User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 05/17/2002 Time: 11:49:38

Sample: AE10901
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 5/17/02 11:30 Ended: 5/17/02 11:30 Analyst: MF
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		< MRL		2.0
2	bis(2-Chloroethyl)ether		< MRL		2.0
3	1,3-Dichlorobenzene		< MRL		2.0
4	1,4-Dichlorobenzene		< MRL		2.0
5	1,2-Dichlorobenzene		< MRL		2.0
6	2,2'-oxybis(1-Chloropropane)		< MRL		2.0
7	n-Nitrosodi-n-propylamine		< MRL		2.0
8	Hexachloroethane		< MRL		2.0
9	Nitrobenzene		< MRL		2.0
10	Isophorone		< MRL		2.0
11	bis(2-Chloroethoxy)methane		< MRL		2.0
12	1,2,4-Trichlorobenzene		< MRL		2.0
13	Naphthalene		< MRL		2.0
14	Hexachlorobutadiene		< MRL		2.0
15	2-Chloronaphthalene		< MRL		2.0
16	Dimethylphthalate		< MRL		2.0
17	2,6-Dinitrotoluene		< MRL		2.0
18	Acenaphthylene		< MRL		2.0
19	Acenaphthene		< MRL		2.0
20	2,4-Dinitrotoluene		< MRL		2.0
21	Diethylphthalate		< MRL		2.0
22	4-Chlorophenylphenylether		< MRL		2.0
23	Fluorene		< MRL		2.0
24	Azobenzene		< MRL		2.0
25	n-Nitrosodiphenylamine		< MRL		2.0
26	4-Bromophenylphenylether		< MRL		2.0
27	Hexachlorobenzene		< MRL		2.0
28	Phenanthrene		< MRL		2.0
29	Anthracene		< MRL		2.0
30	Di-n-butylphthalate		< MRL		2.0
31	Fluoranthene		< MRL		2.0
32	Pyrene		< MRL		2.0
33	Butylbenzylphthalate		< MRL		2.0
34	Benzo(a)anthracene		< MRL		2.0
35	bis(2-Ethylhexyl)phthalate		< MRL		2.0
36	Chrysene		< MRL		2.0
37	Di-n-octylphthalate		< MRL		2.0
38	Benzo(b)fluoranthene		< MRL		2.0
39	Benzo(k)fluoranthene		< MRL		2.0
40	Benzo(a)pyrene		< MRL		2.0
41	Indeno(1,2,3-cd)pyrene		< MRL		2.0
42	Dibenz(a,h)anthracene		< MRL		2.0
43	Benzo(g,h,i)perylene		< MRL		2.0

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 16 14:34:33 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Thu May 16 14:34:28 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.93	152	5738365	40.00 ug/L	0.00
10) Napthalene-d8	13.43	136	22789344	40.00 ug/L	0.00
17) Acenapthalene-d10	18.57	164	12278988	40.00 ug/L	0.00
29) Phenanthranene-d10	22.95	188	17470340	40.00 ug/L	0.00
37) Chrysene-d12	30.81	240	12010307	40.00 ug/L	0.00
43) Perylene-d12	34.71	264	7287670	40.00 ug/L	0.00
System Monitoring Compounds					
27) TCMX	20.55	209	3074858	41.25 ug/L	0.00
Spiked Amount 40.000			Recovery = 103.13%		
Target Compounds					Qvalue
2) n-nitros-dimethyl amine	0.00	74	0	N.D.	
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.	
4) 1,3-Dichlorobenzene	0.00	146	0	N.D.	
5) 1,4-Dichlorobenzene	0.00	146	0	N.D.	
6) 1,2-Dichlorobenzene	0.00	146	0	N.D.	
7) 2,2'oxybis(1-Chloropropane	0.00	45	0	N.D.	
8) N-nitroso-di-propylamine	0.00	70	0	N.D.	
9) Hexachloroethane	0.00	117	0	N.D.	
11) Nitrobenzene	0.00	77	0	N.D.	
12) Isophorone	0.00	82	0	N.D.	
13) bis(2-Chloroethoxy) methan	0.00	93	0	N.D.	
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.	
15) Napthalene	0.00	128	0	N.D.	
16) Hexachlorobutadiene	0.00	225	0	N.D.	
18) 2-Chloronaphthalene	0.00	162	0	N.D.	
19) Dimethylphthalate	0.00	163	0	N.D.	
20) 2,6-Dinitrotoluene	0.00	165	0	N.D.	
21) Acenapthylene	0.00	152	0	N.D.	
22) Acenapthalene	0.00	153	0	N.D.	
23) 2,4-Dinitrotoluene	0.00	165	0	N.D.	
24) Diethylphthalate	0.00	149	0	N.D.	
25) 4-Chlorophenyl-phenylether	20.55	204	18524	N.D.	
26) Fluorene	0.00	166	0	N.D.	
28) Azobenzene	20.55	77	50437	N.D.	
30) Nnitrosodiphenylamine	20.55	169	56950	0.27 ug/L #	60
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.	
32) Hexachlorobenzene	0.00	284	0	N.D.	
33) Phenanthrene	0.00	178	0	N.D.	
34) Anthracene	0.00	178	0	N.D.	
35) Di-n-butylphthalate	0.00	149	0	N.D.	

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

Quantitation Report (QT Reviewed)

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Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 16 14:34:33 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Thu May 16 14:34:28 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
36)	Fluoroanthene	0.00	202	0	N.D.	
38)	Pyrene	0.00	202	0	N.D.	
39)	Butylbenzylphthalate	0.00	149	0	N.D.	
40)	Benzo(a)anthracene	30.81	228	29323	N.D.	
41)	Bis(2-ethylhexyl)phthalate	0.00	149	0	N.D.	
42)	Chrysene	0.00	228	0	N.D.	
44)	Di-n-octylphthalate	0.00	149	0	N.D.	
45)	Benzo(b)fluoranthene	0.00	252	0	N.D.	
46)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
47)	Benzo(a)pyrene	0.00	252	0	N.D.	
48)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
49)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
50)	Benzo(g,h,i)perylene	0.00	276	0	N.D.	

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

Quantitation Report (QT Reviewed)

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Acq On : 14 May 2002 3:07 pm Operator: Mclean
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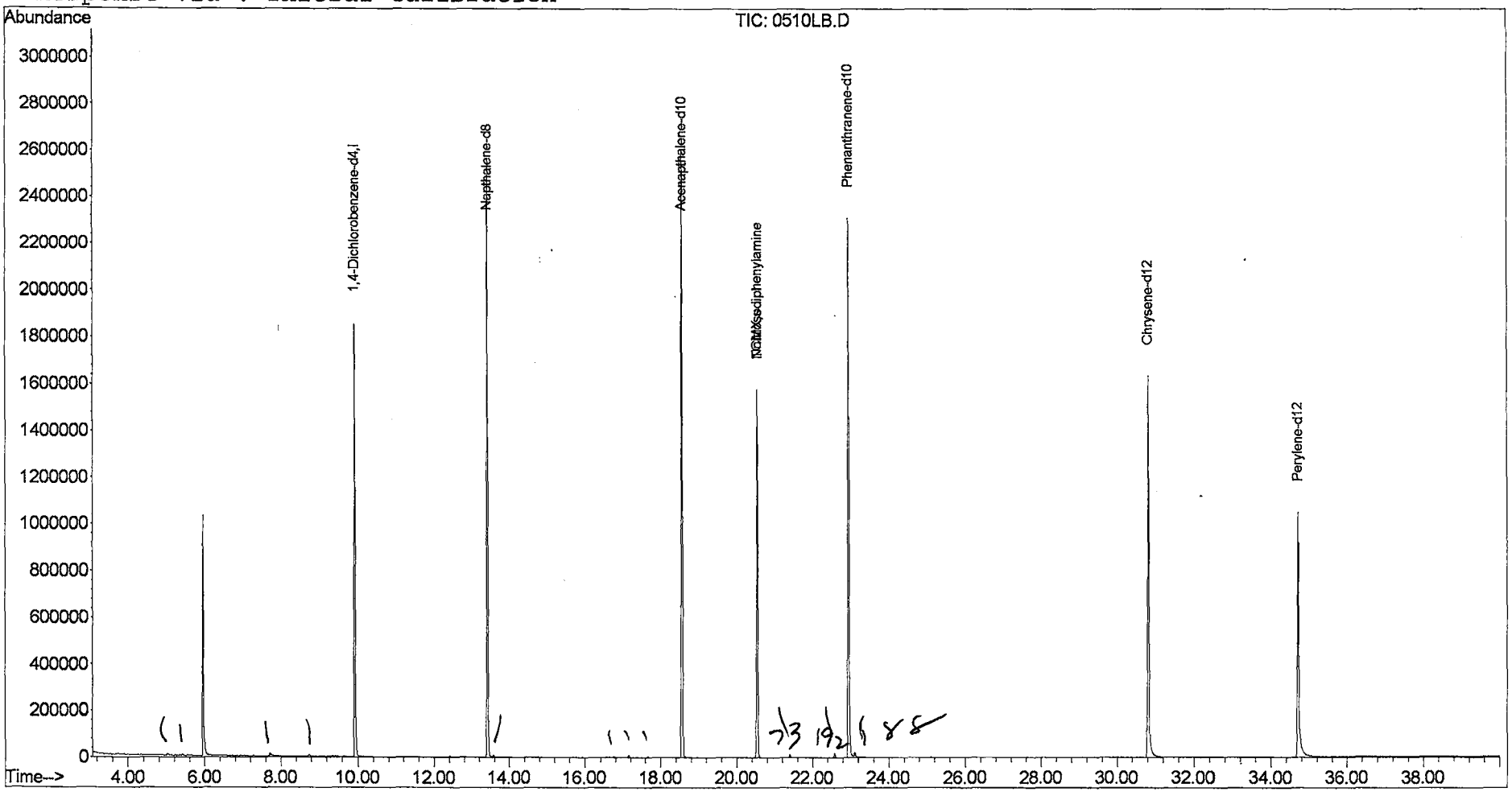
Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Thu May 16 14:34:28 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan

Signal : TIC

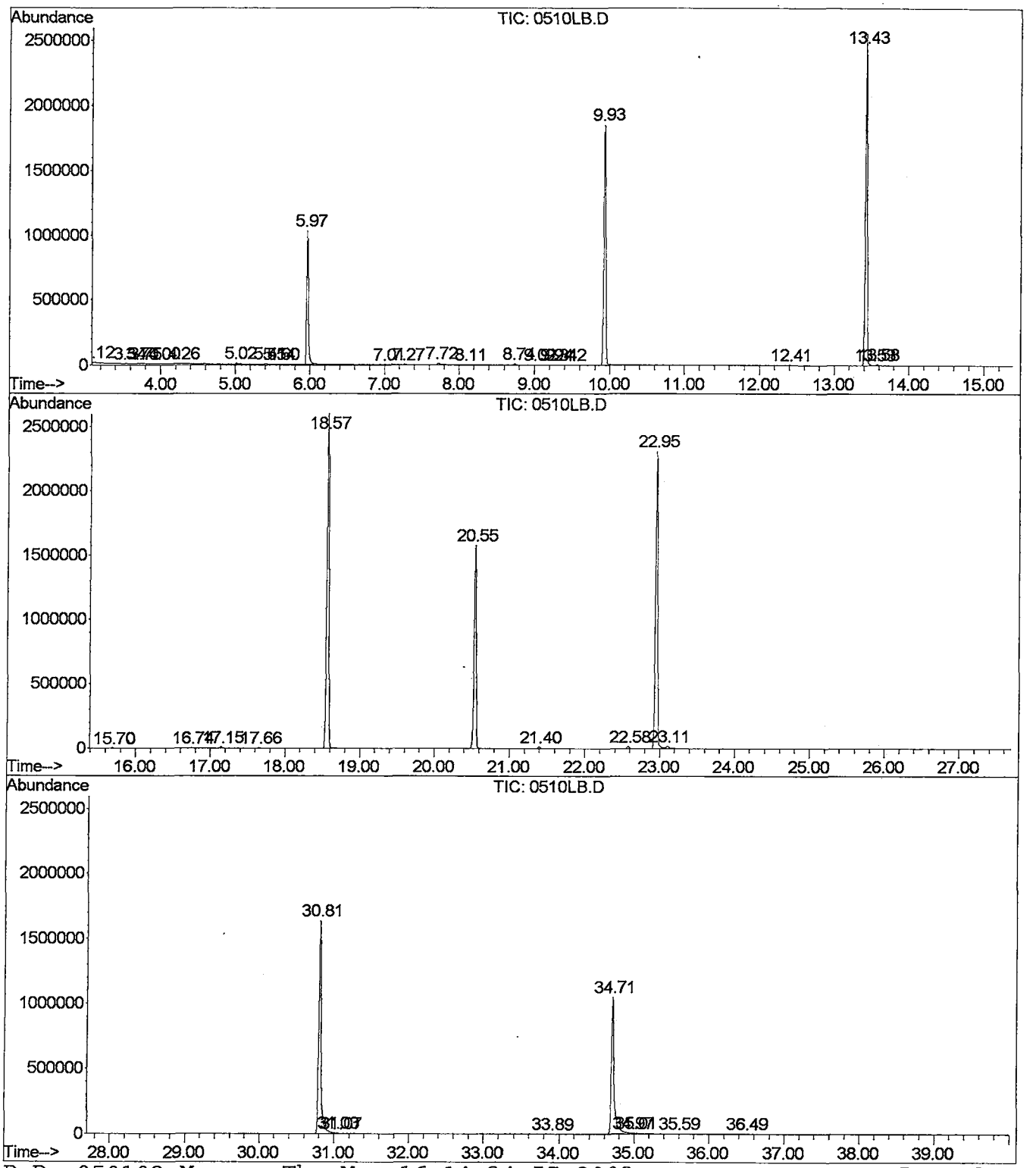
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.120	3	7	23	BV 4	4246	127897	0.25%	0.043%
2	3.543	72	79	88	PV 8	3115	73218	0.15%	0.025%
3	3.702	99	106	109	VV 5	2690	70750	0.14%	0.024%
4	3.749	109	114	118	VV 7	3761	83358	0.17%	0.028%
5	3.996	153	156	162	VV 8	2482	61664	0.12%	0.021%
6	4.254	196	200	206	VV 6	3280	64509	0.13%	0.022%
7	5.018	327	330	346	VV 5	7359	206036	0.41%	0.070%
8	5.412	386	397	411	VV 4	6150	183193	0.36%	0.062%
9	5.541	415	419	423	VV 6	3030	60269	0.12%	0.020%
10	5.594	423	428	434	VV 6	2384	73812	0.15%	0.025%
11	5.964	483	491	518	VV	1011465	17544760	34.83%	5.920%
12	7.010	663	669	676	VV 4	2268	54171	0.11%	0.018%
13	7.274	706	714	721	VV 9	2204	48707	0.10%	0.016%
14	7.721	779	790	812	VV	12368	402233	0.80%	0.136%
15	8.109	849	856	858	PV 6	1789	34457	0.07%	0.012%
16	8.738	955	963	971	VV 2	9016	180743	0.36%	0.061%
17	9.020	1005	1011	1013	VV 5	1940	37098	0.07%	0.013%
18	9.243	1044	1049	1055	VV 4	2602	49066	0.10%	0.017%
19	9.307	1055	1060	1070	VV 4	2839	73191	0.15%	0.025%
20	9.425	1074	1080	1086	PV 5	1685	31487	0.06%	0.011%
21	9.930	1152	1166	1196	PV	1860772	34671714	68.84%	11.699%
22	12.404	1581	1587	1593	VV 2	5010	68315	0.14%	0.023%
23	13.426	1750	1761	1777	PV	2512161	46730819	92.78%	15.768%
24	13.526	1777	1778	1783	VV 3	2458	41400	0.08%	0.014%
25	13.579	1783	1787	1799	VV 2	8463	155630	0.31%	0.053%
26	15.694	2135	2147	2154	PV 3	1661	39682	0.08%	0.013%
27	16.746	2317	2326	2330	PV 3	3514	54658	0.11%	0.018%
28	17.145	2381	2394	2401	PV 3	10157	173567	0.34%	0.059%
29	17.663	2475	2482	2489	VV 2	4309	67741	0.13%	0.023%
30	18.567	2623	2636	2654	PV	2573373	50368192	100.00%	16.996%
31	20.547	2958	2973	2983	BV	1577906	30568290	60.69%	10.315%
32	21.399	3110	3118	3129	PV 3	13244	227373	0.45%	0.077%
33	22.580	3309	3319	3331	VV 2	17958	325698	0.65%	0.110%
34	22.956	3371	3383	3403	PV 2	2289244	47929181	95.16%	16.173%
35	23.109	3403	3409	3421	VV 2	19970	473331	0.94%	0.160%
36	30.812	4706	4720	4756	PV 2	1622877	37428281	74.31%	12.629%
37	31.035	4756	4758	4763	VV 4	4707	86093	0.17%	0.029%

38	31.071	4763	4764	4776	VV	6	2521	63260	0.13%	0.021%
39	33.891	5240	5244	5252	PV	4	1625	24268	0.05%	0.008%
40	34.713	5371	5384	5426	VV	2	1041676	27048324	53.70%	9.127%
41	34.966	5426	5427	5432	VV	5	6956	123006	0.24%	0.042%
42	35.007	5432	5434	5443	VV	6	5307	166088	0.33%	0.056%
43	35.589	5528	5533	5534	VV	4	2085	29511	0.06%	0.010%
44	36.488	5682	5686	5693	VV	5	2225	36449	0.07%	0.012%

Sum of corrected areas: 296361487

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\051402\0510LB.D
Operator : Mclean
Acquired : 14 May 2002 3:07 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 5/10 kb
Misc Info :
Vial Number: 3
Quant File :050102.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

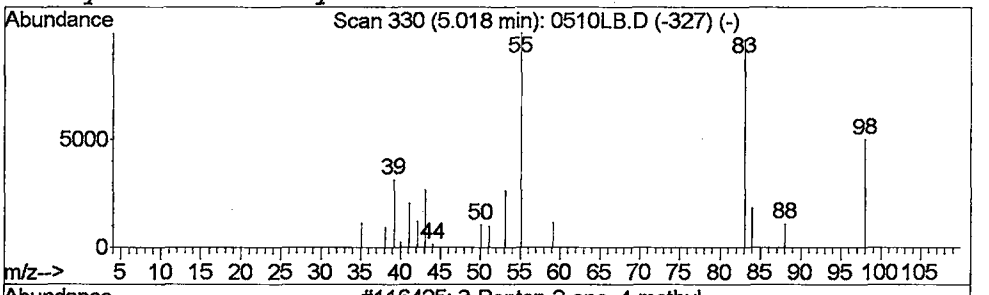
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

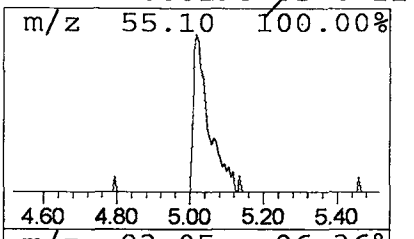
****

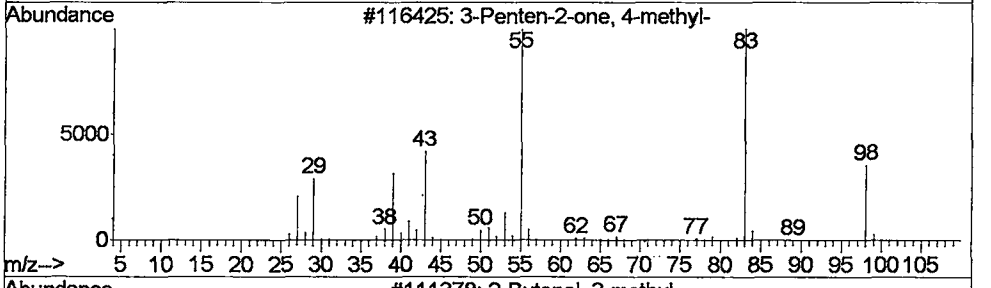
Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 6

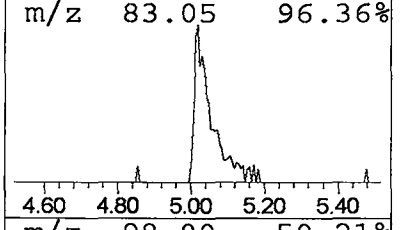
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	0.24 ug/L	206036	1,4-Dichlorobenzene-d4	9.93

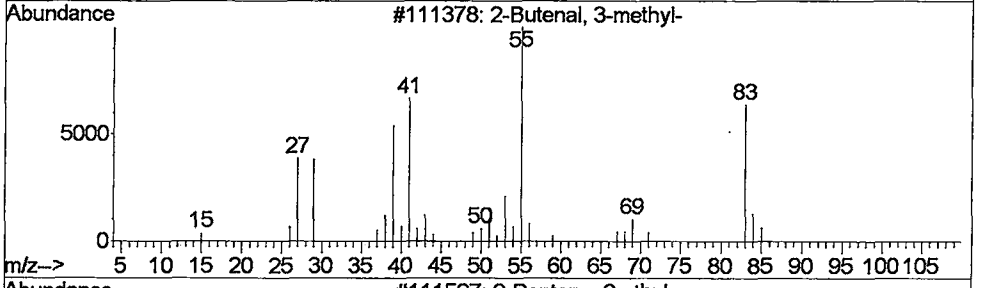
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	59
2	2-Butenal, 3-methyl-	84	C5H8O	000107-86-8	47
3	2-Pentene, 3-ethyl-	98	C7H14	000816-79-5	35
4	Cyclobut-1-enylmethanol	84	C5H8O	1000195-03-0	22

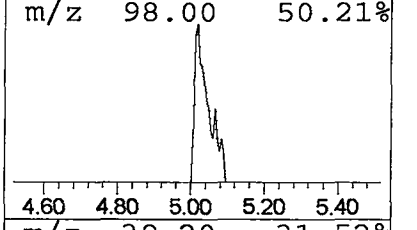


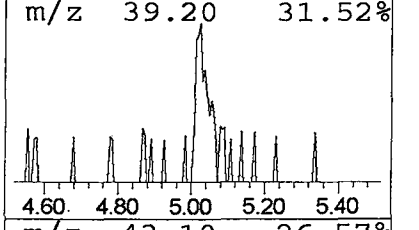


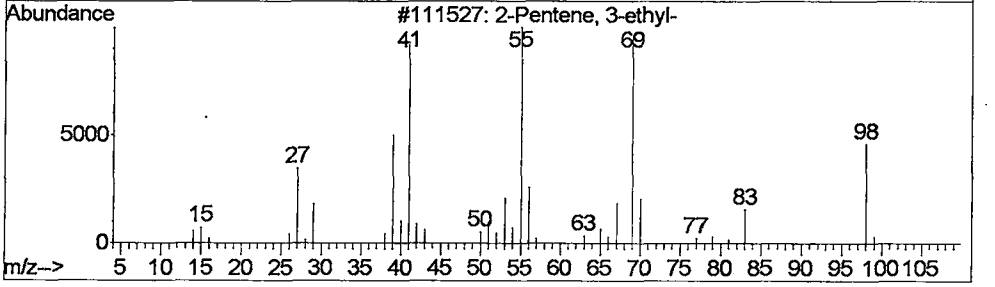


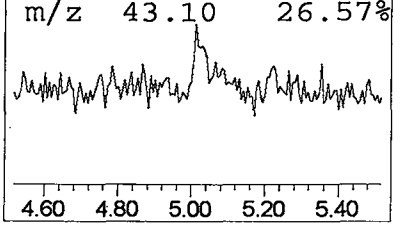












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

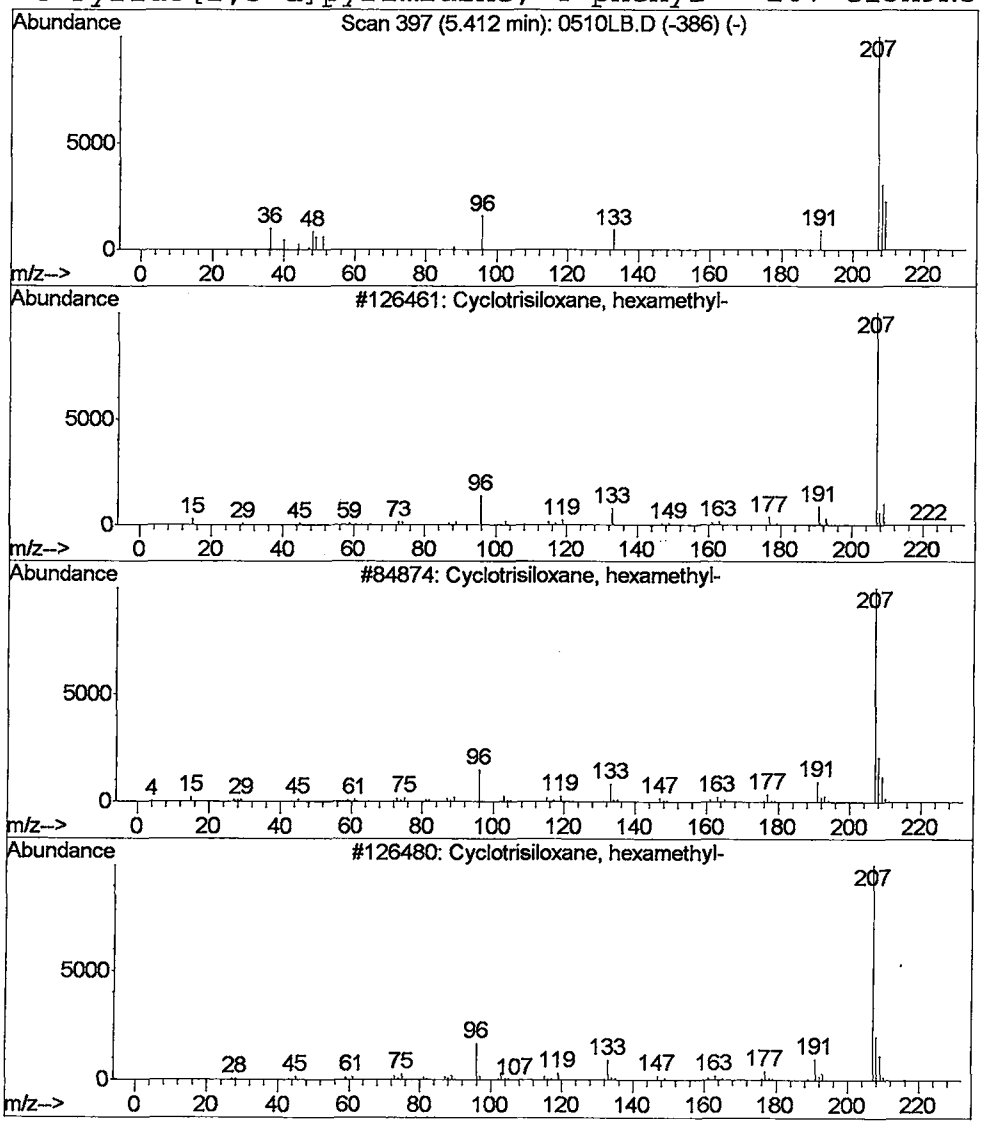
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

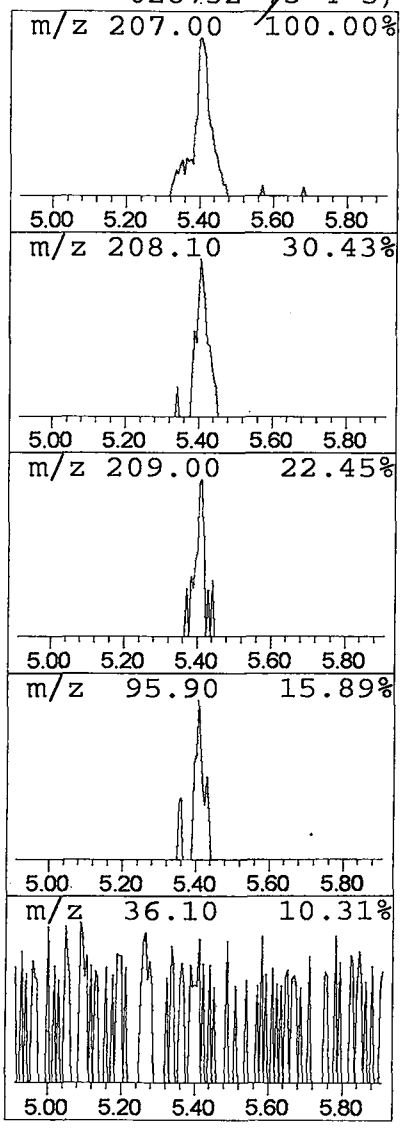
****

Peak Number 2 Cyclotrisiloxane, hexamethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.41	0.21 ug/L	183193	1,4-Dichlorobenzene-d4	9.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	9
2	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	9
3	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	9
4	Pyrido[2,3-d]pyrimidine, 4-phenyl-	207	C13H9N3	028732-75-4	5





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

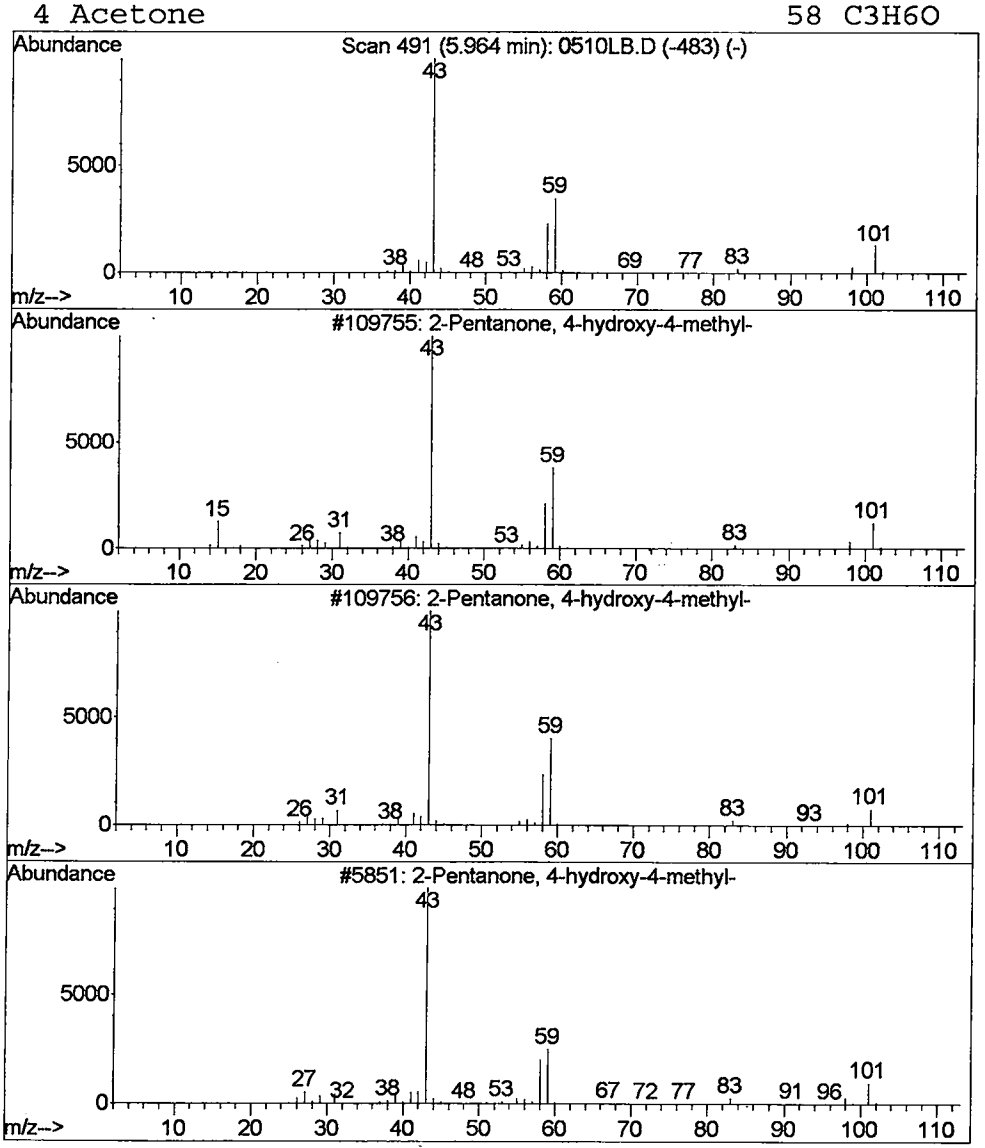
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

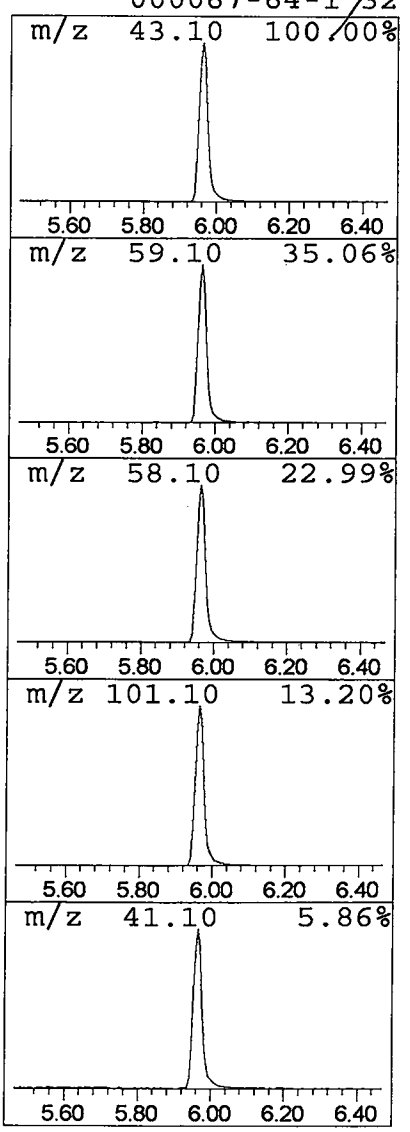
****

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.97	20.24 ug/L	17544800	1,4-Dichlorobenzene-d4	9.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37
4	Acetone	58	C3H6O	000067-64-1	32





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

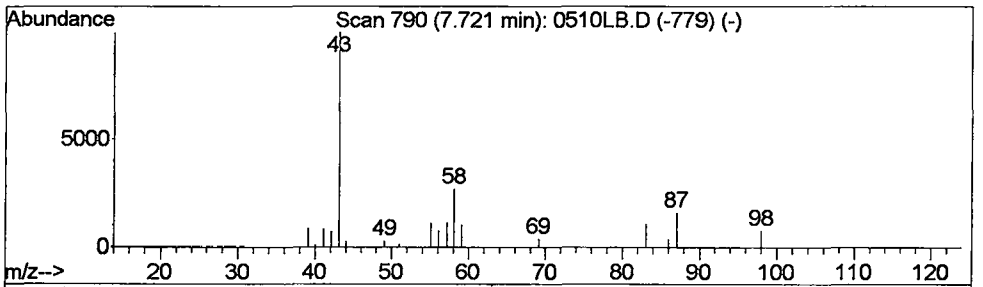
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

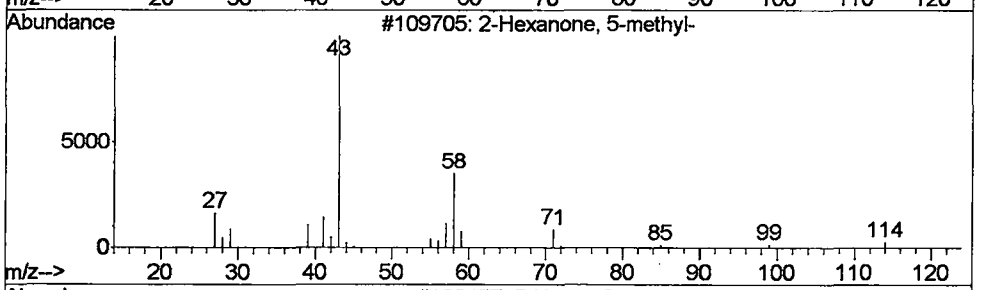
****

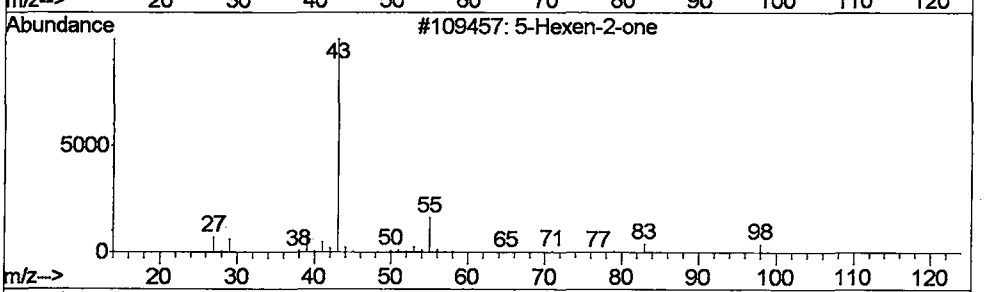
Peak Number 4 2-Hexanone, 5-methyl- Concentration Rank 2

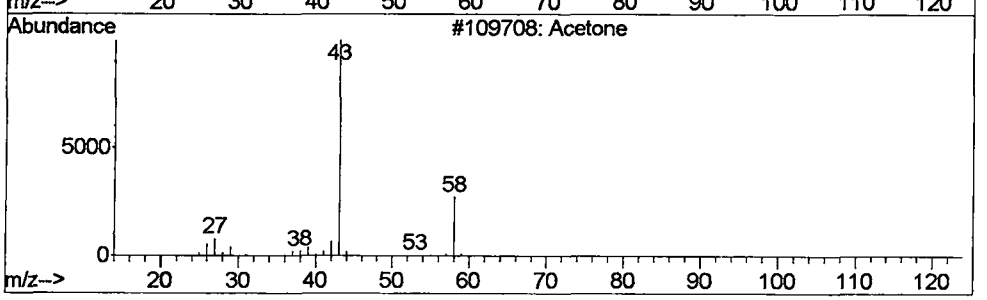
R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	0.46 ug/L	402233	1,4-Dichlorobenzene-d4	9.93

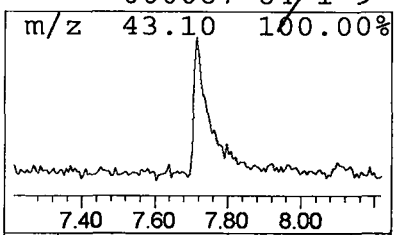
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	28
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9
3	Acetone	58	C3H6O	000067-64-1	9
4	Acetone	58	C3H6O	000067-64-1	9

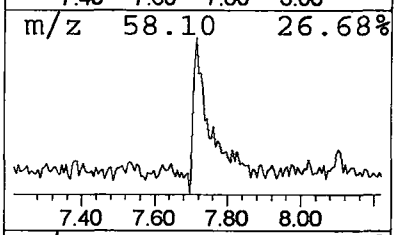


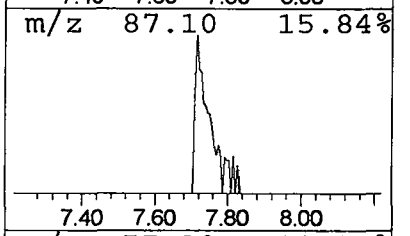


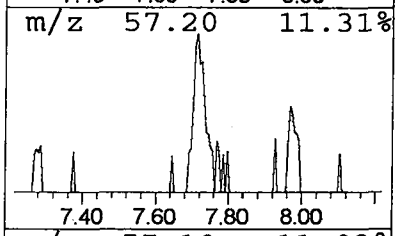


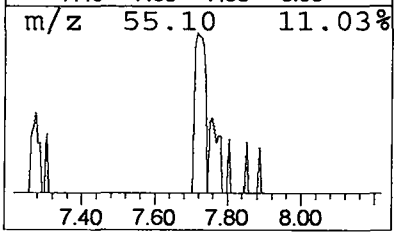












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

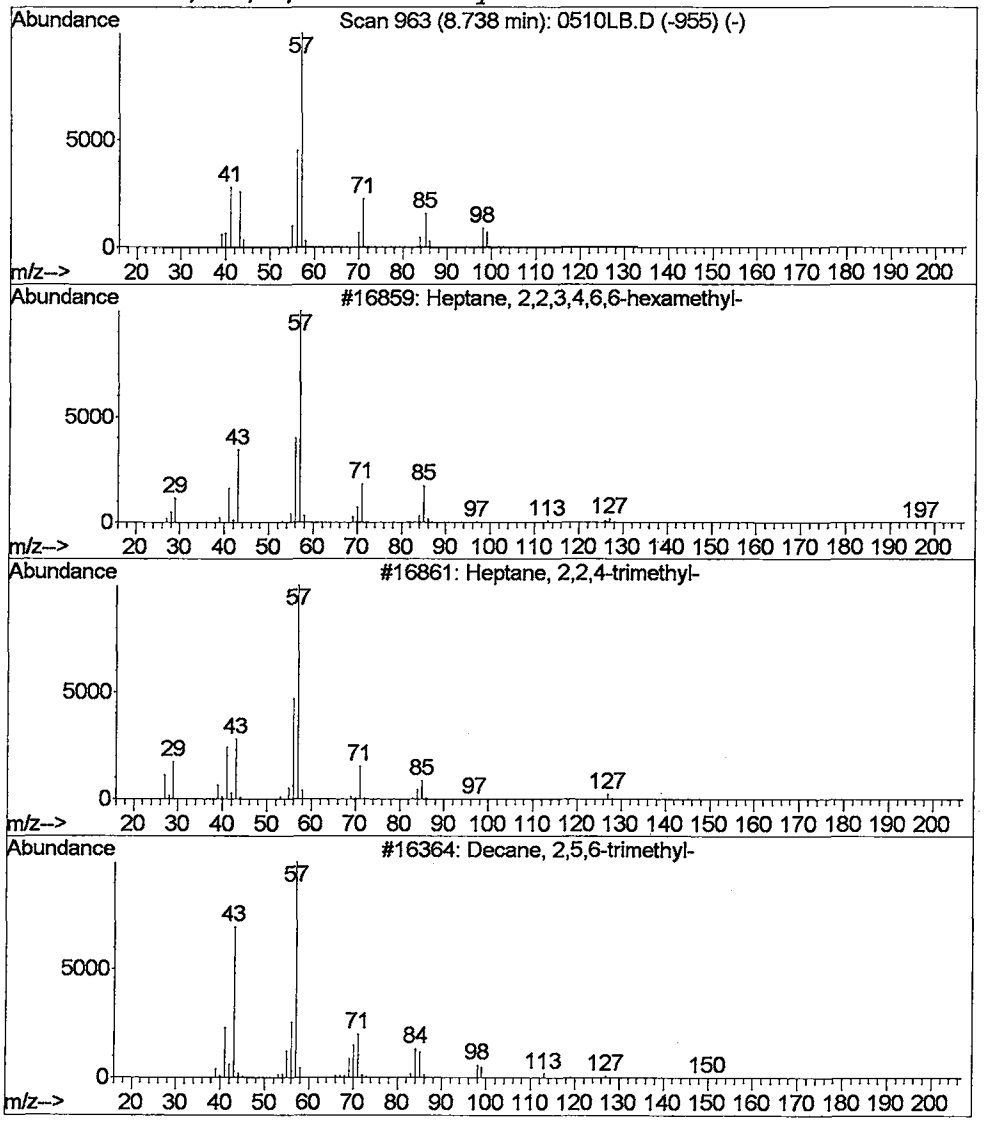
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

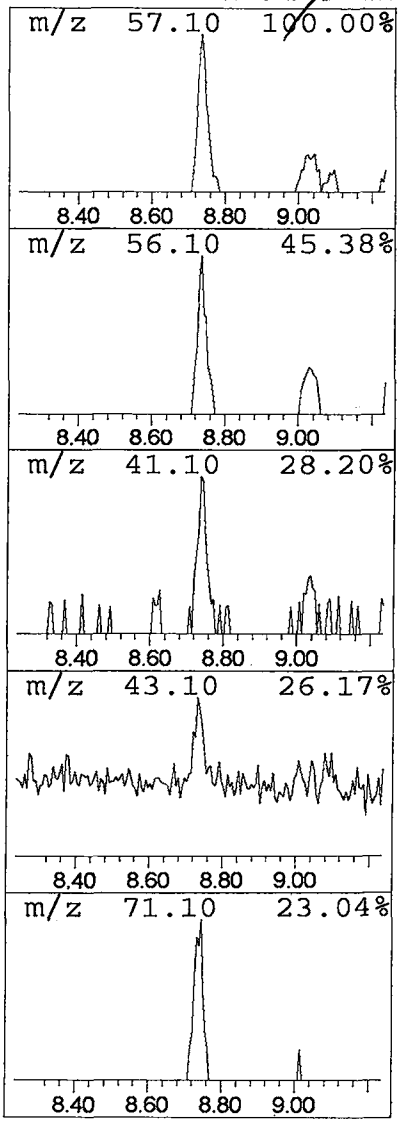
****

Peak Number 5 Heptane, 2,2,3,4,6,6-hexame... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.74	0.21 ug/L	180743	1,4-Dichlorobenzene-d4	9.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,2,3,4,6,6-hexamethyl-	184	C13H28	062108-32-1	64
2	Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	42
3	Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	42
4	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	42





Library Search Compound Report

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Acq On : 14 May 2002 3:07 pm Operator: Mclean
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Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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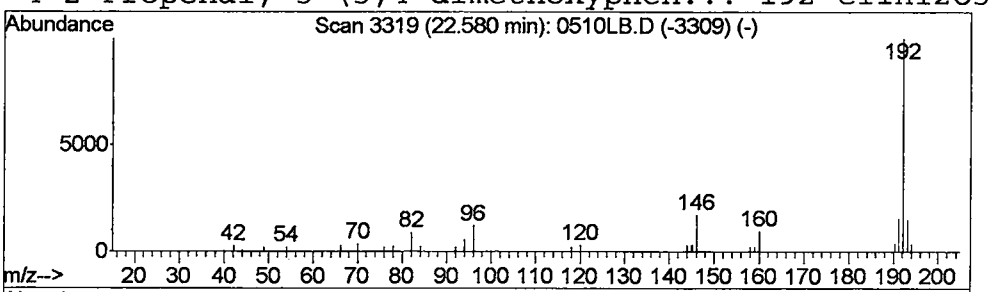
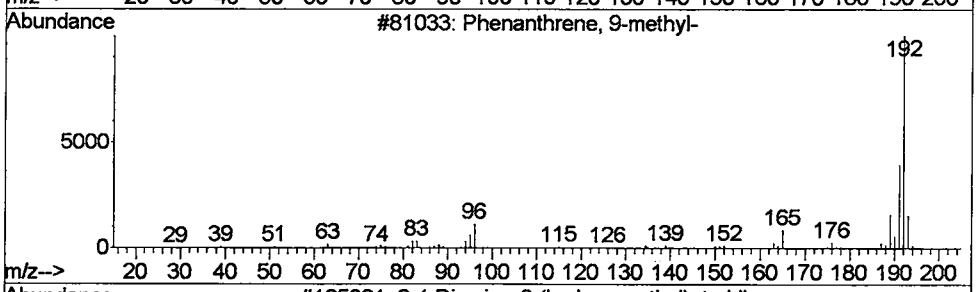
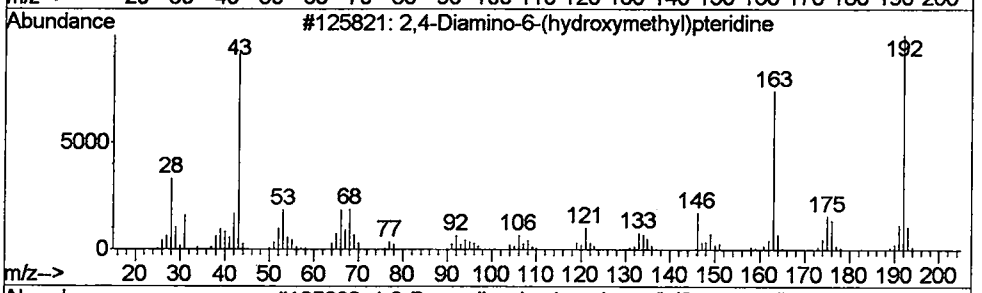
Peak Number 6 Phenanthrene, 9-methyl- Concentration Rank 4

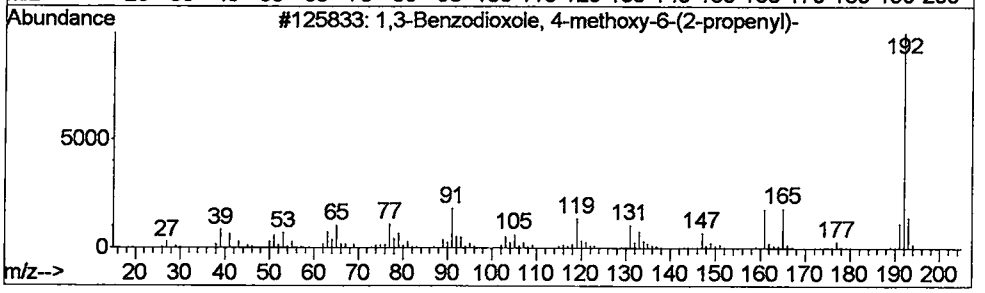
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	0.27 ug/L	325698	Phenanthranene-d10	22.95

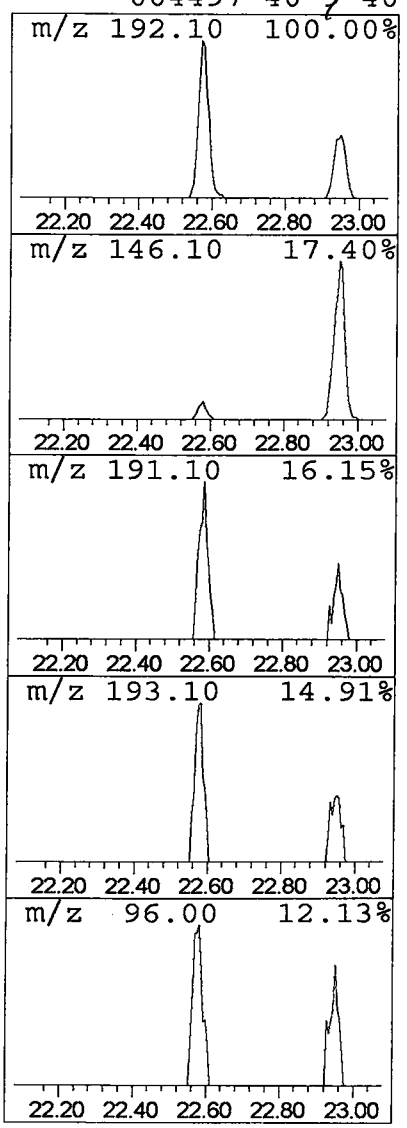
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	53
2	2,4-Diamino-6-(hydroxymethyl)pte...	192	C7H8N6O	000945-24-4	53
3	1,3-Benzodioxole, 4-methoxy-6-(2...	192	C11H12O3	000607-91-0	40
4	2-Propenal, 3-(3,4-dimethoxyphen...	192	C11H12O3	004497-40-9	40











Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

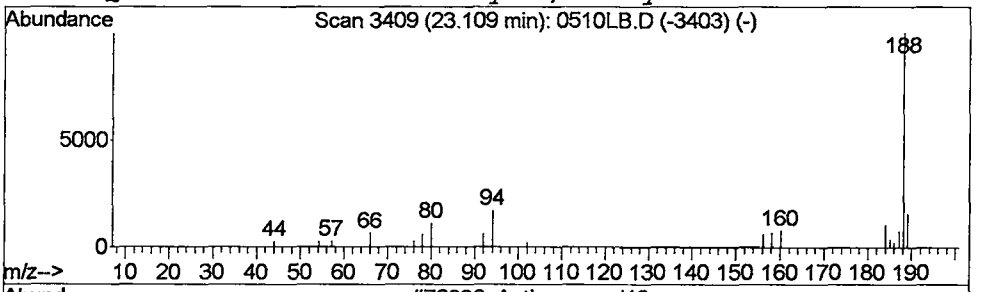
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

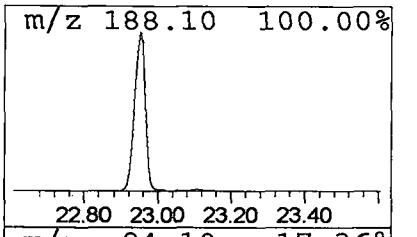
****

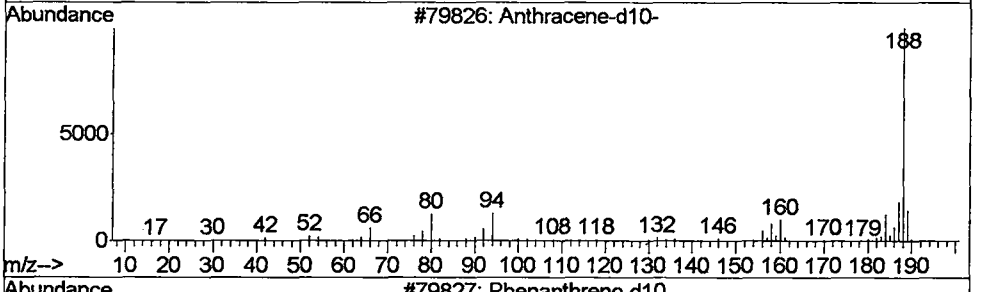
Peak Number 7 Anthracene-d10- Concentration Rank 3

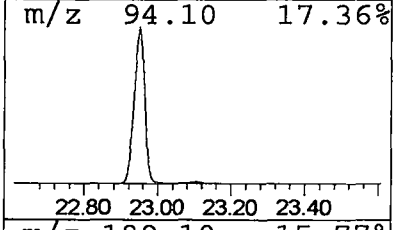
R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	0.40 ug/L	473331	Phenanthranene-d10	22.95

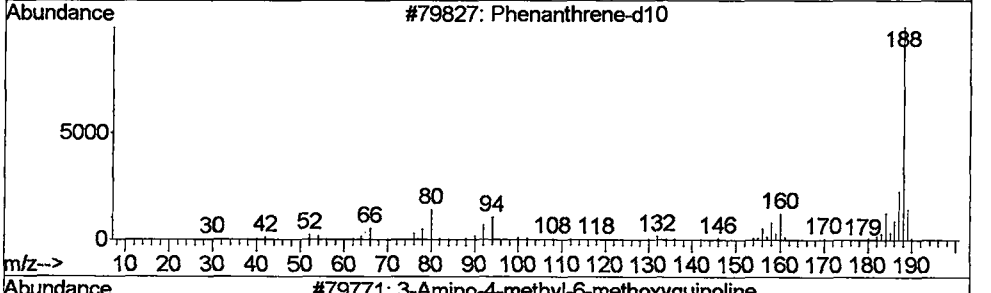
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10-	188	C14D10	001719-06-8	91
2	Phenanthrene-d10	188	C14D10	001517-22-2	86
3	3-Amino-4-methyl-6-methoxyquinoline	188	C11H12N2O	1000213-71-3	42
4	2-Quinolinecarboxaldehyde, 8-hyd...	188	C10H8N2O2	005603-22-5	40

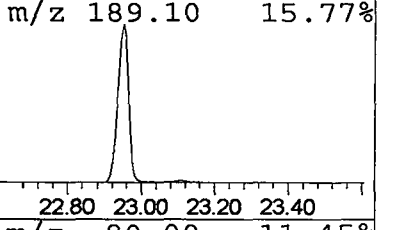


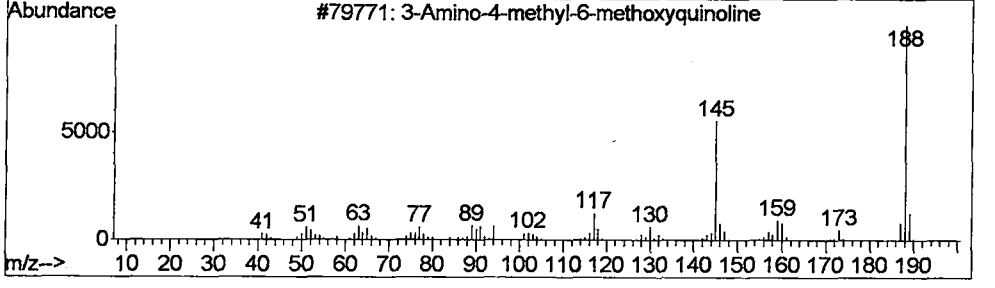


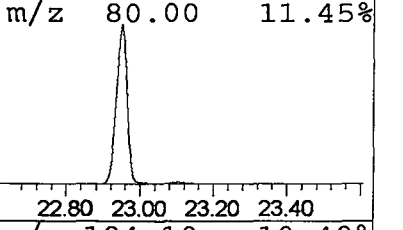


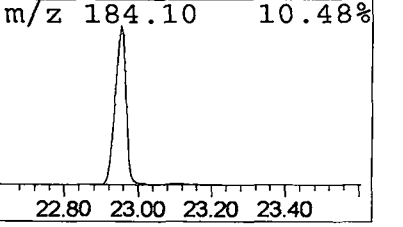












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\051402\0510LB.D Vial: 3
Acq On : 14 May 2002 3:07 pm Operator: Mclean
Sample : 5/10 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

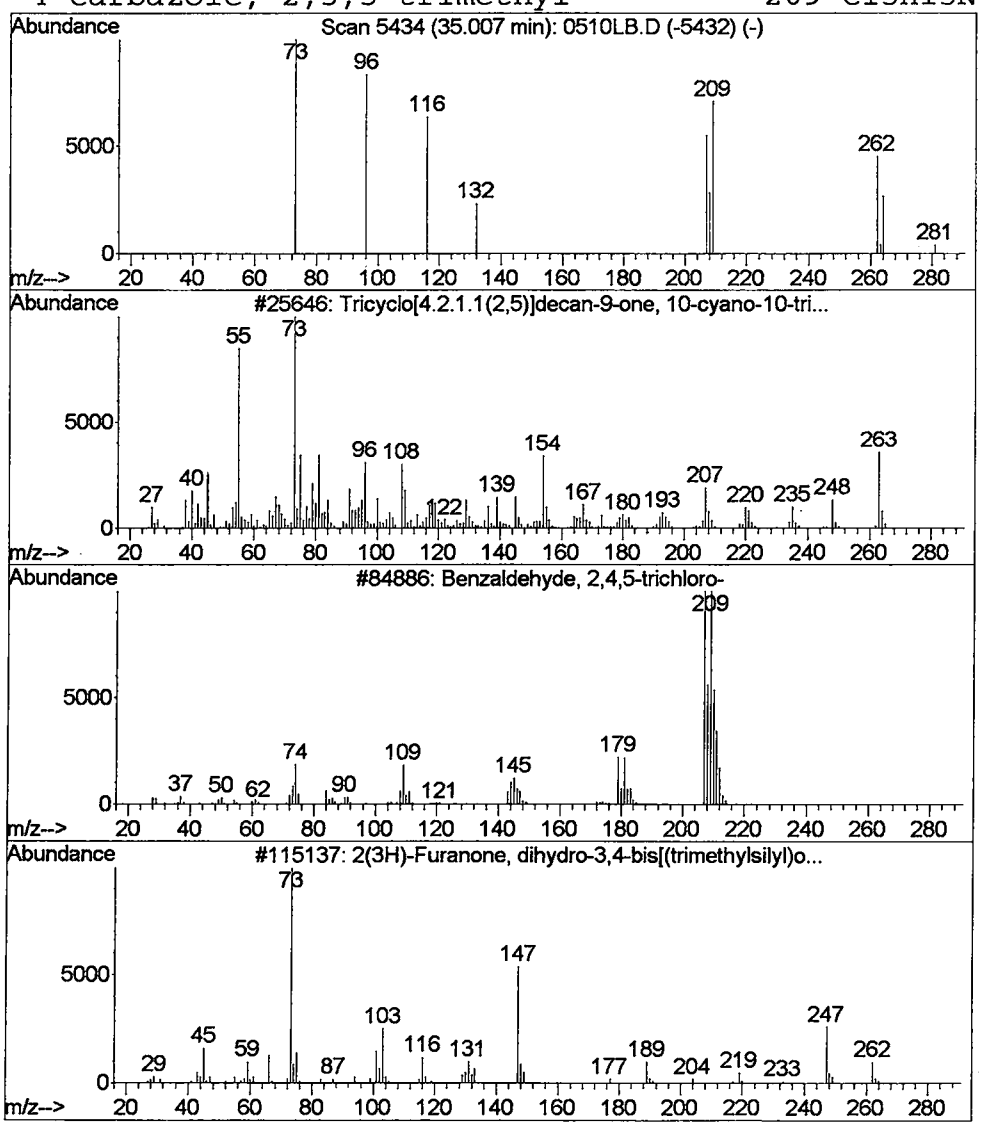
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

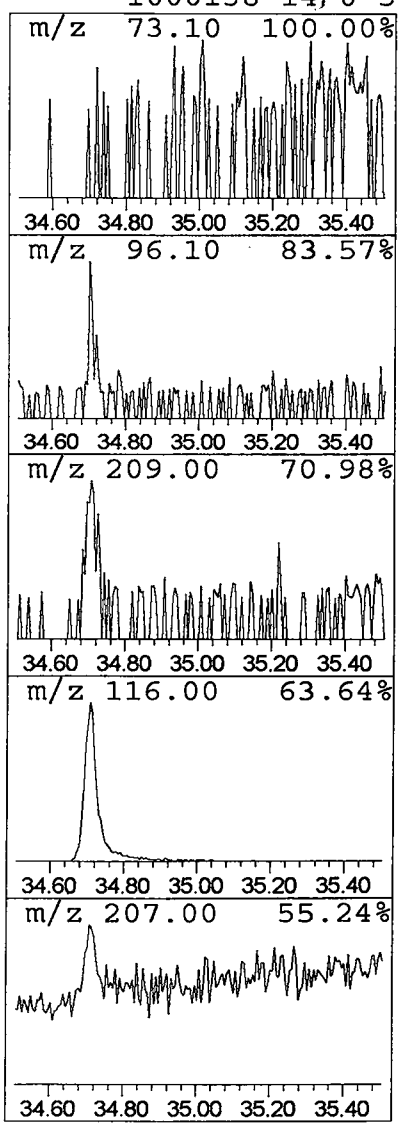
****

Peak Number 8 Tricyclo[4.2.1.1(2,5)]decan... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.01	0.25 ug/L	166088	Perylene-d12	34.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[4.2.1.1(2,5)]decan-9-on...	263	C14H21NO2Si	1000193-78-5	9
2	Benzaldehyde, 2,4,5-trichloro-	208	C7H3Cl3O	035696-87-8	9
3	2(3H)-Furanone, dihydro-3,4-bis[...	262	C10H22O4Si2	055220-75-2	7
4	Carbazole, 2,3,5-trimethyl-	209	C15H15N	1000138-14-0	5





Tentatively Identified Compound (LSC) summary

Operator ID: Mclean Date Acquired: 14 May 2002 3:07 pm
Data File: C:\MSDCHEM\1\DATA\051402\0510LB.D
Name: 5/10 kb
Misc:
Method: C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title: 508 Modified GC/MS scan
Library Searched: C:\DATABASE\NIST98.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--- #	RT	Resp	Conc
3-Penten-2-one, 4...	5.02	0.2	ug/L	206036	1	9.93	34671700	40.0
Cyclotrisiloxane,...	5.41	0.2	ug/L	183193	1	9.93	34671700	40.0
2-Pentanone, 4-hy...	5.97	20.2	ug/L	17544800	1	9.93	34671700	40.0
2-Hexanone, 5-met...	7.72	0.5	ug/L	402233	1	9.93	34671700	40.0
Heptane, 2,2,3,4,...	8.74	0.2	ug/L	180743	1	9.93	34671700	40.0
Phenanthrene, 9-m...	22.58	0.3	ug/L	325698	4	22.95	47929200	40.0
Anthracene-d10-	23.11	0.4	ug/L	473331	4	22.95	47929200	40.0
Tricyclo[4.2.1.1(...	35.01	0.2	ug/L	166088	6	34.71	27048300	40.0