

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 05/15/2002 Time: 15:01:46

Sample: AE09332
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 5/15/02 15:01 Ended: 5/15/02 15:01 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		80		10.0

Comments for Sample AE09332 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 05-15-2002 Time: 15:02:00

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Recoveries for 4 of the 43 compounds in the LCS Standard were low.

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\050902\9332.D

Acq On : 10 May 2002 9:13 am

Sample : 5/8 eh

Misc :

MS Integration Params: EVENTS.E

Quant Time: May 13 11:42:55 2002

Vial: 22

Operator: Mclean

Inst : Instrumen

Multiplr: 1.00

Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)

Title : 508 Modified GC/MS scan

Last Update : Mon May 13 11:40:29 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

*Ne-lyt
buthics*

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.94	152	6466438	40.00	ug/L	0.00
10) Napthalene-d8	13.43	136	25687809	40.00	ug/L	0.00
17) Acenapthalene-d10	18.57	164	13972835	40.00	ug/L	0.00
29) Phenanthranene-d10	22.95	188	20324526	40.00	ug/L	0.00
37) Chrysene-d12	30.81	240	13513526	40.00	ug/L	0.00
43) Perylene-d12	34.70	264	7324137	40.00	ug/L	0.00

System Monitoring Compounds

27) TCMX	20.55	209	2730778	32.19	ug/L	0.00
Spiked Amount	40.000		Recovery	=	80.47%	

Target Compounds

Qvalue

2) n-nitros-dimethyl amine	0.00	74	0	N.D.	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-Chloronapthalene	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	14247	N.D.	
26) Fluorene	0.00	166	0	N.D.	
28) Azobenzene	20.54	77	46210	N.D.	
30) Nnitrosodiphenylamine	20.54	169	51026	N.D.	
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

9332.D 050102.M Mon May 13 11:43:35 2002

Page 1

Quantitation Report

(QT Reviewed)

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Acq On : 10 May 2002 9:13 am

Sample : 5/8 eh

Misc :

MS Integration Params: EVENTS.E

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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	35612	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
9332.D 050102.M Mon May 13 11:43:35 2002 Page 2

Quantitation Report (QT Reviewed)

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Vial: 22

Acq On : 10 May 2002 9:13 am

Operator: Mclean

Sample : 5/8 eh

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: EVENTS.E

Quant Time: May 13 11:43 2002

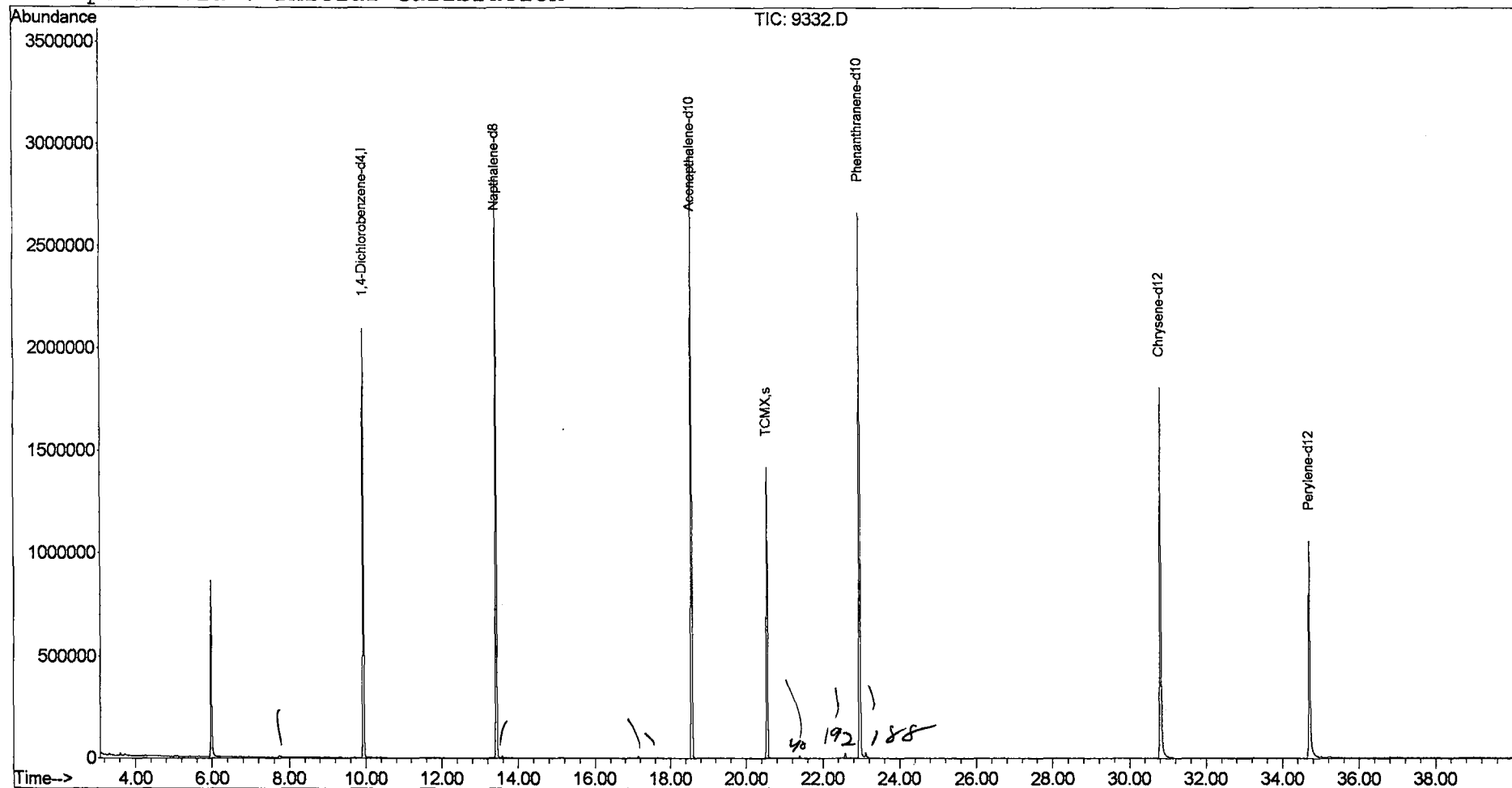
Quant Results File: 050102.RES

Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)

Title : 508 Modified GC/MS scan

Last Update : Mon May 13 11:40:29 2002

Response via : Initial Calibration



9332.D 050102.M

Mon May 13 11:43:35 2002

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LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\050902\9332.D
 Acq On : 10 May 2002 9:13 am
 Sample : 5/8 eh
 Misc :
 MS Integration Params: LSCINT.e

Vial: 22
 Operator: Mclean
 Inst : Instrumen
 Multiplr: 1.00

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

Signal : TIC

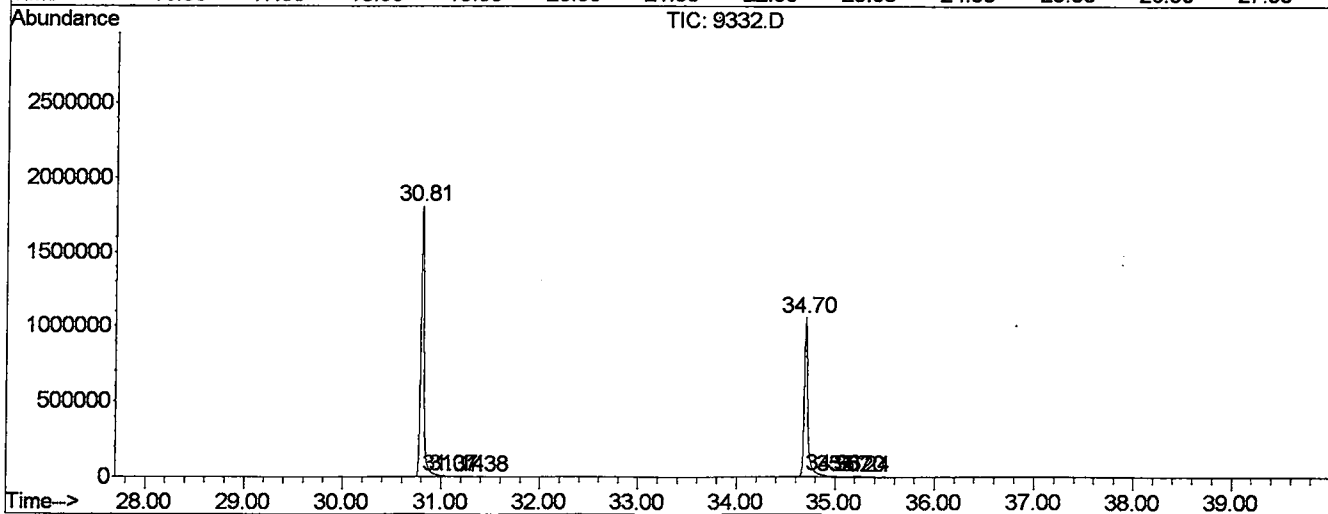
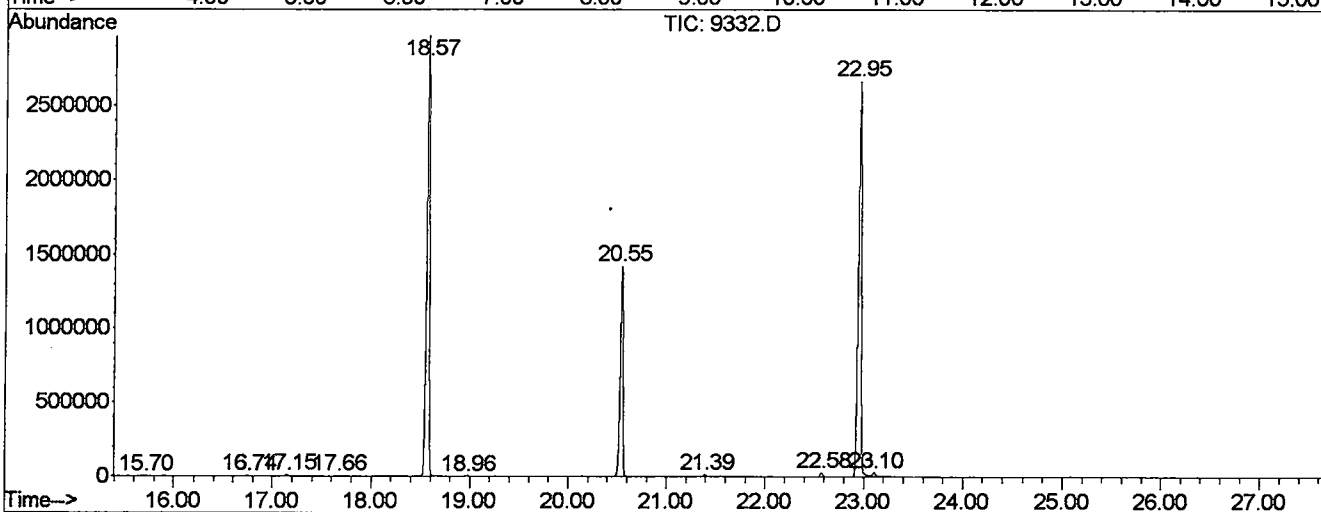
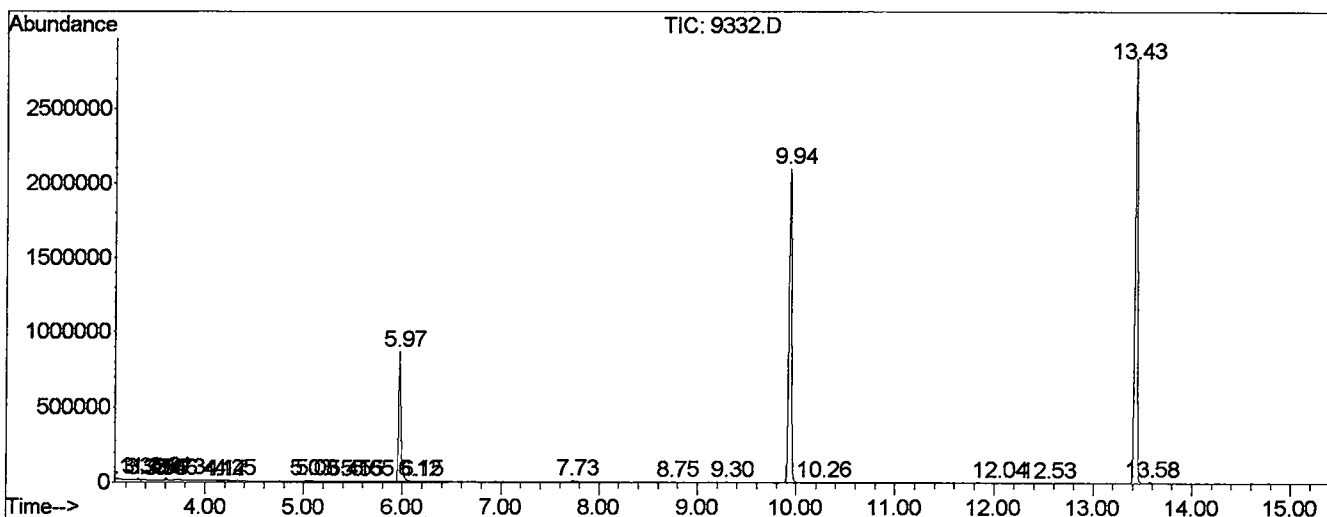
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1	3.114	3	6	16	BV 3	4182	76768	0.13%	0.024%
2	3.326	34	42	47	PV 2	9752	148289	0.26%	0.046%
3	3.379	47	51	58	VV 9	4122	81996	0.14%	0.025%
4	3.538	72	78	79	PV 6	2909	36617	0.06%	0.011%
5	3.561	79	82	84	VV 4	2580	34733	0.06%	0.011%
6	3.608	84	90	97	VV 2	16146	285417	0.50%	0.089%
7	3.661	97	99	100	VV 2	2915	26996	0.05%	0.008%
8	3.731	100	111	126	VV 4	9165	386379	0.67%	0.120%
9	4.143	177	181	183	VV 4	2208	29844	0.05%	0.009%
10	4.166	183	185	191	VV 5	2428	46122	0.08%	0.014%
11	4.254	195	200	207	VV 7	3378	72147	0.13%	0.022%
12	5.030	326	332	338	PV 7	5997	134027	0.23%	0.042%
13	5.083	338	341	363	VV 9	4813	121264	0.21%	0.038%
14	5.412	393	397	405	VV 4	2741	63376	0.11%	0.020%
15	5.547	417	420	422	PV 3	1749	16942	0.03%	0.005%
16	5.653	430	438	448	VV 9	2839	102683	0.18%	0.032%
17	5.970	484	492	517	PV	839786	15010822	26.15%	4.666%
18	6.123	517	518	520	VV 2	2315	21151	0.04%	0.007%
19	6.146	520	522	532	VV 7	2392	38567	0.07%	0.012%
20	7.727	787	791	807	PV 2	8896	249164	0.43%	0.077%
21	8.749	960	965	971	VV 5	2465	41369	0.07%	0.013%
22	9.301	1052	1059	1078	VV 6	2656	53763	0.09%	0.017%
23	9.936	1158	1167	1186	PV	2078859	39273898	68.41%	12.208%
24	10.265	1217	1223	1230	PV 5	2063	37732	0.07%	0.012%
25	12.039	1520	1525	1533	VV 4	2224	47075	0.08%	0.015%
26	12.527	1604	1608	1613	VV 4	2188	37215	0.06%	0.012%
27	13.432	1747	1762	1783	PV	2852381	52834018	92.03%	16.424%
28	13.579	1783	1787	1799	VV 2	9405	176193	0.31%	0.055%
29	15.700	2143	2148	2154	VV 4	2515	48403	0.08%	0.015%
30	16.740	2316	2325	2330	PV	4844	83799	0.15%	0.026%
31	17.151	2382	2395	2400	VV 4	9209	158642	0.28%	0.049%
32	17.657	2475	2481	2487	VV 3	4132	61953	0.11%	0.019%
33	18.567	2625	2636	2659	VV	2926055	57407978	100.00%	17.845%
34	18.967	2698	2704	2708	PV 3	1979	47520	0.08%	0.015%
35	20.547	2960	2973	2984	VV	1407650	27387327	47.71%	8.513%
36	21.393	3109	3117	3124	VV 3	13039	210508	0.37%	0.065%
37	22.580	3312	3319	3336	VV	24356	439453	0.77%	0.137%

38	22.956	3367	3383	3402	PV	2	2639402	55750569	97.11%	17.330%
39	23.103	3402	3408	3421	VV	2	26427	608229	1.06%	0.189%
40	30.806	4707	4719	4761	PV	2	1808026	42658190	74.31%	13.260%
41	31.065	4761	4763	4770	VV	6	4397	100870	0.18%	0.031%
42	31.135	4770	4775	4782	VV	5	2887	92194	0.16%	0.029%
43	31.382	4805	4817	4823	PV	5	1429	30010	0.05%	0.009%
44	34.702	5367	5382	5424	PV		1045860	26819013	46.72%	8.337%
45	34.960	5424	5426	5442	VV	9	4805	179639	0.31%	0.056%
46	35.066	5442	5444	5446	VV	3	1468	13344	0.02%	0.004%
47	35.201	5458	5467	5473	VV	7	2099	77907	0.14%	0.024%
48	35.242	5473	5474	5480	VV	3	2158	34886	0.06%	0.011%

Sum of corrected areas: 321695002

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\050902\9332.D
 Operator : Mclean
 Acquired : 10 May 2002 9:13 am using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: 5/8 eh
 Misc Info :
 Vial Number: 22
 Quant File :050102.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9332.D
Acq On : 10 May 2002 9:13 am
Sample : 5/8 eh
Misc :
MS Integration Params: LSCINT.e

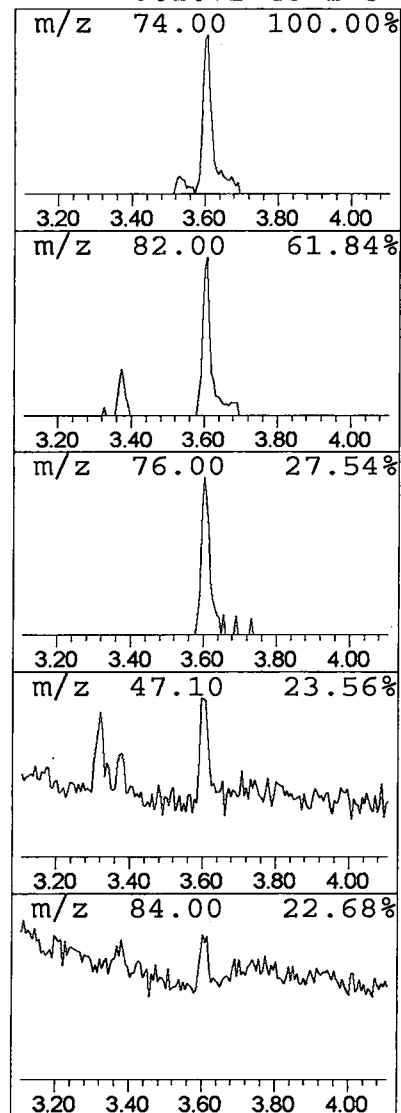
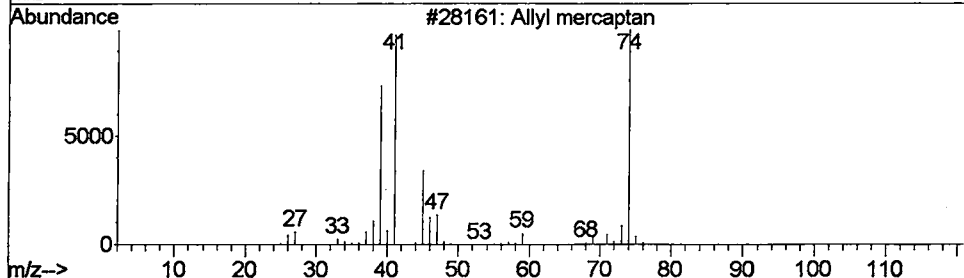
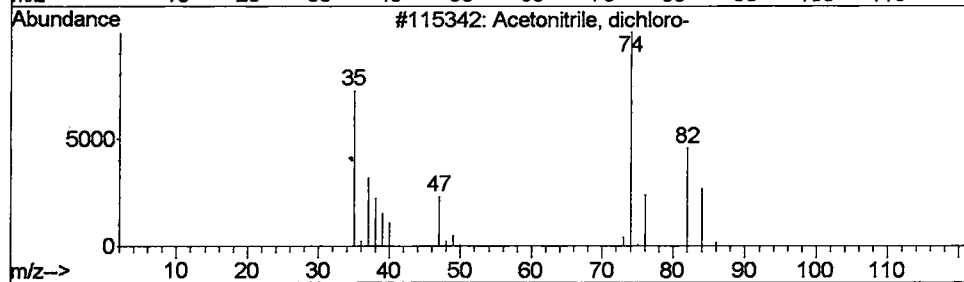
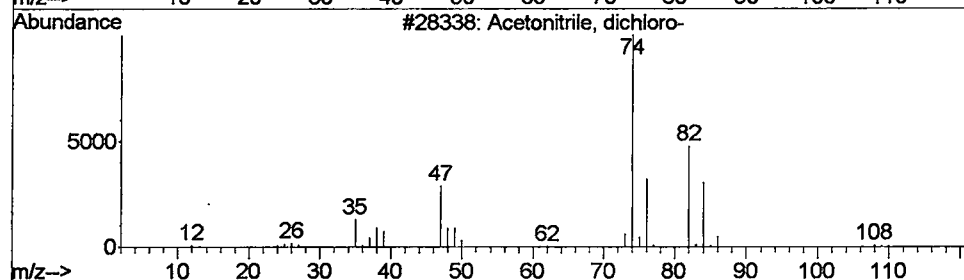
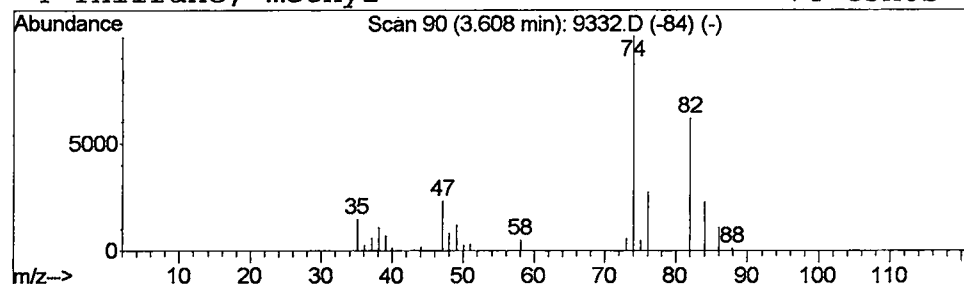
Vial: 22
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 1 Acetonitrile, dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.29 ug/L	285417	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	80
2	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	50
3	Allyl mercaptan	74	C3H6S	000870-23-5	25
4	Thiirane, methyl-	74	C3H6S	001072-43-1	5



Library Search Compound Report

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Acq On : 10 May 2002 9:13 am
Sample : 5/8 eh
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MS Integration Params: LSCINT.e

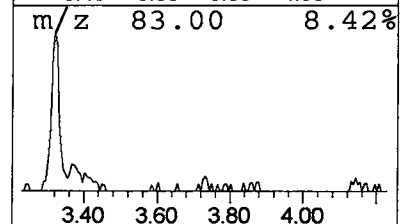
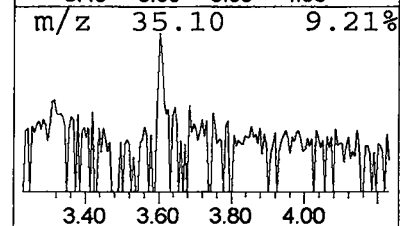
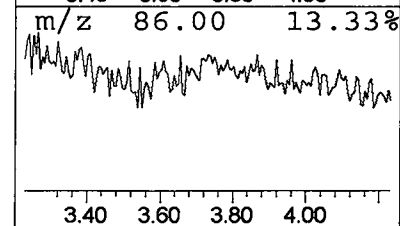
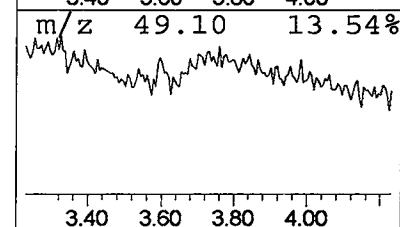
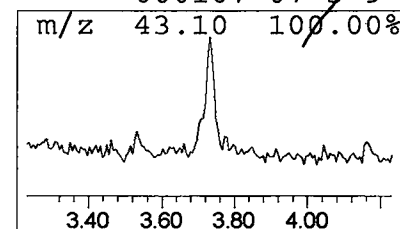
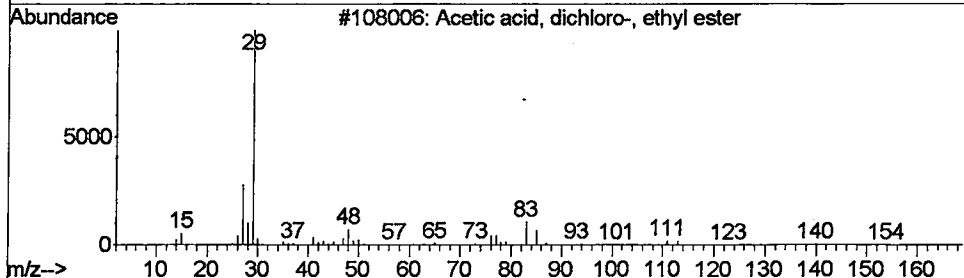
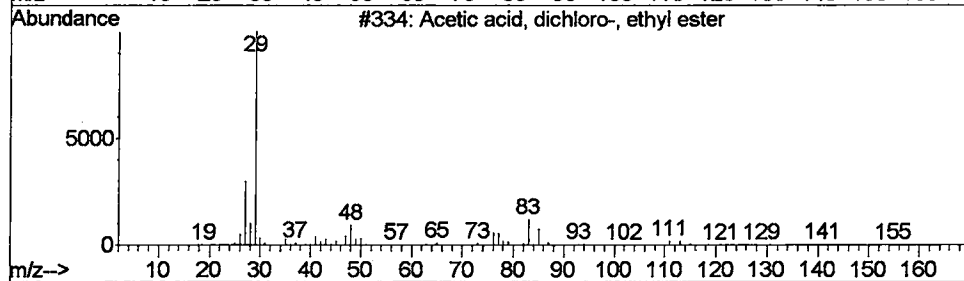
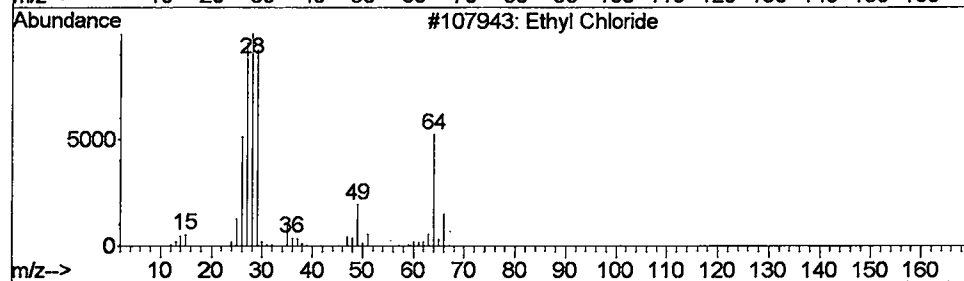
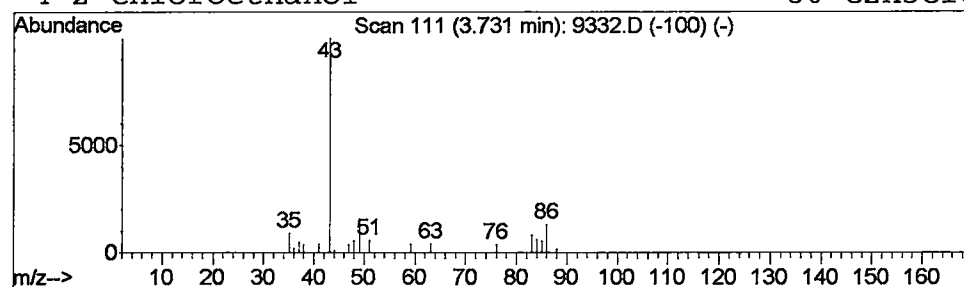
Vial: 22
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 2 Ethyl Chloride Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.73	0.39 ug/L	386379	1,4-Dichlorobenzene-d4	9.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl Chloride	64	C2H5Cl	000075-00-3	12
2			Acetic acid, dichloro-, ethyl ester	156	C4H6Cl2O2	000535-15-9	9
3			Acetic acid, dichloro-, ethyl ester	156	C4H6Cl2O2	000535-15-9	9
4			2-Chloroethanol	80	C2H5ClO	000107-07-3	9



Library Search Compound Report

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Sample : 5/8 eh
Misc :
MS Integration Params: LSCINT.e

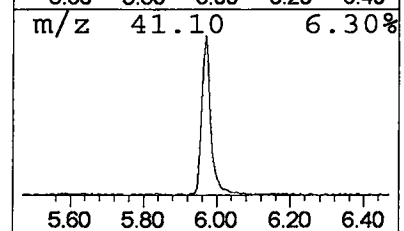
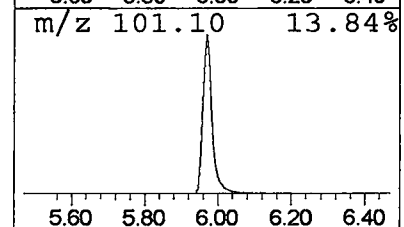
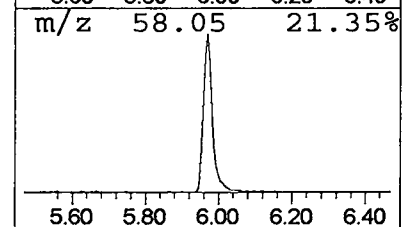
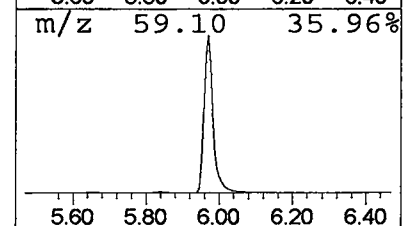
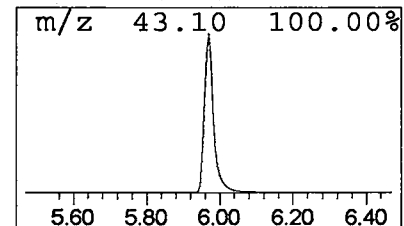
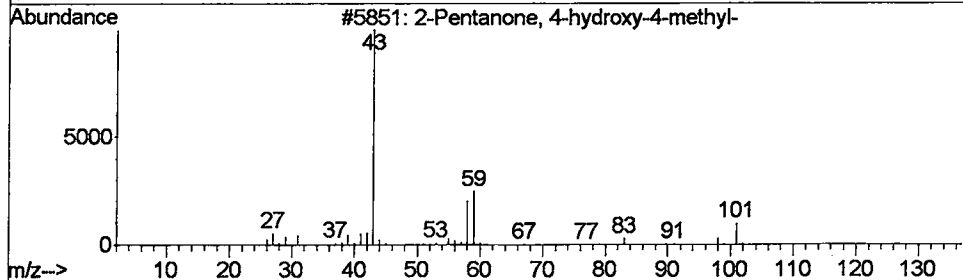
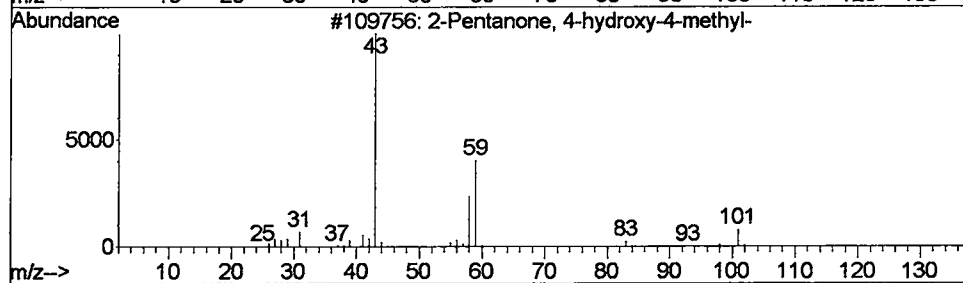
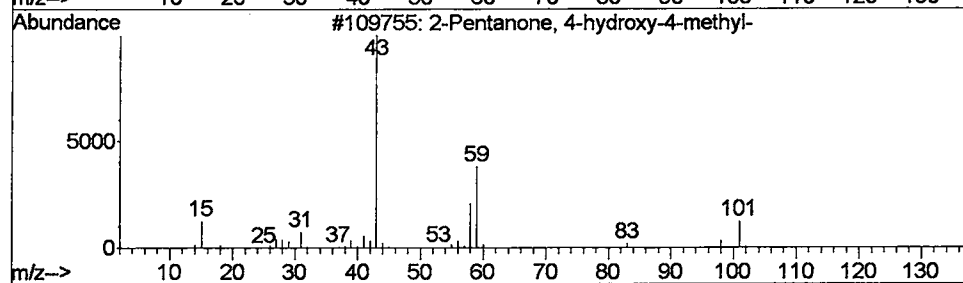
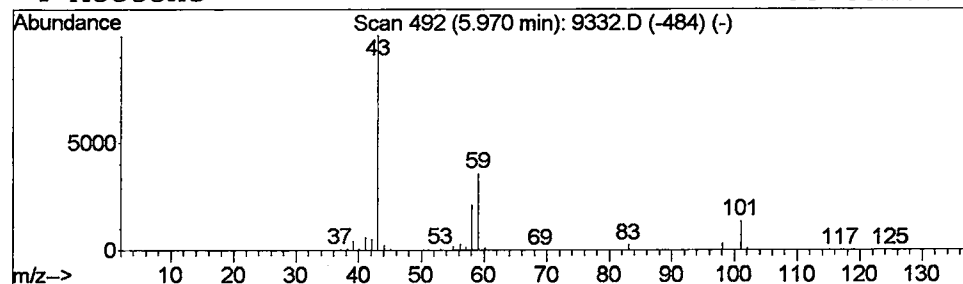
Vial: 22
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

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	15.29 ug/L	15010800	1,4-Dichlorobenzene-d4	9.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	90
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	64
4			Acetone	58	C3H6O	000067-64-1	32



Library Search Compound Report

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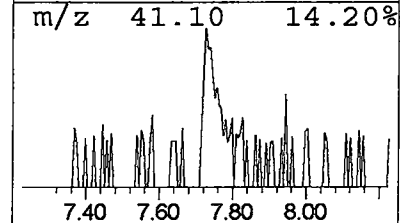
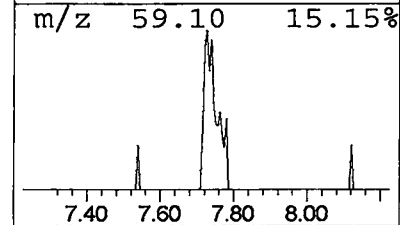
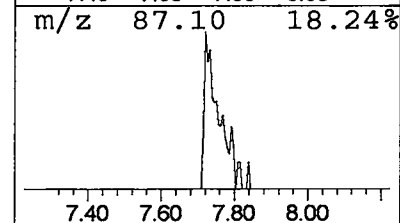
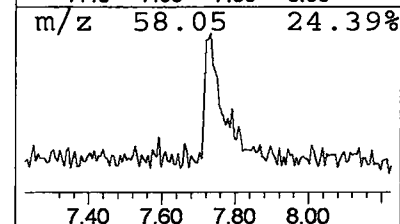
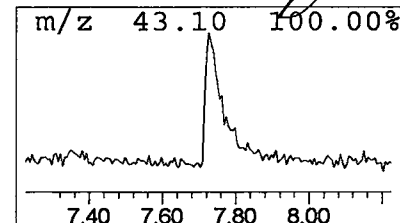
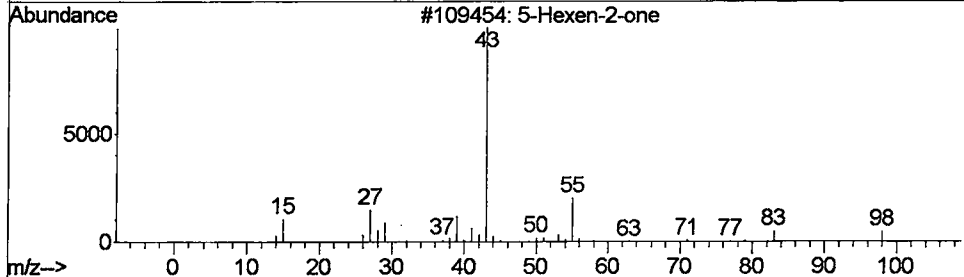
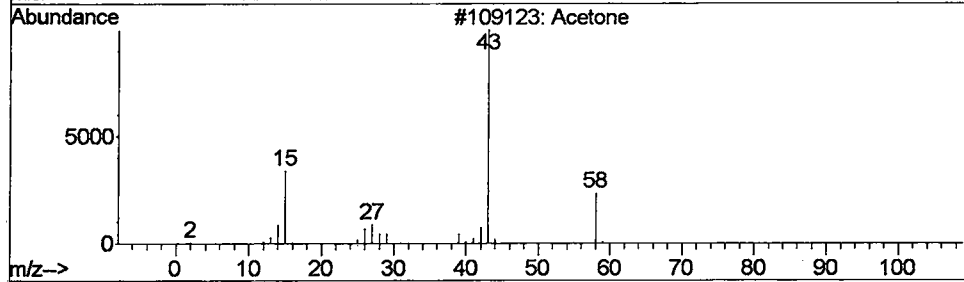
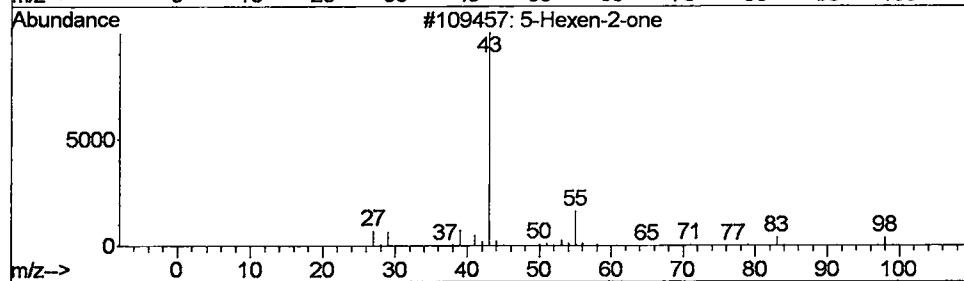
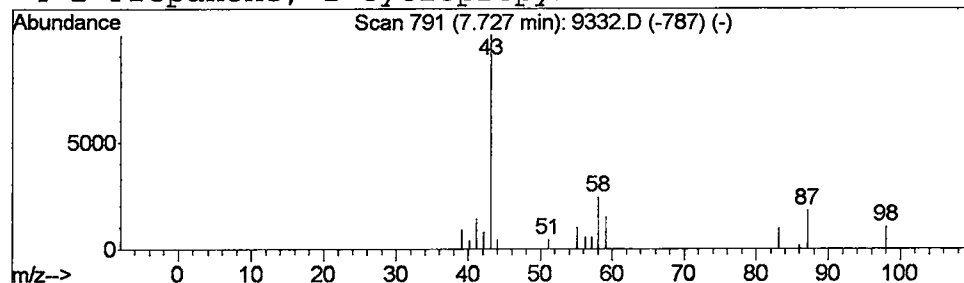
Vial: 22
 Operator: Mclean
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 4 5-Hexen-2-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	0.25 ug/L	249164	1,4-Dichlorobenzene-d4	9.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Hexen-2-one	98	C6H10O	000109-49-9	10
2			Acetone	58	C3H6O	000067-64-1	9
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9332.D
Acq On : 10 May 2002 9:13 am
Sample : 5/8 eh
Misc :
MS Integration Params: LSCINT.e

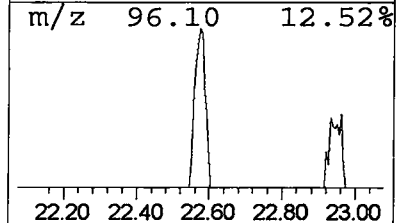
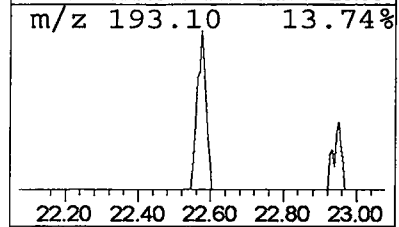
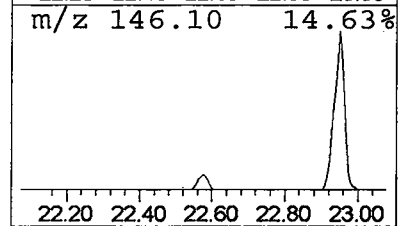
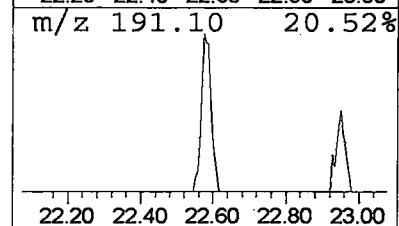
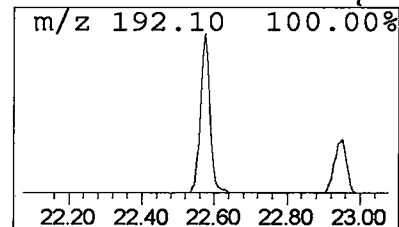
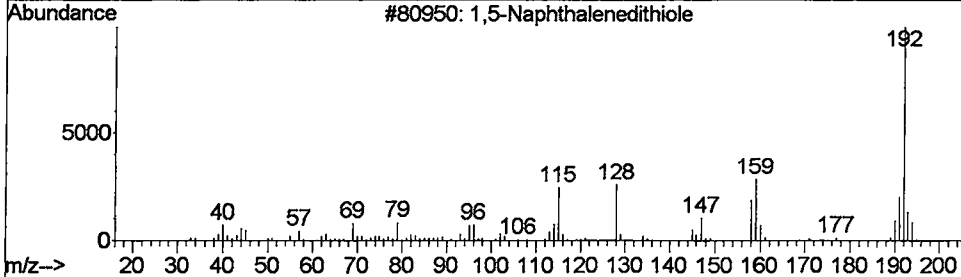
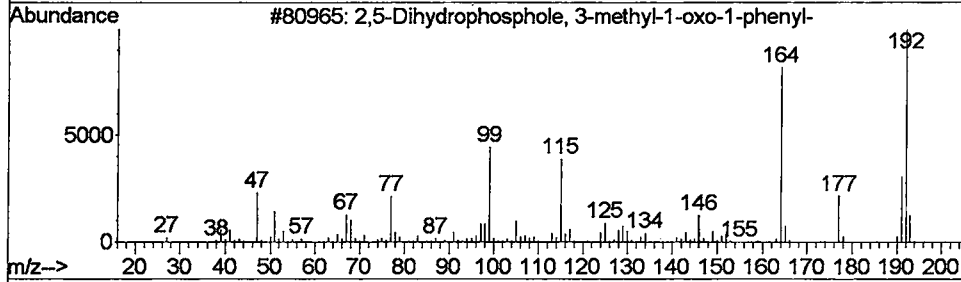
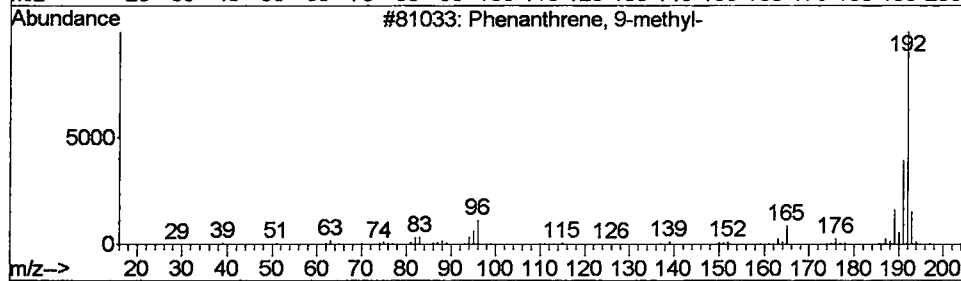
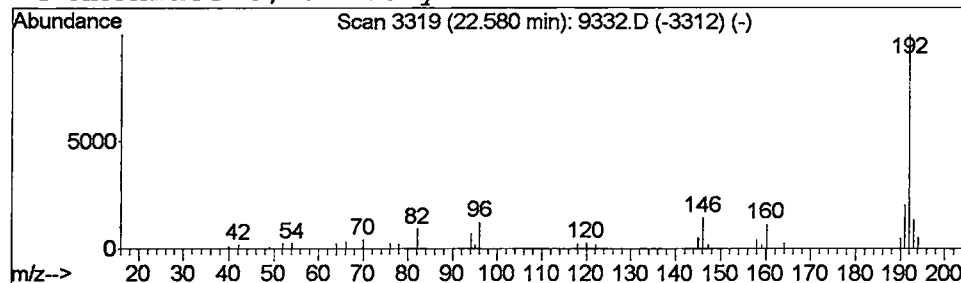
Vial: 22
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 5 Phenanthrene, 9-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	0.32 ug/L	439453	Phenanthranene-d10	22.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	50
2			2,5-Dihydrophosphole, 3-methyl-1...	192	C11H13OP	1000196-97-5	45
3			1,5-Naphthalenedithiole	192	C10H8S2	1000223-69-5	45
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	40



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9332.D
Acq On : 10 May 2002 9:13 am
Sample : 5/8 eh
Misc :
MS Integration Params: LSCINT.e

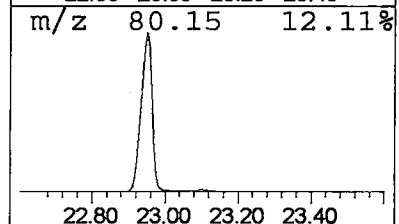
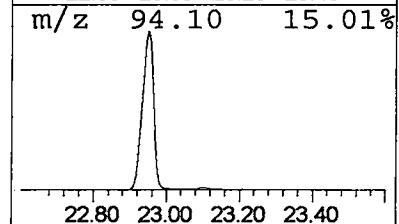
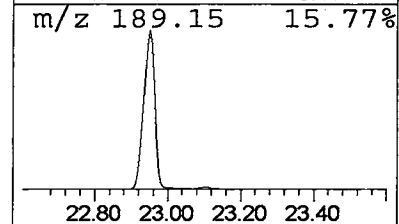
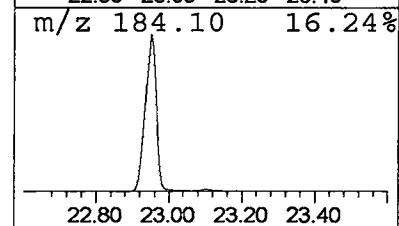
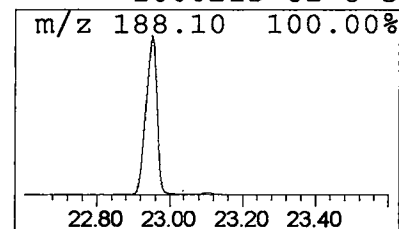
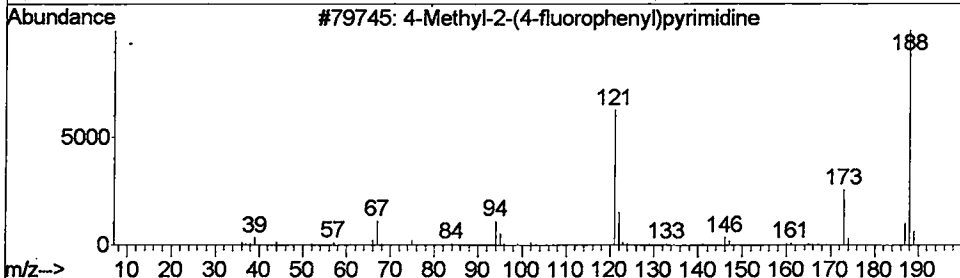
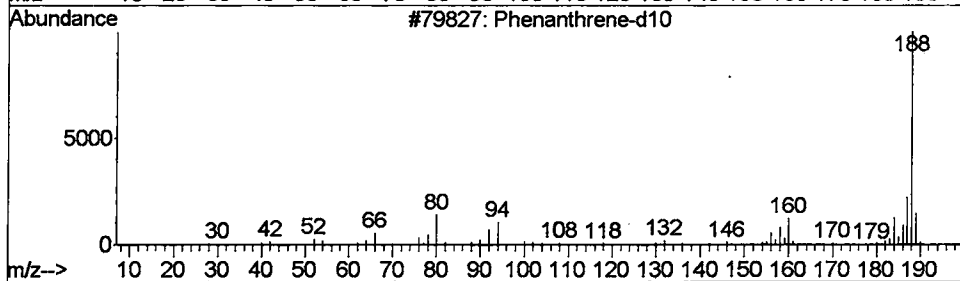
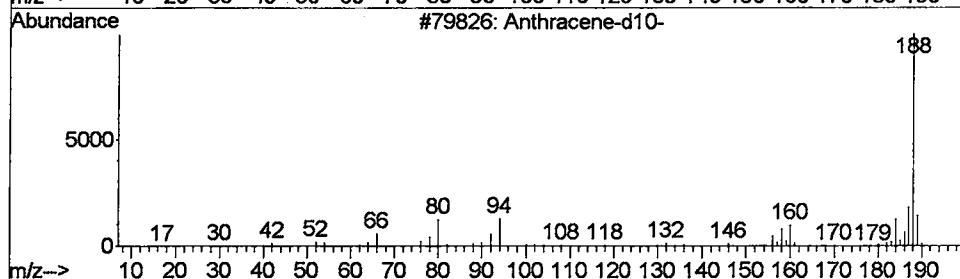
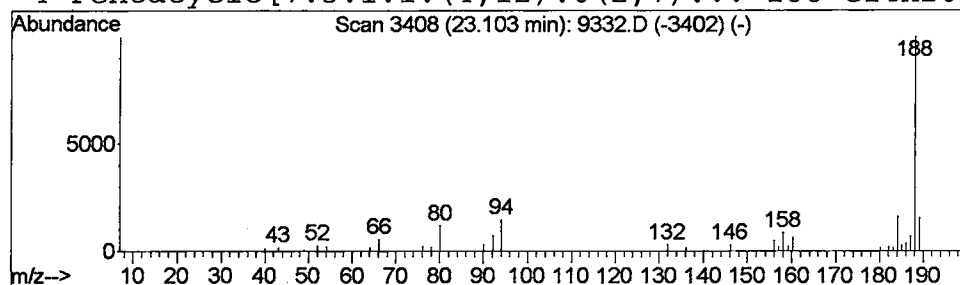
Vial: 22
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 6 Anthracene-d10- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	0.44 ug/L	608229	Phenanthranene-d10	22.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene-d10-	188	C14D10	001719-06-8	93
2			Phenanthrene-d10	188	C14D10	001517-22-2	90
3			4-Methyl-2-(4-fluorophenyl)pyrim...	188	C11H9FN2	076128-67-1	40
4			Pentacyclo[7.3.1.1.(4,12).0(2,7)...]	188	C14H20	1000215-61-8	38



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9332.D
 Acq On : 10 May 2002 9:13 am
 Sample : 5/8 eh
 Misc :
 MS Integration Params: LSCINT.e

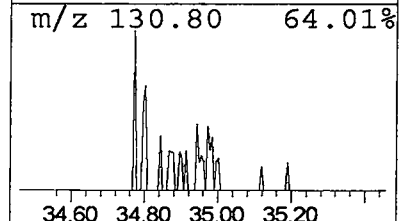
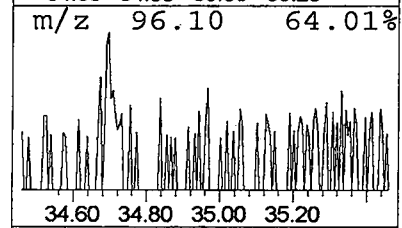
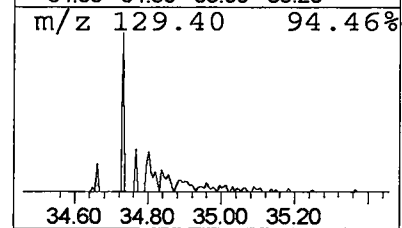
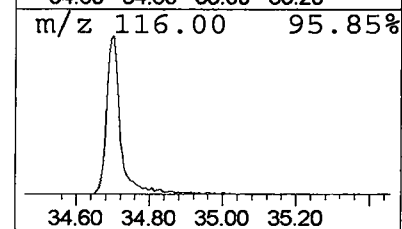
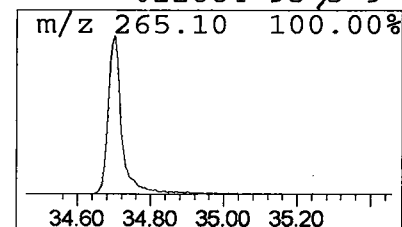
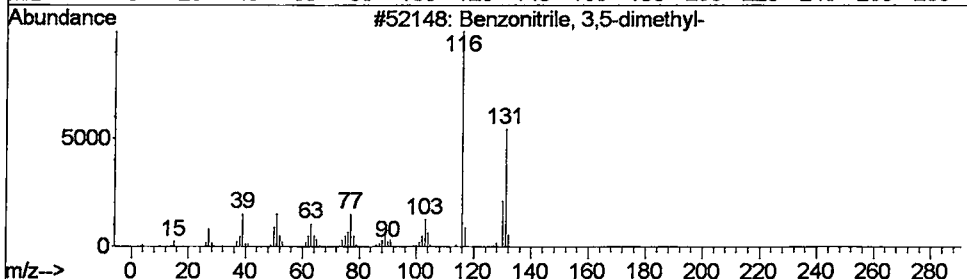
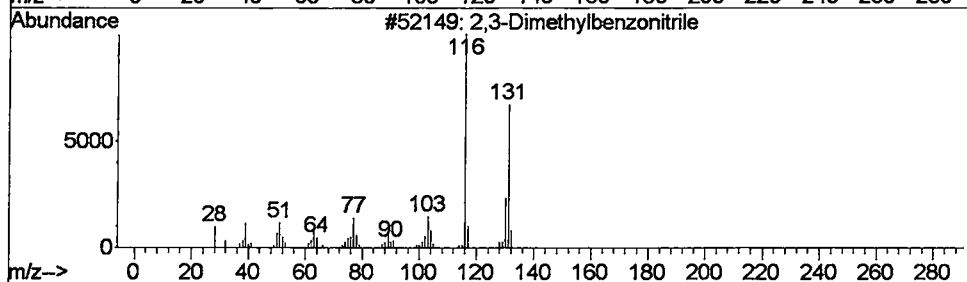
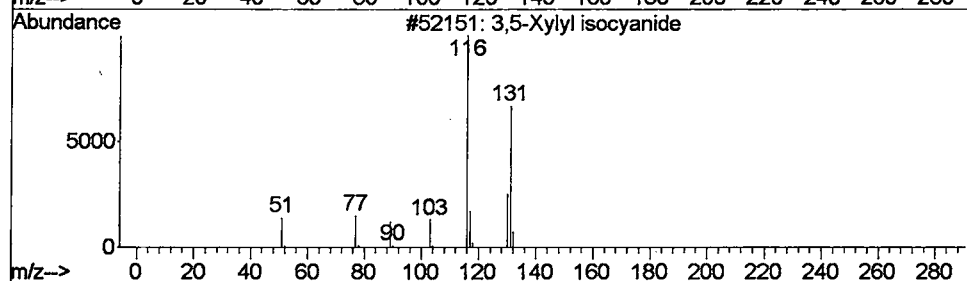
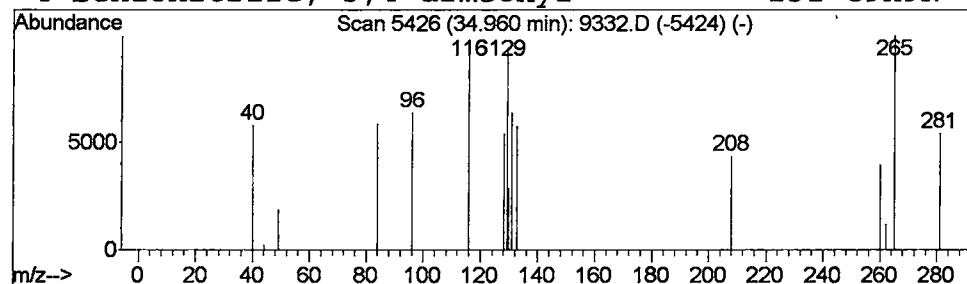
Vial: 22
 Operator: Mclean
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 7 3,5-Xylyl isocyanide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.96	0.27 ug/L	179639	Perylene-d12	34.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Xylyl isocyanide	131	C9H9N	020600-56-0	9
2			2,3-Dimethylbenzonitrile	131	C9H9N	005724-56-1	9
3			Benzonitrile, 3,5-dimethyl-	131	C9H9N	022445-42-7	9
4			Benzonitrile, 3,4-dimethyl-	131	C9H9N	022884-95-3	9



Tentatively Identified Compound (LSC) summary

Operator ID: Mclean Date Acquired: 10 May 2002 9:13 am
Data File: C:\MSDCHEM\1\DATA\050902\9332.D
Name: 5/8.eh
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
Acetonitrile, dic...	3.61	0.3	ug/L	285417	1	9.94	39273900	40.0
Ethyl Chloride	3.73	0.4	ug/L	386379	1	9.94	39273900	40.0
2-Pentanone, 4-hy...	5.97	15.3	ug/L	15010800	1	9.94	39273900	40.0
5-Hexen-2-one	7.73	0.3	ug/L	249164	1	9.94	39273900	40.0
Phenanthrene, 9-m...	22.58	0.3	ug/L	439453	4	22.95	55750600	40.0
Anthracene-d10-	23.11	0.4	ug/L	608229	4	22.95	55750600	40.0
3,5-Xylyl isocyanide	34.96	0.3	ug/L	179639	6	34.70	26819000	40.0