User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/09/2002 Time: 09:29:55

Sample: AE07568
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/9/02 09:29 Ended: 4/9/02 09:29 Analyst: DLV
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		96		5.0

Comments for Sample AE07568 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 04-09-2002 Time: 09:30:10

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

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: Apr 04 08:14:42 2002 Quant Results File: SCAN020403.RE

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Last Update : Thu Apr 04 08:10:12 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN1

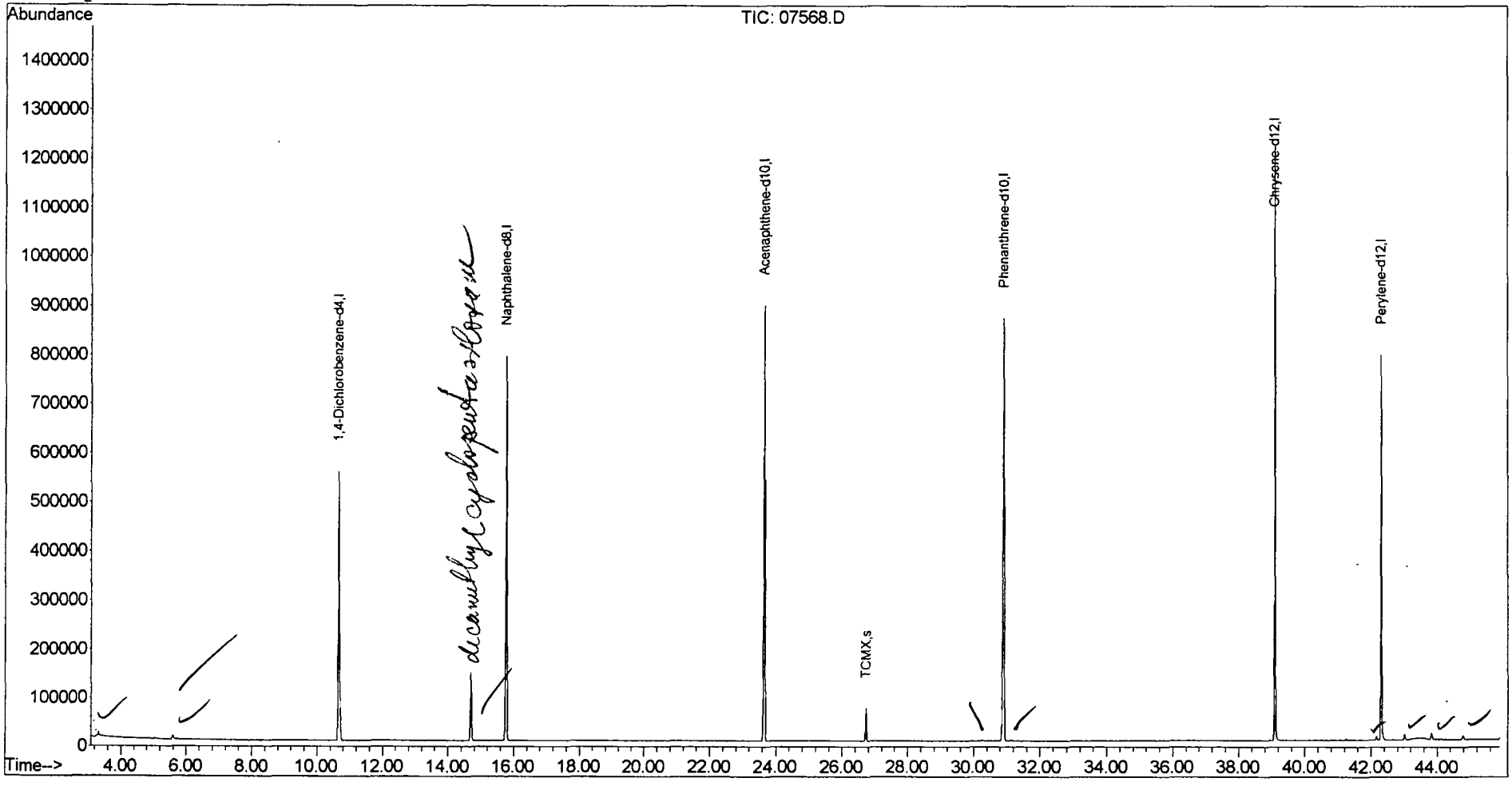
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	10.66	152	2806116	20.00	ug/l	0.00
10) Naphthalene-d8	15.77	136	10301592	20.00	ug/l	0.00
17) Acenaphthene-d10	23.66	164	5927870	20.00	ug/l	-0.01
31) Phenanthrene-d10	30.90	188	9801968	20.00	ug/l	0.00
39) Chrysene-d12	39.10	240	8002225	20.00	ug/l	-0.01
45) Perylene-d12	42.30	264	4532083	20.00	ug/l	0.00
System Monitoring Compounds						
28) TCMX	26.73	207	222224	3.83	ug/l	0.00
Spiked Amount ~~5.000~~ 4.0			Recovery =	~~76.60%~~ 96%		
Target Compounds						Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed
07568.D SCAN020403.M Thu Apr 04 10:29:27 2002 Page 1

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: Apr 4 10:29 2002 Quant Results File: SCAN020403.RES

Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Last Update : Thu Apr 04 08:40:00 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan

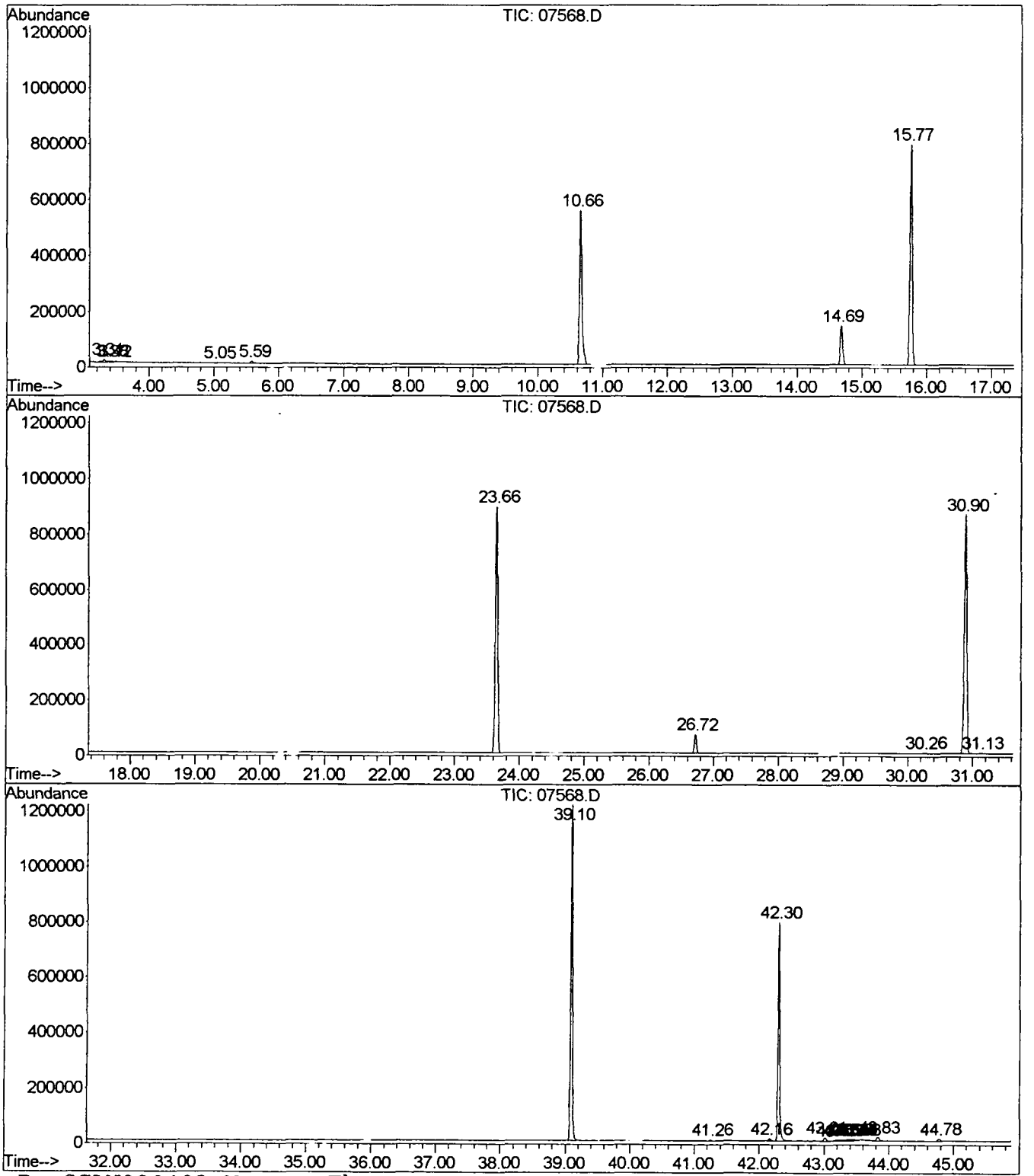
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.314	18	40	50	PV 3	10996	394584	1.65%	0.315%
2	3.385	50	52	55	VV 3	3653	58891	0.25%	0.047%
3	3.420	55	58	65	VV 4	4117	115245	0.48%	0.092%
4	5.047	327	335	338	VV 2	2116	44342	0.18%	0.035%
5	5.588	417	427	438	PV 3	7442	195338	0.81%	0.156%
6	10.659	1279	1290	1314	PV	550008	14598068	60.89%	11.642%
7	14.689	1963	1976	1989	VV	138446	3528208	14.72%	2.814%
8	15.770	2142	2160	2170	PV	784515	19230453	80.21%	15.336%
9	23.655	3486	3502	3518	PV	879074	22770228	94.98%	18.159%
10	26.722	4014	4024	4033	VV 2	66939	1494926	6.24%	1.192%
11	30.260	4620	4626	4641	VV 2	2247	64085	0.27%	0.051%
12	30.894	4720	4734	4753	VV 2	856033	23974721	100.00%	19.119%
13	31.129	4767	4774	4783	PV	2356	61124	0.25%	0.049%
14	39.096	6120	6130	6145	PV 2	1216878	21949145	91.55%	17.504%
15	41.253	6491	6497	6502	PV	3937	69128	0.29%	0.055%
16	42.163	6645	6652	6660	VV	7971	178961	0.75%	0.143%
17	42.305	6665	6676	6701	VV 2	789370	14657052	61.14%	11.689%
18	43.010	6787	6796	6810	VV 3	12907	369332	1.54%	0.295%
19	43.174	6817	6824	6828	VV 3	2646	67154	0.28%	0.054%
20	43.268	6828	6840	6841	VV 3	3883	137446	0.57%	0.110%
21	43.292	6841	6844	6850	VV 3	4361	112683	0.47%	0.090%
22	43.333	6850	6851	6859	VV 4	3829	120278	0.50%	0.096%
23	43.392	6859	6861	6863	VV	4402	53619	0.22%	0.043%
24	43.444	6863	6870	6873	VV 4	4650	144990	0.60%	0.116%
25	43.474	6873	6875	6881	VV 3	5016	139126	0.58%	0.111%
26	43.533	6881	6885	6886	VV 4	4790	79954	0.33%	0.064%
27	43.832	6922	6936	6946	VV 2	13827	487105	2.03%	0.388%
28	44.778	7087	7097	7117	VV 3	8550	299198	1.25%	0.239%

Sum of corrected areas: 125395385

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\020403\07568.D
Operator : Nefliu
Acquired : 3 Apr 2002 10:56 pm using AcqMethod 508SCAN1
Instrument : Instrumen
Sample Name: sample
Misc Info :
Vial Number: 16
Quant File :SCAN020403.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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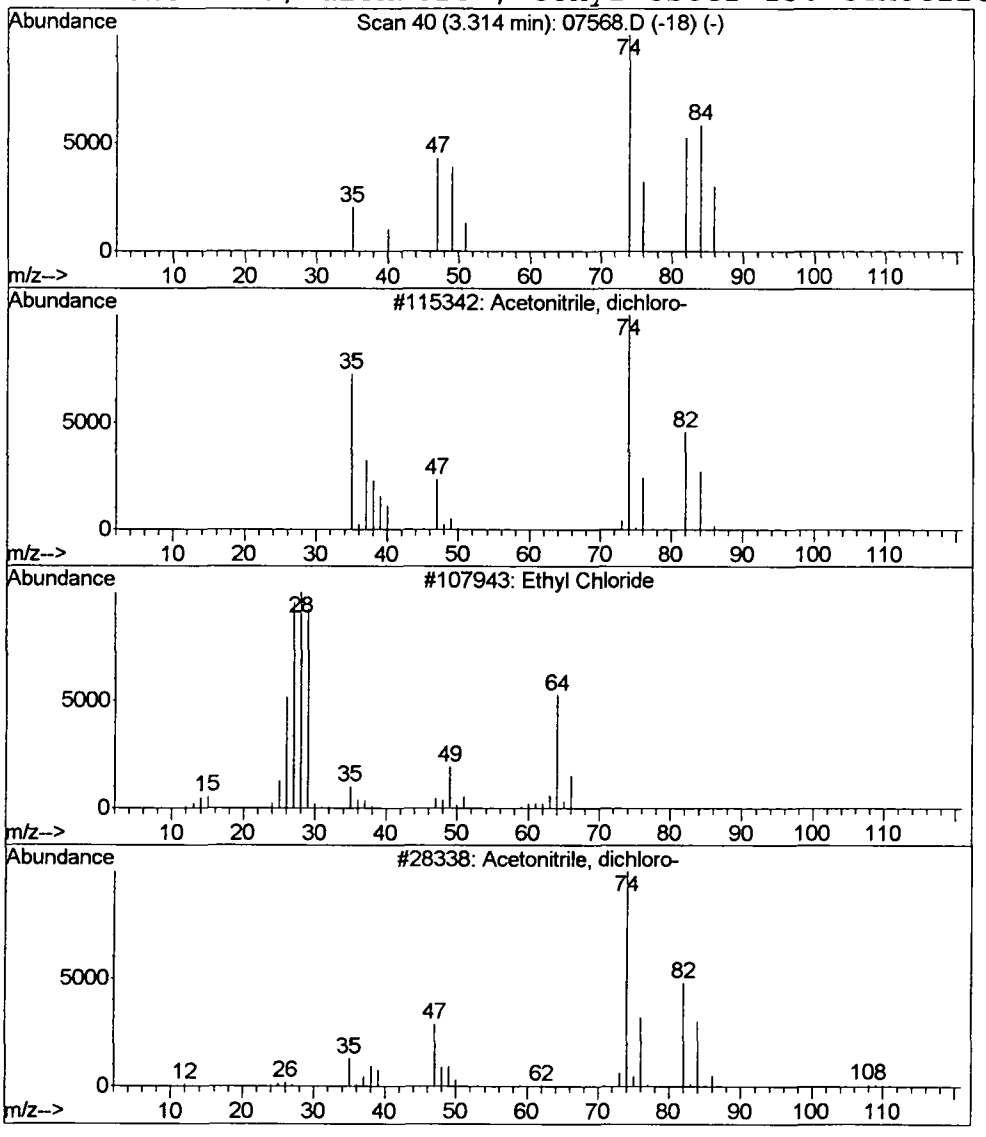
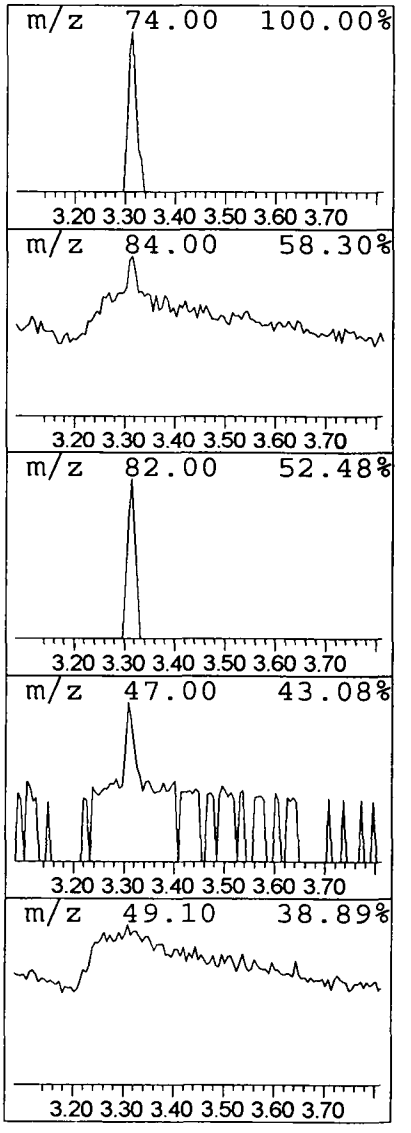
Peak Number 1 Acetonitrile, dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.31	0.54 ug/l	394584	1,4-Dichlorobenzene-d4	10.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	9
2	Ethyl Chloride	64	C2H5Cl	000075-00-3	4
3	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	4
4	Acetic acid, dichloro-, ethyl ester	156	C4H6Cl2O2	000535-15-9	2





Library Search Compound Report

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Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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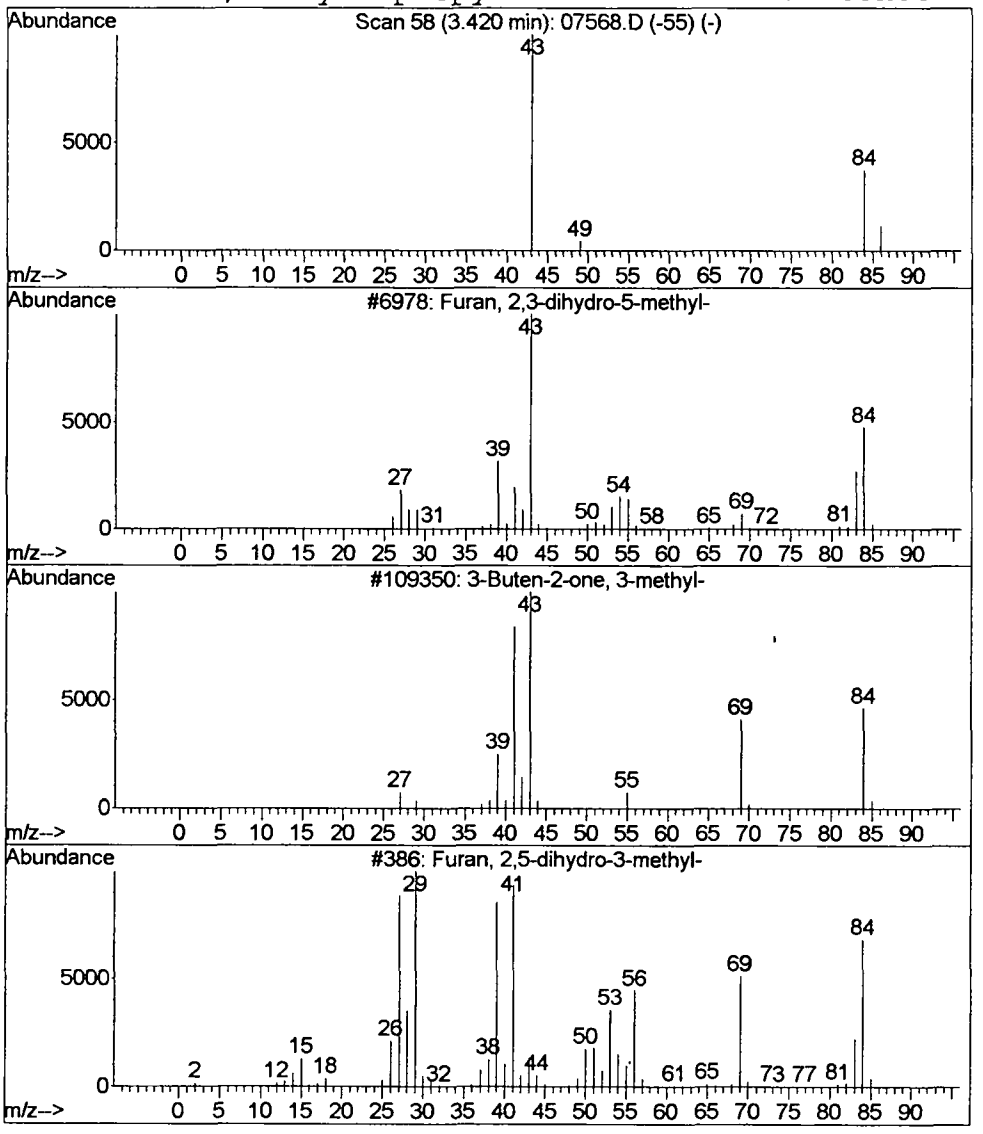
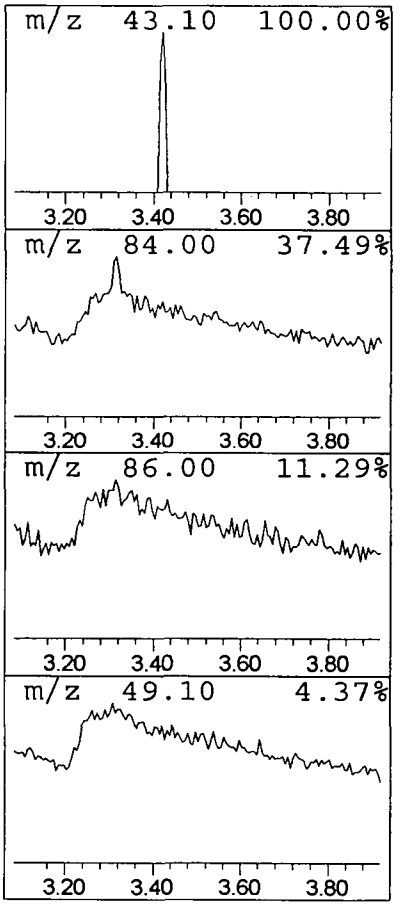
Peak Number 2 Furan, 2,3-dihydro-5-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.42	0.16 ug/l	115245	1,4-Dichlorobenzene-d4	10.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, 2,3-dihydro-5-methyl-	84	C5H8O	001487-15-6	4
2	3-Buten-2-one, 3-methyl-	84	C5H8O	000814-78-8	4
3	Furan, 2,5-dihydro-3-methyl-	84	C5H8O	001708-31-2	4
4	Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	4





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Misc : Multiplr: 1.00
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Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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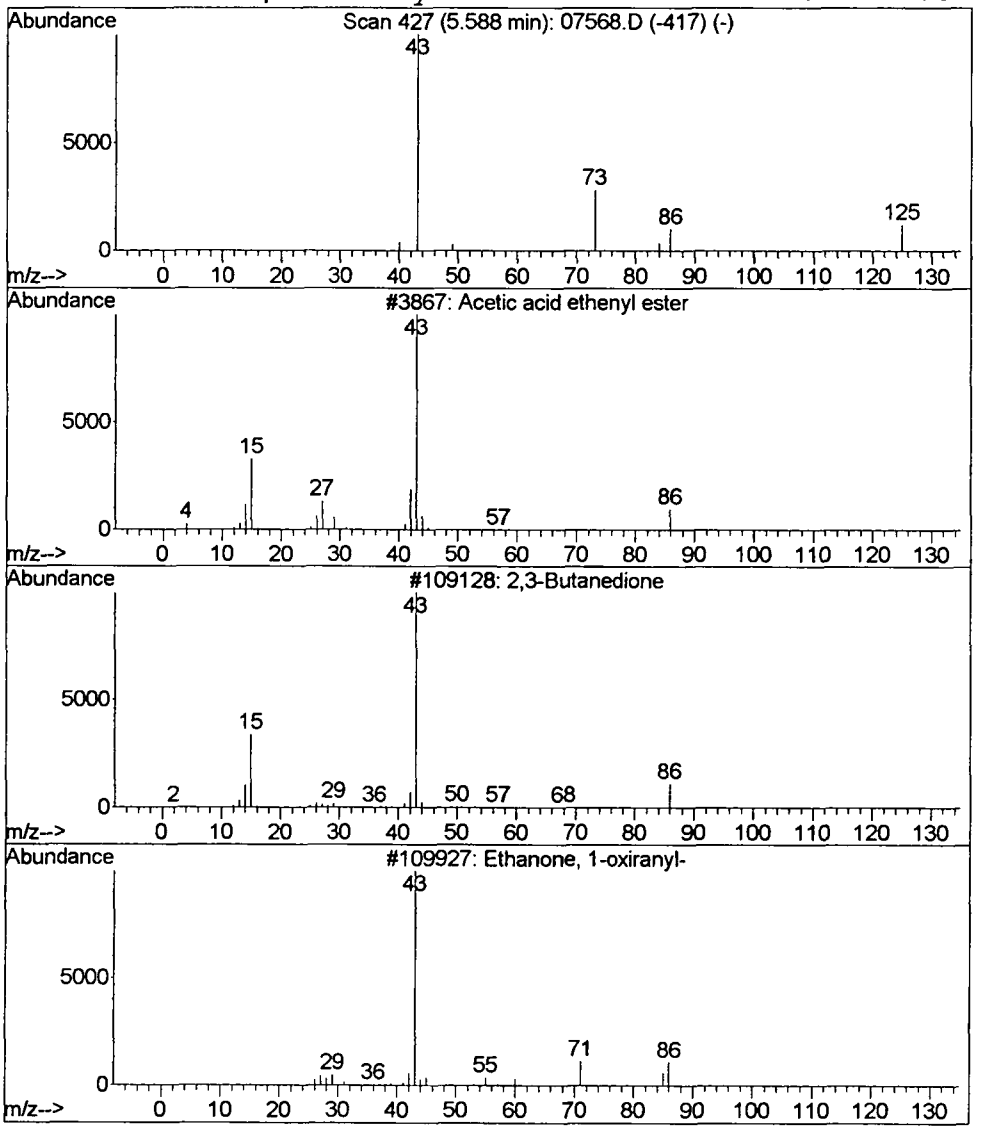
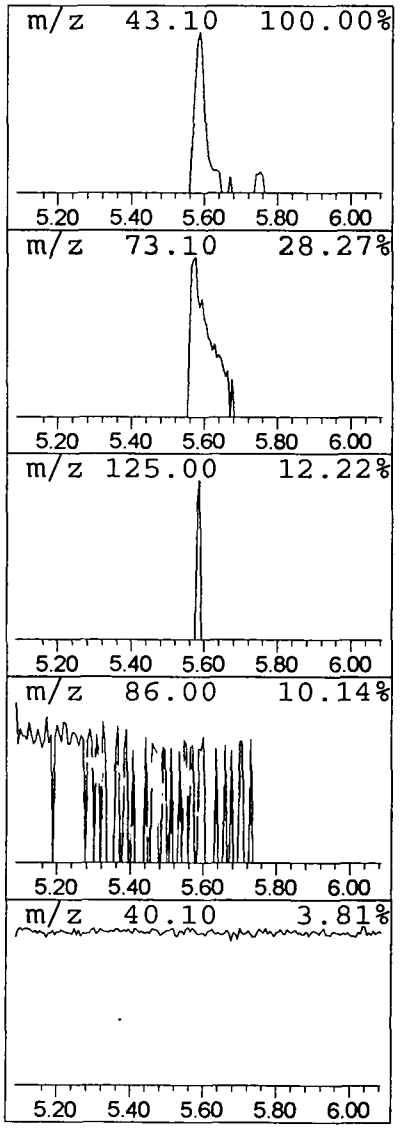
Peak Number 3 Acetic acid ethenyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.59	0.27 ug/l	195338	1,4-Dichlorobenzene-d4	10.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid ethenyl ester	86	C4H6O2	000108-05-4	3
2	2,3-Butanedione	86	C4H6O2	000431-03-8	3
3	Ethanone, 1-oxiranyl-	86	C4H6O2	004401-11-0	3
4	2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	3





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Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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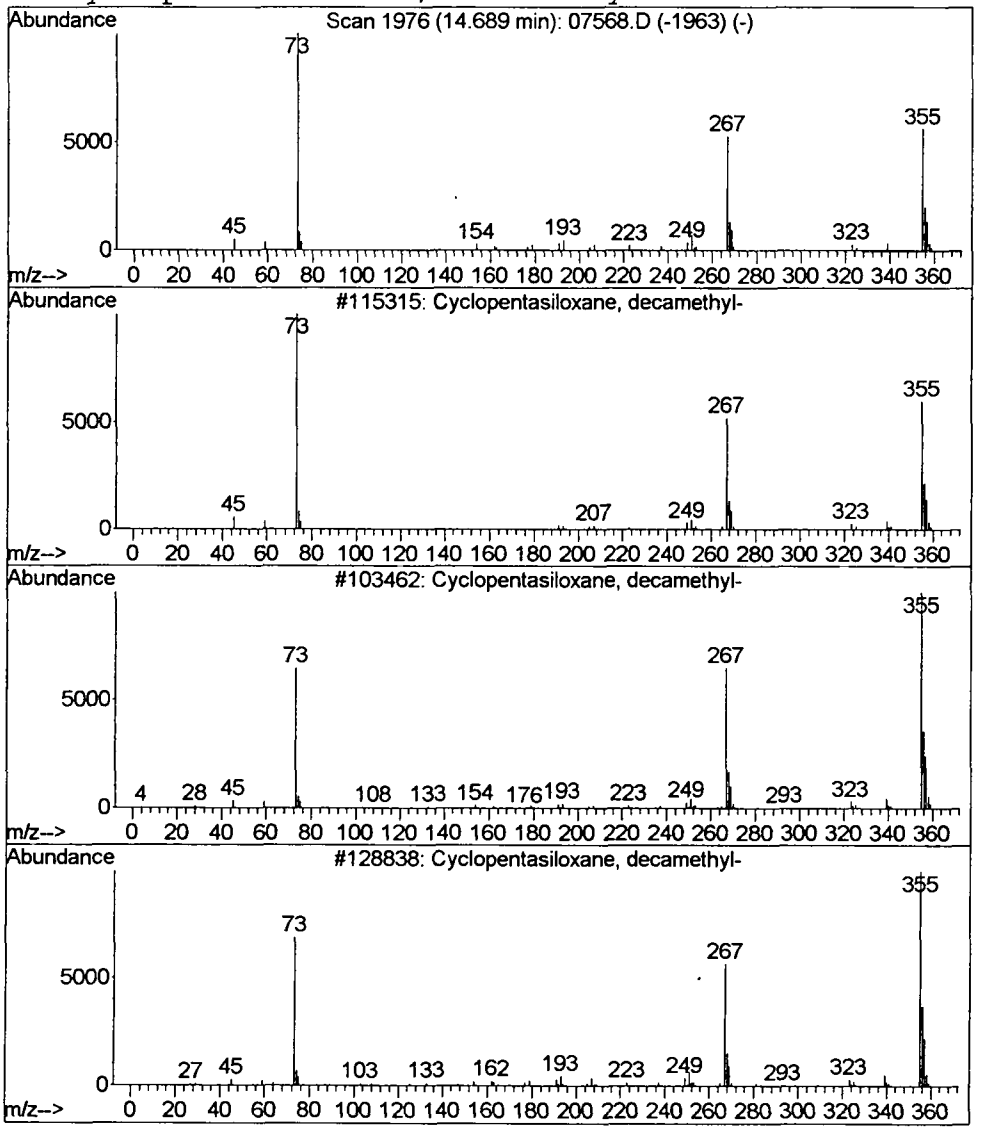
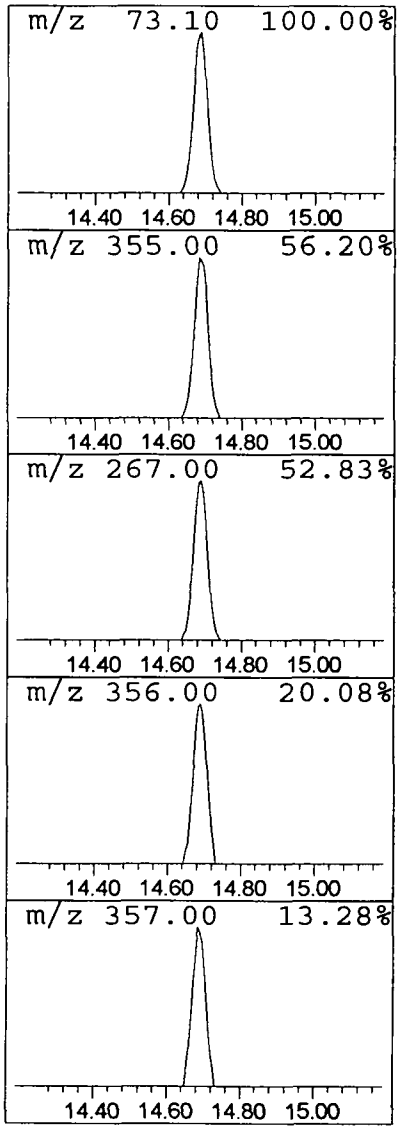
Peak Number 4 Cyclopentasiloxane, decamet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	3.67 ug/l	3528210	Naphthalene-d8	15.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
3	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	83
4	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	68





Library Search Compound Report

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Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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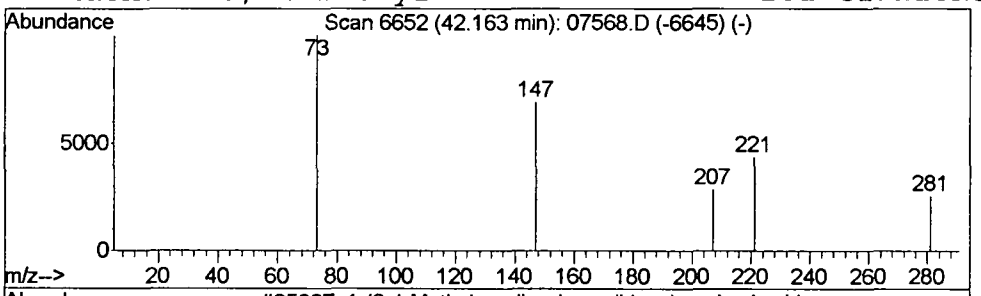
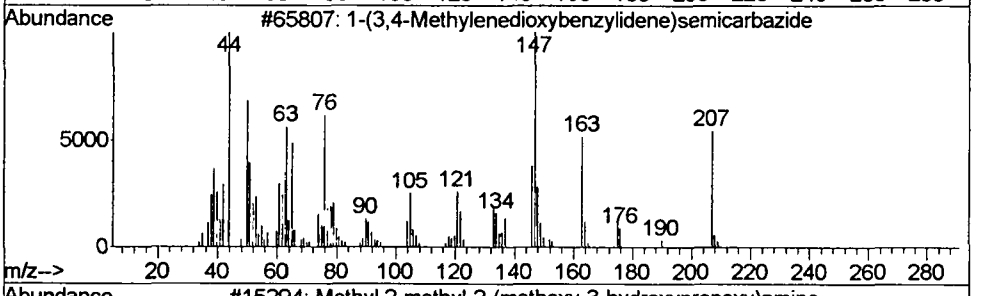
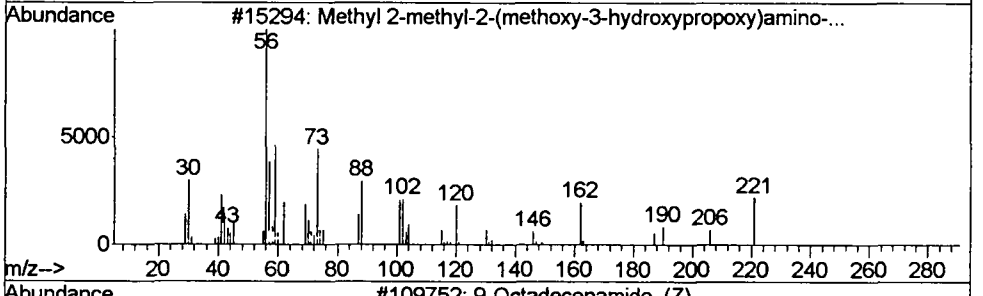
Peak Number 5 1-(3,4-Methylenedioxybenzyl... Concentration Rank 7

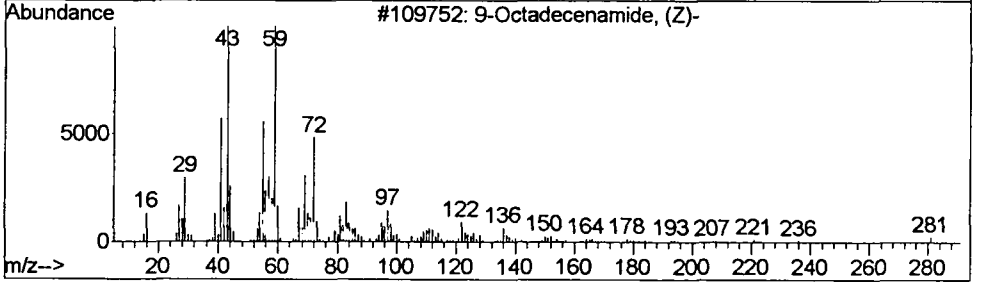
R.T.	EstConc	Area	Relative to ISTD	R.T.
42.16	0.24 ug/l	178961	Perylene-d12	42.30

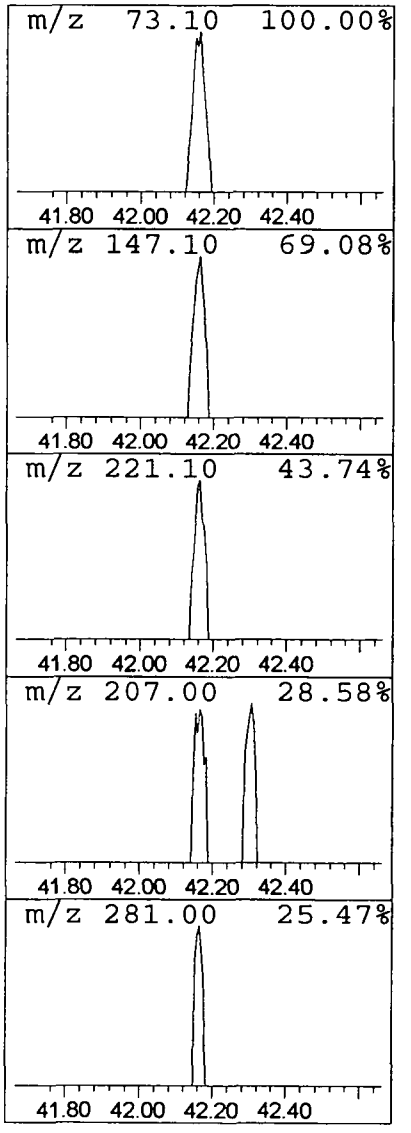
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(3,4-Methylenedioxybenzylidene...	207	C9H9N3O3	1000224-56-9	5
2	Methyl 2-methyl-2-(methoxy-3-hyd...	221	C9H19NO5	1000143-85-8	4
3	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	3
4	Adenosine, 2-methyl-	281	C11H15N5O4	016526-56-0	3











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Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
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Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
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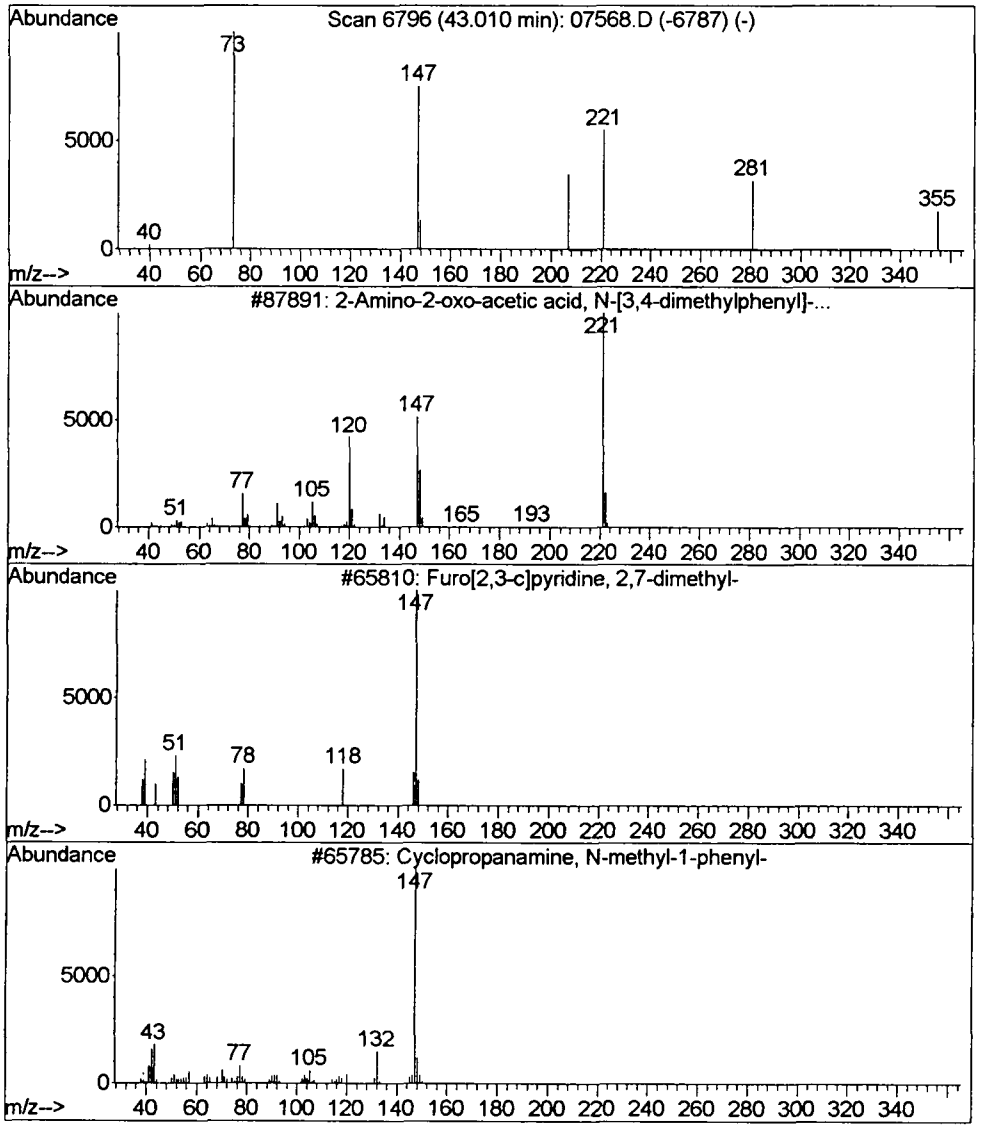
Library : C:\DATABASE\nist98.l

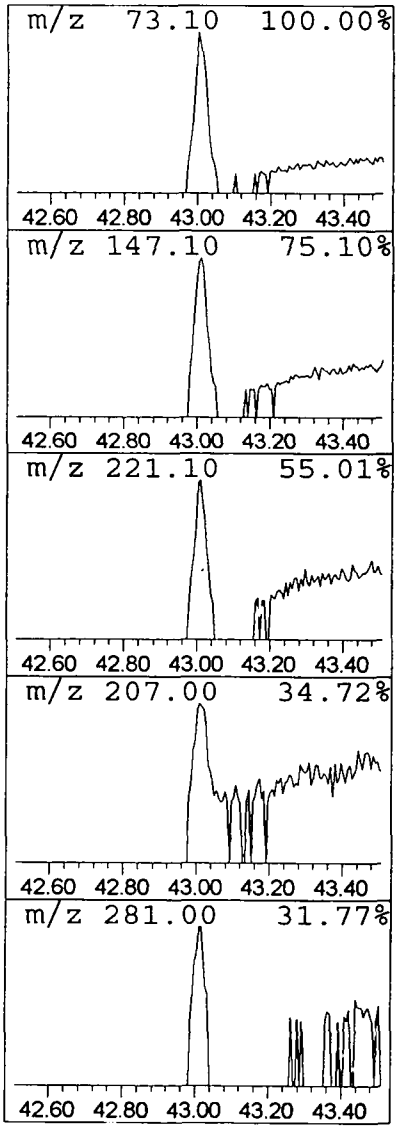
****

Peak Number 6 2-Amino-2-oxo-acetic acid, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.01	0.50 ug/l	369332	Perylene-d12	42.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-2-oxo-acetic acid, N-[3,...	221	C12H15NO3	1000127-85-3	5
2	Furo[2,3-c]pyridine, 2,7-dimethyl-	147	C9H9NO	069022-77-1	4
3	Cyclopropanamine, N-methyl-1-phe...	147	C10H13N	056771-48-3	4
4	2-Indolizinemethanol	147	C9H9NO	022320-21-4	4





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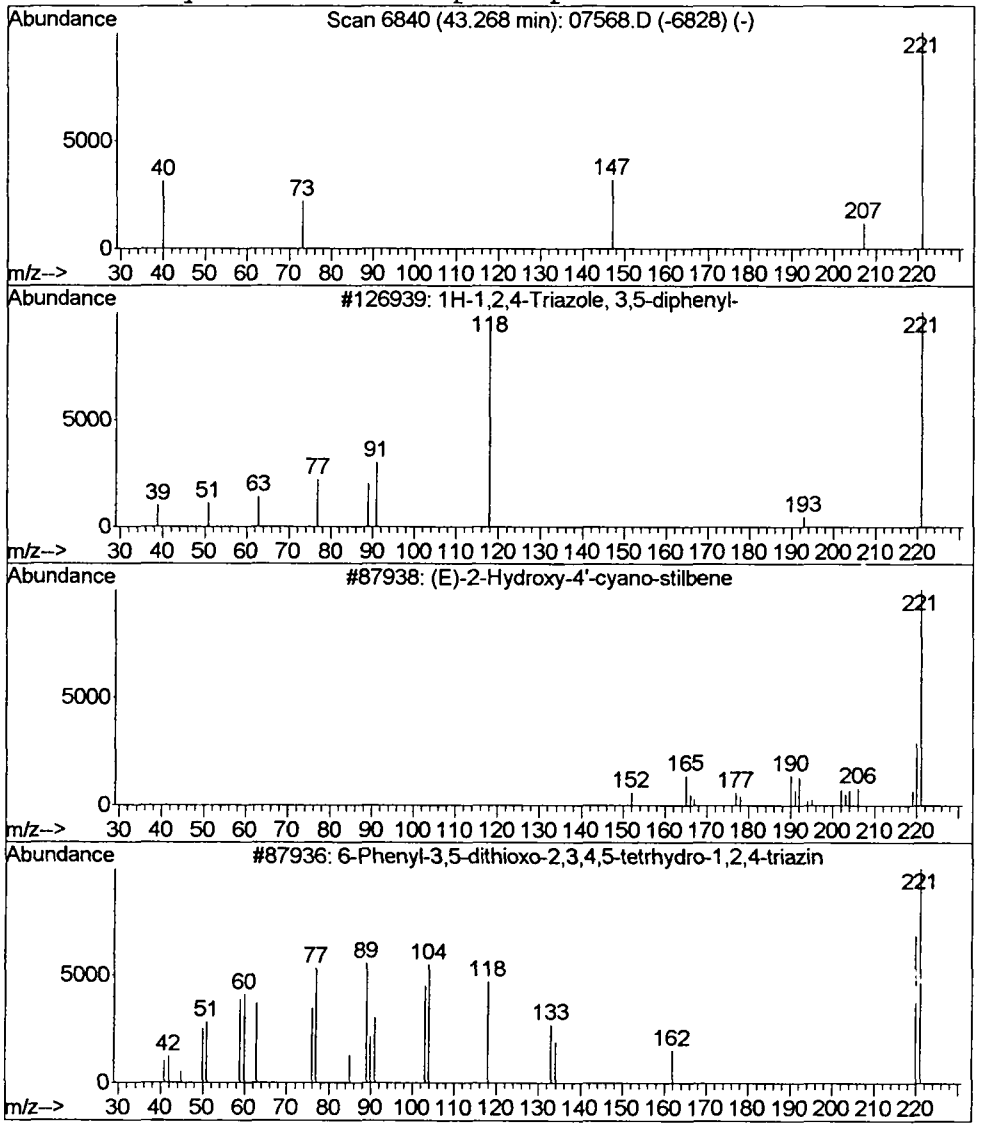
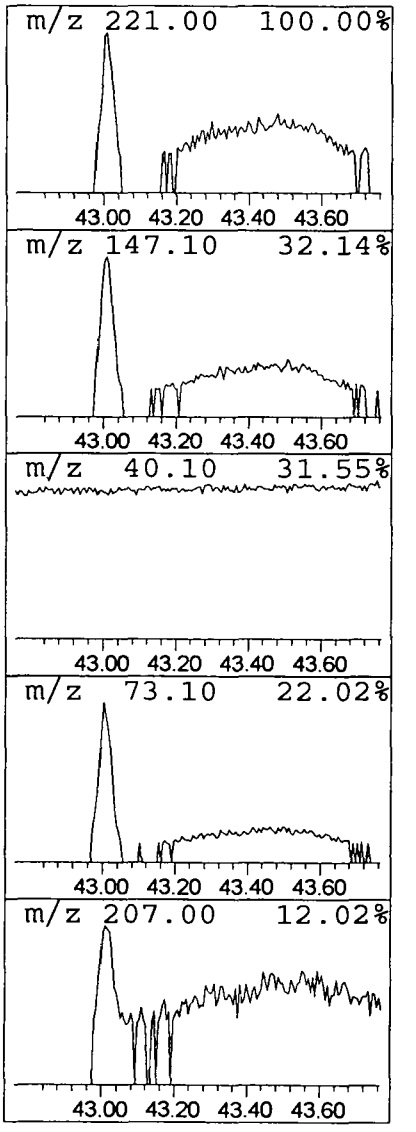
****

Peak Number 7 1H-1,2,4-Triazole, 3,5-diph... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.27	0.19 ug/l	137446	Perylene-d12	42.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-1,2,4-Triazole, 3,5-diphenyl-	221	C14H11N3	002039-06-7	2
2	(E)-2-Hydroxy-4'-cyano-stilbene	221	C15H11NO	1000147-85-5	2
3	6-Phenyl-3,5-dithioxo-2,3,4,5-te...	221	C9H7N3S2	1000147-85-6	2
4	Furfurylideniminosulphur pentafl...	221	C5H4F5NOS	090598-15-5	1





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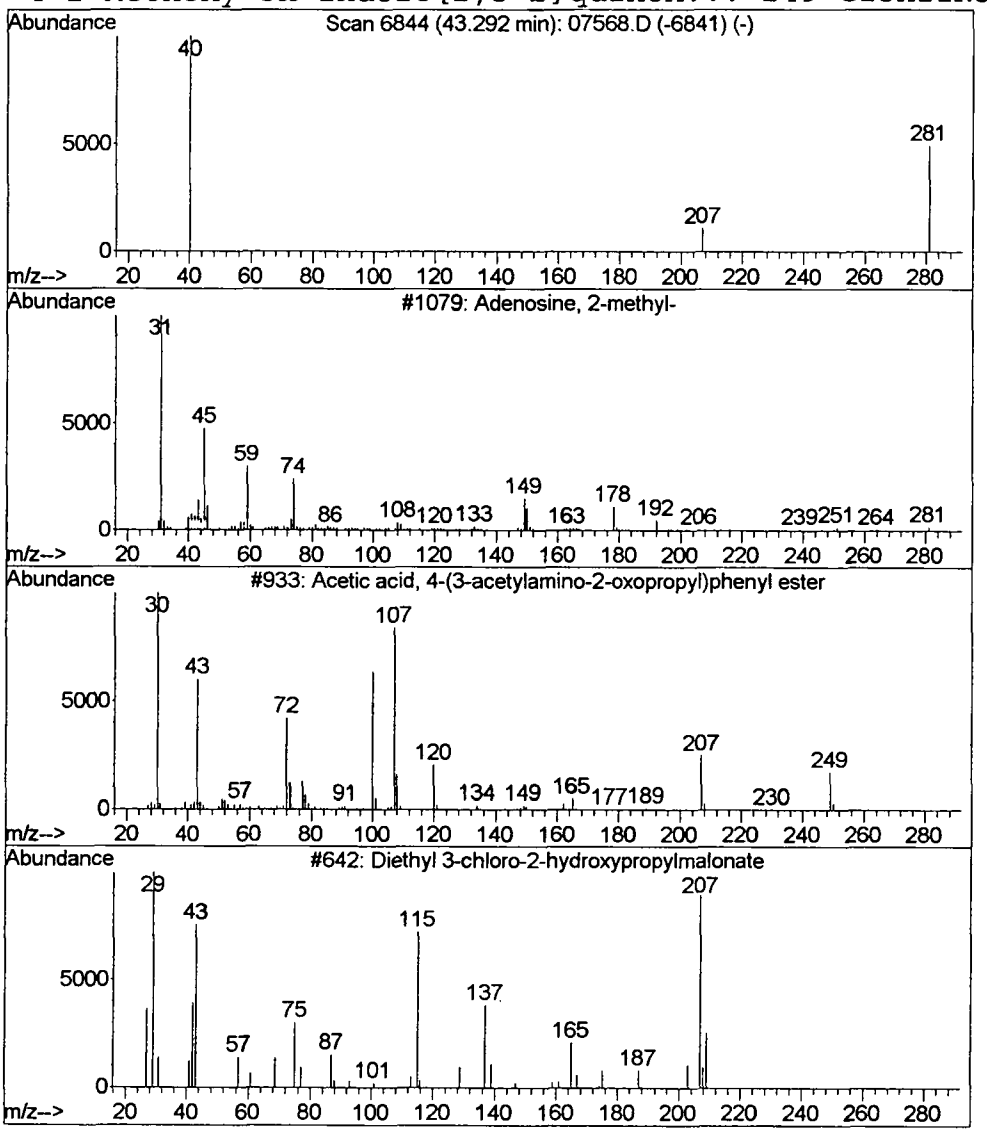
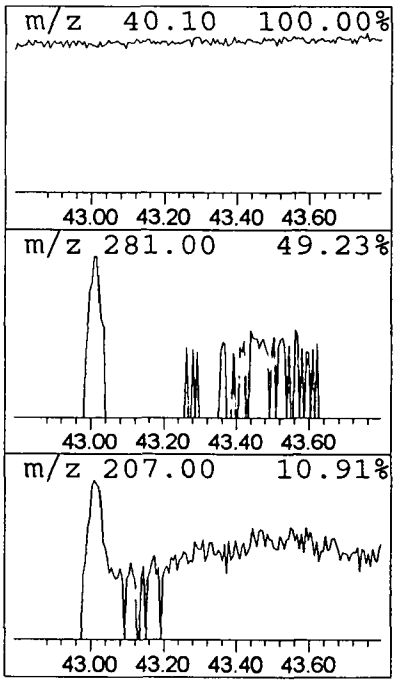
****

Peak Number 8 Adenosine, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.29	0.15 ug/l	112683	Perylene-d12	42.30

Hit# of 4	Tentative ID	MW	MolForm	CAS#	Qual
1	Adenosine, 2-methyl-	281	C11H15N5O4	016526-56-0	1
2	Acetic acid, 4-(3-acetylamino-2-...	249	C13H15NO4	1000194-74-7	1
3	Diethyl 3-chloro-2-hydroxypropyl...	252	C10H17ClO5	1000222-74-8	1
4	2-Methoxy-5H-indolo[2,3-b]quinox...	249	C15H11N3O	1000223-36-6	1





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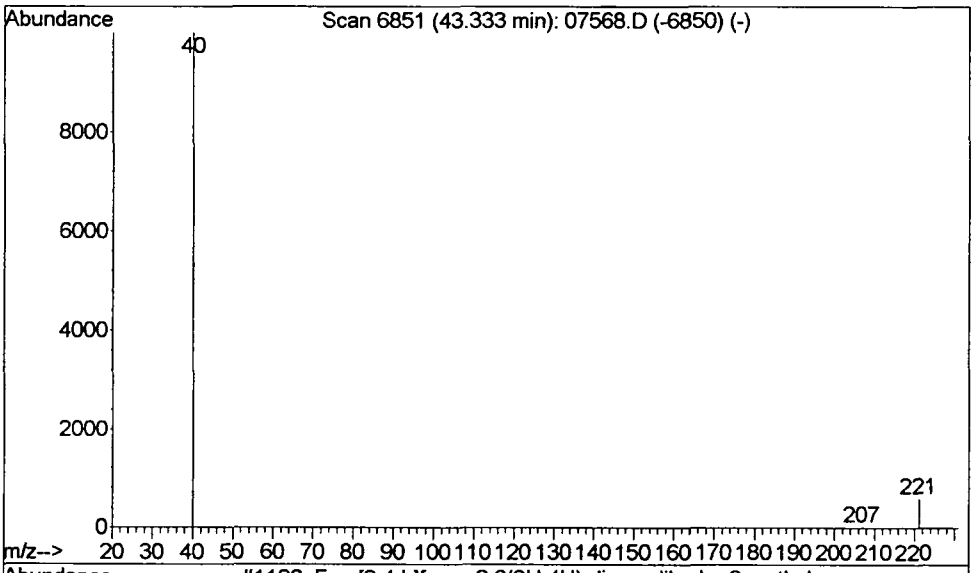
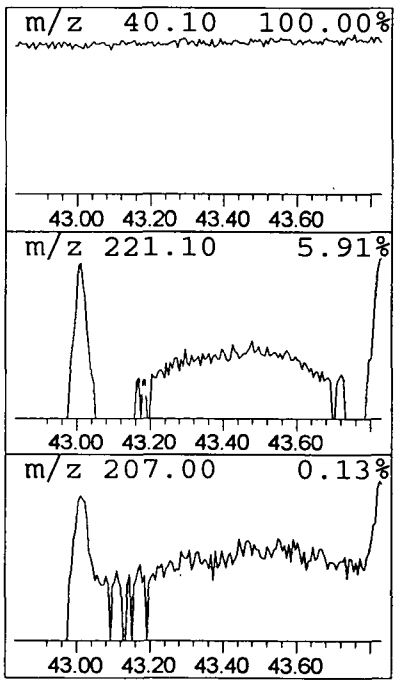
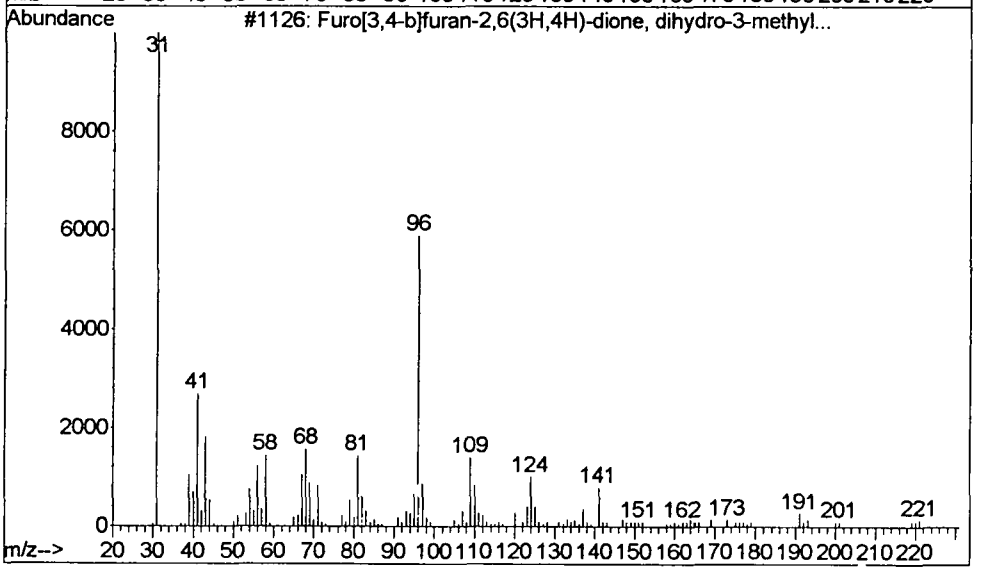
Peak Number 9 Furo[3,4-b]furan-2,6(3H,4H)... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.33	0.16 ug/l	120278	Perylene-d12	42.30

Hit# of 1	Tentative ID	MW	MolForm	CAS#	Qual
1	Furo[3,4-b]furan-2,6(3H,4H)-dion...	266	C15H22O4	020223-76-1	1







Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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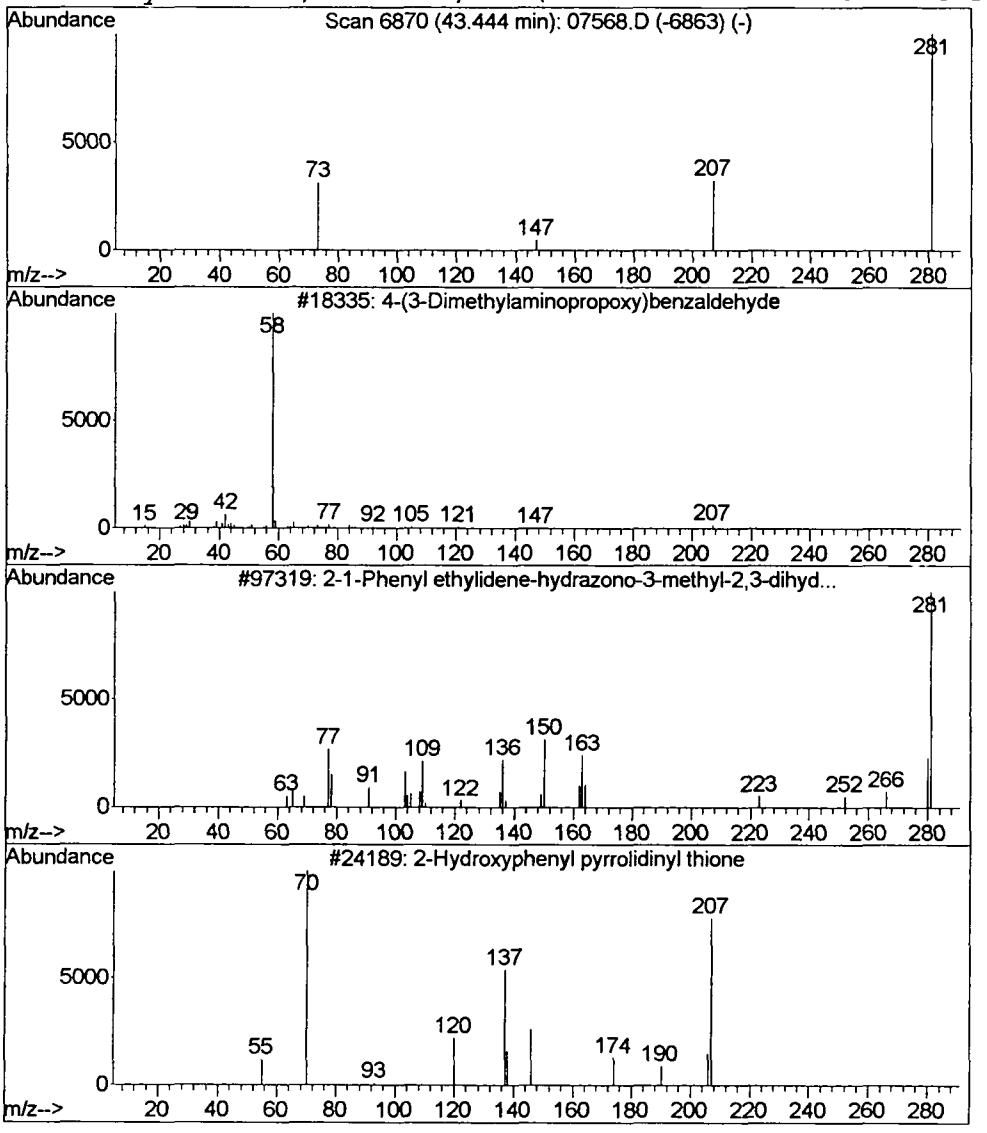
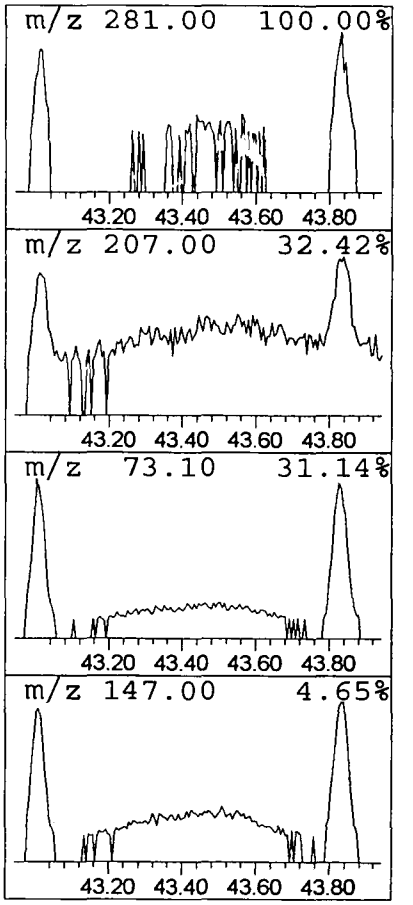
Peak Number 10 4-(3-Dimethylaminopropoxy)b... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.44	0.20 ug/l	144990	Perylene-d12	42.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(3-Dimethylaminopropoxy)benzal...	207	C12H17NO2	026934-35-0	3
2	2-1-Phenyl ethylidene-hydrazono-...	281	C16H15N3S	1000148-55-3	3
3	2-Hydroxyphenyl pyrrolidinyl thione	207	C11H13NOS	084783-01-7	2
4	1H-Pyrrole-2,5-dione, 1-(4-chlor...	207	C10H6ClNO2	001631-29-4	2





Library Search Compound Report

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Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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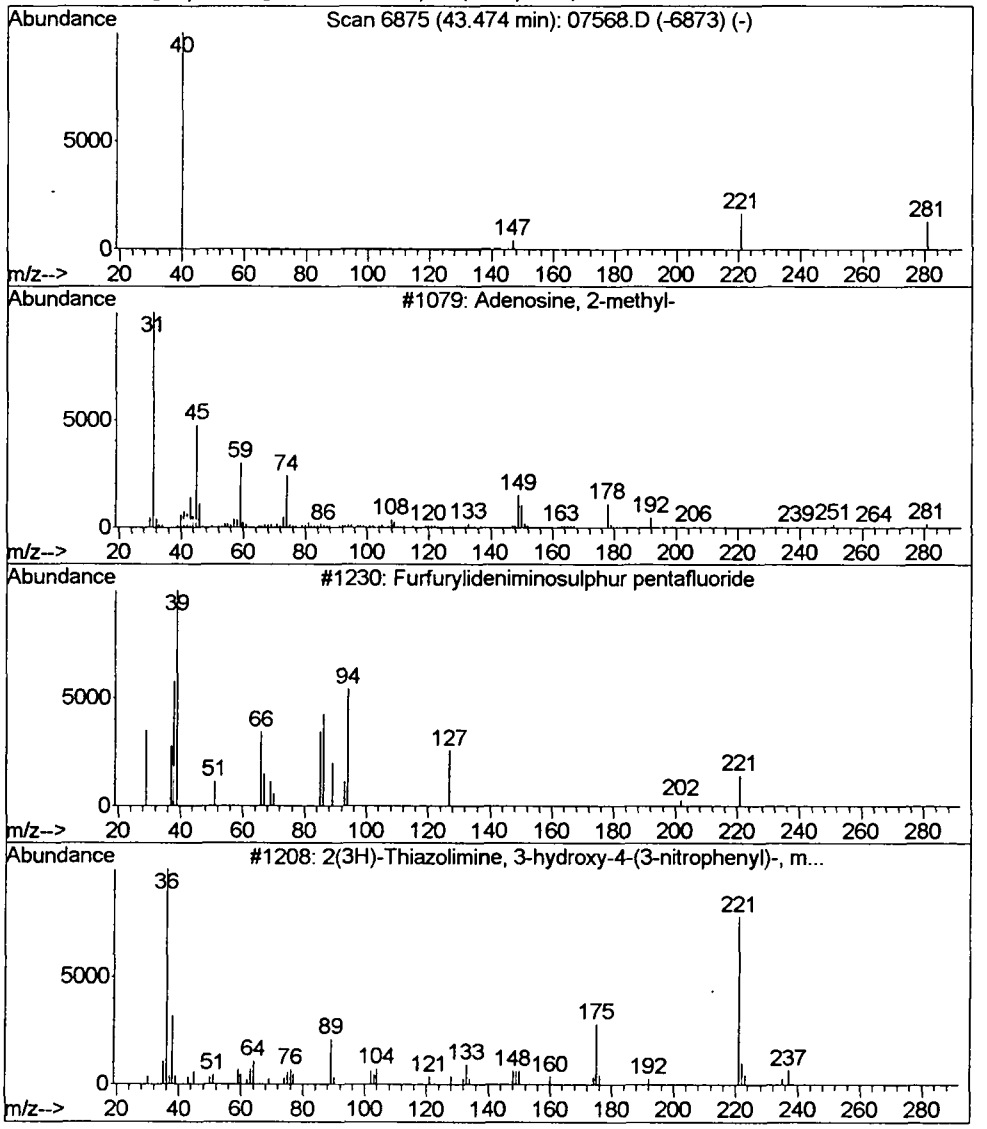
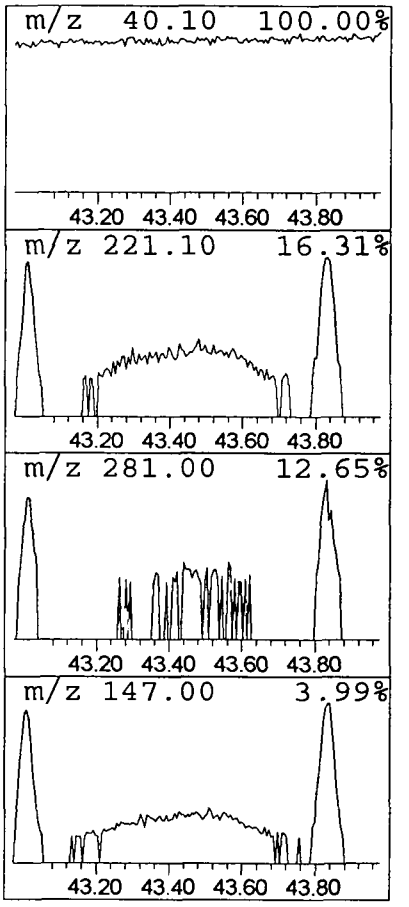
Peak Number 11 Adenosine, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.48	0.19 ug/l	139126	Perylene-d12	42.30

Hit# of 4	Tentative ID	MW	MolForm	CAS#	Qual
1	Adenosine, 2-methyl-	281	C11H15N5O4	016526-56-0	2
2	Furfurylideniminosulphur pentafl...	221	C5H4F5NOS	090598-15-5	1
3	2(3H)-Thiazolimine, 3-hydroxy-4-...	273	C9H8ClN3O3S	058275-70-0	1
4	Furo[3,4-b]furan-2,6(3H,4H)-dion...	266	C15H22O4	020223-76-1	1





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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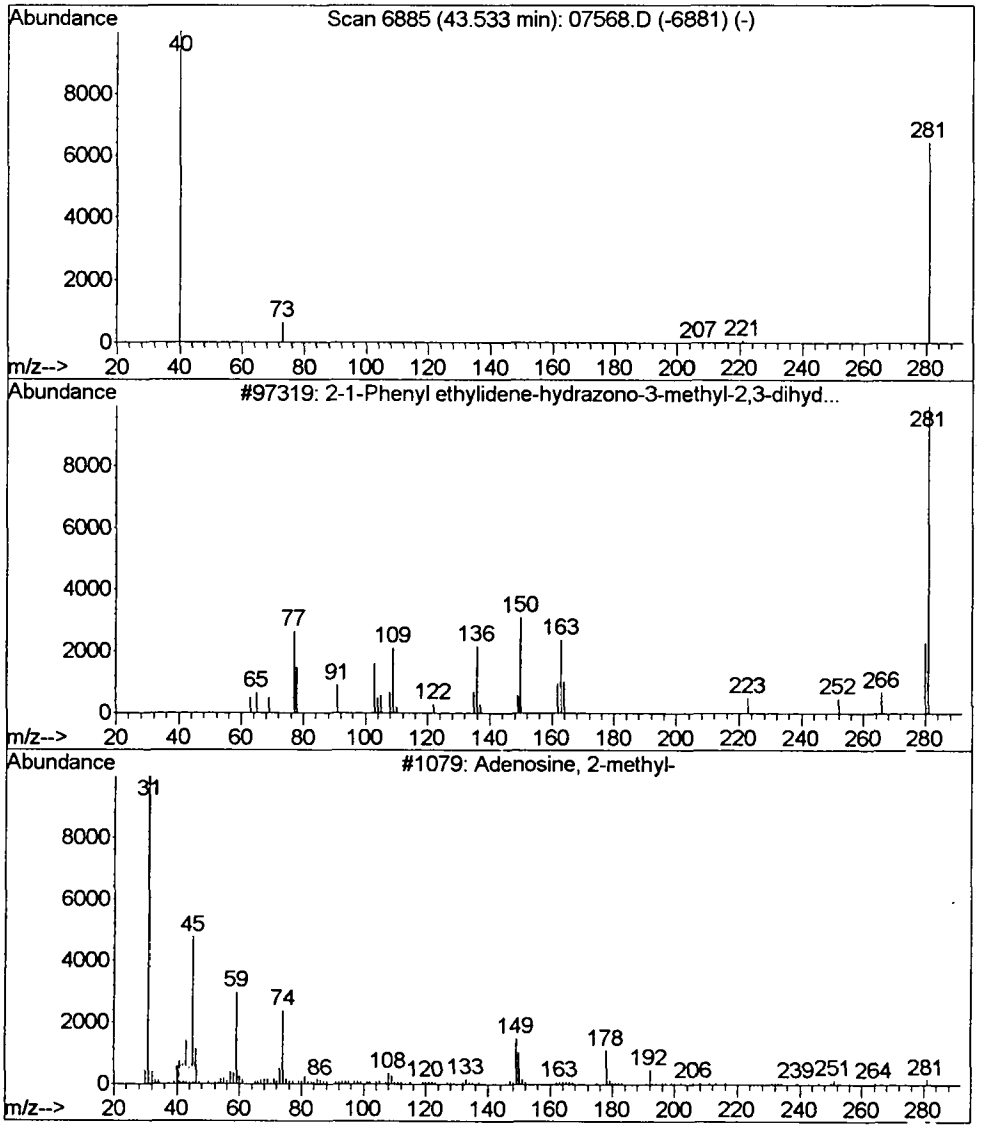
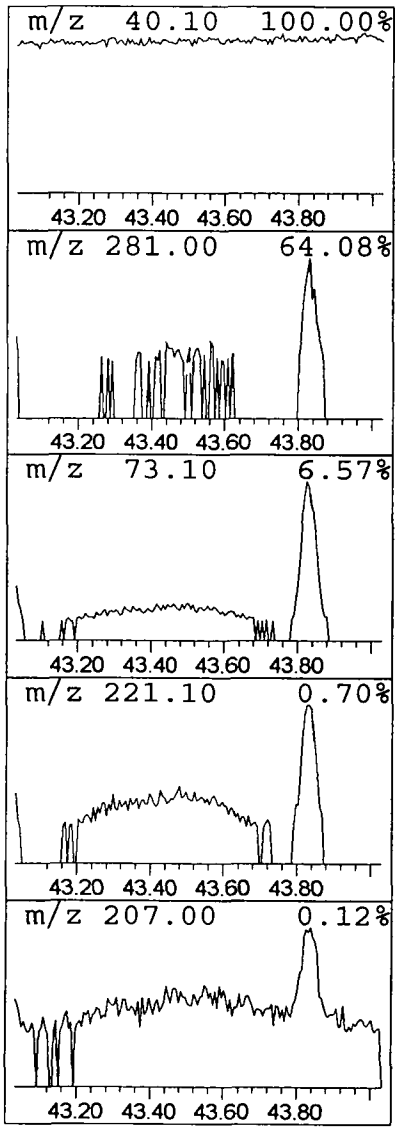
Peak Number 12 2-1-Phenyl ethylidene-hydra... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.53	0.11 ug/l	79954	Perylene-d12	42.30

Hit# of 2	Tentative ID	MW	MolForm	CAS#	Qual
1	2-1-Phenyl ethylidene-hydrazono-...	281	C16H15N3S	1000148-55-3	3
2	Adenosine, 2-methyl-	281	C11H15N5O4	016526-56-0	1





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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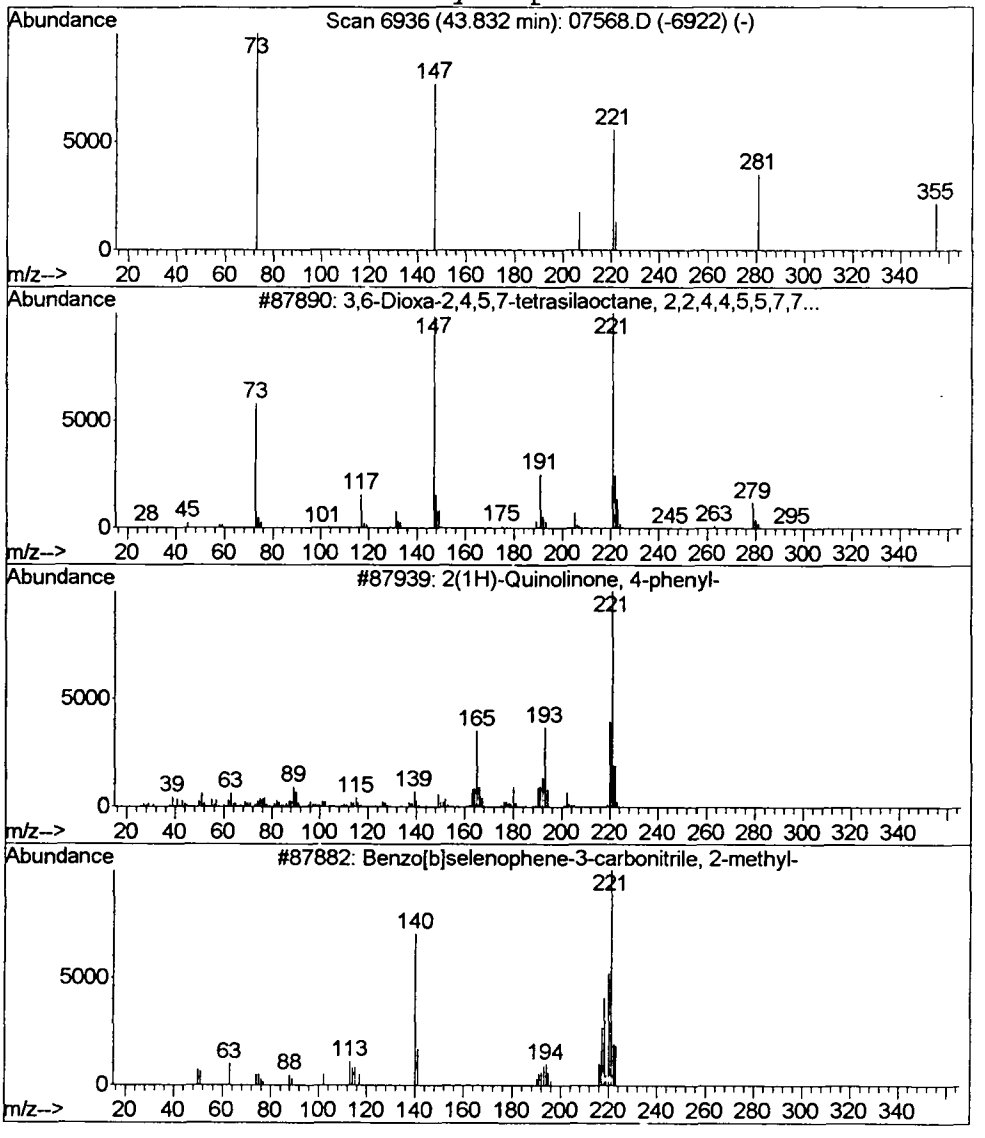
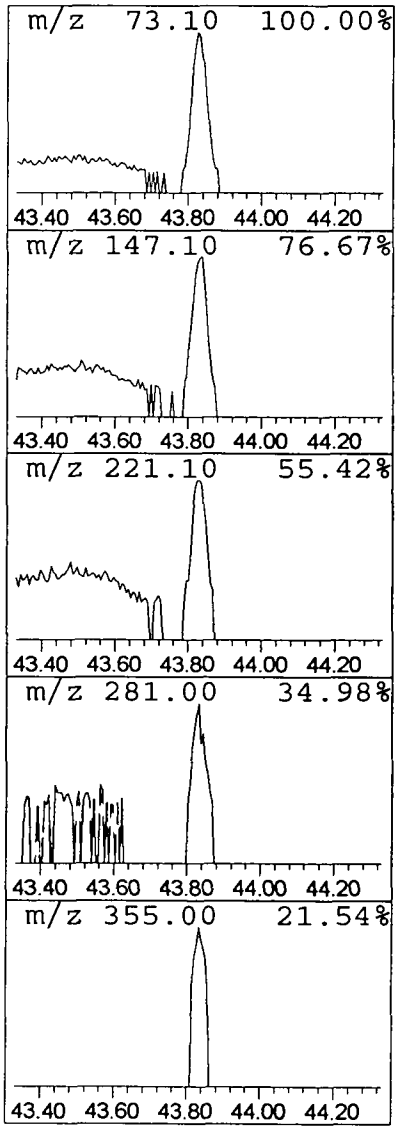
Peak Number 13 3,6-Dioxa-2,4,5,7-tetrasila... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.83	0.66 ug/l	487105	Perylene-d12	42.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	9
2	2(1H)-Quinolinone, 4-phenyl-	221	C15H11NO	005855-57-2	5
3	Benzo[b]selenophene-3-carbonitri...	221	C10H7NSe	037007-54-8	5
4	1-Aza-2-sila-5-boracyclopent-3-e...	221	C12H24BNSi	081620-70-4	5





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07568.D Vial: 16
Acq On : 3 Apr 2002 10:56 pm Operator: Nefliu
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title : Method 508 GC/MS Scan
Library : C:\DATABASE\nist98.l

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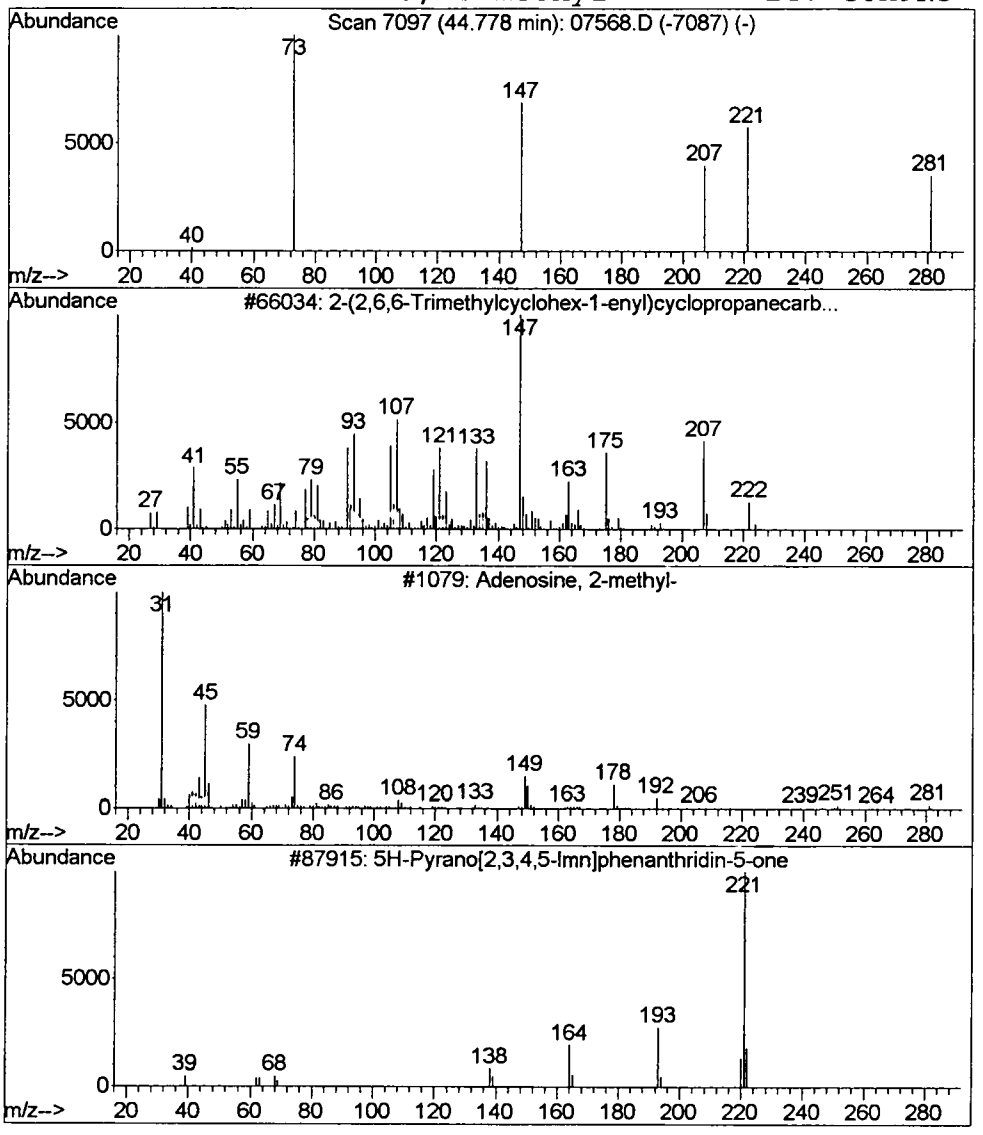
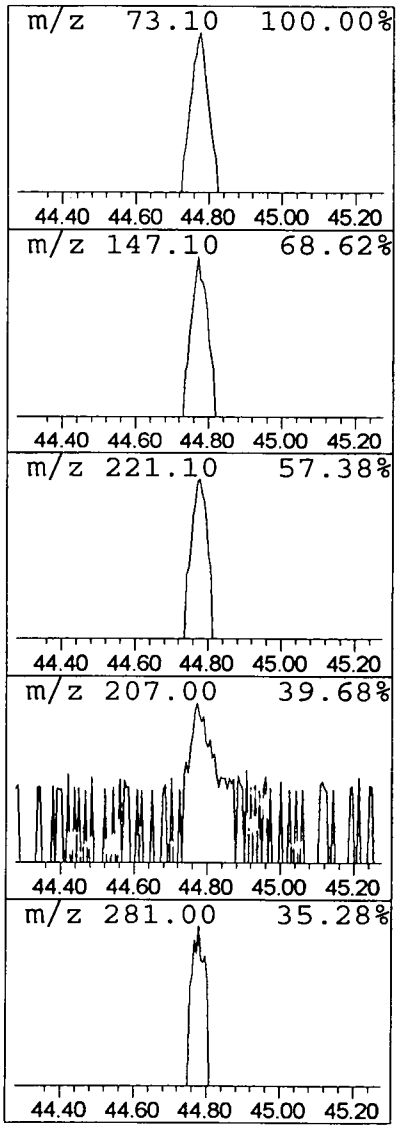
Peak Number 14 2-(2,6,6-Trimethylcyclohex-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
44.78	0.41 ug/l	299198	Perylene-d12	42.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(2,6,6-Trimethylcyclohex-1-eny...	222	C14H22O2	1000185-64-1	9
2	Adenosine, 2-methyl-	281	C11H15N5O4	016526-56-0	3
3	5H-Pyrano[2,3,4,5-lmn]phenanthri...	221	C14H7NO2	004733-28-2	2
4	2H-Indazol-3-amine, 2-methyl-	147	C8H9N3	1000142-18-7	2





Tentatively Identified Compound (LSC) summary

Operator ID: Nefliu Date Acquired: 3 Apr 2002 10:56 pm
Data File: C:\MSDCHEM\1\DATA\020403\07568.D
Name: sample
Misc:
Method: C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)
Title: Method 508 GC/MS Scan
Library Searched: C:\DATABASE\nist98.l

TIC Top Hit name	RT	EstConc	Units	Response	Internal Standard #	Internal Standard RT	Internal Standard Resp	Internal Standard Conc
Acetonitrile, dic...	3.31	0.5	ug/l	394584	1	10.66	14598100	20.0
Furan, 2,3-dihydr...	3.42	0.2	ug/l	115245	1	10.66	14598100	20.0
Acetic acid ethen...	5.59	0.3	ug/l	195338	1	10.66	14598100	20.0
Cyclopentasiloxan...	14.69	3.7	ug/l	3528210	2	15.77	19230500	20.0
1-(3,4-Methylened...	42.16	0.2	ug/l	178961	6	42.30	14657100	20.0
2-Amino-2-oxo-ace...	43.01	0.5	ug/l	369332	6	42.30	14657100	20.0
1H-1,2,4-Triazole...	43.27	0.2	ug/l	137446	6	42.30	14657100	20.0
Adenosine, 2-methyl-	43.29	0.2	ug/l	112683	6	42.30	14657100	20.0
Furo[3,4-b]furan-...	43.33	0.2	ug/l	120278	6	42.30	14657100	20.0
4-(3-Dimethylamin...	43.44	0.2	ug/l	144990	6	42.30	14657100	20.0
Adenosine, 2-methyl-	43.48	0.2	ug/l	139126	6	42.30	14657100	20.0
2-1-Phenyl ethyli...	43.53	0.1	ug/l	79954	6	42.30	14657100	20.0
3,6-Dioxa-2,4,5,7...	43.83	0.7	ug/l	487105	6	42.30	14657100	20.0
2-(2,6,6-Trimethy...	44.78	0.4	ug/l	299198	6	42.30	14657100	20.0