/L	5.0000 U	5.0000 U	5.0000 U	.	.	5.0000 U
I Calcium	UG/L	196.0000 J	170.0000 J	113.0000 J	.	.	2530.0000 J
I Chromium	UG/L	8.0000 U	8.0000 U	8.0000 U	.	.	10.0000 U
I Cobalt	UG/L	8.0000 U	8.0000 U	8.0000 U	.	.	12.0000 U
I Copper	UG/L	4.2000 J	4.0000 U	4.0000 U	.	.	25.0000 U
I Cyanide	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
I Iron	UG/L	55.6000 J	105.0000	61.8000 J	.	.	255.0000
I Lead	UG/L	2.0000 U	2.0000 U	2.0000 U	.	.	2.0000 U
I Magnesium	UG/L	76.4000 J	73.9000 J	80.4000 J	.	.	552.0000 U
I Manganese	UG/L	2.1000 J	2.9000 J	1.3000 J	.	.	6.0000 U
I Mercury	UG/L	0.2000 U	0.2000 U	0.2000 U	.	.	0.2000 U
I Nickel	UG/L	17.0000 U	17.0000 U	17.0000 U	.	.	38.0000 U
I Potassium	UG/L	1350.0000 U	1350.0000 U	1350.0000 U	.	.	3950.0000 U
I Selenium	UG/L	3.0000 U	3.0000 U	3.0000 U	.	.	3.0000 U
I Silver	UG/L	7.0000 U	7.0000 U	7.0000 U	.	.	10.0000 U
I Sodium	UG/L	271.0000 J	319.0000 J	289.0000 J	.	.	483.0000 J
I Thallium	UG/L	4.0000 U	4.0000 U	4.0000 U	.	.	3.0000 U
I Vanadium	UG/L	6.0000 U	6.0000 U	6.0000 U	.	.	15.0000 U
I Zinc	UG/L	5.6000 J	5.9000 J	5.1000 J	.	.	4.5000 J
P 4,4'-DDD	UG/L				.	.	0.1000 U
P 4,4'-DDE	UG/L				.	.	0.1000 U
P 4,4'-DDT	UG/L			.	.	.	0.1000 U
P Aldrin	UG/L			.	.	.	0.0500 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digite.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33S00E30 546724 04/21/93	33S00E31 546745 04/21/93	33S00E32 547109 04/22/93	33S00E32MS 546728 04/22/93	33S00E32MSD 546729 04/22/93	33TANK01 531930 01/14/93
P Aroclor 1016	UG/L	.	.	.		.	1.0000 U
P Aroclor 1221	UG/L	.	.	.		.	2.0000 U
P Aroclor 1232	UG/L	.	.	.		.	1.0000 U
P Aroclor 1242	UG/L	.	.	.		.	1.0000 U
P Aroclor 1248	UG/L	.	.	.		.	1.0000 U
P Aroclor 1254	UG/L	.	.	.		.	1.0000 U
P Aroclor 1260	UG/L	.	.	.		.	1.0000 U
P Dieldrin	UG/L	.	.	.		.	0.1000 U
P Endosulfan I	UG/L	.	.	.		.	0.0500 U
P Endosulfan II	UG/L	.	.	.		.	0.1000 U
P Endosulfan sulfate	UG/L	.	.	.		.	0.1000 U
P Endrin	UG/L	.	.	.		.	0.1000 U
P Endrin aldehyde	UG/L	.	.	.		.	0.1000 U
P Endrin ketone	UG/L						0.1000 U
P Heptachlor	UG/L						0.0500 U
P Heptachlor epoxide	UG/L						0.0500 U
P Methoxychlor	UG/L						0.5000 U
P Toxaphene	UG/L						5.0000 U
P alpha-BHC	UG/L						0.0500 U
P alpha-Chlordane	UG/L						0.0500 U
P beta-BHC	UG/L						0.0500 U
P delta-BHC	UG/L						0.0500 U
P gamma-BHC (Lindane)	UG/L						0.0500 U
P gamma-Chlordane	UG/L						0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	24.0000	29.0000	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	21.0000	25.0000	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S00E30	33S00E31	33S00E32	33S00E32MS	33S00E32MSD	33TANK01
	LAB ID:	546724	546745	547109	546728	546729	531930
	DATE RCVD:	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93	01/14/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	29.0000	34.0000	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	39.0000	42.0000	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	45.0000	SO.0000	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	48.0000	55.0000	25.0000 U
S Acenaphthene	UG/L	10.0000 U	10.0000 U	10.0000 U	30.0000	34.0000	10.0000 U
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S00E30	33S00E31	33S00E32	33S00E32MS	33S00E32MSD	33TANK01
	LAB ID:	546724	546745	547109	546728	546729	531930
	DATE RCVD:	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93	01/14/93
NAME	UNITS						
S Butylbenzylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Carbazole	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Chrysene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	1.0000 U	1.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenz (a,h) anthracene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	10.0000 U	10.0000 u	10.0000 U	1.0000 U	1.0000 u	10.0000 U
S Dimethylphthalate	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Fluorene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	10,0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Isophorone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	10.0000 U	10.0000 U	10.0000 U	29.0000	33.0000	10.0000 U
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 u	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	25.0000 U	25.0000 U	25.0000 U	49.0000	57.0000	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S Phenol	UG/L	10.0000 U	10.0000 U	10.0000 U	39.0000	43.0000	10.0000 U
S Pyrene	UG/L	10.0000 U	10.0000 U	10.0000 U	28.0000	32.0000	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	2.0000 U	2.0000 U	3.0000 U	3.0000 U	2.0000 U	1.0000 U
S bis(2-chloroethyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33S00E30	33S00E31	33S00E32	33S00E32MS	33S00E32MSD	33TANK01
	LAB ID:	546724	546745	547109	546728	546729	531930
	DATE RCVD:	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93	01/14/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,2-Dichloroethane	UG/L	10,0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	37.0000 U	.	.	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Methylene Chloride	UG/L	2.0000 U	3.0000 U	3.0000 U	.	.	2.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits

	ENSAFE ID:	33S00E00	33S00E31	33S00E32	33S00E32MS	33S00E32MSD	33TANK01
	LAB ID:	546724	546745	547109	546728	546729	531930
	DATE RCVD:	04/21/93	04/21/93	04/22/93	04/22/93	04/22/93	01/14/93
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U

V=Volatile, S=Semivolatile P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W00T10	33W00T11	33W00T11MS	33W00T11MSD	33W00T12	33W00T12MS
	LAB ID:	530356	530859	530860	530861	531078	531079
	DATE RCVD:	12/22/92	12/31/92	12/31/92	12/31/92	01/06/93	01/06/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	55.0000	53.0000	10.0000 U	44.0000
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	9.0000 U	7.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	20.0000 U	43.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	53.0000	51.0000	10.0000 U	44.0000
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	51.0000	49.0000	10.0000 U	44.0000
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	1.0000 U	1.0000 U	2.0000 U	2.0000 U	5.0000 U	5.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	52.0000	50.0000	*3.0000* J	47.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W00T10	33W00T11	33W00T11MS	33W00T11MSD	33W00T12	33W00T12MS
	LAB ID:	530356	530859	530860	530861	531078	531079
	DATE RCVD:	12/22/92	12/31/92	12/31/92	12/31/92	01/06/93	01/06/93
NAME	UNITS						
V Trichloroethene	UG/L	10.0000 U	10.0000 U	52.0000	49.0000	10.0000 U	44.0000
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 ANI) SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W00T12MSD	33W00T14	33W00T14	33W00T15	33W00T16	33W00T16
	LAB ID:	531080	531342	531909	532079	532078	532918
	DATE RCVD:	01/06/93	01/08/93	01/14/93	01/15/93	01/15/93	01/20/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	50.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	8.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	63.0000 U	12.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L	48.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	48.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	5.0000 U	2.0000 U	2.0000 U	10.0000 U	2.0000 U	8.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	51.0000	1.0000 J	10.0000 U	10.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W00T12MSD	33W00T14	33W00T14	33W00T15	33W00T16	33W00T16
	LAB ID:	531000	531342	531909	532079	532078	532918
	DATE RCVD:	01/06/93	01/08/93	01/14/93	01/15/93	01/15/93	01/20/93
NAME	UNITS						
V Trichloroethene	UG/L	40.0000	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	2.0000 J
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 u'	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33W00T17 / 532914 / 01/20/93	33W00T18 / 532831 / 01/21/93	33W00T18 / 533343 / 01/22/93	33W01002MS / 532923 / 01/21/93	33W01002MSD / 532924 / 01/21/93	33W02002MS / 532915 / 01/21/93
P 4,4'-DDD	UG/L	.	.	.	0.0068 J	0.0310 J	.
P 4,4'-DDE	UG/L			.	0.0085 J	0.0079 J	
P 4,4'-DDT	UG/L	.		.	1.0000	0.9600	
P Aldrin	UG/L		.	.	0.4300	0.4200	.
P Aroclor 1016	UG/L	.	.	.	1.0000 U	1.0000 U	.
P Aroclor 1221	UG/L	.		.	2.0000 U	2.0000 U	.
P Aroclor 1232	UG/L	.		.	1.0000 U	1.0000 U	.
P Aroclor 1242	UG/L	.	.	.	1.0000 U	1.0000 U	.
P Aroclor 1248	UG/L		.	.	1.0000 U	1.0000 U	.
P Aroclor 1254	UG/L	.		.	1.0000 U	1.0000 U	
P Aroclor 1260	UG/L	.	.		1.0000 U	1.0000 U	.
P Dieldrin	UG/L	.	.	.	0.9200	0.8900	.
P Endosulfan I	UG/L	.	.	.	0.0500 U	0.0500 U	.
P Endosulfan II	UG/L	.	.	.	0.1000 U	0.1000 U	.
P Endosulfan sulfate	UG/L		.	.	0.1000 U	0.1000 U	.
P Endrin	UG/L	.	.	.	1.1000	1.0000	.
P Endrin aldehyde	UG/L	.	.	.	0.1000 U	0.1000 U	.
P Endrin ketone	UG/L	.	.	.	0.0250 J	0.0200 J	.
P Heptachlor	UG/L	.	.	.	0.4900	0.4700	.
P Heptachlor epoxide	UG/L	.	.	.	0.0500 U	0.0500 U	.
P Methoxychlor	UG/L	.	.	.	0.5000 U	0.5000 U	.
P Toxaphene	UG/L	.	.	.	5.0000 U	5.0000 U	.
P alpha-BHC	UG/L	.	.	.	0.0500 U	0.0062 J	
P alpha-Chlordane	UG/L	.	.	.	0.0020 J	0.0020 J	.
P beta-BHC	UG/L	.	.	.	0.0500 U	0.0500 U	
P delta-BHC	UG/L	.	.	.	0.0500 U	0.0500 U	.
P gamma-BHC (Lindane)	UG/L	.	.	.	0.4800	0.4700	.
P gamma-Chlordane	UG/L	.	.	.	0.0062 J	0.0077 J	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W00T17	33W00T18	33W00T18	33W01002MS	33W01002MSD	33W02002MS
	LAB ID:	532914	532831	533343	532923	532924	532915
	DATE RCVD:	01/20/93	01/21/93	01/22/93	01/21/93	01/21/93	01/21/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10,0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,1-Dichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	44.0000
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Acetone	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Benzene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	46.0000
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10,0000 U
V Bromofo m	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Chlorobenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	46.0000
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Methylene Chloride	UG/L	10.0000 U	1.0000 U	5.0000 U	.	.	12.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	10.0000 U
V Toluene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	48.0000
V Trichloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	.	.	45.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W00T17	33W00T18	33W00T18	33W01002MS	33W01002MSD	33W02002MS
	LAB ID:	532914	532831	533343	532923	532924	532915
	DATE RCVD:	01/20/93	01/21/93	01/22/93	01/21/93	01/21/93	01/21/93
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U		.	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U		.	1.0000 J
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U		.	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U		.	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W02002MSD	33W03D02	33W03D02	66WL	NASPAF01	NASPAF02
	LAB ID:	532916	533158	533175	524187	520458	522657
	DATE RCVD:	01/21/93	01/21/93	01/21/93	11/20/92	11/06/92	11/14/92
NAME	UNITS						
I Aluminum	UG/L	.	.	745.0000	.	194.0000 J	128.0000 J
I Antimony	UG/L	.		57.0000 U	.	60.0000 U	60.0000 U
I Arsenic	UG/L	.	.	3.0000 U	.	4.0000 U	4.0000 U
I Barium	UG/L	.	.	12.6000 J	.	2.0000 U	4.7000 J
I Beryllium	UG/L	.	.	3.0000 U	.	1.0000 U	1.0000 U
I Cadmium	UG/L	.	.	5.0000 U	.	5.0000 U	5.0000 U
I Calcium	UG/L	.	.	13300.0000	.	1030.0000 J	424.0000 J
I Chromium	UG/L	.	.	10.0000 U	.	6.0000 U	6.0000 U
I Cobalt	UG/L		.	12.0000 U	.	9.0000 U	9.0000 U
I Copper	UG/L		.	25.0000 U	.	5.0000 U	5.0000 U
I Cyanide	UG/L	.	.	10.0000 U	.	10.0000 U	10.0000 U
I Iron	UG/L	.	.	956.0000	.	57.0000 U	34.0000 U
I Lead	UG/L	.	.	2.0000 U	.	2.0000 U	2.0000 U
I Magnesium	UG/L	.	.	5030.0000	.	665.0000 J	244.0000 J
I Manganese	UG/L	.	.	20.9000	.	3.1000 J	2.0000 U
I Mercury	UG/L	.	.	0.2000 U	.	0.2000 U	0.2000 U
I Nickel	UG/L	.	.	38.0000 U	.	11.0000 U	11.0000 U
I Potaeeium	UG/L	.	.	3950.0000 U	.	1360.0000 U	1360.0000 U
I Selenium	UG/L		.	3.0000 U	.	4.0000 U	4.0000 U
I Silver	UG/L		.	10.0000 U	.	6.0000 U	6.0000 U
I Sodium	UG/L	.	.	21000.0000	.	4880.0000 J	618.0000 J
I Thallium	UG/L		.	3.0000 U	.	2.0000 U	2.0000 U
I Vanadium	UG/L	.	.	15.0000 U	.	5.0000 U	5.0000 U
I Zinc	UG/L	.		38.8000	.	10.1000 J	6.6000 J
P 4,4'-DDD	UG/L	.	0.1300 J		.	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	.	0.0570 J		.	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	.	0.0450 J		.	0.1000 U	0.1000 U
P Aldrin	UG/L	.	0.0500 U		.	0.0500 U	0.0500 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS						
	ENSAFE ID:	33W02002MSD	33W03D02	33W03D02	66WL	NASPAF01	NASPAF02
	LAB ID:	532916	533158	533175	524187	520462	522657
	DATE RCVD:	01/21/93	01/21/93	01/21/93	11/20/92	11/06/92	11/14/92
P Aroclor 1016	UG/L	.	1.0000 U	.	.	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	.	2.0000 U	.	.	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	.	1.0000 U	.	.	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	.	1.0000 U	.	.	1.0000 U	1.0000 U
P Aroclor 1240	UG/L	.	1.0000 U	.	.	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	.	1.0000 U	.		1.0000 U	1.0000 U
P Aroclor 1260	UG/L	.	1.0000 U	.	.	1.0000 U	1.0000 U
P Dieldrin	UG/L	.	0.1000 U	.		0.1000 U	0.1000 U
P Endosulfan I	UG/L	.	0.0500 U	.	.	0.0500 U	0.0500 U
P Endosulfan II	UG/L	.	0.1000 U	.	.	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	.	0.1000 U	.	.	0.1000 U	0.1000 U
P Endrin	UG/L	.	0.1000 U	.	.	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	.	0.1000 U	.	.	0.1000 U	0.1000 U
P Endrin ketone	UG/L	.	0.1000 U	.	.	0.1000 U	0.1000 U
P Heptachlor	UG/L	.	0.0028 J	.	.	0.0500 U	*0.0500* U
P Heptachlor epoxide	UG/L	.	0.0500 U	.	.	0.0500 U	0.0500 U
P Methoxychlor	UG/L	.	0.5000 U	.	.	0.5000 U	0.5000 U
P Toxaphene	UG/L	.	5.0000 U	.	.	5.0000 U	5.0000 U
P alpha-BHC	UG/L	.	0.0030 J	.	.	*0.0500* U	0.0500 U
P alpha-Chlordane	UG/L	.	0.0500 U	.	.	*0.0500* U	0.0500 U
P beta-BHC	UG/L	.	0.0500 U	.	.	*0.0500* U	0.0500 U
P delta-BHC	UG/L	.	0.0500 U	.	.	*0.0500* U	0.0500 U
P gamma-BHC (Lindane)	UG/L	.	0.0500 U	.	.	0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	.	0.0016 J	.	.	0.0500 U	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	.	10,0000 U	.	.	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W02002MSD	33W03D02	33W03D02	66WL	NASPAF01	NASPAF02
	LAB ID:	532916	533158	533175	524187	520462	522657
	DATE RCVD:	01/21/93	01/21/93	01/21/93	11/20/92	11/06/92	11/14/92
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Chloroaniline	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U
S Acenaphthene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Acenaphthylene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Anthracene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits,

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	33W02002MSD 532916 01/21/93	33W03D02 533158 01/21/93	33W03D02 533175 01/21/93	66WL 524187 11/20/92	NASPAF01 520462 11/06/92	NASPAF02 522657 11/14/92
S Butylbenzylphthalate	UG/L	.	4.0000 U	.	.	10.0000 U	10.0000 U
S Carbazole	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Chrysene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Di-n-butylphthalate	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Di-n-octylphthalate	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Dibenz(a,h)anthracene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Dibenzofuran	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Diethylphthalate	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Dimethylphthalate	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Fluoranthene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Fluorene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Hexachloroethane	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Ieophorone	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S N-Nitrosodiphenylamine	UG/L	.	10.0000 U		.	10.0000 U	10.0000 U
S Naphthalene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Nitrobenzene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Pentachlorophenol	UG/L	.	25.0000 U	.	.	25.0000 U	25.0000 U
S Phenanthrene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Phenol	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S Pyrene	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S bis(2-Chloroethoxy)methane	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S bis(2-Ethlyhexy)phthalate	UG/L	.	1.0000 U	.	.	5.0000 U	94.0000
S bis(2-chloroethyl)ether	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U
S bis(2-chloroisopropyl)ether	UG/L	.	10.0000 U	.	.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	WITS	ENSAFE ID:	33W02002MSD	33W03D02	33W03D02	66WL	NASPAF01	NASPAF02
		LAB ID:	532916	533158	533175	524187	520458	522657
		DATE RCVD:	01/21/93	01/21/93	01/21/93	11/20/92	11/06/92	11/14/92
V 1,1,1-Trichloroethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L		44.0000	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L		10.0000 U	10.0000 U		50.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 2-Eutanone	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L		49.0000	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L		46.0000	10.0000 U	.	220.0000	10.0000 U	10.0000 U
V Chloroethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L		1.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L		12.0000 U	12.0000 U	.	42.0000 U	1.0000 U	10.0000 U
V Styrene	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L		10.0000 U	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L		49.0000	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U
V Trichloroethene	UG/L		48.0000	10.0000 U	.	50.0000 U	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	33W02002MSD	33W03D02	33103D02	66WL	NASPAF01	NASPAP02
	LAB ID:	532916	533150	533175	524107	520450	522657
	DATE RCVD:	01/21/93	01/21/93	01/21/93	11/20/92	11/06/92	11/14/92
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U		50.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	1.0000 J		50.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U		50.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U		50.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits

NAME	ENSAFE ID:	NASPAF02DL	NASPAF03	NASPAF04	NASPAF06	NASPAF07	NASPAF07MS
	LAB ID:	522657	524171	525438	528897	529747	529749
	DATE RCVD:	11/14/92	11/20/92	11/28/92	12/12/92	12/18/92	12/18/92
	UNITS						
I Aluminum	UG/L		225.0000 J	139.0000 J	245.0000	206.0000	
I Antimony	UG/L		60.0000 U	60.0000 U	60.0000 U	60.0000 U	
I Arsenic	UG/L		4.0000 U	4.0000 U	4.0000 U	4.0000 U	
I Barium	UG/L		2.2000 J	4.3000 J	3.7000 J	8.9000 J	
I Beryllium	UG/L		1.0000 U	1.0000 U	1.0000 U	1.0000 U	
I Cadmium	UG/L		5.0000 U	5.0000 U	5.0000 U	5.0000 U	
I Calcium	UG/L		944.0000 J	389.0000 J	96.9000 J	253.0000 J	
I Chromium	UG/L		6.0000 U	6.0000 U	6.0000 U	6.0000 U	
I Cobalt	UG/L		9.0000 U	9.0000 U	9.0000 U	9.0000 U	
I Copper	UG/L		7.4000 J	5.0000 U	5.0000 U	5.0000 U	
I Cyanide	UG/L		10.0000 U	10.0000 U	10.0000 U	10.0000 U	
I Iron	UG/L		600.0000	86.6000 J	740.0000	34.0000 U	
I Lead	UG/L		2.0000 U	2.0000 U	2.0000 U	2.0000 U	
I Magnesium	UG/L		592.0000 J	240.0000 J	72.7000 J	219.0000 J	
I Manganese	UG/L		2.7000 J	3.7000 J	3.7000 J	2.0000 U	
I Mercury	UG/L		0.2000 U	0.2000 U	0.2000 U	0.2000 U	
I Nickel	UG/L		11.0000 U	11.0000 U	11.0000 U	11.0000 U	
I Potassium	UG/L		1360.0000 U	1360.0000 U	1360.0000 U	1360.0000 U	
I Selenium	UG/L		4.0000 U	4.0000 U	4.0000 U	4.0000 U	
I Silver	UG/L		6.0000 U	6.0000 U	6.0000 U	6.0000 U	
I Sodium	UG/L		1110.0000 J	717.0000 J	486.0000 J	544.0000 J	
I Thallium	UG/L		2.0000 U	2.0000 U	2.0000 U	2.0000 U	
I Vanadium	UG/L		5.0000 U	5.0000 U	5.0000 U	5.0000 U	
I Zinc	UG/L		22.4000	4.4000 J	7.3000 J	3.0000 U	
P 4,4'-DDD	UG/L					0.1000 U	0.0250 J
P 4,4'-DDE	UG/L					0.1000 U	0.1000 U
P 4,4'-DDT	UG/L					0.1000 U	0.7300
P Aldrin	UG/L					0.0500 U	0.3100

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key preeented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF02DL	NASPAF03	NASPAF04	NASPAF06	NASPAF07	NASPAF07MS
	LAB ID:	522657	524171	525438	528897	529747	529755
	DATE RCVD:	11/14/92	11/20/92	11/28/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
P Aroclor 1016	UG/L		.	.	.	1.0000 U	1.0000 U
P Aroclor 1221	UG/L		.	.	.	2.0000 U	2.0000 U
P Aroclor 1232	UG/L		.	.	.	1.0000 U	1.0000 U
P Aroclor 1242	UG/L		.	.	.	1.0000 U	1.0000 U
P Aroclor 1248	UG/L		.	.	.	1.0000 U	1.0000 U
P Aroclor 1254	UG/L		.	.	.	1.0000 U	1.0000 U
P Aroclor 1260	UG/L		.	.	.	1.0000 U	1.0000 U
P Dieldrin	UG/L		.	.	.	0.1000 U	0.7300
P Endosulfan I	UG/L		.	.	.	0.0500 U	0.0500 U
P Endosulfan II	UG/L		.	.	.	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L		.	.	.	0.1000 U	0.1000 U
P Endrin	UG/L		.	.	.	0.1000 U	0.8300
P Endrin aldehyde	UG/L		.	.	.	0.1000 U	0.1000 U
P Endrin ketone	UG/L		.	.	.	0.1000 U	0.1000 U
P Heptachlor	UG/L		.	.	.	0.0500 U	0.3400
P Heptachlor epoxide	UG/L		.	.	.	0.0500 U	0.0500 U
P Methoxychlor	UG/L		.	.	.	0.5000 U	0.5000 U
P Toxaphene	UG/L		.	.	.	5.0000 U	5.0000 U
P alpha-BHC	UG/L		.	.	.	0.0500 U	0.0500 U
P alpha-Chlordane	UG/L		.	.	.	0.0500 U	0.0500 U
P beta-BHC	UG/L		.	.	.	0.0500 U	0.0500 U
P delta-BHC	UG/L		.	.	.	0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L		.	.	.	0.0500 U	0.3700
P gamma-Chlordane	UG/L		.	.	.	0.0500 U	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	20.0000 U	.	.	.	10.0000 U	35.0000
S 1,2-Dichlorobenzene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	20.0000 U	.	.	.	10.0000 U	32.0000
S 2,4,5-Trichlorophenol	UG/L	50.0000 U	.	.	.	25.0000 U	25.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	NASPAF02DL 522657 11/14/92	NASPAF03 524171 11/20/92	NASPAF04 525438 11/28/92	NASPAF06 528897 12/12/92	NASPAF07 529747 12/18/92	NASPAF07MS 529752 12/18/92
S 2,4,6-Trichlorophenol	UG/L	20.0000 U	.		.	10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	20.0000 U	.		.	10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	50.0000 U			.	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	20.0000 U	.		.	10.0000 U	42.0000
S 2,6-Dinitrotoluene	UG/L	20.0000 U		.	.	10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	20.0000 U	.	.	.	10.0000 U	50.0000
S 2-Methylnaphthalene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	50.0000 U	.	.	.	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	20.0000 U	.		.	10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	50.0000 U	.	.	.	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	50.0000 U	.	.	.	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	20.0000 U	.	.	.	10.0000 U	60.0000
S 4-Chloroaniline	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	20.0000 U	.		.	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	50.0000 U	.	.	.	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	50.0000 U	.	.	.	25.0000 U	62.0000
S Acenaphthene	UG/L	20.0000 U	.	.	.	10.0000 U	36.0000
S Acenaphthylene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S Anthracene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	20.0000 U	.		.	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	20.0000 U	.	.	.	10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	20.0000 U			.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF02DL	NASPAF03	NASPAF04	NASPAF06	NASPAF07	NASPAF07MS
	LAB ID:	522657	524171	525438	528897	529747	529752
	DATE RCVD:	11/14/92	11/20/92	11/28/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
S Butylbenzylphthalate	UG/L	20.0000 ω				10.0000 ω	10.0000 ω
S Carbazole	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Chrysene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Di-n-butylphthalate	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Di-n-octylphthalate	UG/L	20.0000 ω		.	.	10.0000 ω	10.0000 ω
S Dibenz(a,h)anthracene	UG/L	20.0000 ω		.	.	10.0000 ω	10.0000 ω
S Dibenzofuran	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Diethylphthalate	UG/L	20.0000 ω	.	.	.	10.0000 ω	1.0000 ω
S Dimethylphthalate	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Fluoranthene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Fluorene	UG/L	20.0000 ω	.			10.0000 ω	10.0000 ω
S Hexachlorobenzene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Hexachlorobutadiene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Hexachlorocyclopentadiene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Hexachloroethane	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Indeno(1,2,3-cd)pyrene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Isophorone	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S N-Nitroso-di-n-propylamine	UG/L	20.0000 ω	.	.	.	10.0000 ω	35.0000
S N-Nitrosodiphenylamine	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Naphthalene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Nitrobenzene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Pentachlorophenol	UG/L	50.0000 ω	.	.	.	23.0000 ω	61.0000
S Phenanthrene	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S Phenol	UG/L	20.0000 ω	.	.	.	10.0000 ω	46.0000
S Pyrene	UG/L	20.0000 ω	.	.	.	10.0000 ω	38.0000
S bis(2-Chloroethoxy)methane	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S bis(2-Ethlyhexy)phthalate	UG/L	83.0000	.	.	.	10.0000 ω	27.0000 ω
S bis(2-chloroethyl)ether	UG/L	20.0000 ω	.	.	.	10.0000 ω	10.0000 ω
S bis(2-chloroisopropyl)ether	UG/L	20.0000 ω				10.0000 ω	10.0000 ω

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF02DL	NASPAF03	NASPAF04	NASPAF06	NASPAF07	NASPAF07MS
	LAB ID:	522657	524171	525438	528897	529747	529749
	DATE RCVD:	11/14/92	11/20/92	11/28/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	63.0000
V 1,2-Dichloroethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 2-Butanone	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	9.0000 U
V 2-Hexanone	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Acetone	UG/L	.	10.0000 U	10.0000 U	.	20.0000	22.0000
V Benzene	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	59.0000
V Bromodichloromethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Bromofo m	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Bromomethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	52.0000
V Chloroethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Chloroform	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Chloromethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	.	16.0000 U	5.0000 U	.	4.0000 U	5.0000 U
V Styrene	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	10.0000 U
V Toluene	UG/L	.	10.0000 U	10.0000 U	.	2.0000 J	55.0000
V Trichloroethene	UG/L	.	10.0000 U	10.0000 U	.	10.0000 U	55.0000

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF02DL	NASPAF03	NASPAP04	NASPAF06	NASPAF07	NASPAF07MS
	LAB ID:	522657	524171	525430	520097	529747	529749
	DATE RCVD:	11/14/92	11/20/92	11/28/92	12/12/92	12/18/92	12/18/92
NAME	UNITS						
V Vinyl Chloride	UG/L		10.0000 U	10.0000 U		10.0000 U	10.0000 U
V Xylene (total)	UG/L		10.0000 U	10.0000 U		10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L		10.0000 U	10.0000 U		10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L		10.0000 U	10.0000 U		10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF07MSD	NASPAF08	NASPAF08MS	NASPAF08MSD	NASPAF09	NASPAF10
	LAB ID:	529750	530842	530843	530844	531338	531346
	DATE RCVD:	12/18/92	12/31/92	12/31/92	12/31/92	01/08/93	01/08/93
NAME	UNITS						
I Aluminum	UG/L		211.0000	.	.	80.3000 J	101.0000 J
I Antimony	UG/L		57.0000 U	.	.	57.0000 U	52.0000 U
I Arsenic	UG/L		4.0000 U	.	.	4.0000 U	4.5000 J
I Barium	UG/L		3.0000 U	.	.	3.0000 U	2.0000 U
I Beryllium	UG/L		2.0000 U	.	.	2.0000 U	1.0000 U
I Cadmium	UG/L		5.0000 U	.	.	5.0000 U	5.0000 U
I Calcium	UG/L		924.0000 U	.	.	924.0000 U	55.7000 J
I Chromium	UG/L		10.0000 U	.	.	10.0000 U	6.0000 U
I Cobalt	UG/L		11.0000 U	.	.	11.0000 U	9.0000 U
I Copper	UG/L		25.0000 U	.	.	25.0000 U	4.0000 U
I Cyanide	UG/L		10.0000 U	.	.	10.0000 U	10.0000 U
I Iron	UG/L		28.0000 U	.	.	69.3000 J	33.5000 J
I Lead	UG/L		2.0000 U	.	.	2.0000 U	2.0000 U
I Magnesium	UG/L		425.0000 U	.	.	425.0000 U	53.9000 J
I Manganese	UG/L		7.0000 U	.	.	7.0000 U	1.0000 U
I Mercury	UG/L		0.2000 U	.	.	0.2000 U	0.2000 U
I Nickel	UG/L		40.0000 U	.	.	40.0000 U	13.0000 U
I Potassium	UG/L		4860.0000 U	.	.	4860.0000 U	1440.0000 U
I Selenium	UG/L		4.0000 U	.	.	4.0000 U	3.0000 U
I Silver	UG/L		10.0000 U	.	.	37.8000	5.0000 U
I sodium	UG/L		418.0000 J	.	.	255.0000 U	345.0000 J
I Thallium	UG/L		2.0000 U	.	.	2.0000 U	3.0000 U
I Vanadium	UG/L		19.0000 U	.	.	19.0000 U	4.0000 U
I Zinc	UG/L		5.0000 U	.	.	47.7000	18.7000 J
P 4,4'-DDD	UG/L	0.0330 J	0.1000 U	.	.	0.1000 U	0.1000 U
P 4,4'-DDE	UG/L	0.1000 U	0.1000 U	.	.	0.1000 U	0.1000 U
P 4,4'-DDT	UG/L	0.8700	0.1000 U	.	.	0.1000 U	0.1000 U
P Aldrin	UG/L	0.3600	0.0500 U	.	.	0.0500 U	0.0500 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	NASPAF07MSD	NASPAF08	NASPAF08MS	NASPAF08MSD	NASPAF09	NASPAF10
	LAB ID:	529756	530842	530843	530844	531338	531882
	DATE RCVD:	12/18/92	12/31/92	12/31/92	12/31/92	01/08/93	01/14/93
P Aroclor 1016	UG/L	1.0000 U	1.0000 U		.	1.0000 U	1.0000 U
P Aroclor 1221	UG/L	2.0000 U	2.0000 U		.	2.0000 U	2.0000 U
P Aroclor 1232	UG/L	1.0000 U	1.0000 U	.	.	1.0000 U	1.0000 U
P Aroclor 1242	UG/L	1.0000 U	1.0000 U		.	1.0000 U	1.0000 U
P Aroclor 1248	UG/L	1.0000 U	1.0000 U	.	.	1.0000 U	1.0000 U
P Aroclor 1254	UG/L	1.0000 U	1.0000 U	.	.	1.0000 U	1.0000 U
P Aroclor 1260	UG/L	1.0000 U	1.0000 U	.	.	1.0000 U	1.0000 U
P Dieldrin	UG/L	0.8400	0.1000 U	.	.	0.1000 U	0.1000 U
P Endosulfan I	UG/L	0.0500 U	0.0500 U	.	.	0.0500 U	0.0500 U
P Endosulfan II	UG/L	0.1000 U	0.1000 U	.	.	0.1000 U	0.1000 U
P Endosulfan sulfate	UG/L	0.1000 U	0.1000 U	.	.	0.1000 U	0.1000 U
P Endrin	UG/L	0.9700	0.1000 U		.	0.1000 U	0.1000 U
P Endrin aldehyde	UG/L	0.1000 U	0.1000 U	.	.	0.1000 U	0.1000 U
P Endrin ketone	UG/L	0.1000 U	0.1000 U		.	0.1000 U	0.1000 U
P Heptachlor	UG/L	0.3800	0.0500 U	.	.	0.0500 U	0.0500 U
P Heptachlor epoxide	UG/L	0.0500 U	0.0500 U	.	.	0.0500 U	0.0500 U
P Methoxychlor	UG/L	0.5000 U	0.5000 U	.	.	0.5000 U	0.5000 U
P Toxaphene	UG/L	5.0000 U	5.0000 U	.	.	5.0000 U	5.0000 U
P alpha-BHC	UG/L	0.0500 U	0.0500 U	.	.	0.0500 U	0.0500 U
P alpha-Chlordane	UG/L	0.0500 U	0.0500 U	.	.	0.0500 U	0.0500 U
P beta-BHC	UG/L	0.0500 U	0.0500 U	.	.	0.0500 U	0.0500 U
P delta-BHC	UG/L	0.0500 U	0.0500 U	.	.	0.0500 U	0.0500 U
P gamma-BHC (Lindane)	UG/L	0.4100	0.0500 U	.	.	0.0500 U	0.0500 U
P gamma-Chlordane	UG/L	0.0500 U	0.0500 U	.	.	0.0500 U	0.0500 U
S 1,2,4-Trichlorobenzene	UG/L	41.0000	10.0000 U	.	.	10.0000 U	10.0000 U
S 1,2-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S 1,3-Dichlorobenzene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S 1,4-Dichlorobenzene	UG/L	39.0000	10.0000 U	.	.	10.0000 U	10.0000 U
S 2,4,5-Trichlorophenol	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	25.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	NASPAF07MSD 529753 12/18/92	NASPAF08 530842 12/31/92	NASPAF08MS 530843 12/31/92	NASPAF08MSD 530844 12/31/92	NASPAF09 531338 01/08/93	NASPAF10 531882 01/14/93
S 2,4,6-Trichlorophenol	UG/L	10.0000 U	10.0000 U			10.0000 U	10.0000 U
S 2,4-Dichlorophenol	UG/L	10.0000 U	10.0000 U	.		10.0000 U	10.0000 U
S 2,4-Dimethylphenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S 2,4-Dinitrophenol	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	25.0000 U
S 2,4-Dinitrotoluene	UG/L	49.0000	10.0000 U	.	.	10.0000 U	10.0000 U
S 2,6-Dinitrotoluene	UG/L	10.0000 U	10.0000 U	.		10.0000 U	10.0000 U
S 2-Chloronaphthalene	UG/L	10.0000 U	10.00.00 U	.		10.0000 U	10.0000 U
S 2-Chlorophenol	UG/L	55.0000	10.0000 U	.	.	10.0000 U	10.0000 U
S 2-Methylnaphthalene	UG/L	10.0000 U	10.0000 U	.		10.0000 U	10.0000 U
S 2-Methylphenol	UG/L	10.0000 U	10.0000 U	.		10.0000 U	10.0000 U
S 2-Nitroaniline	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	25.0000 U
S 2-Nitrophenol	UG/L	10.0000 U	10.0000 U	.		10.0000 U	10.0000 U
S 3,3'-Dichlorobenzidine	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S 3-Nitroaniline	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	25.0000 U
S 4,6-Dinitro-2-methylphenol	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	25.0000 U
S 4-Bromophenyl-phenylether	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Chloro-3-methylphenol	UG/L	66.0000	10.0000 U		.	10.0000 u	10.0000 U
S 4-Chloroaniline	UG/L	10.0000 U	10.0000 U	.	.	10.0000 u	10.0000 U
S 4-Chlorophenyl-phenylether	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Methylphenol	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S 4-Nitroaniline	UG/L	25.0000 U	25.0000 U	.	.	25.0000 U	25.0000 U
S 4-Nitrophenol	UG/L	67.0000	25.0000 U		.	25.0000 U	25.0000 U
S Acenaphthene	UG/L	40.0000	10.0000 U		.	10.0000 U	10.0000 U
S Acenaphthylene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S Anthracene	UG/L	10.0000 U	10.0000 U		.	10.0000 U	10.0000 U
S Benzo(a)anthracene	UG/L	10.0000 U	10.0000 U		.	10.0000 U	10.0000 U
S Benzo(a)pyrene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S Benzo(b,k)fluoranthene	UG/L	10.0000 U	10.0000 U	.		10.0000 U	10.0000 U
S Benzo(g,h,i)perylene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF07MSD	NASPAF08	NASPAF08MS	NASPAF08MSD	NASPAF09	NASPAF10
	LAB ID:	529753	530842	530843	530844	531338	531882
	DATE RCVD:	12/18/92	12/31/92	12/31/92	12/31/92	01/08/93	01/14/93
NAME	WITS						
S Butylbenzylphthalate	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	10.0000 u
S Carbazole	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	10.0000 u
S Chrysene	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	10.0000 u
S Di-n-butylphthalate	UG/L	10.0000 u	1.0000 U	.	.	10.0000 U	10.0000 u
S Di-n-octylphthalate	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	10,0000 u
S Dibenz (a,h) anthracene	UG/L	10.0000 u	10.0000 U	.	.	10.0000 U	10.0000 u
S Dibenzofuran	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 u
S Diethylphthalate	UG/L	1.0000 U	2.0000 U	.	.	10.0000 U	1.0000 u
S Dimethylphthalate	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	10.0000 U
S Fluoranthene	UG/L	10.0000 u	10.0000 U	.	.	10.0000 U	10.0000 u
S Fluorene	UG/L	10.0000 u	10.0000 U	.	.	10.0000 U	10.0000 U
S Hexachlorobenzene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S Hexachlorobutadiene	UG/L	10.0000 u	10.0000 U	.	.	10.0000 U	10.0000 U
S Hexachlorocyclopentadiene	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	10.0000 u
S Hexachloroethane	UG/L	10.0000 u	10.0000 U	.	.	10.0000 U	10.0000 U
S Indeno(1,2,3-cd)pyrene	UG/L	10.0000 u	10.0000 u	.	.	10.0000 U	10.0000 u
S Isophorone	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S N-Nitroso-di-n-propylamine	UG/L	42.0000	10.0000 U	.	.	10.0000 U	10.0000 u
S N-Nitrosodiphenylamine	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S Naphthalene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 u
S Nitrobenzene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 u
S Pentachlorophenol	UG/L	72.0000	25.0000 U	.	.	25.0000 U	25.0000 U
S Phenanthrene	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 U
S Phenol	UG/L	51.0000	5.0000 U	.	.	10.0000 U	10.0000 U
S Pyrene	UG/L	47.0000	10.0000 U	.	.	10.0000 U	10.0000 U
S bis (2-Chloroethoxy) methane	UG/L	10.0000 U	10.0000 U	.	.	10.0000 U	10.0000 u
S bis (2-Ethlyhexy) phthalate	UG/L	18.0000 u	2.0000 U	.	.	11.0000 U	2.0000 u
S bis (2-chloroethyl) ether	UG/L	10.0000 u	10.0000 U	.	.	10.0000 U	10.0000 u
S bis(2-chloroisopropyl) ether	UG/L	10.0000 u	10.0000 U	.	.	10.0000 U	10.0000 u

NAS I'ENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF07MSD	NASPAF08	NASPAF08MS	NASPAF08MSD	NASPAF09	NASPAF10
	LAB ID:	529750	530842	530843	530844	531338	531882
	DATE RCVD:	12/18/92	12/31/92	12/31/92	12/31/92	01/08/93	01/14/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	61.0000	10.0000 U	53.0000	54.0000	10.0000 U	10.0000 U
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	23.0000	47.0000	58.0000	35.0000	10.0000 U	6.0000 J
V Benzene	UG/L	58.0000	10.0000 U	53.0000	53.0000	10.0000 U	10.0000 U
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	*10.0000* U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	51.0000	10.0000 U	51.0000	51.0000	10.0000 U	10.0000 U
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	*10.0000* U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	5.0000 U	10.0000 U	2.0000 U	2.0000 U	1.0000 U	2.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	54.0000	10.0000 U	52.0000	52.0000	1.0000 J	10.0000 U
V Trichloroethene	UG/L	54.0000	10.0000 U	50.0000	50.0000	10.0000 U	10.0000 U

NAS YENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digite.

	ENSAFE ID:	NASPAF07MSD	NASPAF08	NASPAF08MS	NASPAF08MSD	NASPAF09	NASPAF10
	LAB ID:	529750	530842	530843	530844	531338	531882
	DATE RCVD:	12/18/92	12/31/92	12/31/92	12/31/92	01/08/93	01/14/93
NAME	UNITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	ENSAFE ID: / LAB ID: / DATE RCVD: / UNITS	NASPAF10MS 531003 01/14/93	NASPAF10MSD 531004 01/14/93	NASPAF11 532030 01/21/93	NASPAF12 534044 01/29/93	NASPAF12MS 534052 01/29/93	NASPAF12MSD 534053 01/29/93
		.	.			.	.
I Aluminum	UG/L	.	.	241.0000	351.0000	.	.
I Antimony	UG/L	.	.	52.0000 U	52.0000 U	.	.
I Arsenic	UG/L	.	.	3.0000 U	3.0000 U	.	.
I Barium	UG/L	.	.	3.8000 J	S.7000 J	.	.
I Beryllium	UG/L	.	.	1.0000 U	1.0000 U	.	.
I Cadmium	UG/L	.	.	5.0000 U	5.0000 U	.	.
I Calcium	UG/L	.	.	82.0000 J	76.5000 J	.	.
I Chromium	UG/L	.	.	6.0000 U	6.0000 U	.	.
I Cobalt	UG/L	.	.	9.0000 U	9.0000 U	.	.
I Copper	UG/L	.	.	4.0000 U	4.0000 U	.	.
I Cyanide	UG/L	.	.	10.0000 U	10.0000 U	.	.
I Iron	UG/L	.	.	56.4000 J	191.0000	.	.
I Lead	UG/L	.	.	2.0000 U	2.0000 U	.	.
I Magnesium	UG/L	.	.	104.0000 J	251.0000 J	.	.
I Manganese	UG/L	.	.	1.1000 J	2.6000 J	.	.
I Mercury	UG/L	.	.	0.2000 U	0.2000 U	.	.
I Nickel	UG/L	.	.	13.0000 U	13.0000 U	.	.
I Potassium	UG/L	.	.	1440.0000 U	1440.0000 U	.	.
I Selenium	UG/L	.	.	3.0000 U	3.0000 U	.	.
I Silver	UG/L	.	.	5.0000 U	5.0000 U	.	.
I Sodium	UG/L	.	.	453.0000 J	486.0000 J	.	.
I Thallium	UG/L	.	.	3.0000 U	3.0000 U	.	.
I Vanadium	UG/L	.	.	4.0000 U	4.0000 U	.	.
I Zinc	UG/L	.	.	5.4000 J	5.0000 U	.	.
P 4,4'-DDD	UG/L	.	.	0.1000 U	0.1000 U	.	.
P 4,4'-DDE	UG/L	.	.	0.1000 U	0.1000 U	.	.
P 4,4'-DDT	UG/L	.	.	0.1000 U	0.1000 U	.	.
P Aldrin	UG/L	.	.	0.0500 U	0.0500 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in eignificant digits.

	ENSAFE ID:	NASPAF10MS	NASPAF10MSD	NASPAF11	NASPAF12	NASPAF12MS	NASPAF12MSD
	LAB ID:	531883	531884	532830	534045	534052	534053
	DATE RCVD:	01/14/93	01/14/93	01/21/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
P Aroclor 1016	UG/L	.	.	1.0000 U	1.0000 U	.	.
P Aroclor 1221	UG/L	.	.	2.0000 U	2.0000 U	.	.
P Aroclor 1232	UG/L	.	.	1.0000 U	1.0000 U	.	.
P Aroclor 1242	UG/L	.	.	1.0000 U	1.0000 U	.	.
P Aroclor 1248	UG/L	.	.	1.0000 U	1.0000 U	.	.
P Aroclor 1254	UG/L	.	.	1.0000 U	1.0000 U	.	.
P Aroclor 1260	UG/L	.	.	1.0000 U	1.0000 U	.	.
P Dieldrin	UG/L	.	.	0.1000 U	0.1000 U	.	.
P Endosulfan I	UG/L	.	.	0.0500 U	0.0500 U	.	.
P Endosulfan II	UG/L	.	.	0.1000 U	0.1000 U	.	.
P Endosulfan sulfate	UG/L	.	.	0.1000 U	0.1000 U	.	.
P Endrin	UG/L	.	.	0.1000 U	0.1000 U	.	.
P Endrin aldehyde	UG/L	.	.	0.1000 U	0.1000 U	.	.
P Endrin ketone	UG/L	.	.	0.1000 U	0.1000 U	.	.
P Heptachlor	UG/L	.	.	0.0016 J	0.0500 U	.	.
P Heptachlor epoxide	UG/L	.	.	0.0500 U	0.0500 U	.	.
P Methoxychlor	UG/L	.	.	0.5000 U	0.5000 U	.	.
P Toxaphene	UG/L	.	.	5.0000 U	5.0000 U	.	.
P alpha-BHC	UG/L	.	.	0.0500 U	0.0500 U	.	.
P alpha-Chlordane	UG/L	.	.	0.0500 U	0.0500 U	.	.
P beta-BHC	UG/L	.	.	0.0500 U	0.0500 U	.	.
P de1ta-BHC	UG/L	.	.	0.0500 U	0.0500 U	.	.
P gamma-BHC (Lindane)	UG/L	.	.	0.0500 U	0.0500 U	.	.
P gamma-Chlordane	UG/L	.	.	0.0016 J	0.0500 U	.	.
S 1,2,4-Trichlorobenzene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S 1,2-Dichlorobenzene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S 1,3-Dichlorobenzene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S 1,4-Dichlorobenzene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S 2,4,5-Trichlorophenol	UG/L	.	.	25.0000 U	25.0000 U	.	.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF10MS	NASPAF10MSD	NASPAF11	NASPAF12	NASPAF12MS	NASPAF12MSD
	LAB ID:	531883	531884	532830	534045	534052	534053
	DATE RCVD:	01/14/93	01/14/93	01/21/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
S 2,4,6-Trichlorophenol	UG/L	.	.	10.0000 U	10.0000 U		.
S 2,4-Dichlorophenol	UG/L	.	.	10.0000 U	10.0000 U		.
S 2,4-Dimethylphenol	UG/L	.	.	10.0000 U	10.0000 U		.
S 2,4-Dinitrophenol	UG/L	.	.	25.0000 U	25.0000 U		.
S 2,4-Dinitrotoluene	UG/L	.	.	10.0000 U	10.0000 U		.
S 2,6-Dinitrotoluene	UG/L	.	.	10.0000 U	10.0000 U		.
S 2-Chloronaphthalene	UG/L	.		10.0000 U	10.0000 U		
S 2-Chlorophenol	UG/L		.	10.0000 U	10.0000 U		.
S 2-Methylnaphthalene	UG/L			10.0000 U	10.0000 U		.
S 2-Methylphenol	UG/L	.	.	10.0000 U	10.0000 U		.
S 2-Nitroaniline	UG/L	.	.	25.0000 U	25.0000 U		.
S 2-Nitrophenol	UG/L	.	.	10.0000 U	10.0000 U		
S 3,3'-Dichlorobenzidine	UG/L	.	.	10.0000 U	10.0000 U		.
S 3-Nitroaniline	UG/L	.	.	25.0000 U	25.0000 U		.
S 4,6-Dinitro-2-methylphenol	UG/L		.	25.0000 U	25.0000 U		.
S 4-Brmophenyl-phenylether	UG/L	.	.	10.0000 U	10.0000 U		.
S 4-Chloro-3-methylphenol	UG/L	.	.	10.0000 U	10.0000 U		.
S 4-Chloroaniline	UG/L			10.0000 U	10.0000 U		.
S 4-Chlorophenyl-phenylether	UG/L	.	.	10.0000 U	10.0000 U		.
S 4-Methylphenol	UG/L	.		10.0000 U	10.0000 U		.
S 4-Nitroaniline	UG/L	.		25.0000 U	25.0000 U		.
S 4-Nitrophenol	UG/L	.	.	25.0000 U	25.0000 U		.
S Acenaphthene	UG/L	.		10.0000 U	10.0000 U		.
S Acenaphthylene	UG/L	.	.	10.0000 U	10.0000 U		.
S Anthracene	UG/L			10.0000 U	10.0000 U		.
S Benzo(a)anthracene	UG/L	.	.	10.0000 U	10.0000 U		.
S Benzo(a)pyrene	UG/L		.	10.0000 U	10.0000 U		.
S Benzo(b,k)fluoranthene	UG/L	.		10.0000 U	10.0000 U		.
S Benzo(g,h,i)perylene	UG/L	.	.	10.0000 U	10.0000 U		.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

NAME	UNITS	NASPAF10MS	NASPAF10MSD	NASPAF11	NASPAF12	NASPAF12MS	NASPAF12MSD
ENSAFE ID:		NASPAF10MS	NASPAF10MSD	NASPAF11	NASPAF12	NASPAF12MS	NASPAF12MSD
LAB ID:		531883	531884	532830	534045	534052	534053
DATE RCVD:		01/14/93	01/14/93	01/28/93	01/29/93	01/29/93	01/29/93
S Butylbenzylphthalate	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Carbazole	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Chrysene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Di-n-butylphthalate	UG/L	.	.	2.0000 U	10.0000 U	.	.
S Di-n-octylphthalate	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Dibenz(a,h)anthracene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Dibenzofuran	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Diethylphthalate	UG/L	.	.	1.0000 U	10.0000 U	.	.
S Dimethylphthalate	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Fluoranthene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Fluorene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Hexachlorobenzene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Hexachlorobutadiene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Hexachlorocyclopentadiene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Hexachloroethane	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Indeno(1,2,3-cd)pyrene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Isophorone	UG/L	.	.	10.0000 U	10.0000 U	.	.
S N-Nitroso-di-n-propylamine	UG/L	.	.	10.0000 U	10.0000 U	.	.
S N-Nitrosodiphenylamine	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Naphthalene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Nitrobenzene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Pentachlorophenol	UG/L	.	.	25.0000 U	25.0000 U	.	.
S Phenanthrene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Phenol	UG/L	.	.	10.0000 U	10.0000 U	.	.
S Pyrene	UG/L	.	.	10.0000 U	10.0000 U	.	.
S bis(2-Chloroethoxy)methane	UG/L	.	.	10.0000 U	10.0000 U	.	.
S bis(2-Ethlyhexy)phthalate	UG/L	.	.	11.0000 U	10.0000 U	.	.
S bis(2-chloroethyl)ether	UG/L	.	.	10.0000 U	10.0000 U	.	.
S bis(2-chloroisopropyl)ether	UG/L			10.0000 U	10.0000 U		.

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF10MS	NASPAF10MSD	NASPAF11	NASPAF12	NASPAF12MS	NASPAF12MSD
	LAB ID:	531883	531884	532830	534045	534052	534053
	DATE RCVD:	01/14/93	01/14/93	01/21/93	01/29/93	01/29/93	01/29/93
NAME	UNITS						
V 1,1,1-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2,2-Tetrachloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1,2-Trichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,1-Dichloroethene	UG/L	53.0000	54.0000	10.0000 U	10.0000 U	46.0000	42.0000
V 1,2-Dichloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloroethene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 1,2-Dichloropropane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Butanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 2-Hexanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V 4-Methyl-2-Pentanone	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Acetone	UG/L	9.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Benzene	UG/L	54.0000	54.0000	10.0000 U	10.0000 U	47.0000	47.0000
V Bromodichloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromoform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Bromomethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Disulfide	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Carbon Tetrachloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chlorobenzene	UG/L	49.0000	49.0000	10.0000 U	10.0000 U	45.0000	43.0000
V Chloroethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloroform	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Chloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Dibromochloromethane	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Ethylbenzene	UG/L	10.0000 U	10,0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Methylene Chloride	UG/L	3.0000 U	9.0000 U	1.0000 U	10.0000 U	1.0000 U	10.0000 U
V Styrene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Tetrachloroethene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Toluene	UG/L	49.0000	50.0000	10.0000 U	10.0000 U	47.0000	44.0000
V Trichloroethene	UG/L	51.0000	51.0000	10.0000 U	10.0000 U	44.0000	43.0000

NAS PENSACOLA OU-10 AND SITE 13
QAQC SAMPLES

Key presented at end of Appendix. Data not shown in significant digits.

	ENSAFE ID:	NASPAF10MS	NASPAF10MSD	NASPAF11	NASPAF12	NASPAF12MS	NASPAF12MSD
	LAB ID:	531883	531884	532830	534045	534052	534053
	DATE RCVD:	01/14/93	01/14/93	01/21/93	01/29/93	01/29/93	01/29/93
NAME	WITS						
V Vinyl Chloride	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V Xylene (total)	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V cis-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U
V trans-1,3-Dichloropropene	UG/L	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U	10.0000 U

V=Volatile, S=Semivolatile, P=Pesticide, I=Inorganic, R=Remediation