User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/28/2002 Time: 17:05:31

Sample: AE03510
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/28/02 17:05 Ended: 3/28/02 17:05 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		84		5.0

Comments for Sample AE03510 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 03-28-2002 Time: 17:05:38

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 6 of the 43 compounds in the LCS Standard were high.

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Mar 15 10:05:11 2002 Quant Results File: HP022102.RES

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Last Update : Mon Feb 25 10:47:59 2002
Response via : Initial Calibration
DataAcq Meth : HP022102

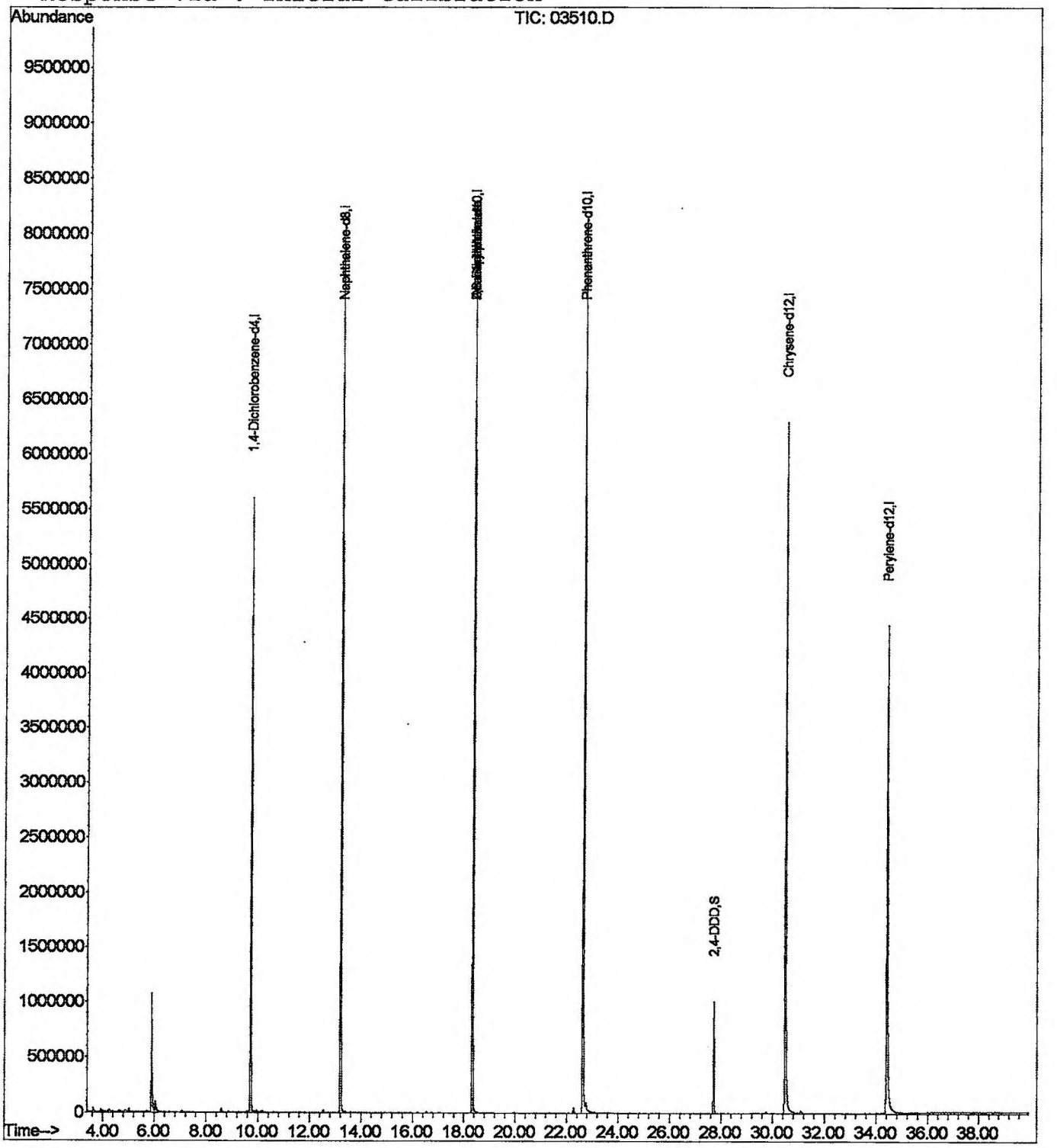
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.71	152	1866527	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	8236577	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	4094716	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	6940644	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	5228806	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	3324914	20.00	ug/L	-0.04
System Monitoring Compounds						
37) 2,4-DDD	27.72	235	388790	4.20	ug/L	0.00
Spiked Amount 5.000			Recovery = 84.00%			
Target Compounds						Qvalue
20) Dimethylphthalate	18.33	163	656477	2.64	ug/L #	59
21) 2,6-Dinitrotoluene	18.33	165	528287	9.21	ug/L #	27

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

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Mar 15 10:05 2002 Quant Results File: HP022102.R

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title : 508 scan method
Last Update : Mon Feb 25 10:47:59 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.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

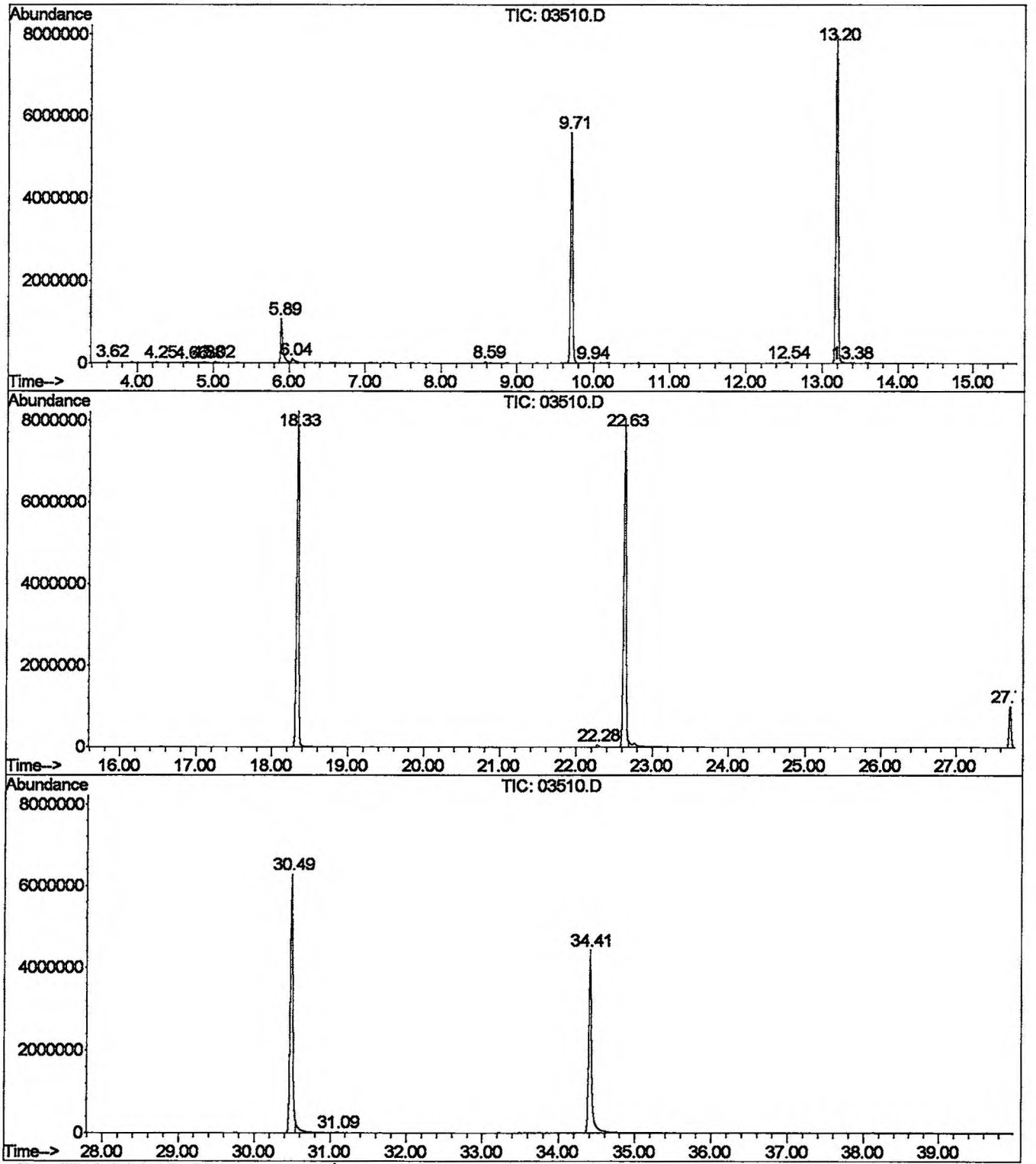
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.622	19	22	29	rBV2	44970	75716	0.43%	0.082%
2	4.254	75	78	85	rVB4	22084	55631	0.32%	0.060%
3	4.661	107	114	121	rVB6	12619	57111	0.33%	0.062%
4	4.875	127	133	143	rBV3	16797	59784	0.34%	0.065%
5	5.022	143	146	153	rVB	31548	55629	0.32%	0.060%
6	5.891	219	223	235	rVV	1073031	2124734	12.16%	2.307%
7	6.038	235	236	261	rVB	97115	311909	1.78%	0.339%
8	8.589	459	462	465	rVV	38305	67094	0.38%	0.073%
9	9.707	557	561	567	rBV	5593600	10702645	61.25%	11.621%
10	9.944	579	582	593	rVB2	20169	64050	0.37%	0.070%
11	12.540	809	812	817	rVB	24954	43692	0.25%	0.047%
12	13.195	865	870	875	rBV	7950868	15847715	90.69%	17.207%
13	13.376	883	886	901	rVB2	16677	48154	0.28%	0.052%
14	18.332	1319	1325	1331	rBV	8219639	16837646	96.36%	18.282%
15	22.283	1671	1675	1685	rVB2	47421	99358	0.57%	0.108%
16	22.633	1701	1706	1711	rBV2	8037929	17474214	100.00%	18.973%
17	27.724	2153	2157	2161	rBV	1006281	1921553	11.00%	2.086%
18	30.490	2395	2402	2407	rBV	6289217	14647696	83.82%	15.904%
19	31.088	2451	2455	2465	rVB	22828	67646	0.39%	0.073%
20	34.407	2743	2749	2791	rVV	4430949	11537370	66.03%	12.527%

Sum of corrected areas: 92099347

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02MAR08\03510.D
Operator :
Acquired : 8 Mar 2002 2:46 pm using AcqMethod HP022102
Instrument : Instrumen
Sample Name: 508 scan
Misc Info : March 8, 2002
Vial Number: 6
Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

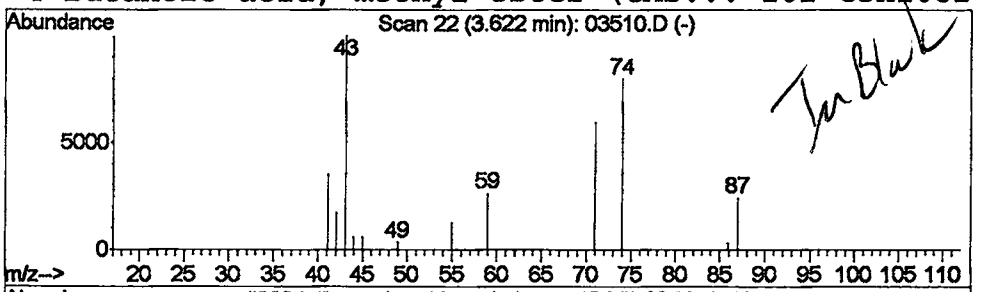
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

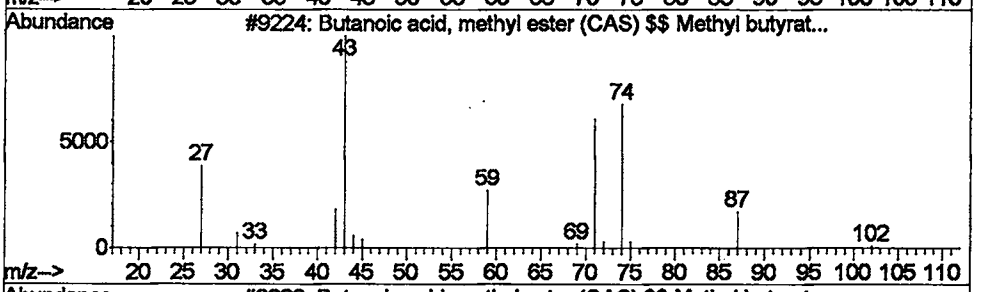
****

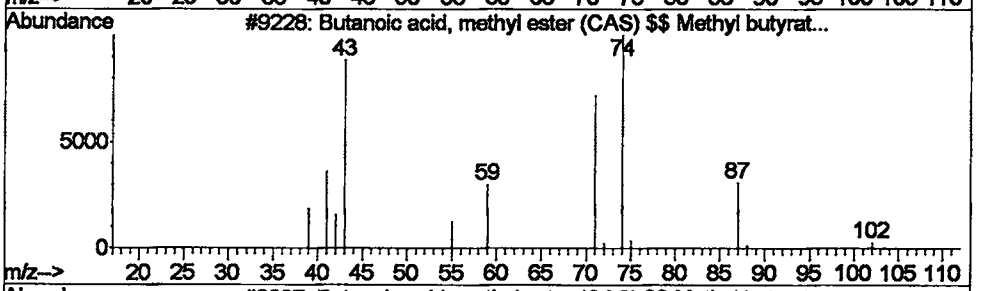
Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 3

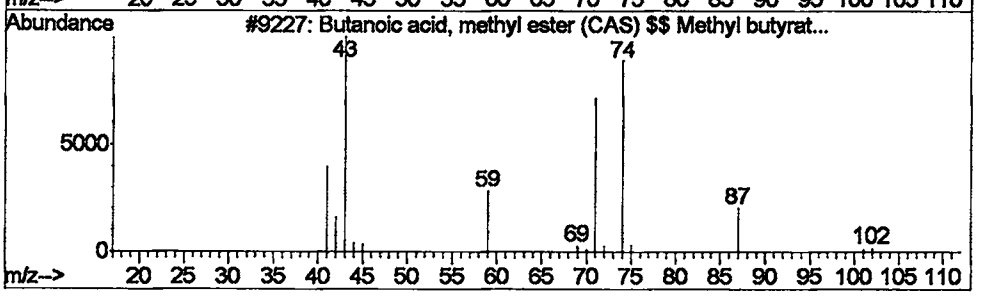
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	0.14 ug/L	75716	1,4-Dichlorobenzene-d4	9.71

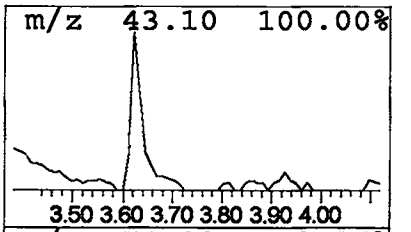
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
2	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
3	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
4	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83

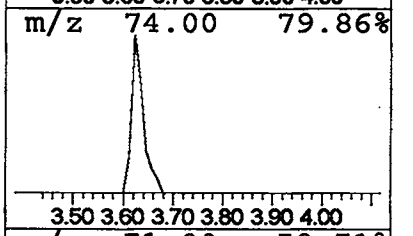


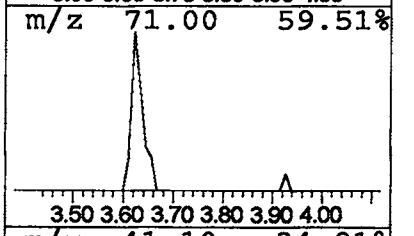


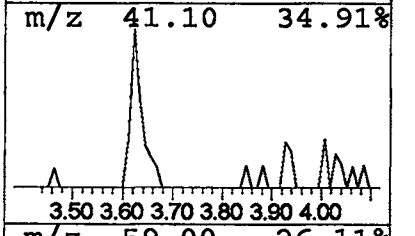


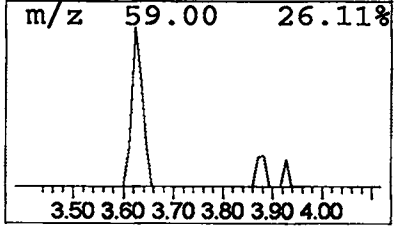












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

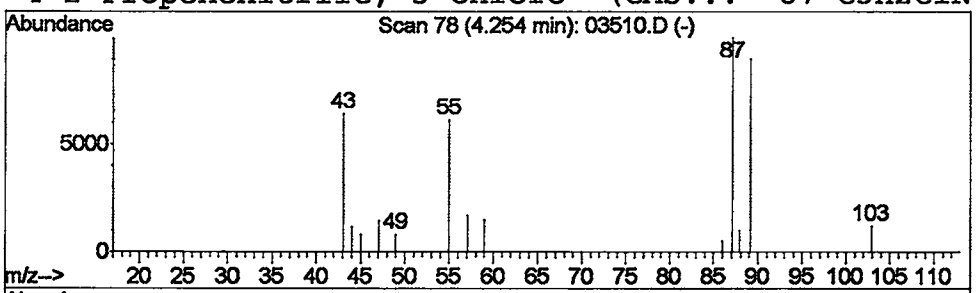
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

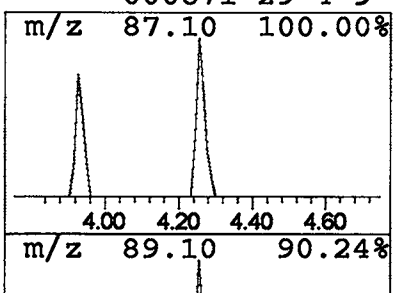
****

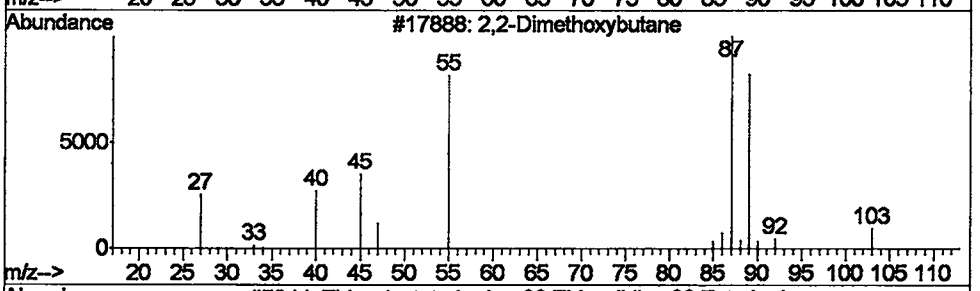
Peak Number 2 2,2-Dimethoxybutane Concentration Rank 9

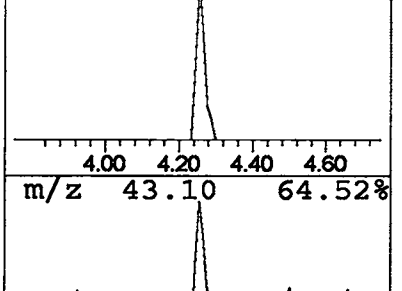
R.T.	EstConc	Area	Relative to ISTD	R.T.
4.25	0.10 ug/L	55631	1,4-Dichlorobenzene-d4	9.71

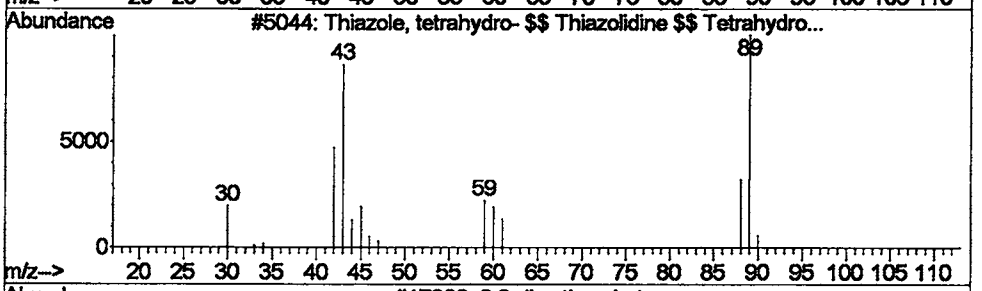
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethoxybutane	118	C6H14O2	003453-99-4	64
2	Thiazole, tetrahydro- $$ Thiazol...	89	C3H7NS	000504-78-9	36
3	2,2-dimethoxybutane	118	C6H14O2	003453-99-4	10
4	2-Propenenitrile, 3-chloro- (CAS...	87	C3H2ClN	000871-29-4	9

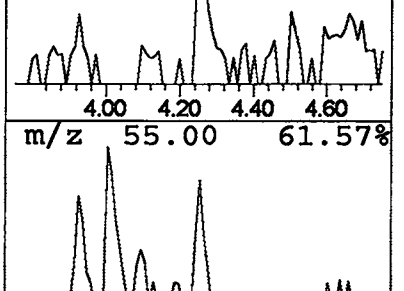


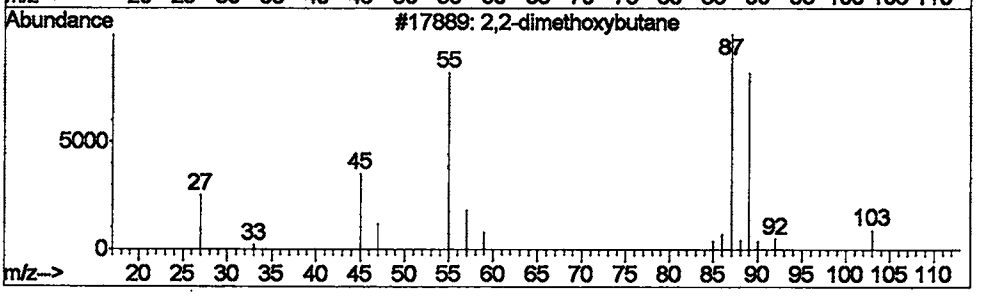


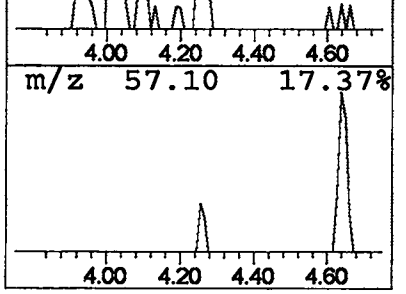












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Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

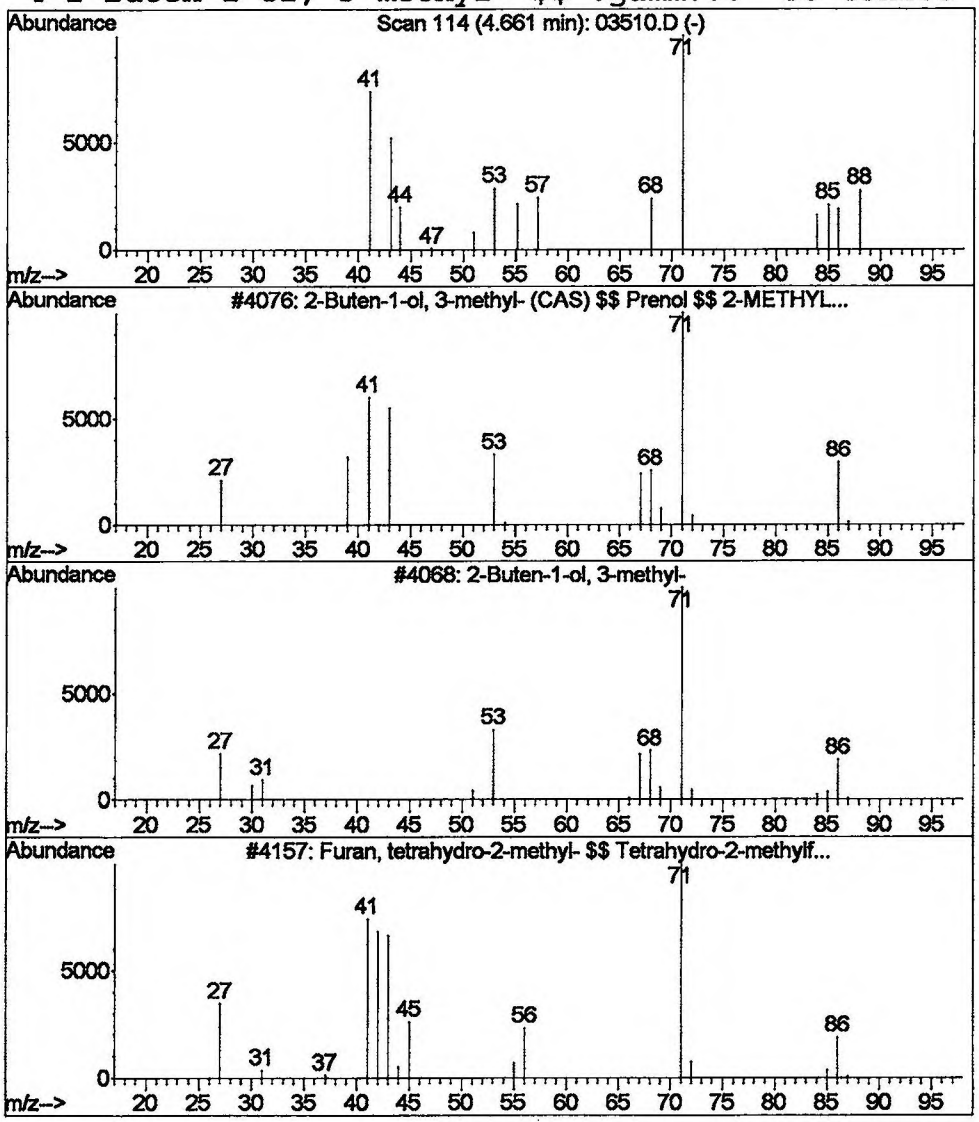
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

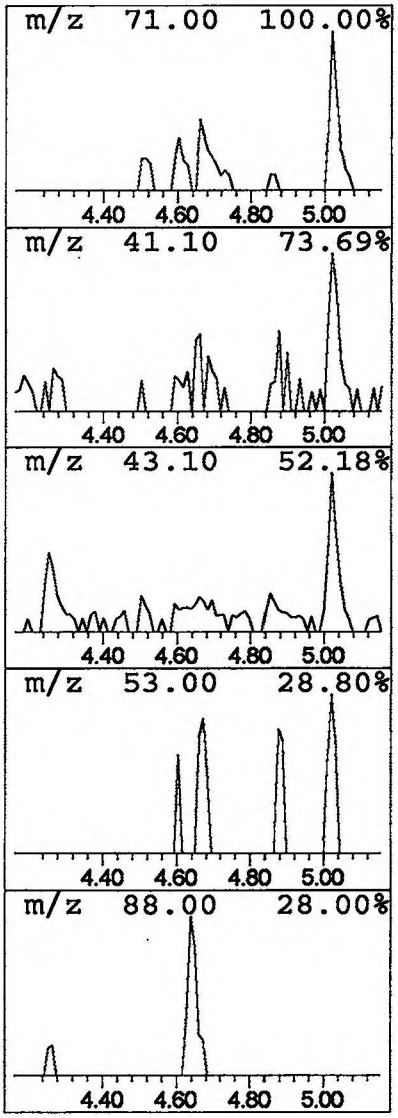
****

Peak Number 3 2-Buten-1-ol, 3-methyl- (CA... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	0.11 ug/L	57111	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Buten-1-ol, 3-methyl- (CAS) $$...	86	C5H10O	000556-82-1	52
2	2-Buten-1-ol, 3-methyl-	86	C5H10O	000556-82-1	47
3	Furan, tetrahydro-2-methyl- $$ T...	86	C5H10O	000096-47-9	37
4	2-Buten-1-ol, 3-methyl- $$ .gamm...	86	C5H10O	000556-82-1	9





Library Search Compound Report

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Acq On : 8 Mar 2002 2:46 pm Operator:
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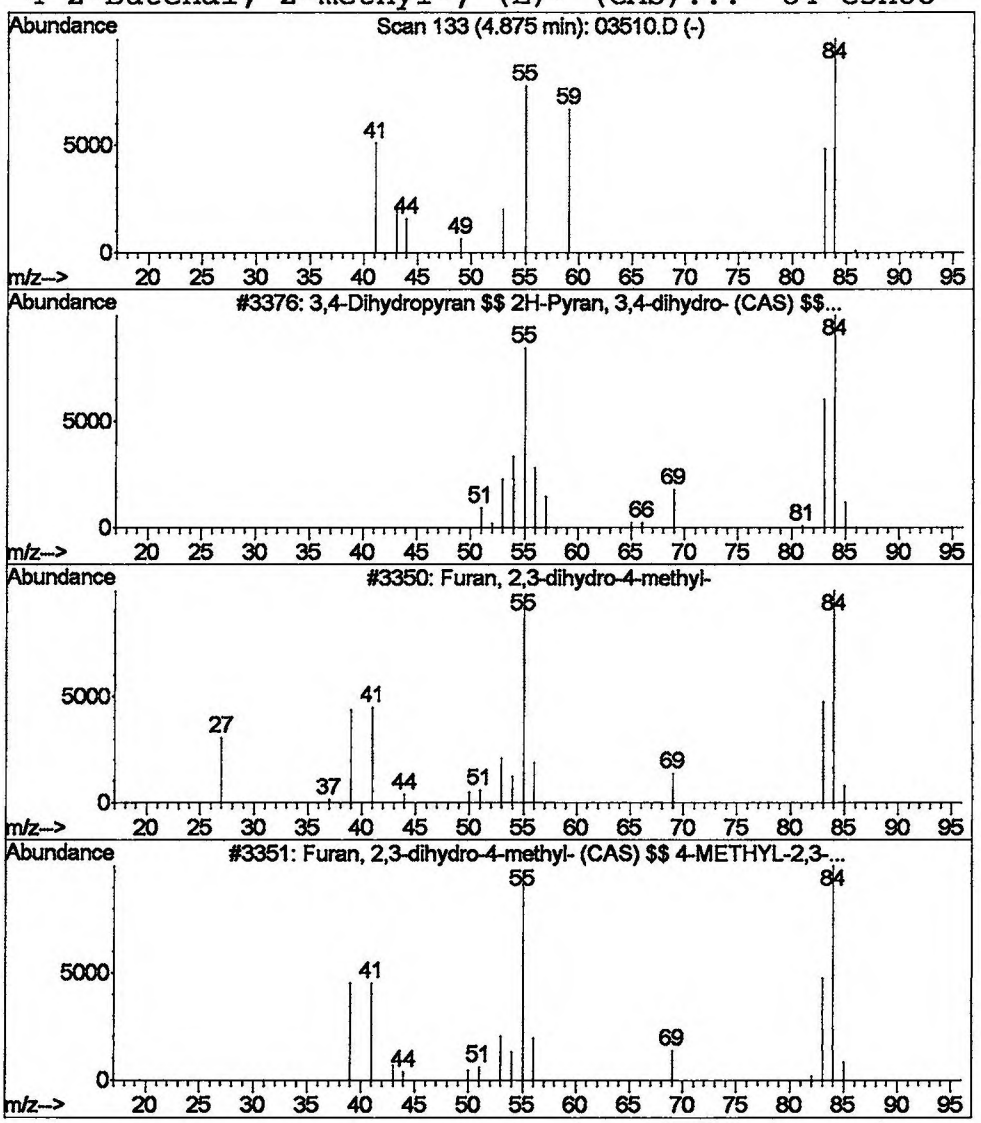
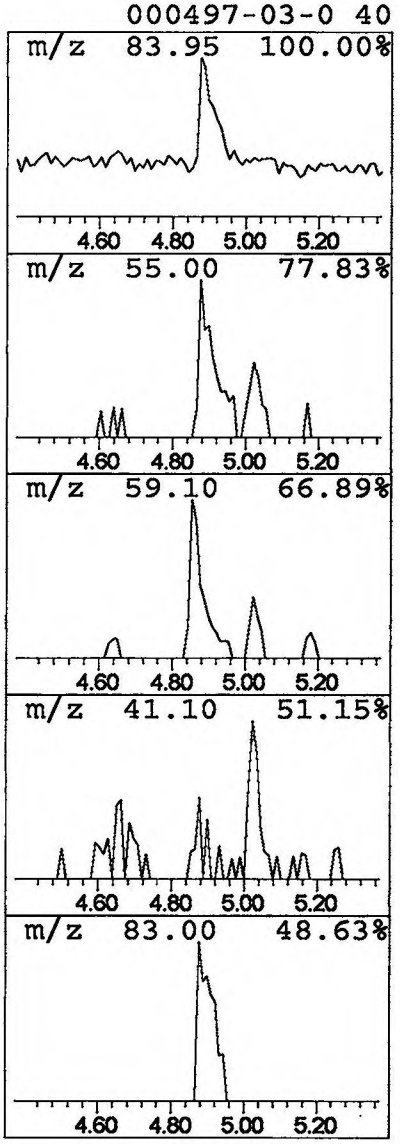
****

Peak Number 4 3,4-Dihydropyran $$ 2H-Pyra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	0.11 ug/L	59784	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dihydropyran $$ 2H-Pyran, 3,...	84	C5H8O	000110-87-2	64
2	Furan, 2,3-dihydro-4-methyl-	84	C5H8O	034314-83-5	59
3	Furan, 2,3-dihydro-4-methyl- (CA...	84	C5H8O	034314-83-5	59
4	2-Butenal, 2-methyl-, (E)- (CAS)...	84	C5H8O	000497-03-0	40





Library Search Compound Report

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Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

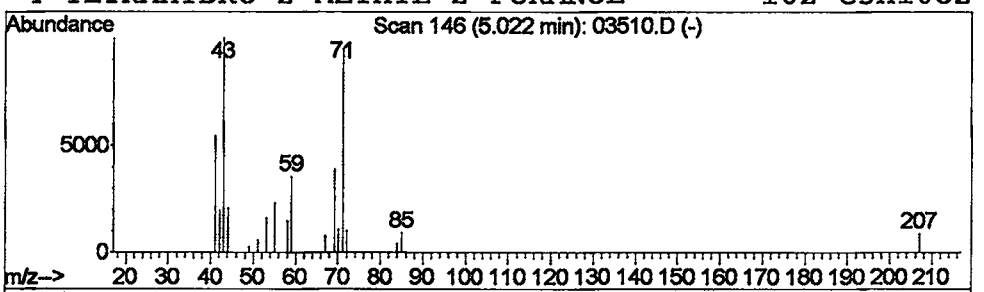
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

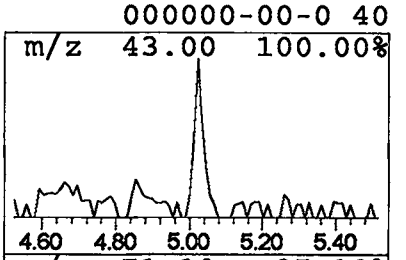
****

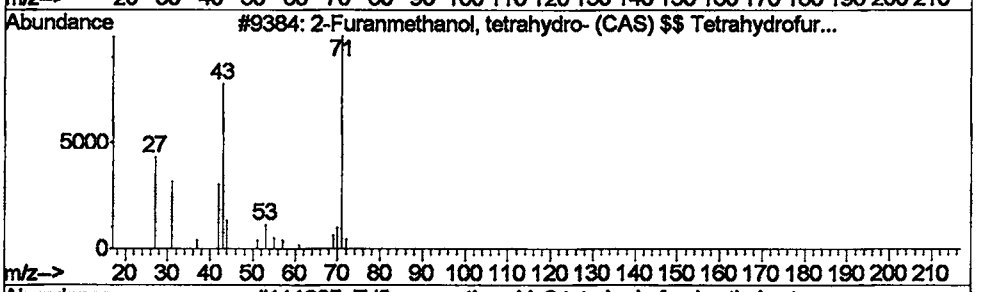
Peak Number 5 2-Furanmethanol, tetrahydro... Concentration Rank 10

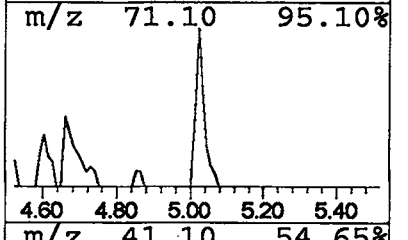
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	0.10 ug/L	55629	1,4-Dichlorobenzene-d4	9.71

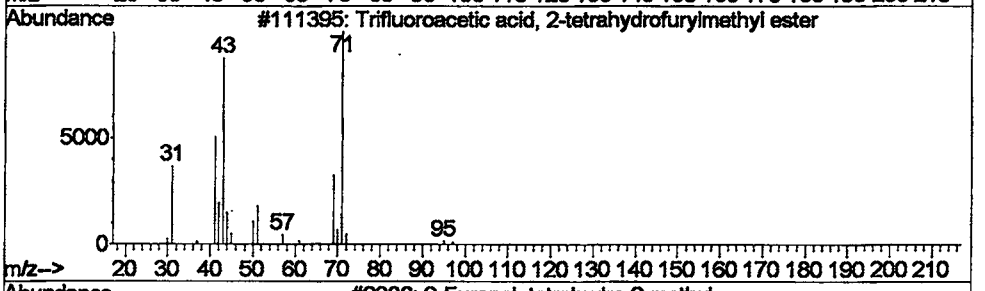
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furanmethanol, tetrahydro- (CA...	102	C5H10O2	000097-99-4	47
2	Trifluoroacetic acid, 2-tetrahyd...	198	C7H9F3O3	000000-00-0	45
3	2-Furanol, tetrahydro-2-methyl-	102	C5H10O2	007326-46-7	40
4	TETRAHYDRO-2-METHYL-2-FURANOL	102	C5H10O2	000000-00-0	40

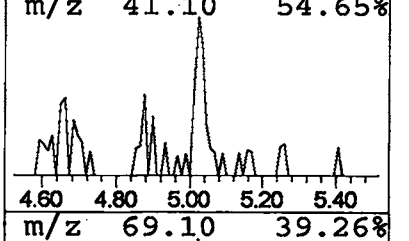


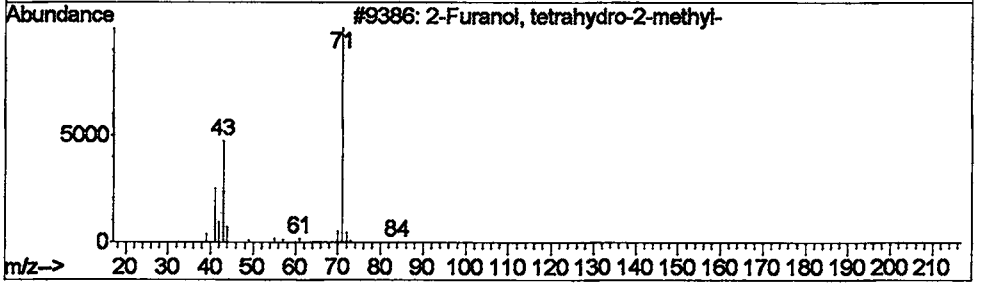


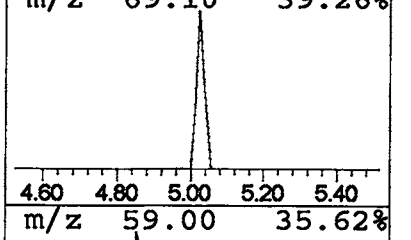


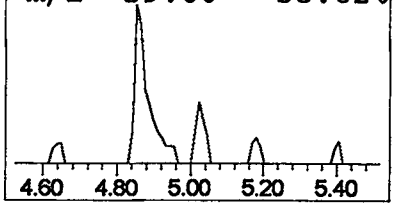












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

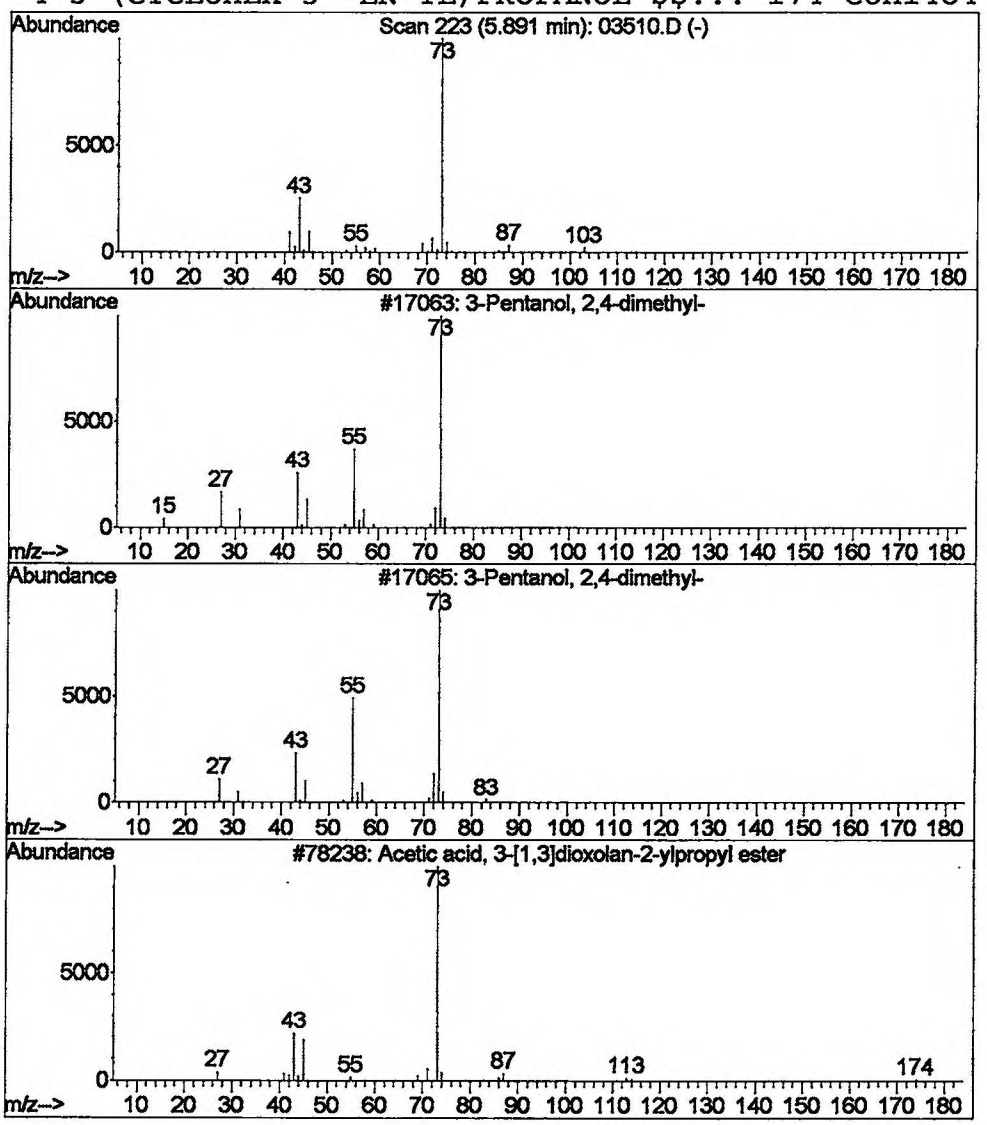
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

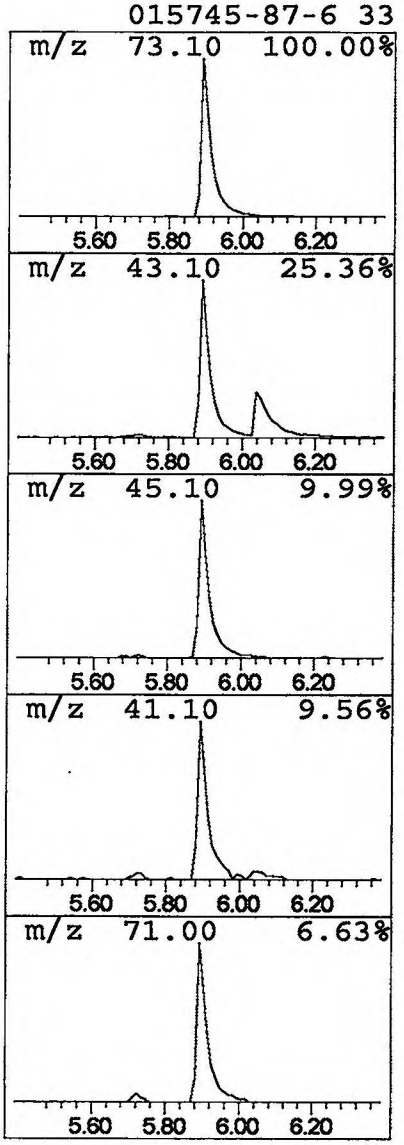
****

Peak Number 6 3-Pentanol, 2,4-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	3.97 ug/L	2124730	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	86
2	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	86
3	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33
4	3-(CYCLOHEX-3'-EN-YL) PROPANOL $$...	174	C8H14O4	015745-87-6	33





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
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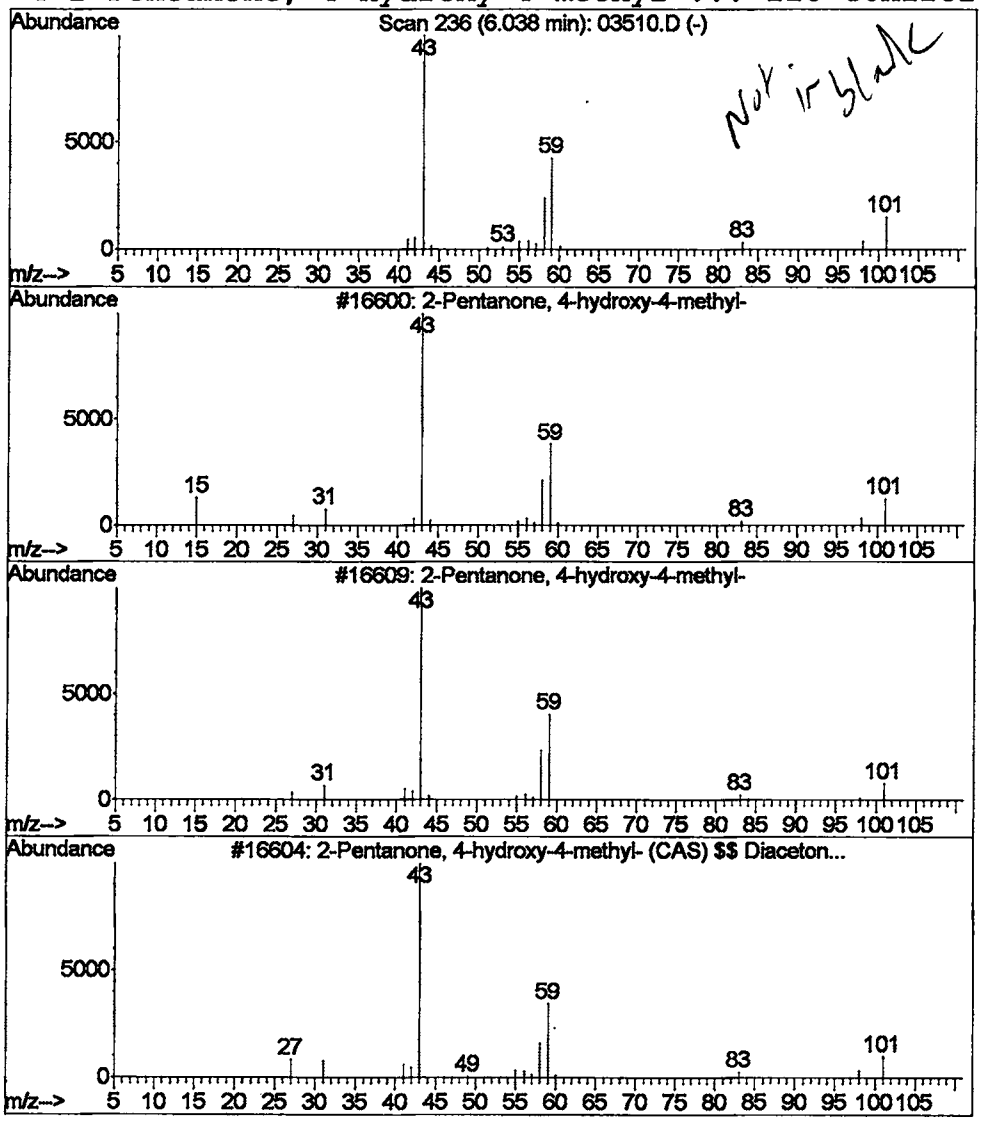
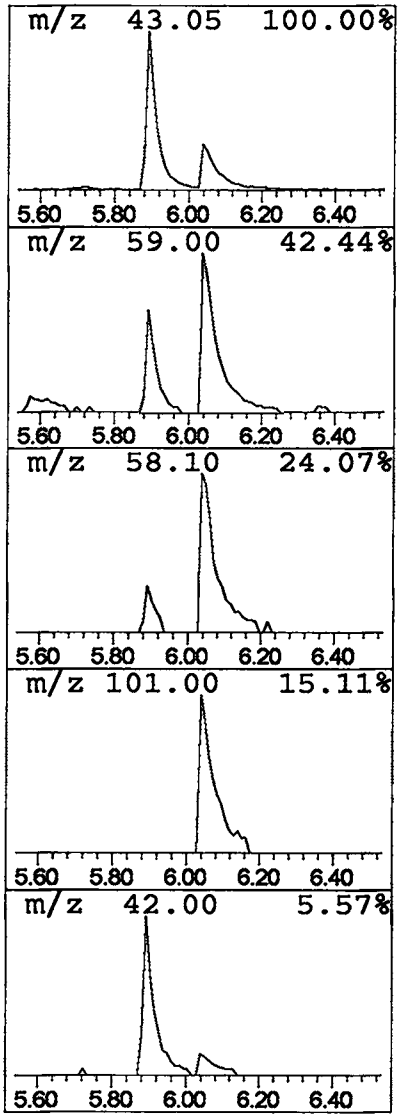
****

Peak Number 7 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.04	0.58 ug/L	311909	1,4-Dichlorobenzene-d4	9.71

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	83
3	2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	83
4	2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	83





Library Search Compound Report

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Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

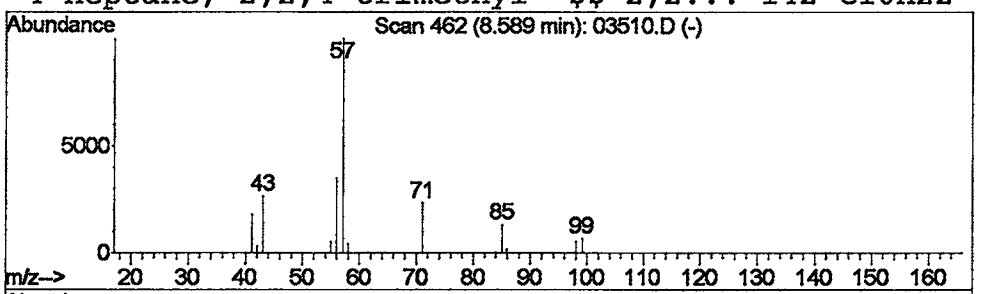
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

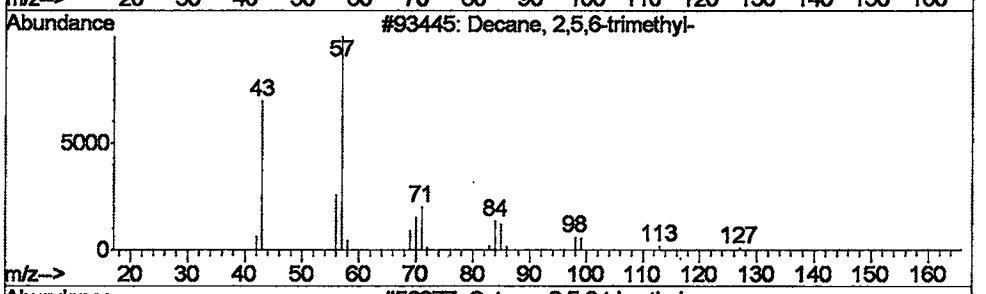
****

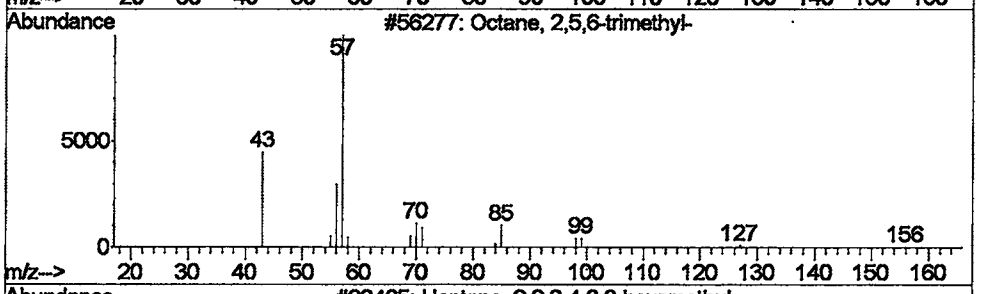
Peak Number 8 Decane, 2,5,6-trimethyl- Concentration Rank 4

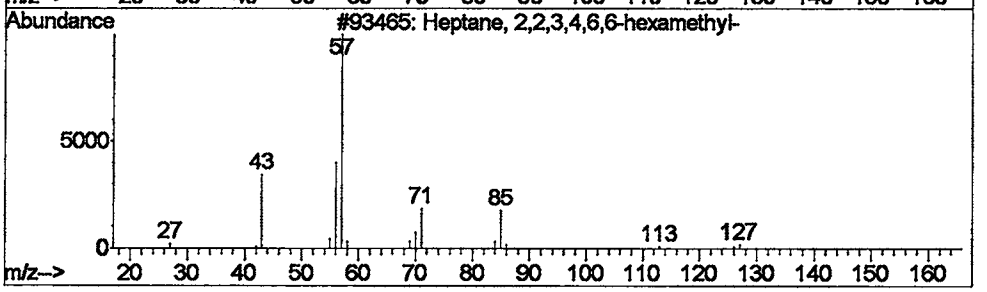
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.59	0.13 ug/L	67094	1,4-Dichlorobenzene-d4	9.71

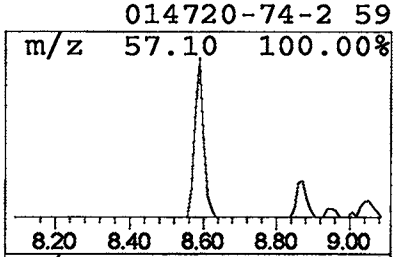
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	74
2	Octane, 2,5,6-trimethyl-	156	C11H24	062016-14-2	72
3	Heptane, 2,2,3,4,6,6-hexamethyl-	184	C13H28	062108-32-1	64
4	Heptane, 2,2,4-trimethyl- $$ 2,2...	142	C10H22	014720-74-2	59

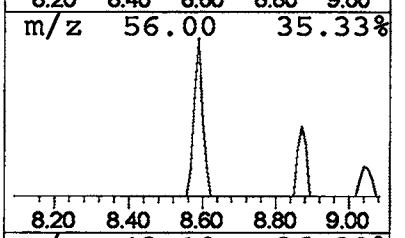


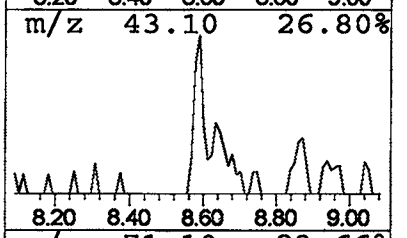


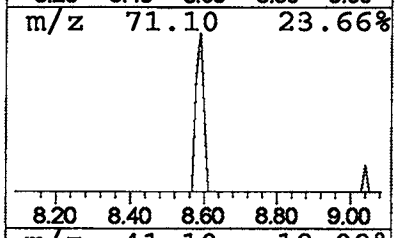


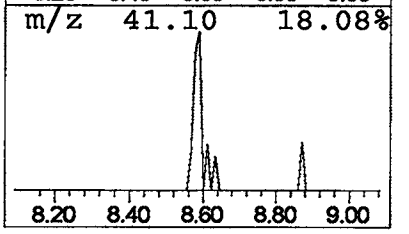












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

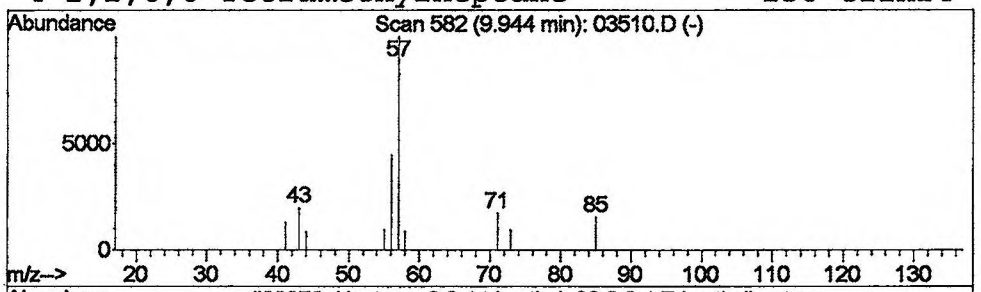
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

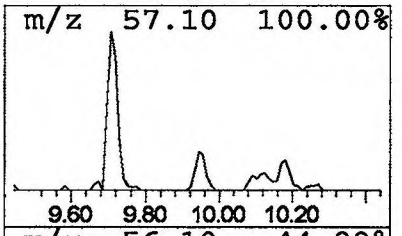
****

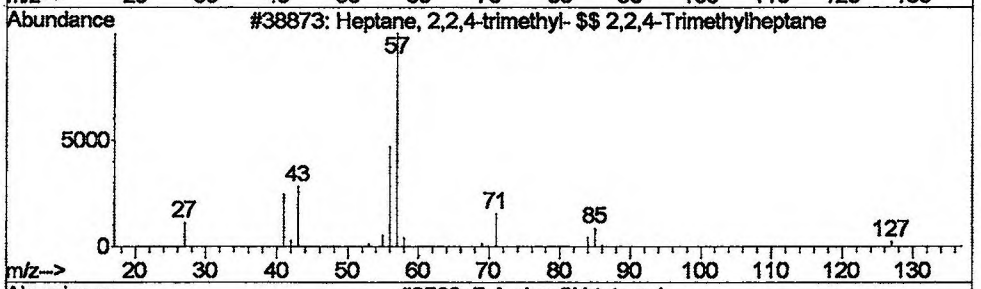
Peak Number 9 Heptane, 2,2,4-trimethyl- $... Concentration Rank 5

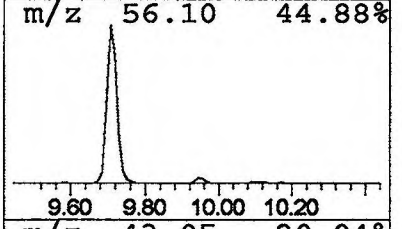
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.94	0.12 ug/L	64050	1,4-Dichlorobenzene-d4	9.71

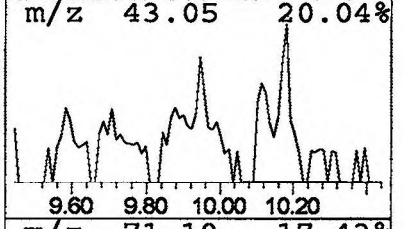
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,2,4-trimethyl- $$ 2,2...	142	C10H22	014720-74-2	43
2	5-Amino-2H-tetrazole	85	CH3N5	000000-00-0	9
3	5-Amino-2H-tetrazole	85	CH3N5	000000-00-0	9
4	2,2,6,6-Tetramethylheptane	156	C11H24	040117-45-1	9

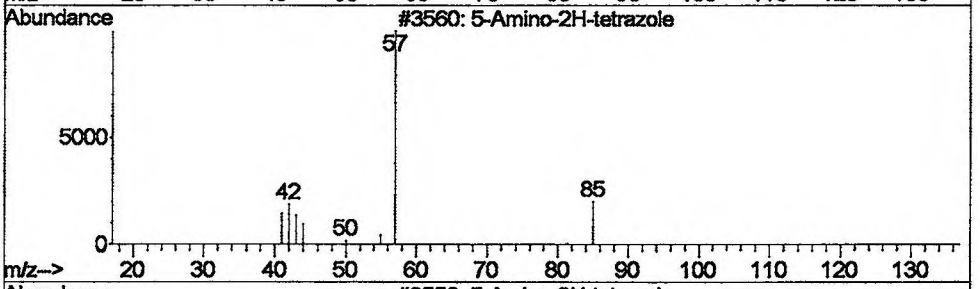


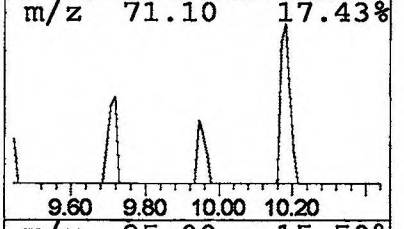


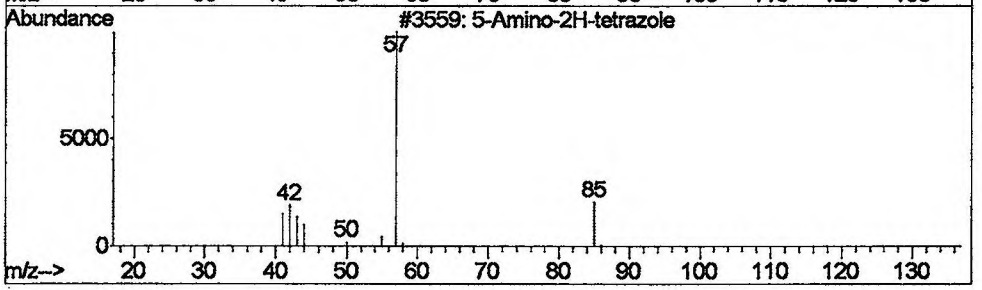


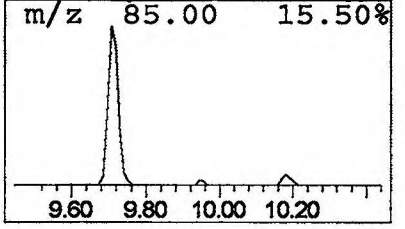












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03510.D Vial: 6
Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

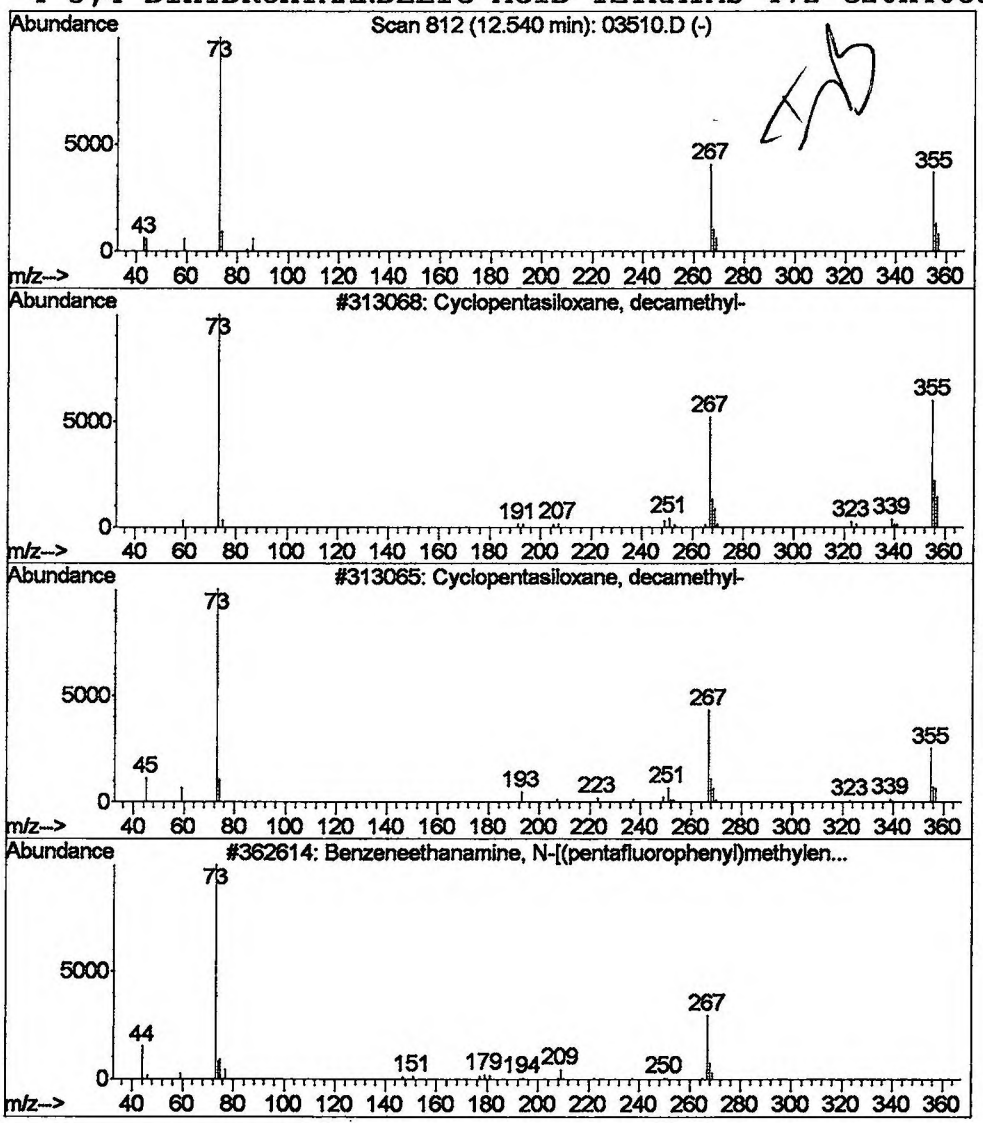
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

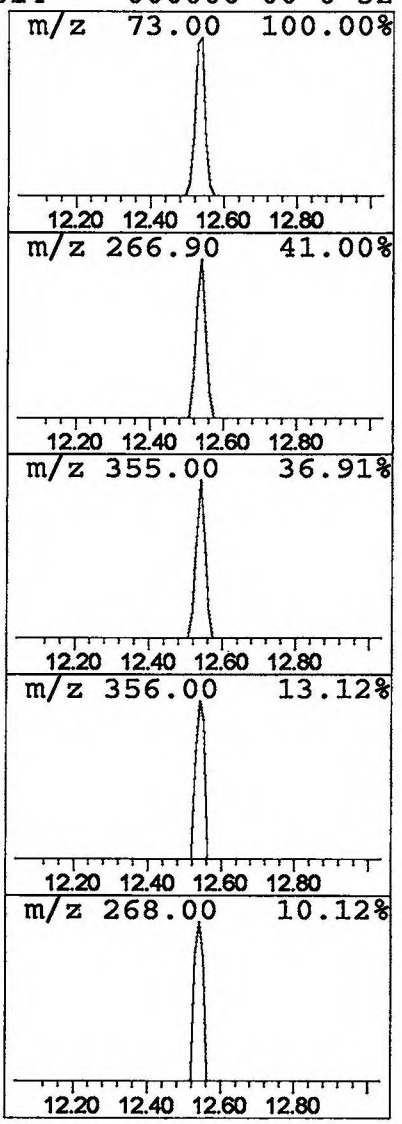
****

Peak Number 10 Cyclopentasiloxane, decamet... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.06 ug/L	43692	Naphthalene-d8	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	83
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	83
3	Benzeneethanamine, N-[(pentafluo...	475	C21H26F5NO2Si2	055429-85-1	33
4	3,4-DIHYDROXYMANDELIC ACID-TETRATMS	472	C20H40O5Si4	000000-00-0	32





Library Search Compound Report

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Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
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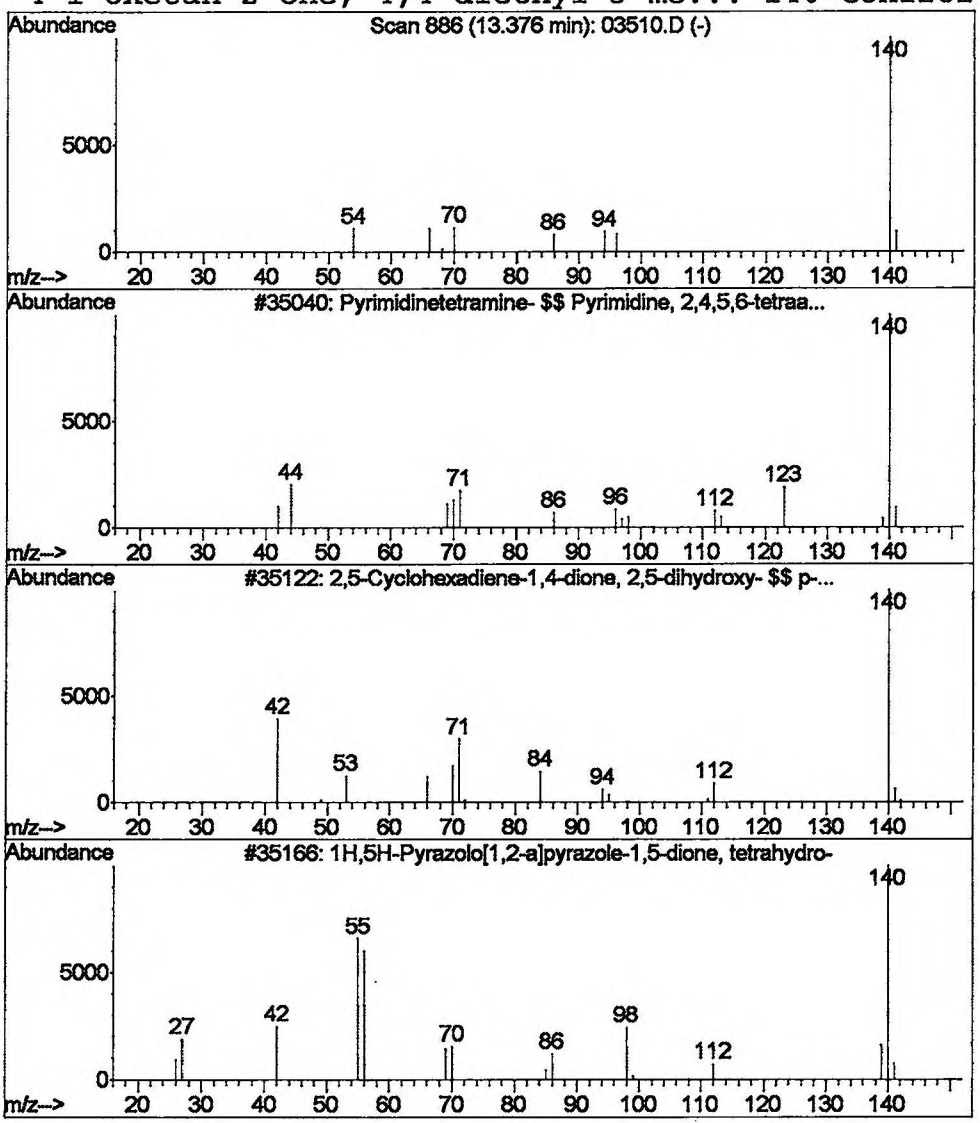
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

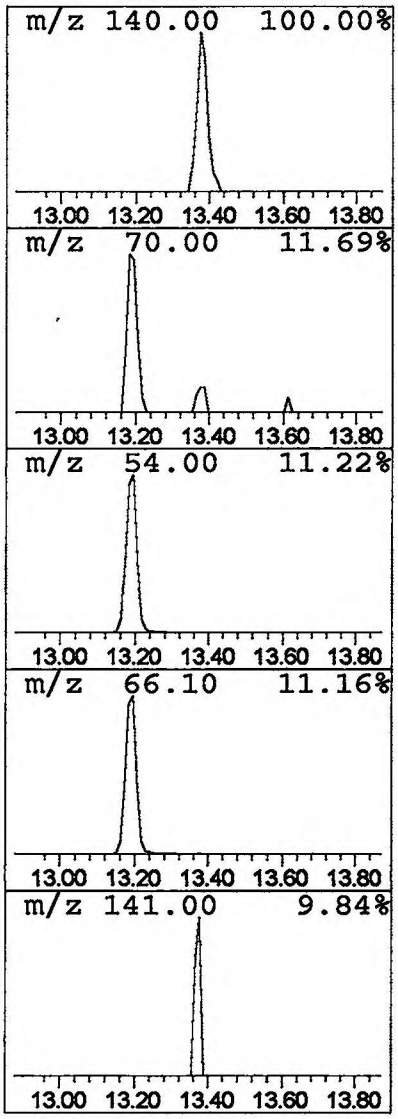
****

Peak Number 11 Pyrimidinetetramine- $$ Pyr... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	0.06 ug/L	48154	Naphthalene-d8	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrimidinetetramine- $$ Pyrimidi...	140	C4H8N6	001004-74-6	64
2	2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	56
3	1H,5H-Pyrazolo[1,2-a]pyrazole-1,...	140	C6H8N2O2	019720-72-0	50
4	1-Oxetan-2-one, 4,4-diethyl-3-me...	140	C8H12O2	000000-00-0	50





Library Search Compound Report

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Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
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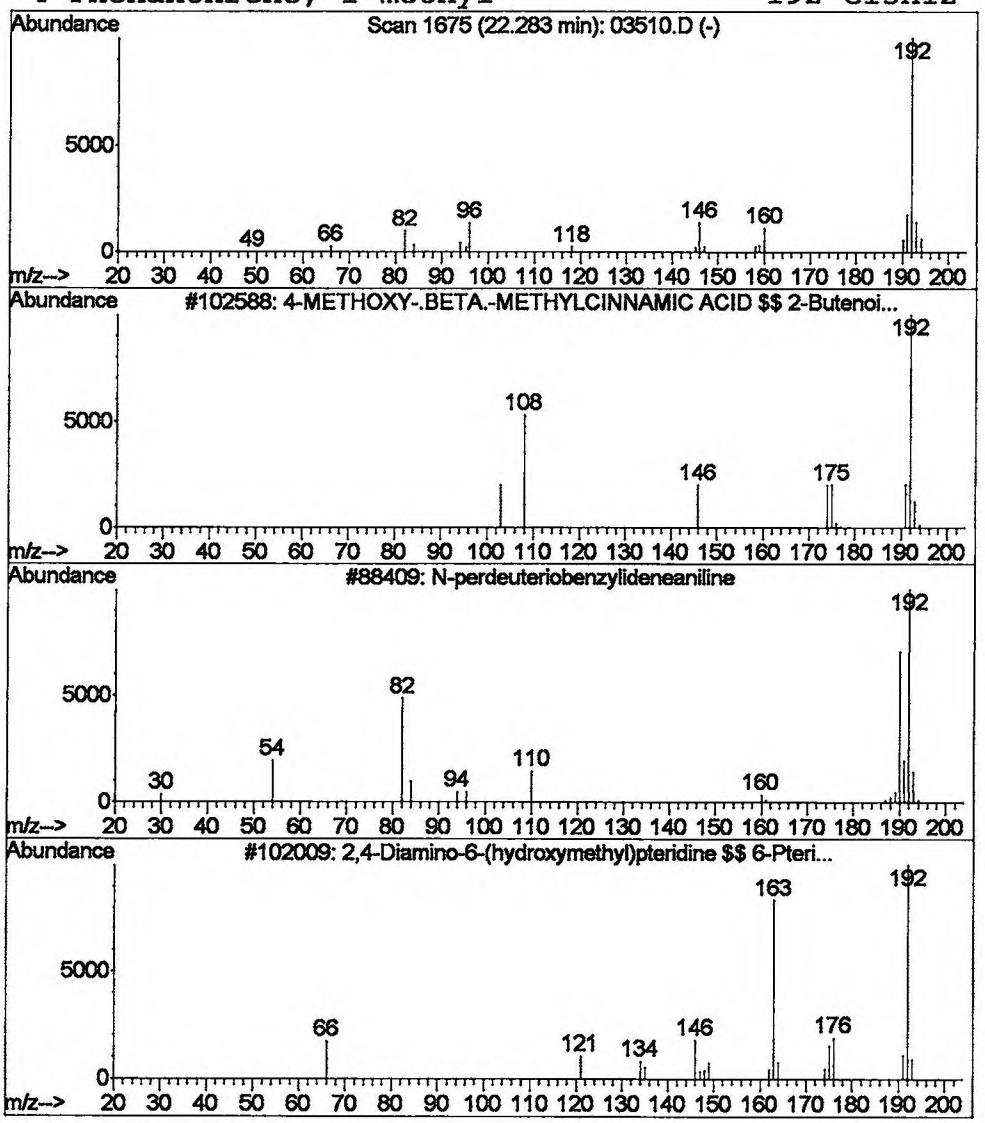
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

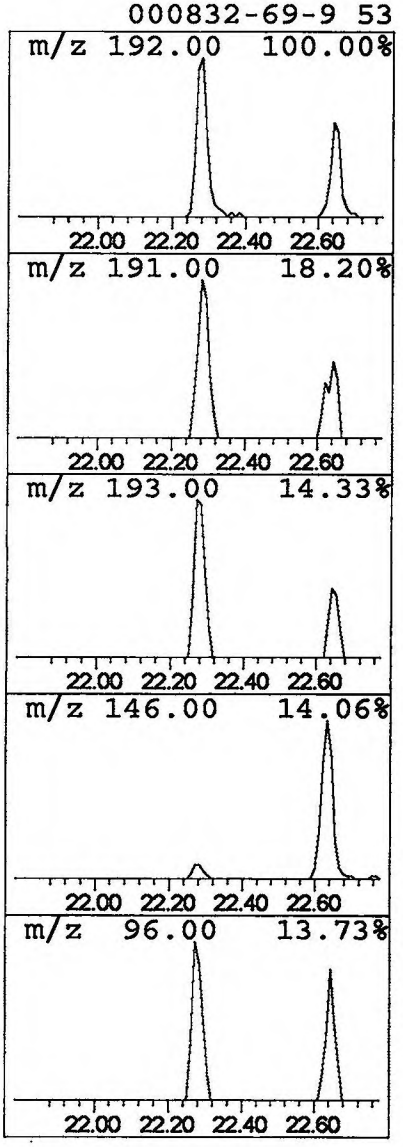
****

Peak Number 12 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.11 ug/L	99358	Phenanthrene-d10	22.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	72
2	N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	59
3	2,4-Diamino-6-(hydroxymethyl)pte...	192	C7H8N6O	000945-24-4	53
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	53





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Acq On : 8 Mar 2002 2:46 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
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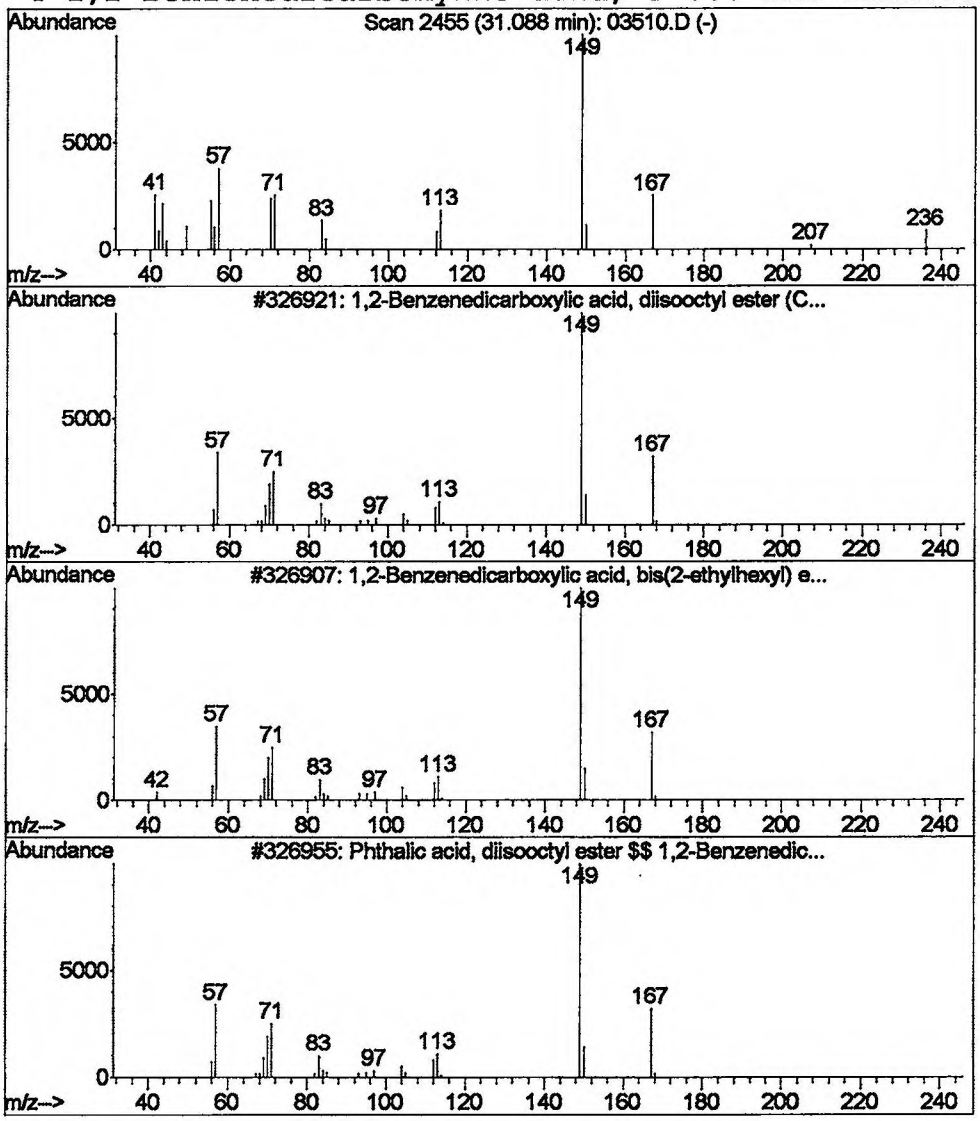
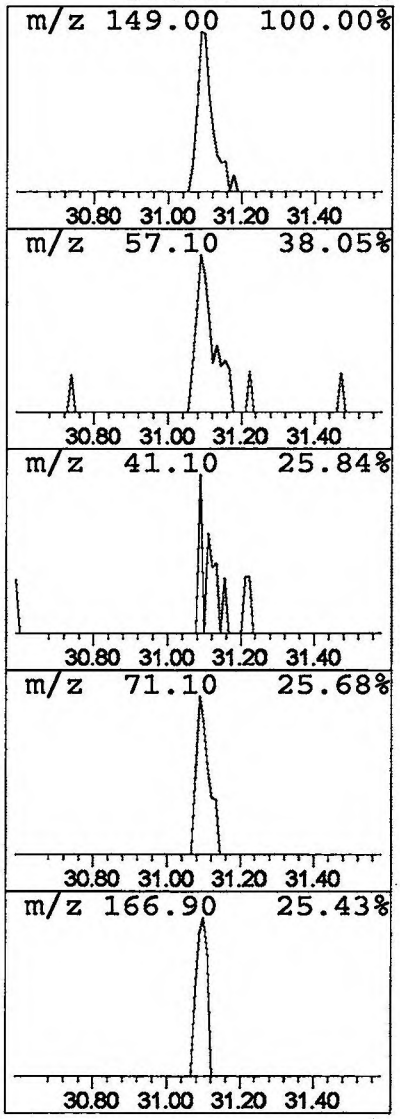
****

Peak Number 13 1,2-Benzenedicarboxylic aci... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.09	0.09 ug/L	67646	Chrysene-d12	30.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, di...	390	C24H38O4	027554-26-3	72
2	1,2-Benzenedicarboxylic acid, bi...	390	C24H38O4	000117-81-7	72
3	Phthalic acid, diisooctyl ester ...	390	C24H38O4	001330-91-2	72
4	1,2-Benzenedicarboxylic acid, 3-...	211	C8H5NO6	000603-11-2	64





Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 8 Mar 2002 2:46 pm
Data File: C:\MSDCHEM\1\5973N\02MAR08\03510.D
Name: 508 scan
Misc: March 8, 2002
Method: C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title: 508 scan method
Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--- #	RT	Resp	Conc
Butanoic acid, me...	3.62	0.1	ug/L	75716	1	9.71	10702600	20.0
2,2-Dimethoxybutane	4.25	0.1	ug/L	55631	1	9.71	10702600	20.0
2-Buten-1-ol, 3-m...	4.66	0.1	ug/L	57111	1	9.71	10702600	20.0
3,4-Dihydropyran ...	4.88	0.1	ug/L	59784	1	9.71	10702600	20.0
2-Furanmethanol, ...	5.02	0.1	ug/L	55629	1	9.71	10702600	20.0
3-Pentanol, 2,4-d...	5.89	4.0	ug/L	2124730	1	9.71	10702600	20.0
2-Pentanone, 4-hy...	6.04	0.6	ug/L	311909	1	9.71	10702600	20.0
Decane, 2,5,6-tri...	8.59	0.1	ug/L	67094	1	9.71	10702600	20.0
Heptane, 2,2,4-tr...	9.94	0.1	ug/L	64050	1	9.71	10702600	20.0
Cyclopentasiloxan...	12.54	0.1	ug/L	43692	2	13.20	15847700	20.0
Pyrimidinetetrami...	13.38	0.1	ug/L	48154	2	13.20	15847700	20.0
4-METHOXY-.BETA.-...	22.28	0.1	ug/L	99358	4	22.63	17474200	20.0
1,2-Benzenedicarb...	31.09	0.1	ug/L	67646	5	30.49	14647700	20.0