User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 05/28/2002 Time: 14:43:08

Sample: AE11799
Analysis: $REGCMS Reference for GCMS Scan
Units: ug
Started: 5/28/02 14:21 Ended: 5/28/02 14:21 Analyst: MF
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		7.45		2.0
2	bis(2-Chloroethyl)ether		15.32		2.0
3	1,3-Dichlorobenzene		13.43		2.0
4	1,4-Dichlorobenzene		12.98		2.0
5	1,2-Dichlorobenzene		14.45		2.0
6	2,2-oxybis(1-Chloropropane)		16.04		2.0
7	n-Nitrosodi-n-propylamine		15.61		2.0
8	Hexachloroethane		10.11		2.0
9	Nitrobenzene		15.70		2.0
10	Isophorone		17.68		2.0
11	bis(2-Chloroethoxy)methane		16.05		2.0
12	1,2,4-Trichlorobenzene		14.55		2.0
13	Naphthalene		17.11		2.0
14	Hexachlorobutadiene		11.11		2.0
15	2-Chloronaphthalene		18.17		2.0
16	Dimethylphthalate		16.58		2.0
17	2,6-Dinitrotoluene		16.49		2.0
18	Acenaphthylene		16.51		2.0
19	Acenaphthene		17.70		2.0
20	2,4-Dinitrotoluene		16.37		2.0
21	Diethylphthalate		17.52		2.0
22	4-Chlorophenylphenylether		16.63		2.0
23	Fluorene		17.89		2.0
24	Azobenzene		16.52		2.0
25	n-Nitrosodiphenylamine		19.15		2.0
26	4-Bromophenylphenylether		17.26		2.0
27	Hexachlorobenzene		17.71		2.0
28	Phenanthrene		17.53		2.0
29	Anthracene		15.95		2.0
30	Di-n-butylphthalate		17.27		2.0
31	Fluoranthene		16.38		2.0
32	Pyrene		19.47		2.0
33	Butylbenzylphthalate		12.90		2.0
34	Benzo(a)anthracene		16.82		2.0
35	bis(2-Ethylhexyl)phthalate		10.46		2.0
36	Chrysene		17.22		2.0
37	Di-n-octylphthalate		8.66		2.0
38	Benzo(b)fluoranthene		16.97		2.0
39	Benzo(k)fluoranthene		17.68		2.0
40	Benzo(a)pyrene		12.12		2.0
41	Indeno(1,2,3-cd)pyrene		12.53		2.0
42	Dibenz(a,h)anthracene		15.14		2.0
43	Benzo(g,h,i)perylene		14.99		2.0

User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 05/28/2002 Time: 14:43:17

Sample: AE11799
Analysis: $Q_GCMS Ref calc for GCMS Scan
Units: ug
Started: 5/28/02 14:21 Ended: 5/28/02 14:21 Analyst: MF
Result source: calculation
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		37		2.0
2	bis(2-Chloroethyl)ether		77		2.0
3	1,3-Dichlorobenzene		67		2.0
4	1,4-Dichlorobenzene		65		2.0
5	1,2-Dichlorobenzene		72		2.0
6	2,2-oxybis(1-Chloropropane)		80		2.0
7	n-Nitrosodi-n-propylamine		78		2.0
8	Hexachloroethane		51		2.0
9	Nitrobenzene		78		2.0
10	Isophorone		88		2.0
11	bis(2-Chloroethoxy)methane		80		2.0
12	1,2,4-Trichlorobenzene		73		2.0
13	Naphthalene		86		2.0
14	Hexachlorobutadiene		56		2.0
15	2-Chloronaphthalene		91		2.0
16	Dimethylphthalate		83		2.0
17	2,6-Dinitrotoluene		82		2.0
18	Acenaphthylene		83		2.0
19	Acenaphthene		88		2.0
20	2,4-Dinitrotoluene		82		2.0
21	Diethylphthalate		88		2.0
22	4-Chlorophenylphenylether		83		2.0
23	Fluorene		89		2.0
24	Azobenzene		83		2.0
25	n-Nitrosodiphenylamine		96		2.0
26	4-Bromophenylphenylether		86		2.0
27	Hexachlorobenzene		89		2.0
28	Phenanthrene		88		2.0
29	Anthracene		80		2.0
30	Di-n-butylphthalate		86		2.0
31	Fluoranthene		82		2.0
32	Pyrene		97		2.0
33					
34	Benzo(a)anthracene		84		2.0
35					
36	Chrysene		86		2.0
37					
38	Benzo(b)fluoranthene		85		2.0
39	Benzo(k)fluoranthene		88		2.0
40	Benzo(a)pyrene		61		2.0
41	Indeno(1,2,3-cd)pyrene		63		2.0
42	Dibenz(a,h)anthracene		76		2.0
43	Benzo(g,h,i)perylene		75		2.0

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF1.D Vial: 4
Acq On : 22 May 2002 4:16 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:36:28 2002 Quant Results File: 052202.RES

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.90	152	4935975	40.00	ug/L	0.00
10) Napthalene-d8	13.40	136	19575446	40.00	ug/L	-0.01
17) Acenapthalene-d10	18.53	164	10866616	40.00	ug/L	0.00
29) Phenanthranene-d10	22.92	188	16461139	40.00	ug/L	0.00
37) Chrysene-d12	30.77	240	11464744	40.00	ug/L	-0.04
43) Perylene-d12	34.66	264	6637249	40.00	ug/L	-0.02
System Monitoring Compounds						
27) TCMX	20.51	209	2382371	38.05	ug/L	0.00
Spiked Amount 40.000			Recovery = 95.13%			
Target Compounds						Qvalue
2) n-nitros-dimethyl amine	3.87	74	686299	7.45	ug/L #	81
3) bis(2-Chloroethyl)ether	9.31	93	2800493	15.32	ug/L	98
4) 1,3-Dichlorobenzene	9.73	146	2407556	13.07	ug/L	99
5) 1,4-Dichlorobenzene	9.94	146	2558495	12.78	ug/L	99
6) 1,2-Dichlorobenzene	10.32	146	2546625	14.45	ug/l	99
7) 2,2'oxybis(1-Chloropropane	10.75	45	4494638m	16.04	ug/L	
8) N-nitroso-di-propylamine	11.10	70	1857123	15.61	ug/L	97
9) Hexachloroethane	11.22	117	729485	10.05	ug/L	97
11) Nitrobenzene	11.46	77	2697740	15.70	ug/L	98
12) Isophorone	12.17	82	5505613	17.68	ug/L	99
13) bis(2-Chloroethoxy) methan	12.91	93	3501273	16.05	ug/L	98
14) 1,2,4-Trichlorobenzene	13.27	180	2151370	14.55	ug/L	100
15) Napthalene	13.45	128	8831360	17.11	ug/L	100
16) Hexachlorobutadiene	13.90	225	784147	11.11	ug/L	97
18) 2-Chloronaphthalene	16.89	162	5056444	18.17	ug/L	99
19) Dimethylphthalate	17.97	163	5663723	16.58	ug/L	99
20) 2,6-Dinitrotoluene	18.08	165	904752	16.49	ug/L	98
21) Acenapthylene	18.09	152	7663758	16.51	ug/L	99
22) Acenapthalene	18.62	153	5187449	17.70	ug/L	99
23) 2,4-Dinitrotoluene	19.24	165	1174605	16.37	ug/L	95
24) Diethylphthalate	20.11	149	5592315	17.52	ug/L	99
25) 4-Chlorophenyl-phenylether	20.29	204	2831377	16.63	ug/L	97
26) Fluorene	20.16	166	6181011	17.89	ug/L	100
28) Azobenzene	20.74	77	6211729	16.52	ug/L	98
30) Nnitrosodiphenylamine	20.66	169	3180273	19.15	ug/L	100
31) 4-Bromophenyl-phenylether	21.72	248	1455411	17.26	ug/L	99
32) Hexachlorobenzene	21.75	284	1451047	17.71	ug/L	98
33) Phenanthrene	22.98	178	8681059	17.53	ug/L	99
34) Anthracene	23.13	178	7772950	15.95	ug/L	99
35) Di-n-butylphthalate	25.05	149	8493781	17.27	ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF1.D Vial: 4
Acq On : 22 May 2002 4:16 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:36:28 2002 Quant Results File: 052202.RES

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

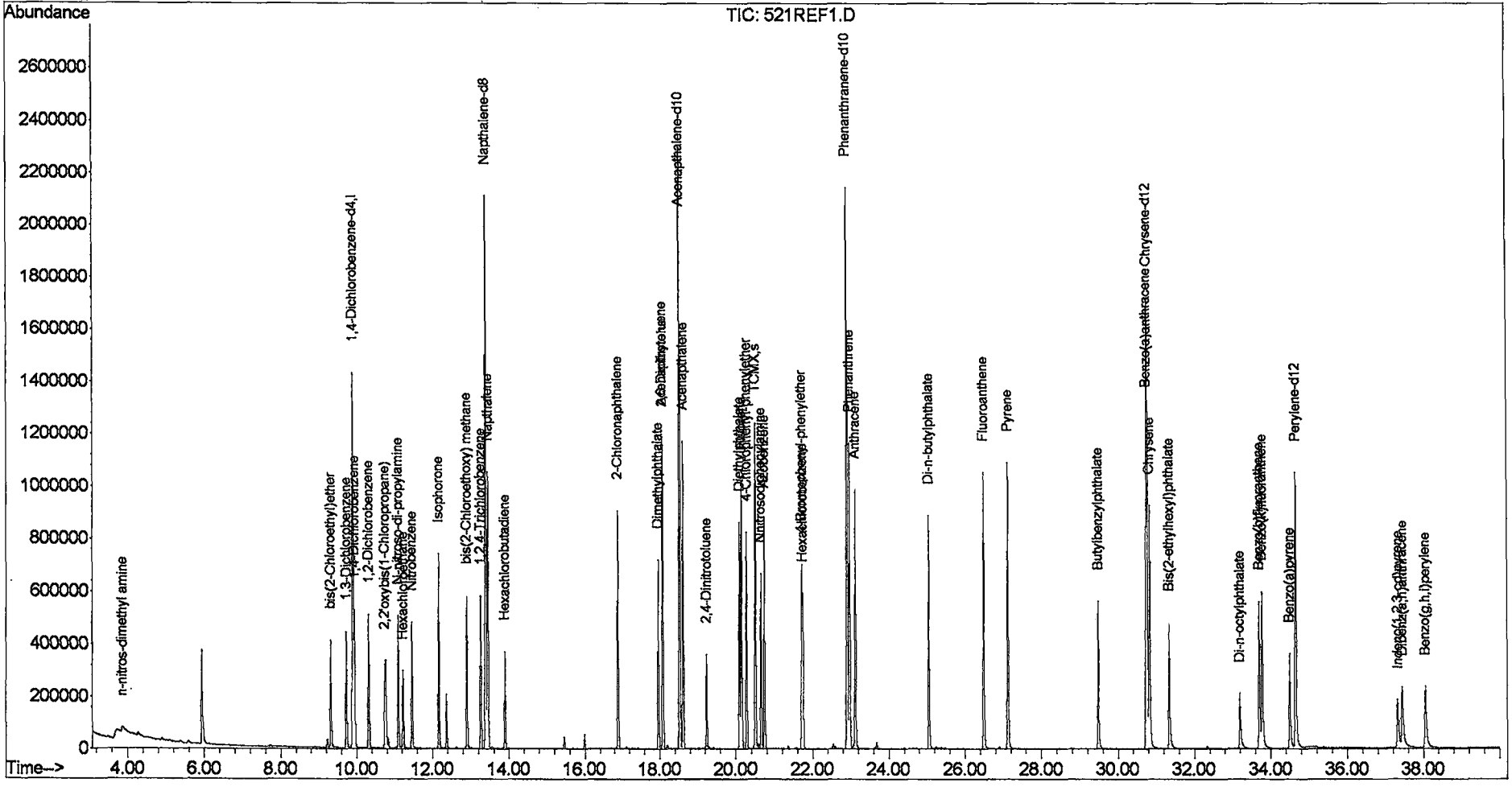
	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
36)	Fluoroanthene	26.50	202	8519902	16.38	ug/L	99
38)	Pyrene	27.13	202	8867504	19.47	ug/L	99
39)	Butylbenzylphthalate	29.48	149	2513007	12.90	ug/L	96
40)	Benzo(a)anthracene	30.74	228	5782730	16.82	ug/L	99
41)	Bis(2-ethylhexyl)phthalate	31.35	149	2651734	10.46	ug/L	98
42)	Chrysene	30.84	228	6802658m	17.22	ug/L	
44)	Di-n-octylphthalate	33.20	149	2477992	8.66	ug/L	100
45)	Benzo(b)fluoranthene	33.70	252	4474230	16.97	ug/L	99
46)	Benzo(k)fluoranthene	33.78	252	5221851	17.68	ug/L	99
47)	Benzo(a)pyrene	34.51	252	2857820	12.12	ug/L	100
48)	Indeno(1,2,3-cd)pyrene	37.32	276	1969979	12.53	ug/L	96
49)	Dibenz(a,h)anthracene	37.45	278	2618109	15.14	ug/L	100
50)	Benzo(g,h,i)perylene	38.06	276	2751330	14.99	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF1.D Vial: 4
Acq On : 22 May 2002 4:16 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:36 2002 Quant Results File: 052202.RES

Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF2.D Vial: 5
Acq On : 22 May 2002 5:05 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:37:23 2002 Quant Results File: 052202.RES

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.90	152	3820779	40.00	ug/L	0.00
10) Napthalene-d8	13.39	136	15277958	40.00	ug/L	-0.01
17) Acenapthalene-d10	18.53	164	8518901	40.00	ug/L	0.00
29) Phenanthranene-d10	22.91	188	12919843	40.00	ug/L	0.00
37) Chrysene-d12	30.76	240	8646037	40.00	ug/L	-0.04
43) Perylene-d12	34.66	264	5033337	40.00	ug/L	-0.02
System Monitoring Compounds						
27) TCMX	20.51	209	1773550	36.13	ug/L	0.00
Spiked Amount 40.000			Recovery = 90.33%			
Target Compounds						Qvalue
2) n-nitros-dimethyl amine	3.89	74	481285	6.75	ug/L #	78
3) bis(2-Chloroethyl)ether	9.31	93	2101353	14.85	ug/L	98
4) 1,3-Dichlorobenzene	9.73	146	1915752	13.43	ug/L	99
5) 1,4-Dichlorobenzene	9.94	146	2011016	12.98	ug/L	99
6) 1,2-Dichlorobenzene	10.32	146	1969131	14.44	ug/l	98
7) 2,2'oxybis(1-Chloropropane	10.77	45	3358686m	15.49	ug/L	
8) N-nitroso-di-propylamine	11.10	70	1398722	15.19	ug/L	98
9) Hexachloroethane	11.22	117	567978	10.11	ug/L	99
11) Nitrobenzene	11.46	77	1969687	14.68	ug/L	99
12) Isophorone	12.17	82	4135136	17.02	ug/L	100
13) bis(2-Chloroethoxy) methan	12.91	93	2622897	15.41	ug/L	99
14) 1,2,4-Trichlorobenzene	13.27	180	1632853	14.15	ug/L	98
15) Napthalene	13.45	128	6618861	16.43	ug/L	99
16) Hexachlorobutadiene	13.90	225	586266	10.64	ug/L	97
18) 2-Chloronaphthalene	16.89	162	3786725	17.35	ug/L	99
19) Dimethylphthalate	17.97	163	4295368	16.04	ug/L	99
20) 2,6-Dinitrotoluene	18.08	165	656596	15.27	ug/L	97
21) Acenapthylene	18.09	152	5801427	15.94	ug/L	98
22) Acenapthalene	18.62	153	3928356	17.09	ug/L	98
23) 2,4-Dinitrotoluene	19.24	165	837436	14.89	ug/L	98
24) Diethylphthalate	20.11	149	4327636	17.30	ug/L	100
25) 4-Chlorophenyl-phenylether	20.29	204	2170867	16.26	ug/L	99
26) Fluorene	20.16	166	4731624	17.47	ug/L	99
28) Azobenzene	20.74	77	4695422	15.93	ug/L	99
30) Nnitrosodiphenylamine	20.65	169	2535595	19.45	ug/L	98
31) 4-Bromophenyl-phenylether	21.72	248	1095524	16.56	ug/L	99
32) Hexachlorobenzene	21.75	284	1094270	17.02	ug/L	95
33) Phenanthrene	22.98	178	6694490	17.23	ug/L	99
34) Anthracene	23.13	178	5877630	15.36	ug/L	99
35) Di-n-butylphthalate	25.05	149	6335464	16.42	ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF2.D Vial: 5
Acq On : 22 May 2002 5:05 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:37:23 2002 Quant Results File: 052202.RES

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

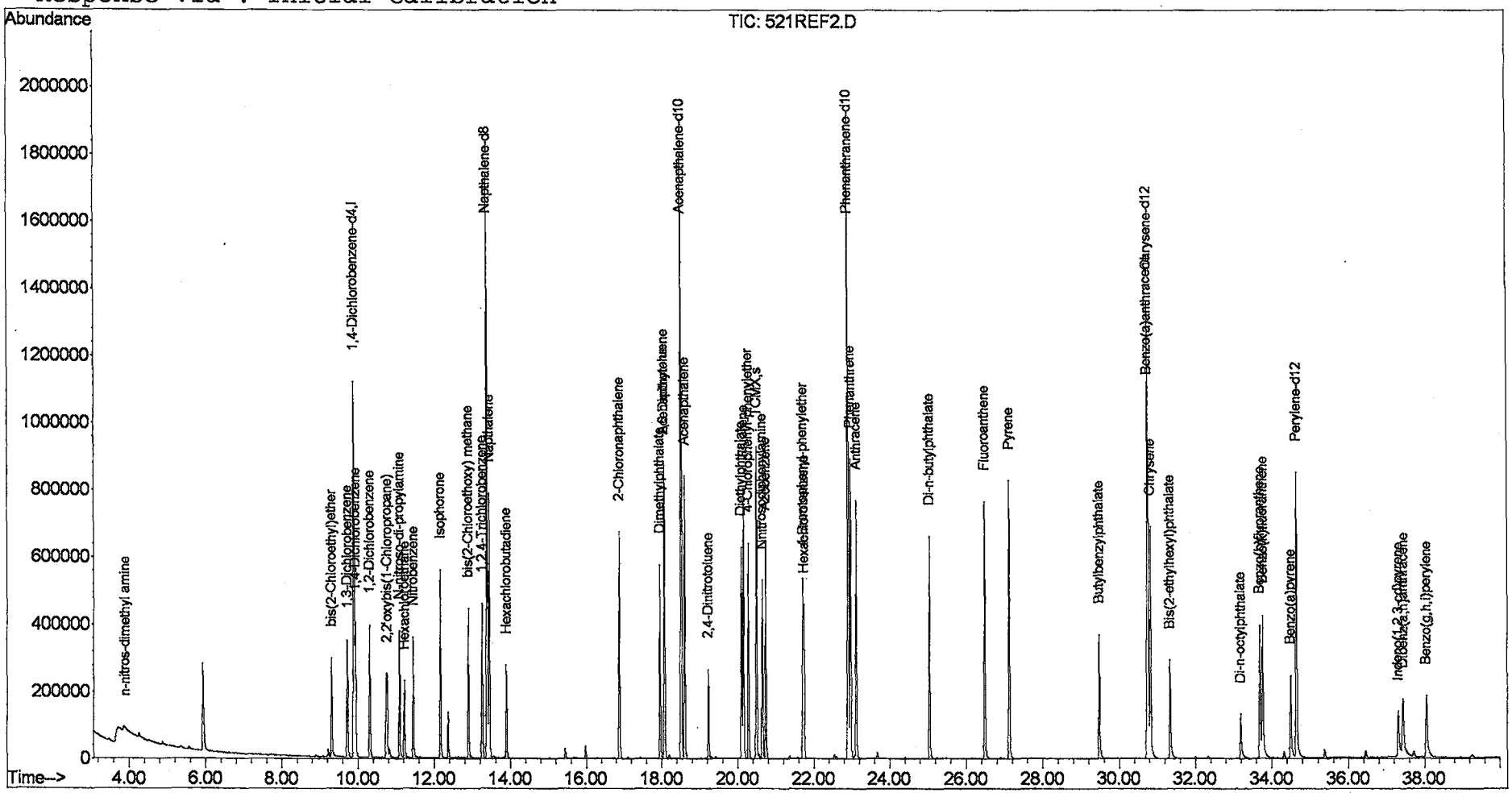
	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
36)	Fluoroanthene	26.49	202	6329723	15.51	ug/L	99
38)	Pyrene	27.13	202	6650358	19.37	ug/L	99
39)	Butylbenzylphthalate	29.48	149	1684438	11.46	ug/L	95
40)	Benzo(a)anthracene	30.74	228	4127024	15.92	ug/L	100
41)	Bis(2-ethylhexyl)phthalate	31.34	149	1660056	8.69	ug/L	99
42)	Chrysene	30.83	228	4965155m	16.67	ug/L	
44)	Di-n-octylphthalate	33.20	149	1477077	6.81	ug/L	99
45)	Benzo(b)fluoranthene	33.70	252	3185507	15.94	ug/L	98
46)	Benzo(k)fluoranthene	33.78	252	3861250	17.24	ug/L	100
47)	Benzo(a)pyrene	34.50	252	2058703	11.51	ug/L	98
48)	Indeno(1,2,3-cd)pyrene	37.32	276	1492028	12.52	ug/L	97
49)	Dibenz(a,h)anthracene	37.45	278	2021457	15.42	ug/L	98
50)	Benzo(g,h,i)perylene	38.06	276	2221318	15.96	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF2.D Vial: 5
Acq On : 22 May 2002 5:05 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:37 2002 Quant Results File: 052202.RES

Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF2.D Vial: 5
Acq On : 22 May 2002 5:05 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:37:23 2002 Quant Results File: 052202.RES

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.90	152	3820779	40.00	ug/L	0.00
10) Napthalene-d8	13.39	136	15277958	40.00	ug/L	-0.01
17) Acenapthalene-d10	18.53	164	8518901	40.00	ug/L	0.00
29) Phenanthranene-d10	22.91	188	12919843	40.00	ug/L	0.00
37) Chrysene-d12	30.76	240	8646037	40.00	ug/L	-0.04
43) Perylene-d12	34.66	264	5033337	40.00	ug/L	-0.02
System Monitoring Compounds						
27) TCMX	20.51	209	1773550	36.13	ug/L	0.00
Spiked Amount 40.000			Recovery = 90.33%			
Target Compounds						Qvalue
2) n-nitros-dimethyl amine	3.89	74	481285	6.75	ug/L #	78
3) bis(2-Chloroethyl)ether	9.31	93	2101353	14.85	ug/L	98
4) 1,3-Dichlorobenzene	9.73	146	1915752	13.43	ug/L	99
5) 1,4-Dichlorobenzene	9.94	146	2011016	12.98	ug/L	99
6) 1,2-Dichlorobenzene	10.32	146	1969131	14.44	ug/l	98
7) 2,2'oxybis(1-Chloropropane	10.77	45	3358686m	15.49	ug/L	
8) N-nitroso-di-propylamine	11.10	70	1398722	15.19	ug/L	98
9) Hexachloroethane	11.22	117	567978	10.11	ug/L	99
11) Nitrobenzene	11.46	77	1969687	14.68	ug/L	99
12) Isophorone	12.17	82	4135136	17.02	ug/L	100
13) bis(2-Chloroethoxy) methan	12.91	93	2622897	15.41	ug/L	99
14) 1,2,4-Trichlorobenzene	13.27	180	1632853	14.15	ug/L	98
15) Napthalene	13.45	128	6618861	16.43	ug/L	99
16) Hexachlorobutadiene	13.90	225	586266	10.64	ug/L	97
18) 2-Chloronaphthalene	16.89	162	3786725	17.35	ug/L	99
19) Dimethylphthalate	17.97	163	4295368	16.04	ug/L	99
20) 2,6-Dinitrotoluene	18.08	165	656596	15.27	ug/L	97
21) Acenapthylene	18.09	152	5801427	15.94	ug/L	98
22) Acenapthalene	18.62	153	3928356	17.09	ug/L	98
23) 2,4-Dinitrotoluene	19.24	165	837436	14.89	ug/L	98
24) Diethylphthalate	20.11	149	4327636	17.30	ug/L	100
25) 4-Chlorophenyl-phenylether	20.29	204	2170867	16.26	ug/L	99
26) Fluorene	20.16	166	4731624	17.47	ug/L	99
28) Azobenzene	20.74	77	4695422	15.93	ug/L	99
30) Nnitrosodiphenylamine	20.65	169	2535595	19.45	ug/L	98
31) 4-Bromophenyl-phenylether	21.72	248	1095524	16.56	ug/L	99
32) Hexachlorobenzene	21.75	284	1094270	17.02	ug/L	95
33) Phenanthrene	22.98	178	6694490	17.23	ug/L	99
34) Anthracene	23.13	178	5877630	15.36	ug/L	99
35) Di-n-butylphthalate	25.05	149	6335464	16.42	ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF2.D Vial: 5
Acq On : 22 May 2002 5:05 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:37:23 2002 Quant Results File: 052202.RES

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
36)	Fluoroanthene	26.49	202	6329723	15.51	ug/L	99
38)	Pyrene	27.13	202	6650358	19.37	ug/L	99
39)	Butylbenzylphthalate	29.48	149	1684438	11.46	ug/L	95
40)	Benzo(a)anthracene	30.74	228	4127024	15.92	ug/L	100
41)	Bis(2-ethylhexyl)phthalate	31.34	149	1660056	8.69	ug/L	99
42)	Chrysene	30.83	228	4965155m	16.67	ug/L	
44)	Di-n-octylphthalate	33.20	149	1477077	6.81	ug/L	99
45)	Benzo(b)fluoranthene	33.70	252	3185507	15.94	ug/L	98
46)	Benzo(k)fluoranthene	33.78	252	3861250	17.24	ug/L	100
47)	Benzo(a)pyrene	34.50	252	2058703	11.51	ug/L	98
48)	Indeno(1,2,3-cd)pyrene	37.32	276	1492028	12.52	ug/L	97
49)	Dibenz(a,h)anthracene	37.45	278	2021457	15.42	ug/L	98
50)	Benzo(g,h,i)perylene	38.06	276	2221318	15.96	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052202\521REF2.D Vial: 5
Acq On : 22 May 2002 5:05 pm Operator: Mclean
Sample : 5/21 ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 28 14:37 2002 Quant Results File: 052202.RES

Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 28 14:25:08 2002
Response via : Initial Calibration