User: Baglioni, Walter
Westchester Dept. of Labs & Research
Date: 02/01/2002 Time: 14:19:17

Sample: AE00830
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 1/31/02 14:18 Ended: 2/1/02 14:18 Analyst: WB
Result source: manual entry
Page: 1

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

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Feb 01 08:46:16 2002 Quant Results File: SCAN508.RES

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Last Update : Fri Jan 11 13:20:07 2002
Response via : Initial Calibration
DataAcq Meth : SCAN508

POSSIBLE ADDITION OF 2x NORMAL AMT. OF [redacted]. INT. STD.

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.72	152	2941985	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	12275652	20.00	ug/L	-0.01
17) Acenaphthene-d10	18.35	164	6355278	20.00	ug/L	0.00
29) Phenanthrene-d10	22.66	188	10306726	20.00	ug/L	-0.01
38) Chrysene-d12	30.51	240	8972198	20.00	ug/L	0.00
44) Perylene-d12	34.44	264	6914690	20.00	ug/L	0.02
System Monitoring Compounds						
37) 2,4-DDD	27.73	235	303662	1.86	ug/L	0.00
Spiked Amount 5.000			Recovery =	37.20%		
Target Compounds						Qvalue
20) Dimethylphthalate	18.35	163	1043558	2.49	ug/L #	56
21) 2,6-Dinitrotoluene	18.35	165	805008	9.83	ug/L #	17
42) Bis(2-ethylhexyl)phthalate	31.09	149	25410	0.61	ug/L	99

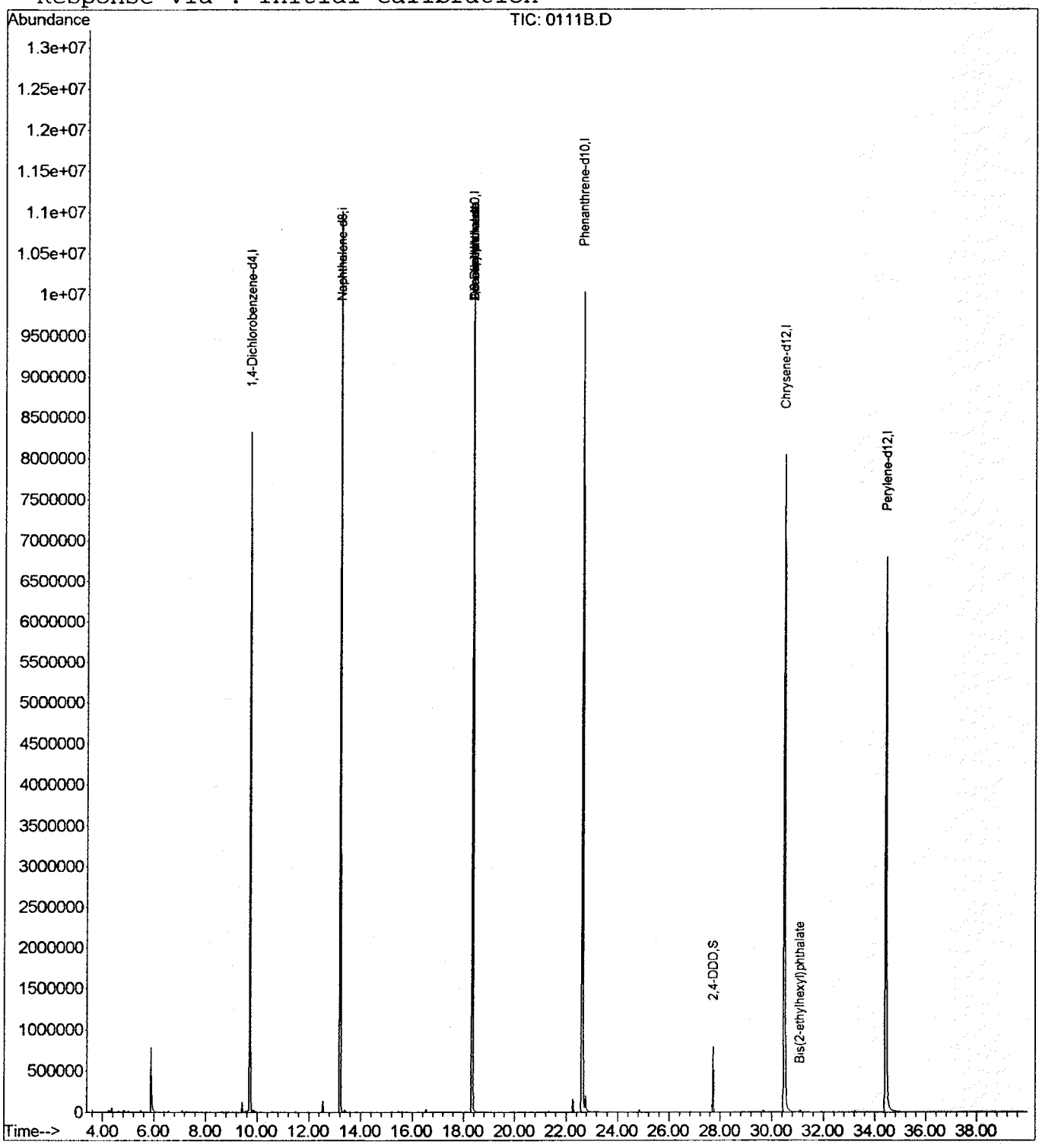
∴ REALLY SHOULD BE ≈ 74%

(#) = qualifier out of range (m) = manual integration
0111B.D SCAN508.M Fri Feb 01 08:46:17 2002 Page 1

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Feb 1 8:46 2002 Quant Results File: SCAN508.RE

Method : C:\MSDCHEM\1\METHODS\METHODS\SCAN508.M (RTE Integrator)
Title : 508 scan method
Last Update : Fri Jan 11 13:20:07 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\METHODS\METHODS\SCAN508.M (RTE Integrator)
Title : 508 scan method
Smoothing : OFF Filtering: 5
Sampling : 2 Min Area: 100 Area counts
Start Thrs: 0.2 Max Peaks: 20
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

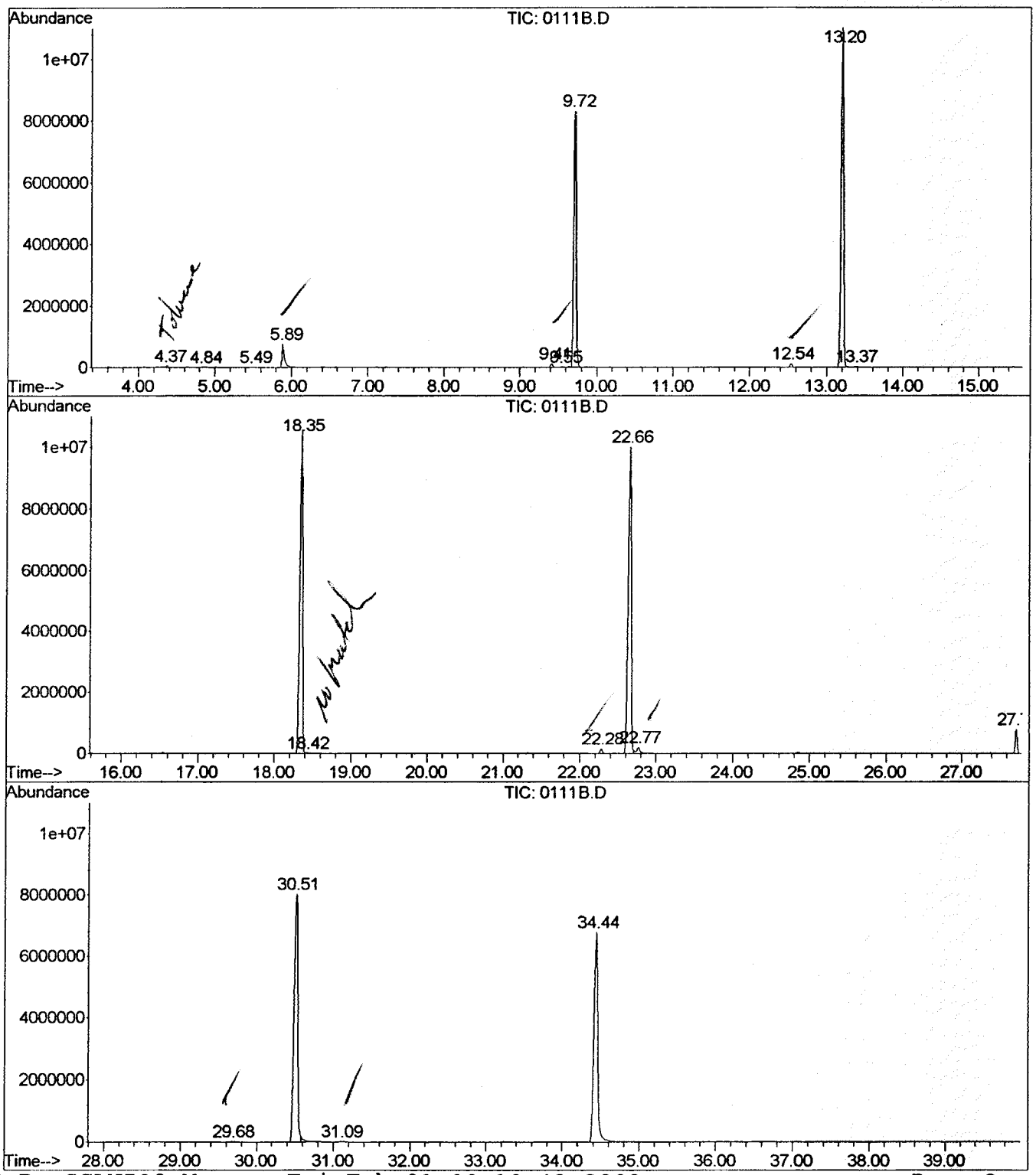
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.367	83	87	91	rVB	55196	105665	0.38%	0.070%
2	4.835	125	128	139	rBV2	27571	75439	0.27%	0.050%
3	5.486	181	185	197	rBV3	19434	68914	0.25%	0.046%
4	5.885	215	220	245	rVV	778357	1621072	5.87%	1.077%
5	9.413	525	529	539	rVB	125293	227568	0.82%	0.151%
6	9.550	539	541	547	rBV	26606	67434	0.24%	0.045%
7	9.722	551	556	561	rVV	8307141	17347657	62.77%	11.525%
8	12.541	797	803	807	rVV	134023	225191	0.81%	0.150%
9	13.204	855	861	865	rBV	11017835	24341311	88.08%	16.172%
10	13.375	873	876	891	rVB	36881	87608	0.32%	0.058%
11	18.353	1305	1312	1317	rBV	10577812	26049455	94.26%	17.306%
12	18.421	1317	1318	1331	rVB	19914	66656	0.24%	0.044%
13	22.280	1651	1656	1661	rBV2	159874	294127	1.06%	0.195%
14	22.657	1681	1689	1693	rBV2	10018495	27636100	100.00%	18.360%
15	22.771	1693	1699	1711	rVB	199802	606415	2.19%	0.403%
16	27.726	2127	2133	2137	rBV	792606	1479814	5.35%	0.983%
17	29.678	2299	2304	2313	rVV2	24088	70209	0.25%	0.047%
18	30.512	2369	2377	2383	rBV2	8035811	26447695	95.70%	17.571%
19	31.094	2421	2428	2443	rVB6	25821	111083	0.40%	0.074%
20	34.439	2713	2721	2755	rBV2	6782933	23589968	85.36%	15.672%

Sum of corrected areas: 150519381

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D
Operator :
Acquired : 31 Jan 2002 10:31 am using AcqMethod SCAN508
Instrument : Instrumen
Sample Name: lab blank 1/11
Misc Info : January 31, 2002
Vial Number: 1
Quant File :SCAN508.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

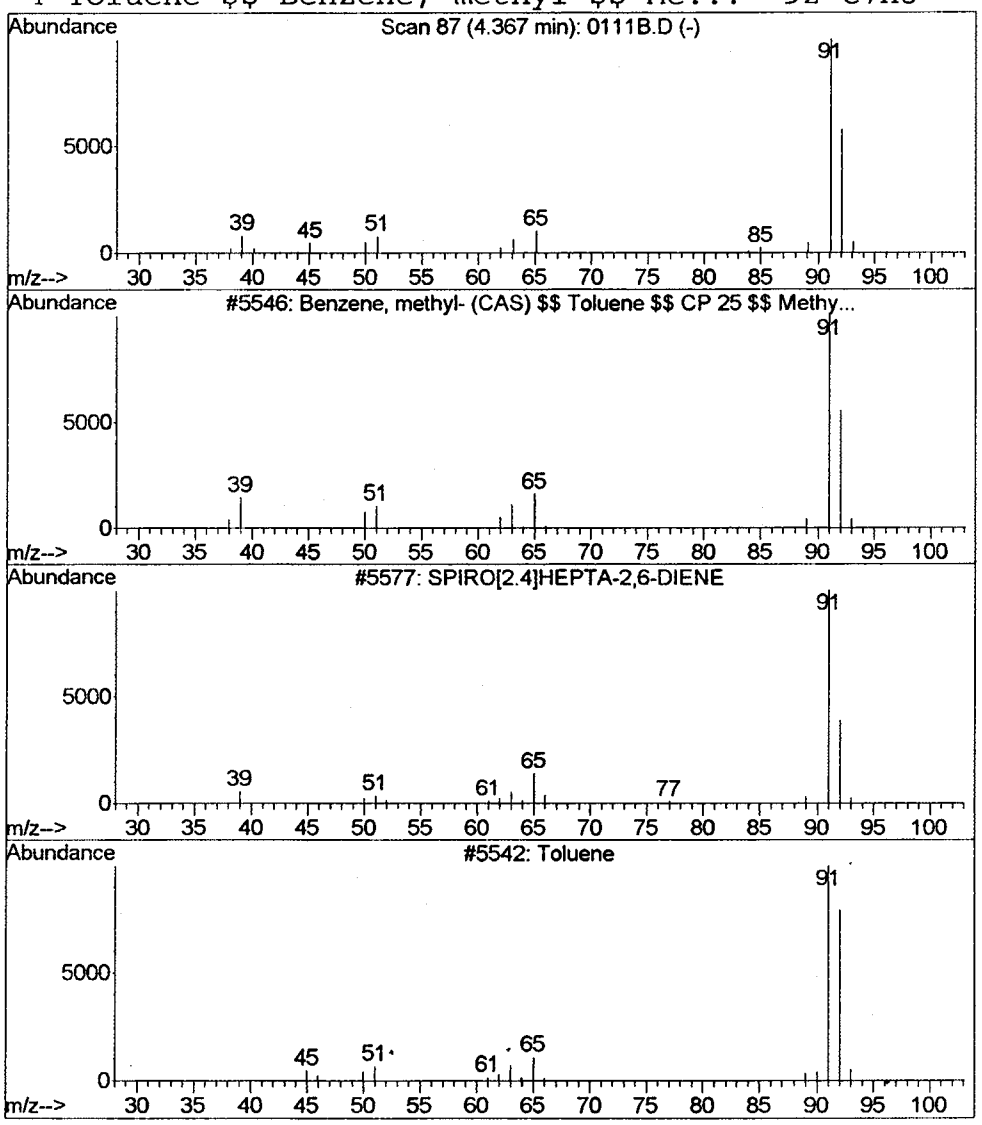
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

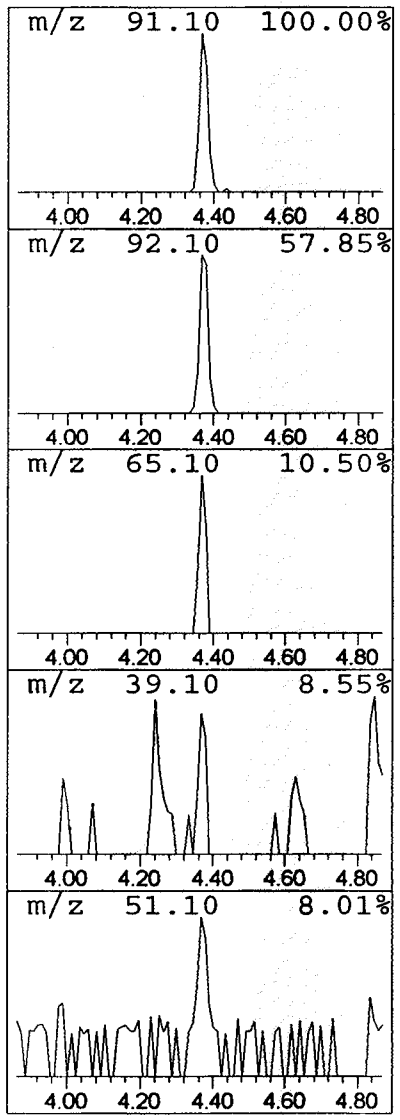
****

Peak Number 1 Benzene, methyl- (CAS) $$ T... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	0.12 ug/L	105665	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, methyl- (CAS) $$ Toluen...	92	C7H8	000108-88-3	90
2	SPIRO[2.4]HEPTA-2,6-DIENE	92	C7H8	000000-00-0	90
3	Toluene	92	C7H8	000108-88-3	90
4	Toluene $$ Benzene, methyl $$ Me...	92	C7H8	000108-88-3	90





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

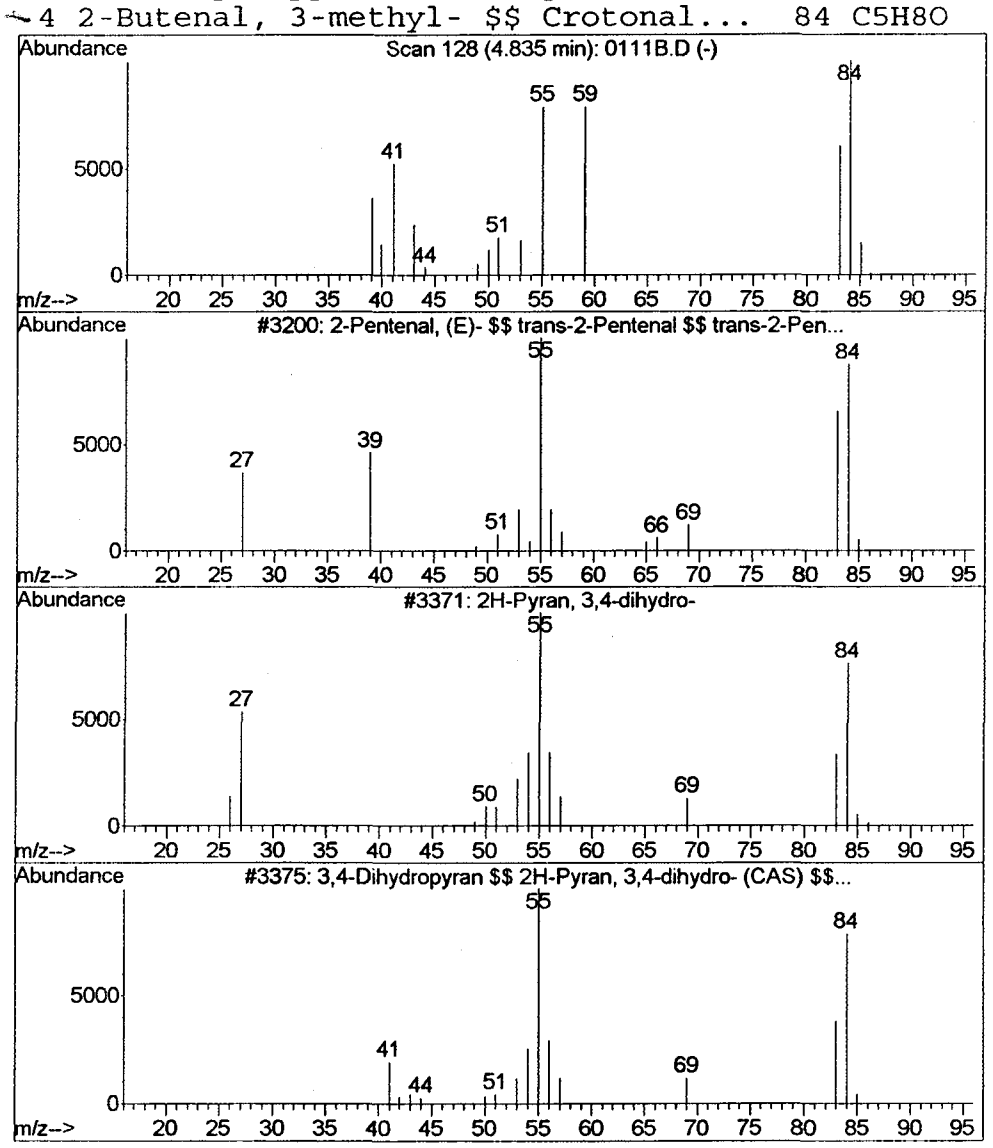
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

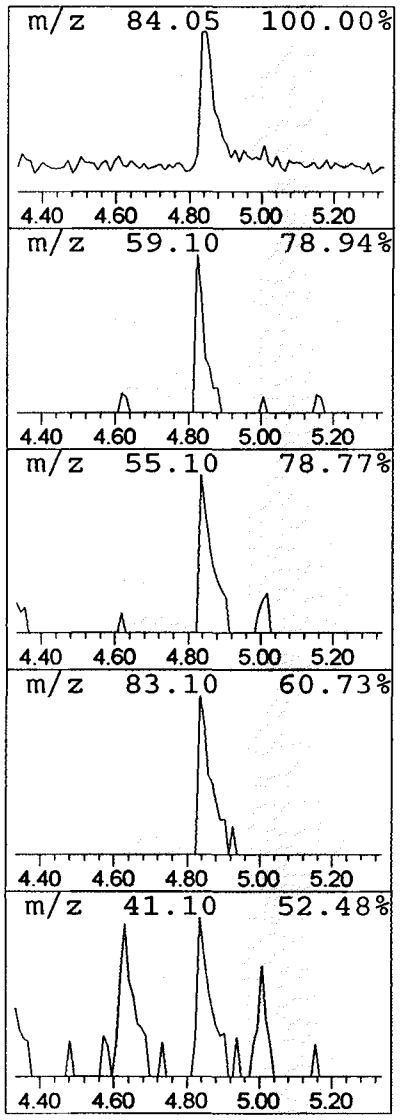
****

Peak Number 2 2-Pentenal, (E)- $$ trans-2... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	0.09 ug/L	75439	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentenal, (E)- $$ trans-2-Pent...	84	C5H8O	001576-87-0	49
2	2H-Pyran, 3,4-dihydro-	84	C5H8O	000110-87-2	47
3	3,4-Dihydropyran $$ 2H-Pyran, 3,...	84	C5H8O	000110-87-2	47
4	2-Butenal, 3-methyl- $$ Crotonal...	84	C5H8O	000107-86-8	47





Library Search Compound Report

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Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

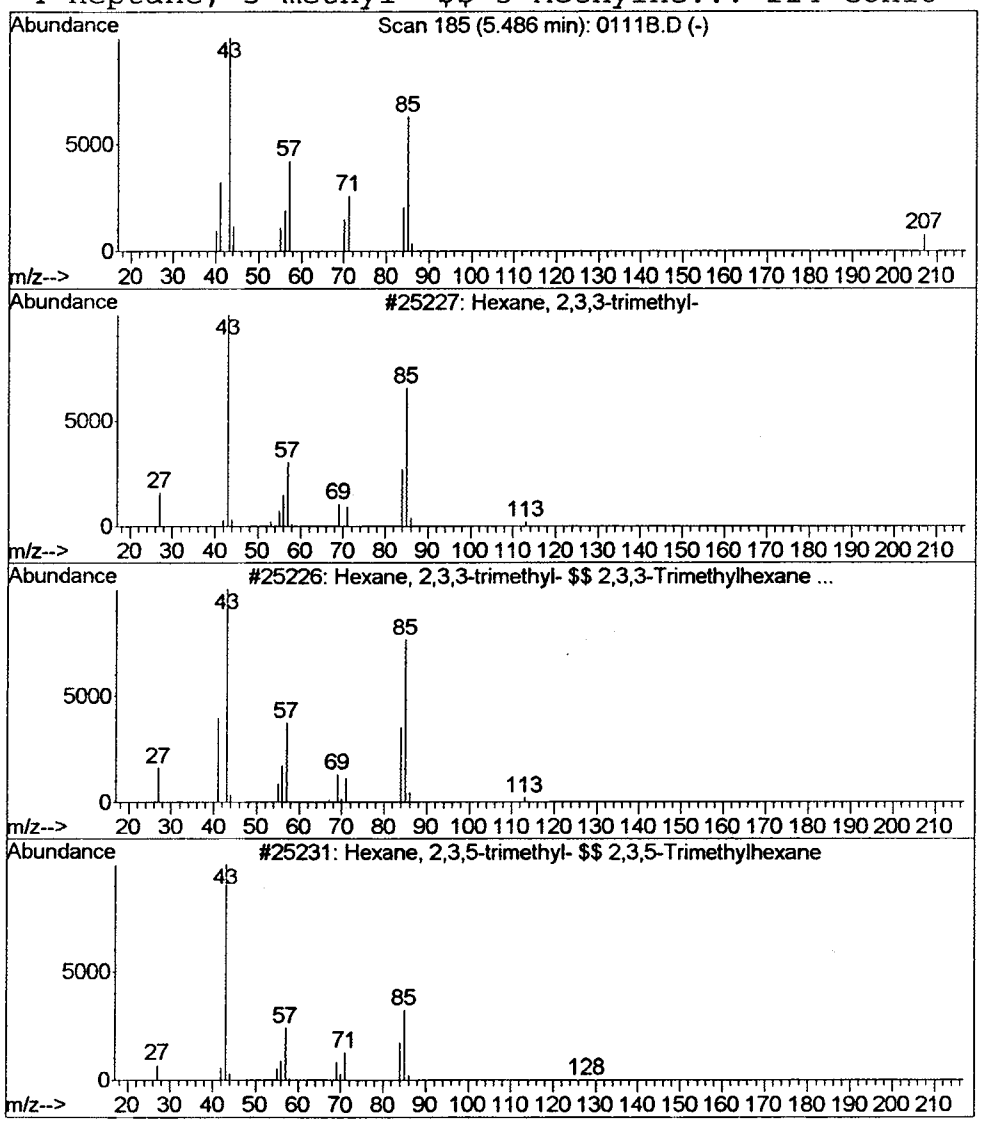
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

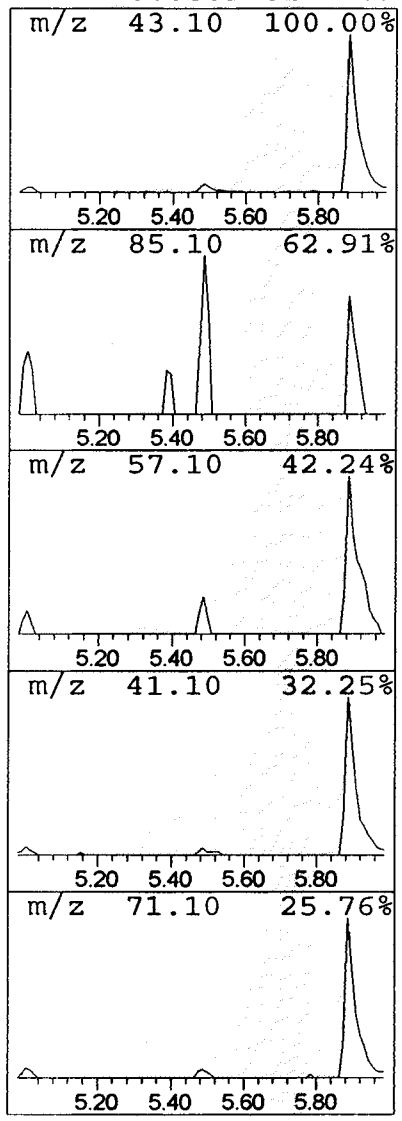
****

Peak Number 3 Hexane, 2,3,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.49	0.08 ug/L	68914	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,3,3-trimethyl-	128	C9H20	016747-28-7	64
2	Hexane, 2,3,3-trimethyl- $$ 2,3,...	128	C9H20	016747-28-7	59
3	Hexane, 2,3,5-trimethyl- $$ 2,3,...	128	C9H20	001069-53-0	59
4	Heptane, 3-methyl- $$ 3-Methylhe...	114	C8H18	000589-81-1	59





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

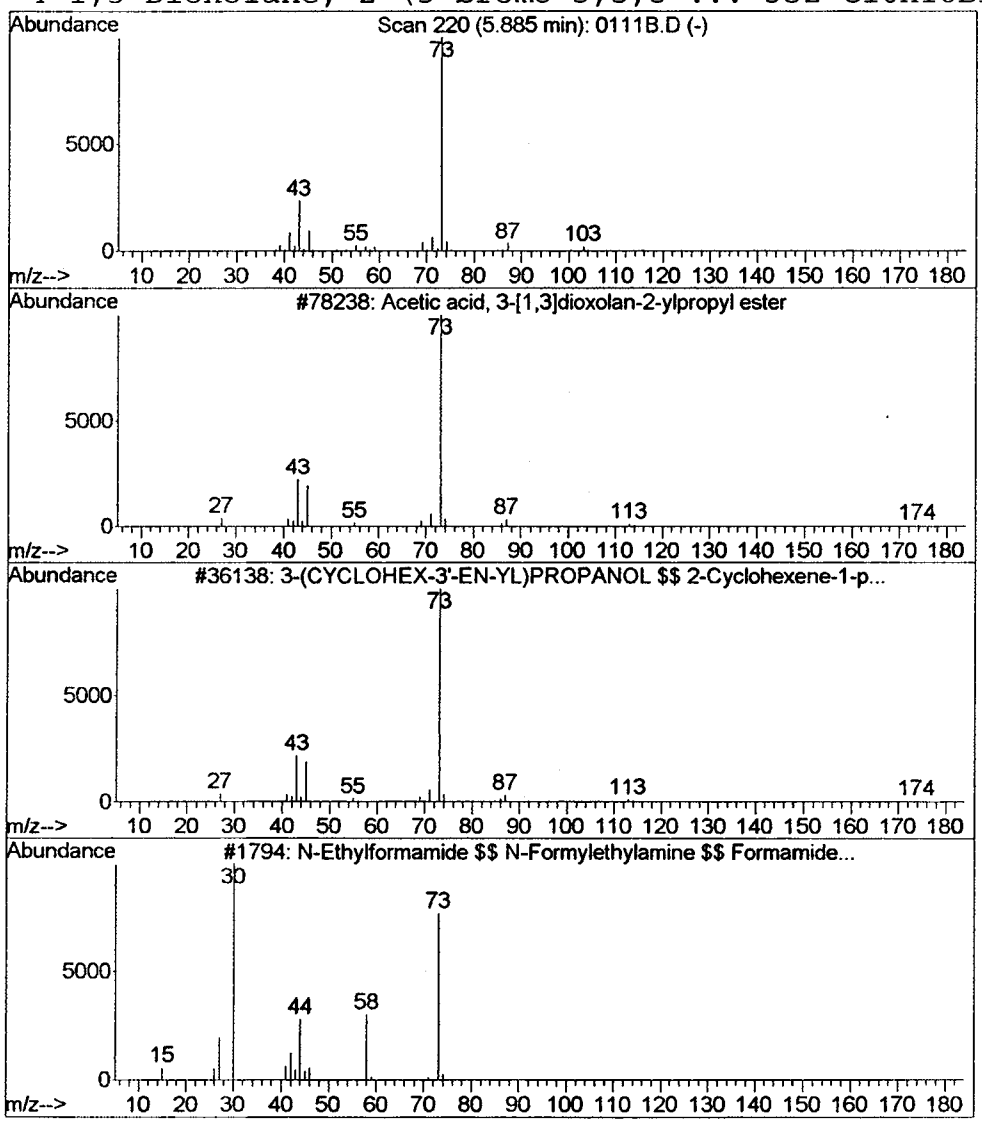
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

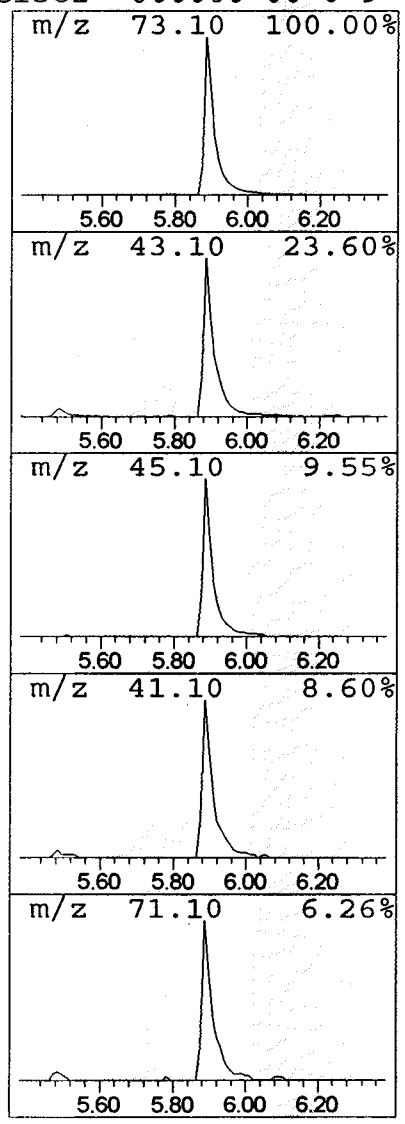
****

Peak Number 4 Acetic acid, 3-[1,3]dioxola... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	1.87 ug/L	1621070	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33
2	3-(CYCLOHEX-3'-EN-YL)PROPANOL $$...	174	C8H14O4	015745-87-6	33
3	N-Ethylformamide $$ N-Formylethy...	73	C3H7NO	000627-45-2	9
4	1,3-Dioxolane, 2-(3-bromo-5,5,5-...	352	C10H16BrCl3O2	000000-00-0	9





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

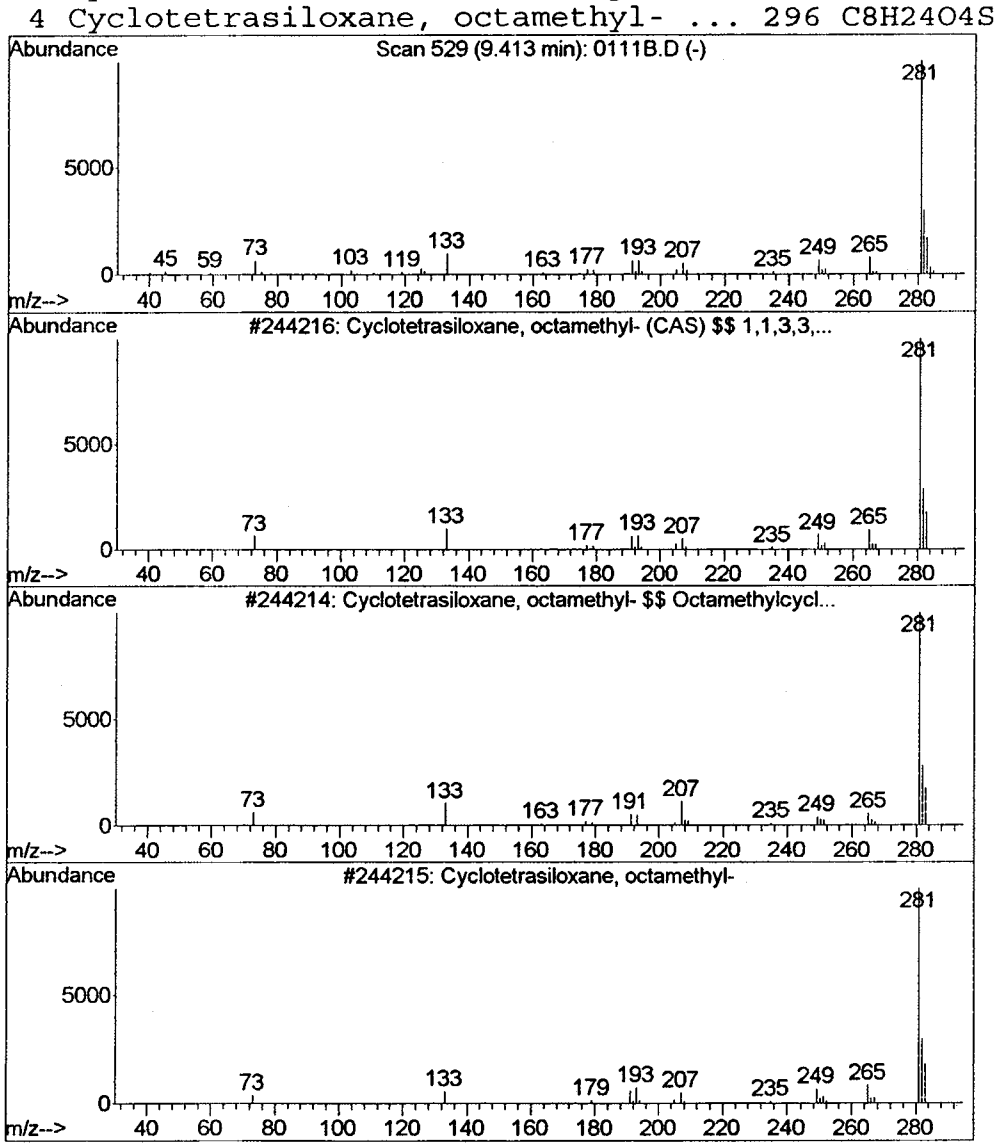
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

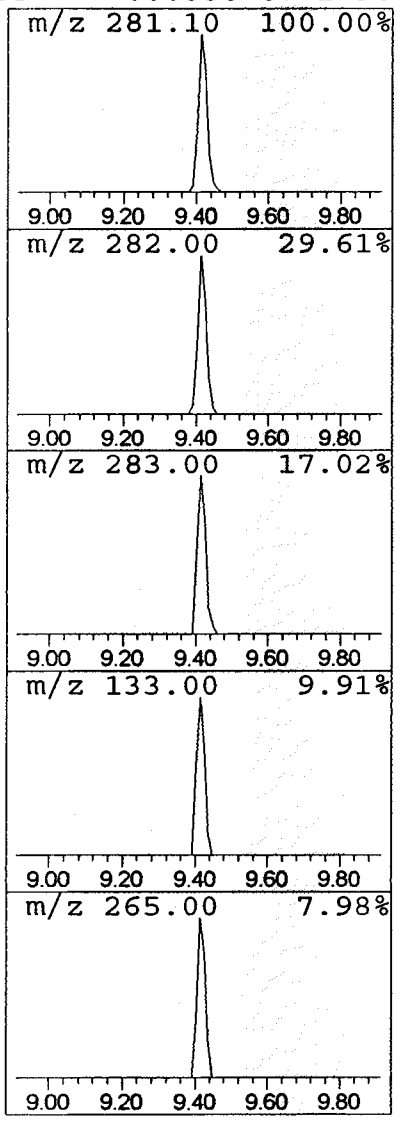
****

Peak Number 5 Cyclotetrasiloxane, octamet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.41	0.26 ug/L	227568	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetrasiloxane, octamethyl- ...	296	C8H24O4Si4	000556-67-2	91
2	Cyclotetrasiloxane, octamethyl- ...	296	C8H24O4Si4	000556-67-2	91
3	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	90
4	Cyclotetrasiloxane, octamethyl- ...	296	C8H24O4Si4	000556-67-2	90





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

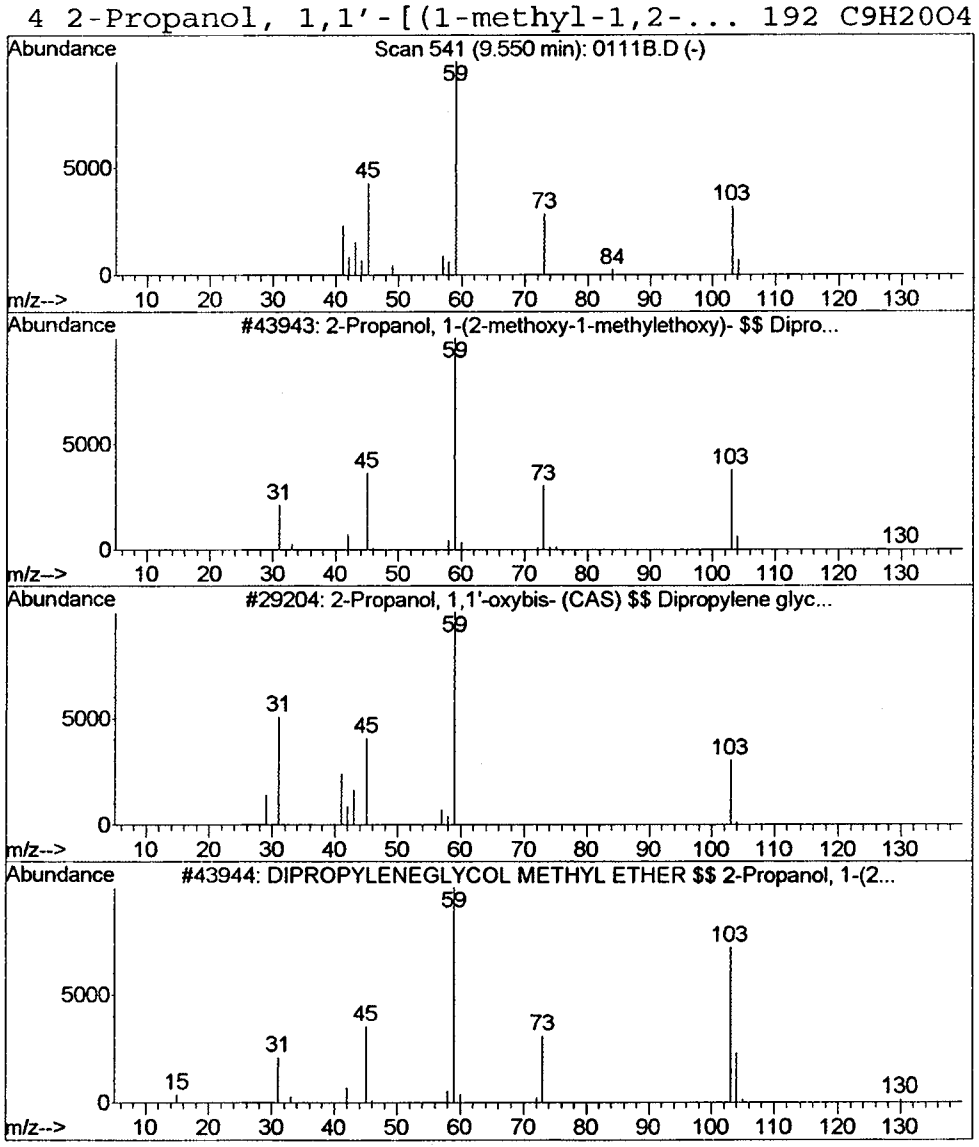
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

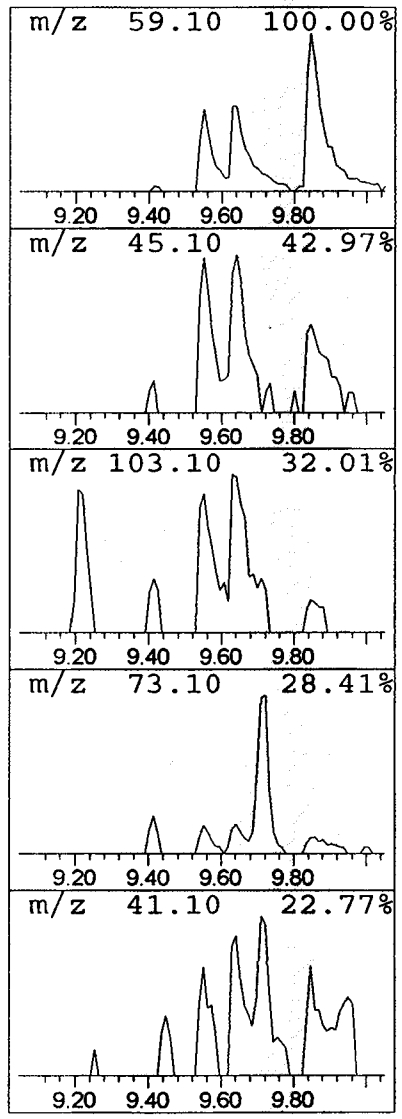
****

Peak Number 6 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.55	0.08 ug/L	67434	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	50
2	2-Propanol, 1,1'-oxybis- (CAS) $...	134	C6H14O3	000110-98-5	50
3	DIPROPYLENEGLYCOL METHYL ETHER $...	148	C7H16O3	020324-32-7	50
4	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	42





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

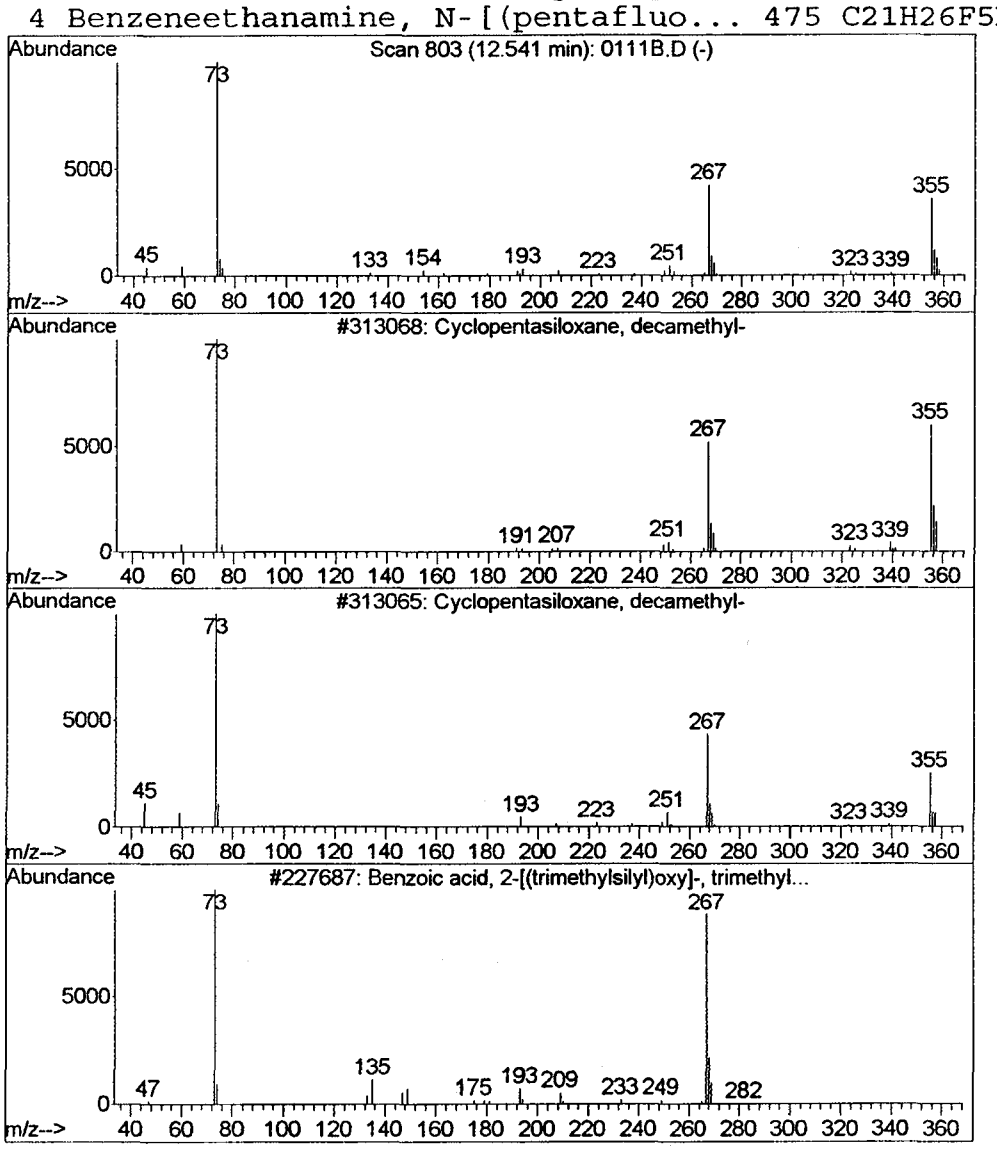
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

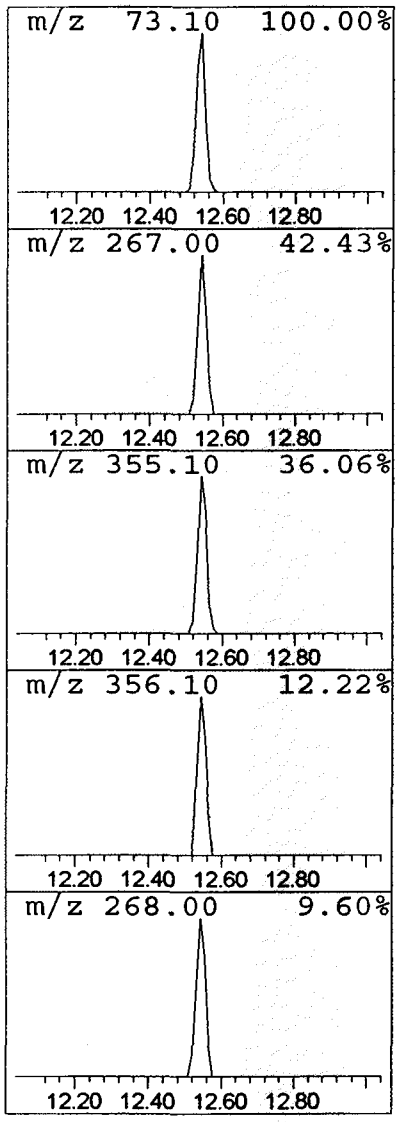
****

Peak Number 7 Cyclopentasiloxane, decamet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.19 ug/L	225191	Naphthalene-d8	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	80
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	43
3	Benzoic acid, 2-[(trimethylsilyl...	282	C13H22O3Si2	003789-85-3	38
4	Benzeneethanamine, N-[(pentafluo...	475	C21H26F5NO2Si2	055429-85-1	38





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

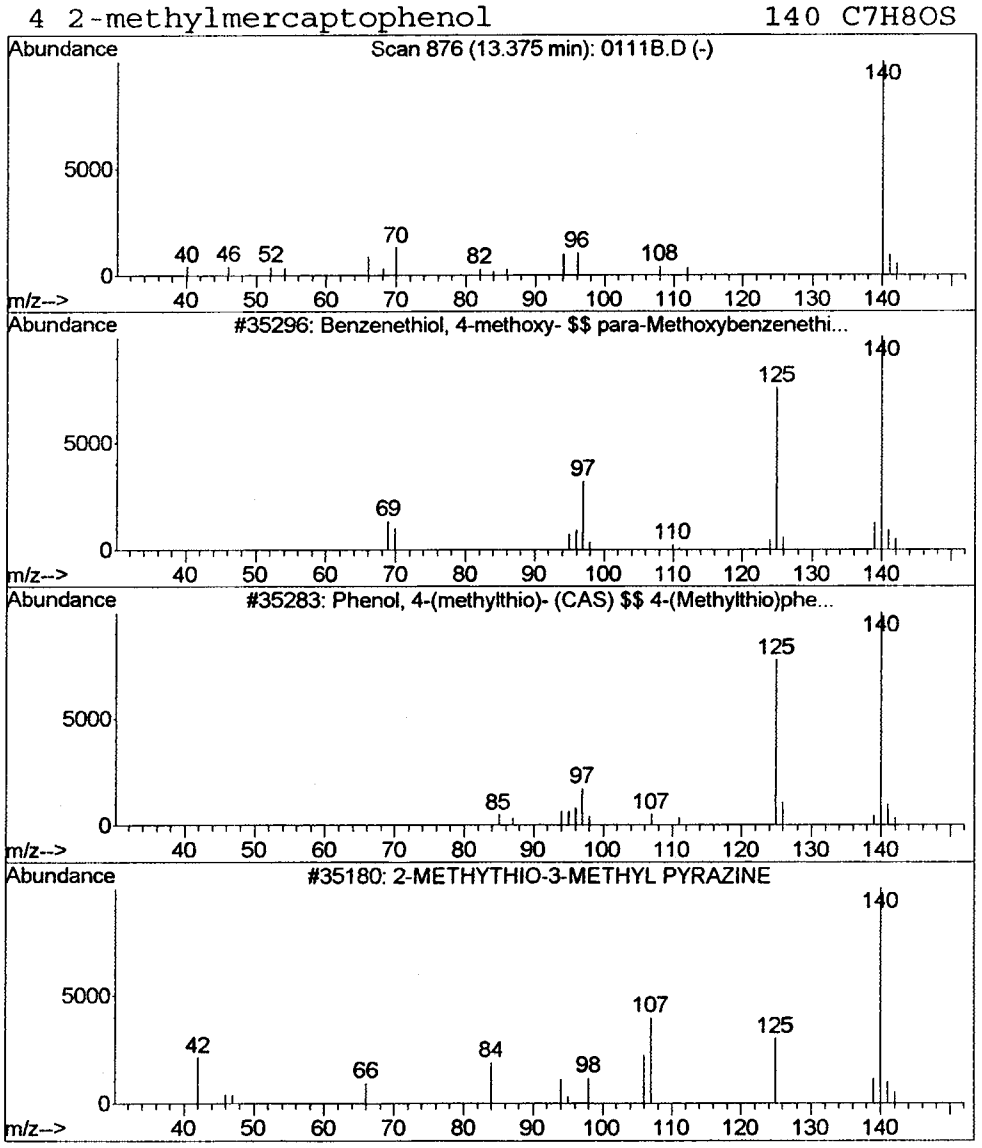
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

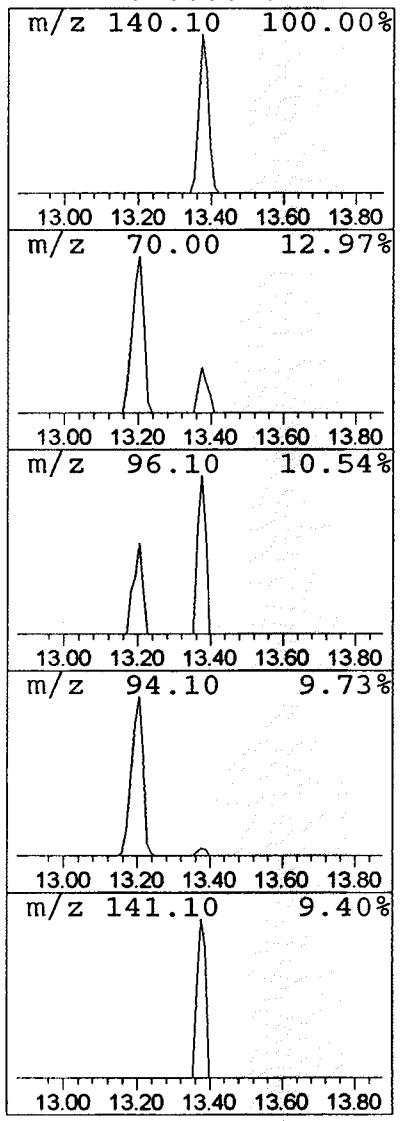
****

Peak Number 8 Benzenethiol, 4-methoxy- $$... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	0.07 ug/L	87608	Naphthalene-d8	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenethiol, 4-methoxy- $$ para...	140	C7H8OS	000696-63-9	64
2	Phenol, 4-(methylthio)- (CAS) $$...	140	C7H8OS	001073-72-9	64
3	2-METHYTHIO-3-METHYL PYRAZINE	140	C6H8N2S	002882-20-4	64
4	2-methylmercaptophenol	140	C7H8OS	000000-00-0	59





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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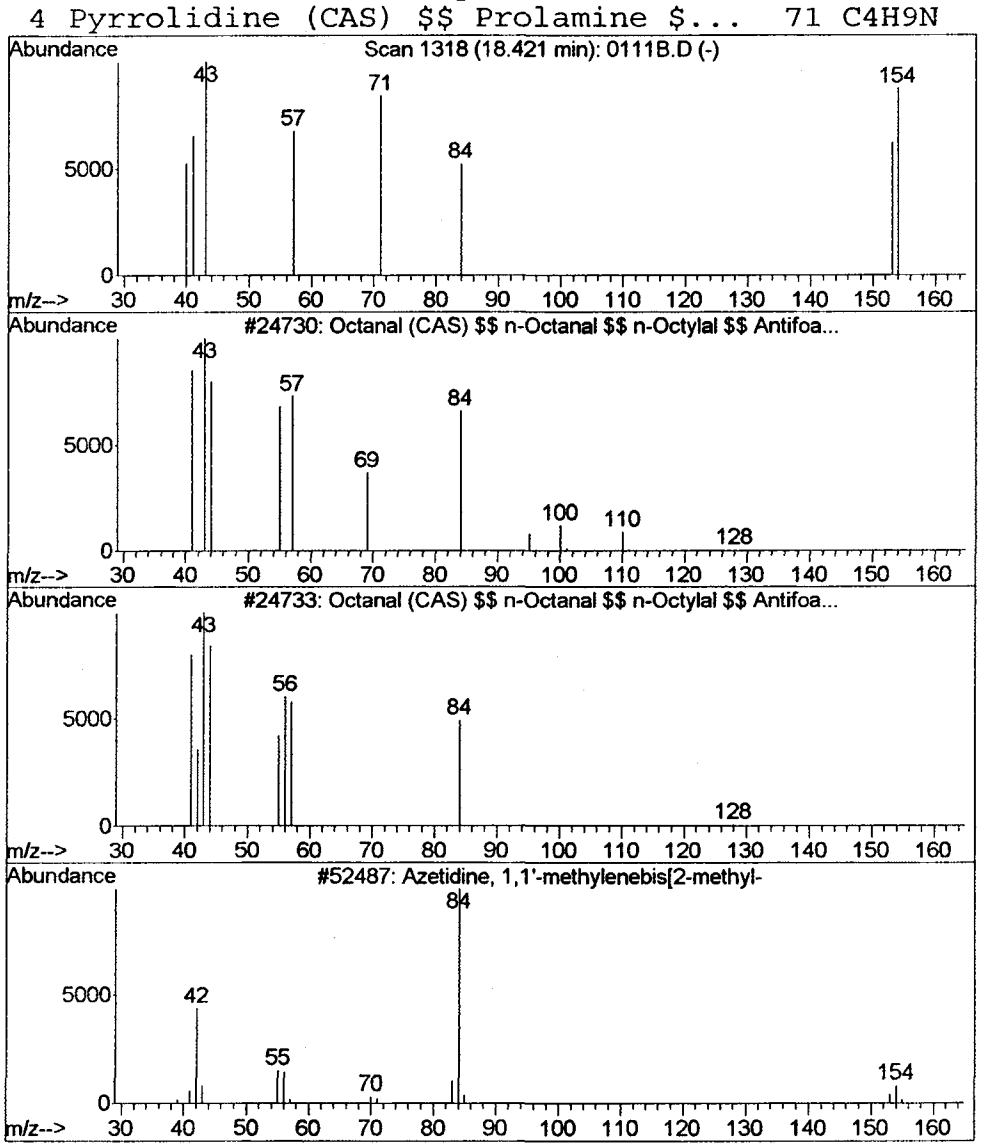
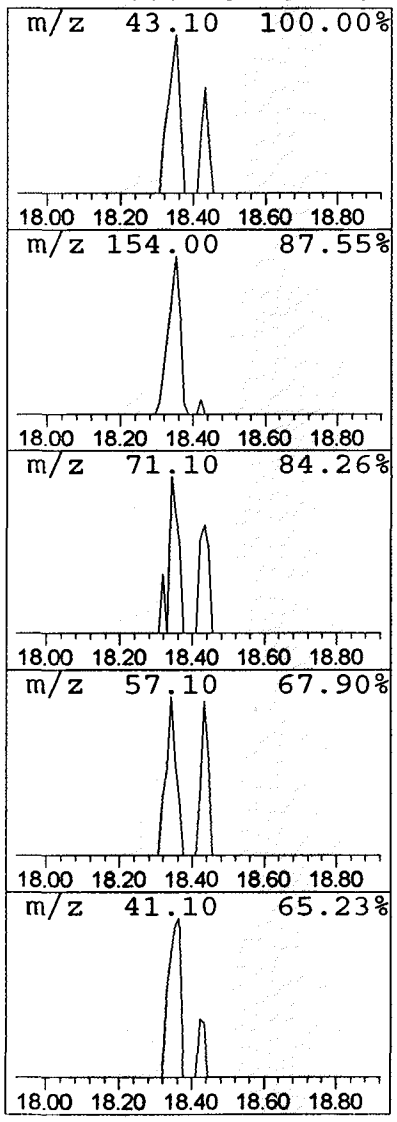
Peak Number 9 Octanal (CAS) $$ n-Octanal ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.42	0.05 ug/L	66656	Acenaphthene-d10	18.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanal (CAS) $$ n-Octanal $$ n-...	128	C8H16O	000124-13-0	9
2	Octanal (CAS) $$ n-Octanal $$ n-...	128	C8H16O	000124-13-0	9
3	Azetidine, 1,1'-methylenebis[2-m...	154	C9H18N2	038455-30-0	7
4	Pyrrolidine (CAS) $$ Prolamine $...	71	C4H9N	000123-75-1	5





Library Search Compound Report

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Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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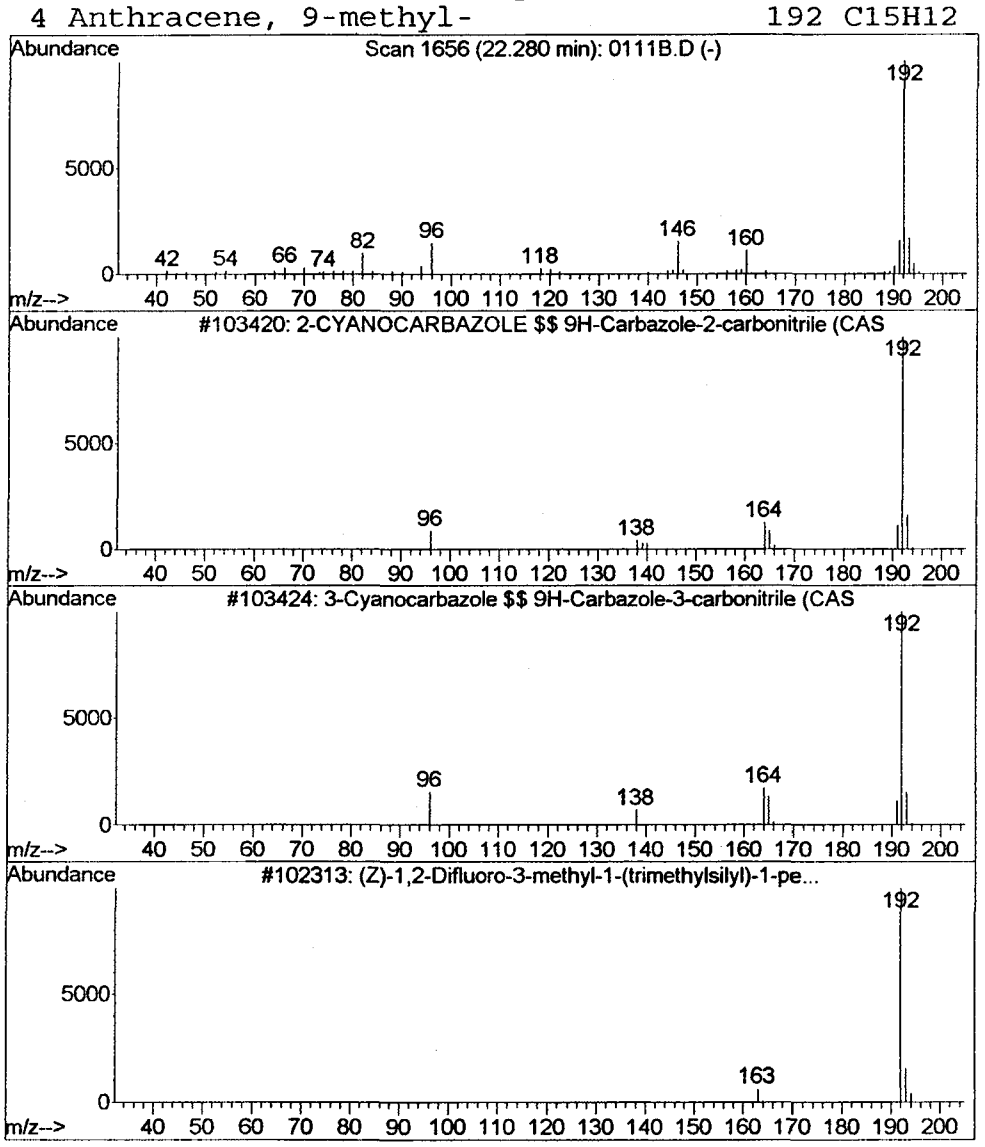
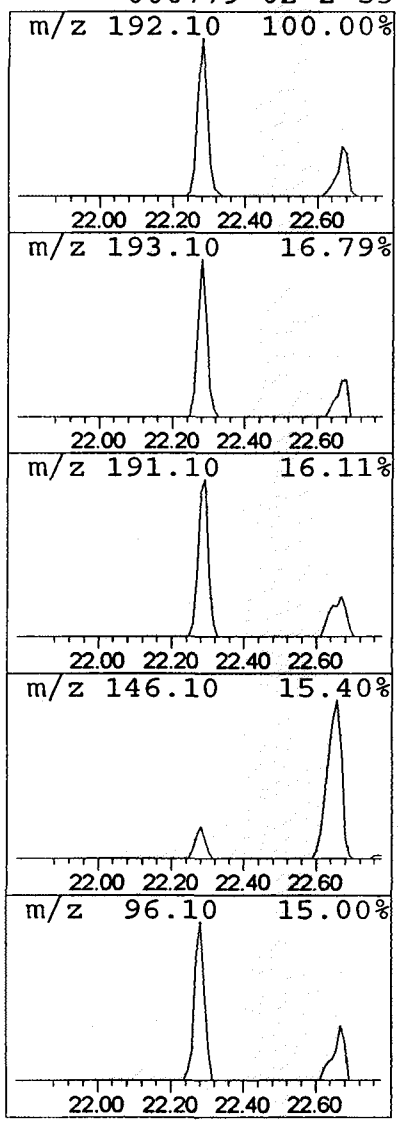
Peak Number 10 2-CYANOCARBAZOLE $$ 9H-Carb... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.21 ug/L	294127	Phenanthrene-d10	22.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-CYANOCARBAZOLE $$ 9H-Carbazole...	192	C13H8N2	057955-18-7	59
2	3-Cyanocarbazole $$ 9H-Carbazole...	192	C13H8N2	057102-93-9	59
3	(Z)-1,2-Difluoro-3-methyl-1-(tri...	192	C9H18F2Si	000000-00-0	56
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	53





Library Search Compound Report

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Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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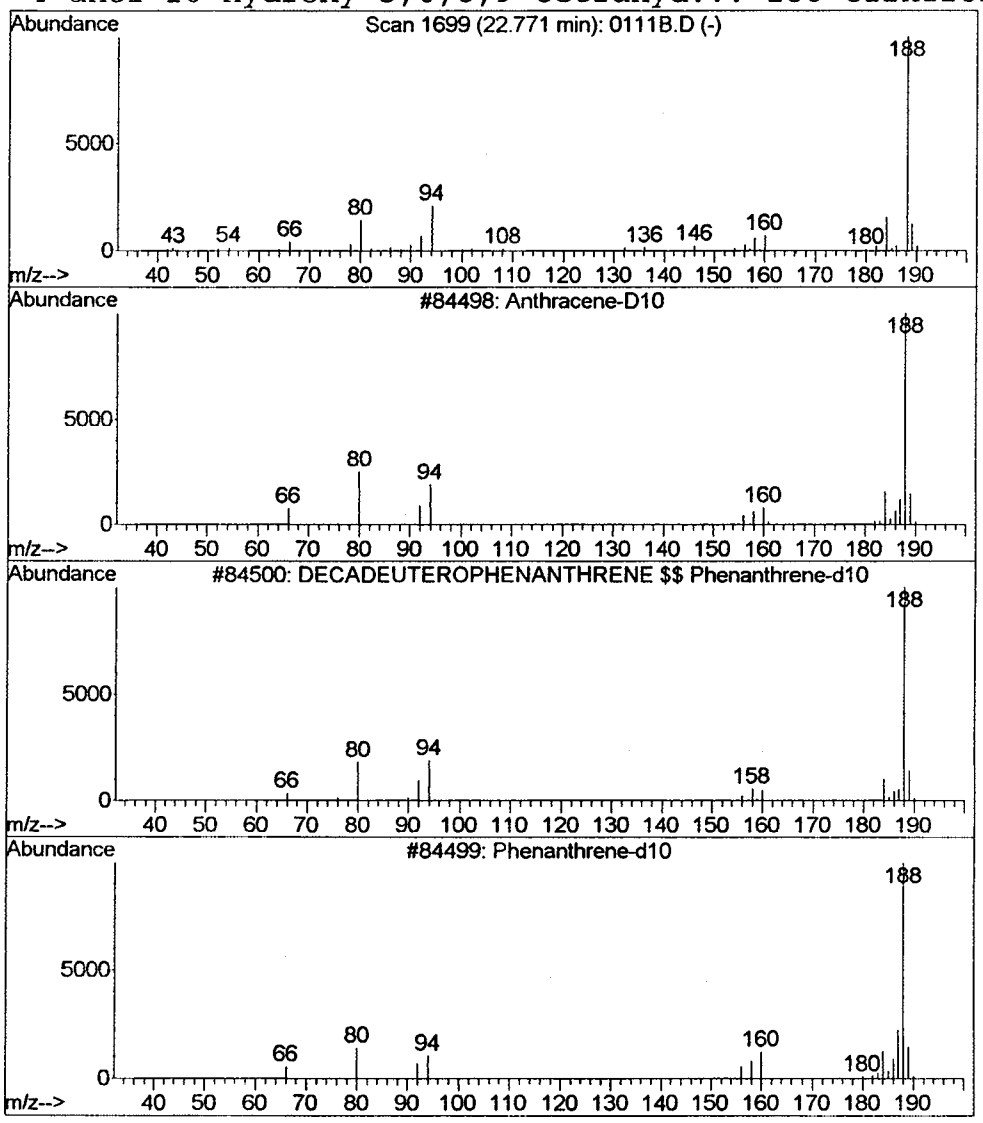
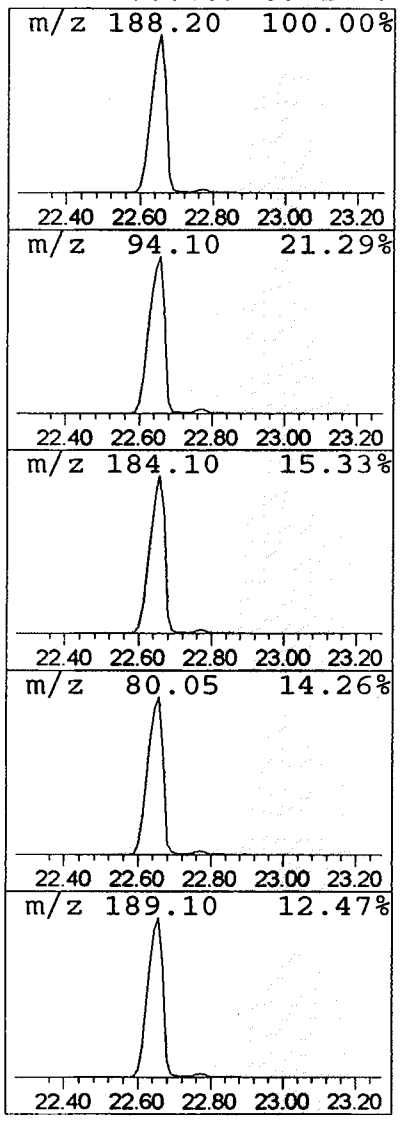
Peak Number 11 Anthracene-D10 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.44 ug/L	606415	Phenanthrene-d10	22.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-D10	188	C14D10	000000-00-0	86
2	DECADEUTEROPHENANTHRENE $$ Phena...	188	C14D10	001517-22-2	72
3	Phenanthrene-d10	188	C14D10	001517-22-2	64
4	anti-10-hydroxy-5,6,8,9-tetrahyd...	188	C12H12O2	055139-53-2	49





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\0111B.D Vial: 1
Acq On : 31 Jan 2002 10:31 am Operator:
Sample : lab blank 1/11 Inst : Instrumen
Misc : January 31, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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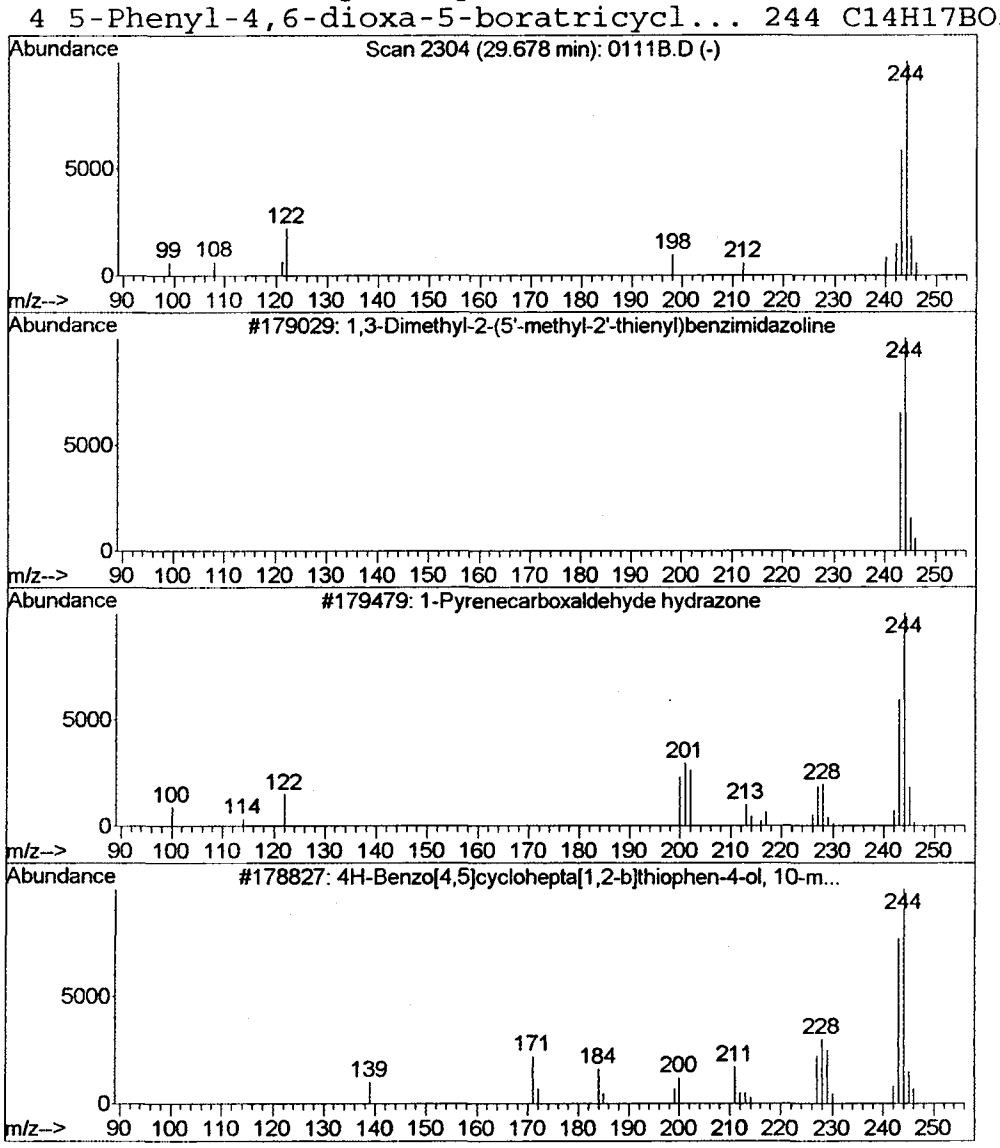
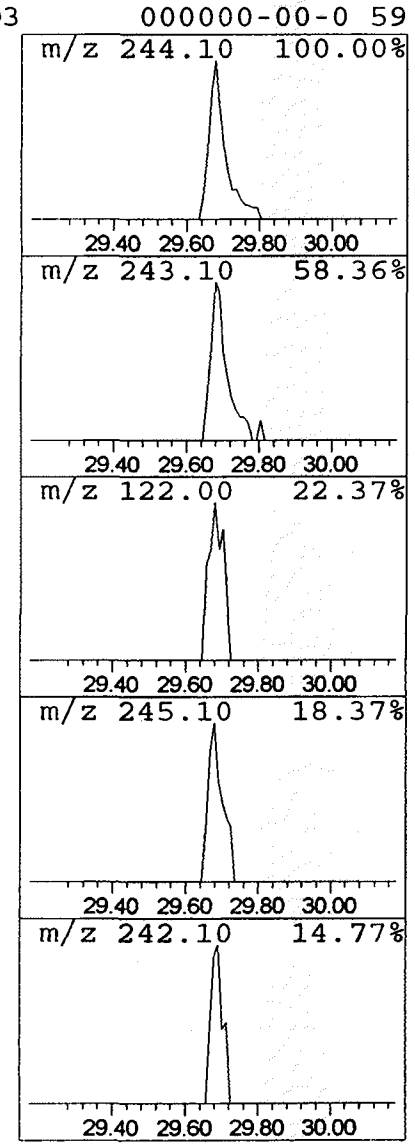
Peak Number 12 1,3-Dimethyl-2-(5'-methyl-2... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.68	0.05 ug/L	70209	Chrysene-d12	30.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dimethyl-2-(5'-methyl-2'-thi...	244	C14H16N2S	000000-00-0	83
2	1-Pyrenecarboxaldehyde hydrazone	244	C17H12N2	076465-53-7	72
3	4H-Benzo[4,5]cyclohepta[1,2-b]th...	244	C14H12O2S	000000-00-0	64
4	5-Phenyl-4,6-dioxa-5-boratricycl...	244	C14H17BO3	000000-00-0	59





Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 31 Jan 2002 10:31 am
Data File: C:\MSDCHEM\1\5973N\02JAN31\0111B.D
Name: lab blank 1/11
Misc: January 31, 2002
Method: C:\MSDCHEM\1\METHODS\METHODS\SCAN508.M (RTE Integrator)
Title: 508 scan method
Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	Internal Standard #	Internal Standard RT	Internal Standard Resp	Internal Standard Conc
Benzene, methyl- ...	4.37	0.1	ug/L	105665	1	9.72	17347700	20.0
2-Pentenal, (E)- ...	4.84	0.1	ug/L	75439	1	9.72	17347700	20.0
Hexane, 2,3,3-tri...	5.49	0.1	ug/L	68914	1	9.72	17347700	20.0
Acetic acid, 3-[1...	5.89	1.9	ug/L	1621070	1	9.72	17347700	20.0
Cyclotetrasiloxan...	9.41	0.3	ug/L	227568	1	9.72	17347700	20.0
2-Propanol, 1-(2-...	9.55	0.1	ug/L	67434	1	9.72	17347700	20.0
Cyclopentasiloxan...	12.54	0.2	ug/L	225191	2	13.20	24341300	20.0
Benzenethiol, 4-m...	13.37	0.1	ug/L	87608	2	13.20	24341300	20.0
Octanal (CAS) $$ ...	18.42	0.1	ug/L	66656	3	18.35	26049500	20.0
2-CYANOCARBAZOLE ...	22.28	0.2	ug/L	294127	4	22.66	27636100	20.0
Anthracene-D10	22.77	0.4	ug/L	606415	4	22.66	27636100	20.0
1,3-Dimethyl-2-(5...	29.68	0.1	ug/L	70209	5	30.51	26447700	20.0