User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/05/2002 Time: 11:43:04

Sample: AE02713
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/5/02 11:42 Ended: 3/5/02 11:42 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		80		5.0

Comments for Sample AE02713 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 03-05-2002 Time: 11:44:23

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 14 of the 43 compounds in the LCS Standard were above their acceptance ranges.

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02713.D Vial: 3
Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

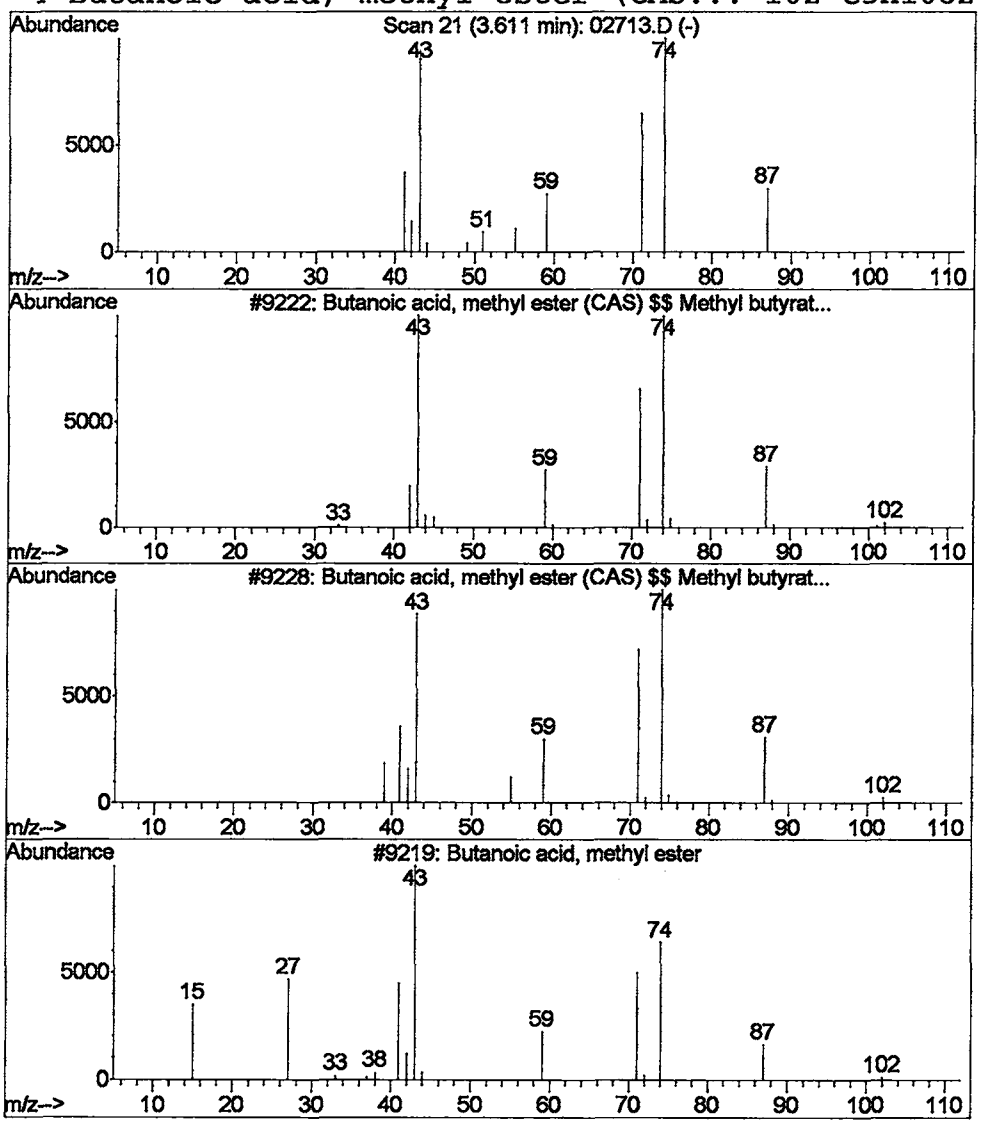
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

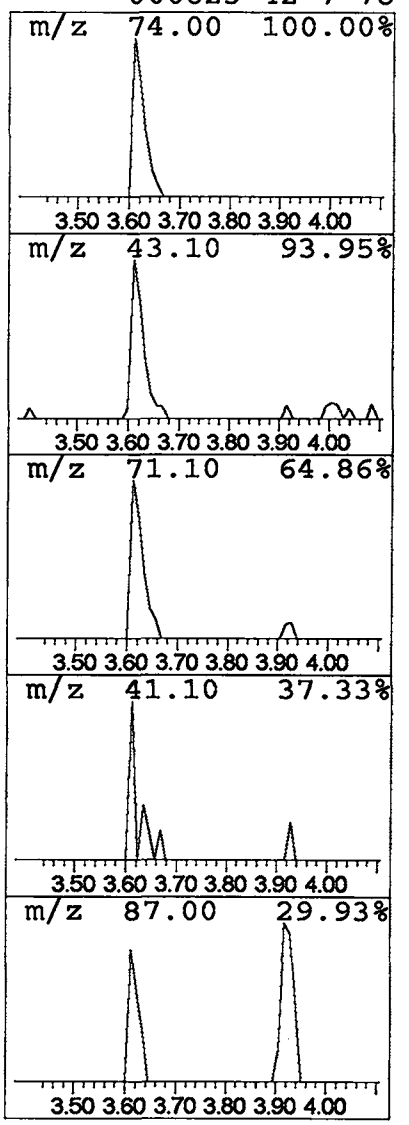
****

Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.16 ug/L	58932	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	78
2	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	78
3	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	78
4	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	78





Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02713.D Vial: 3
Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Mar 01 09:02:48 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 : HP020702

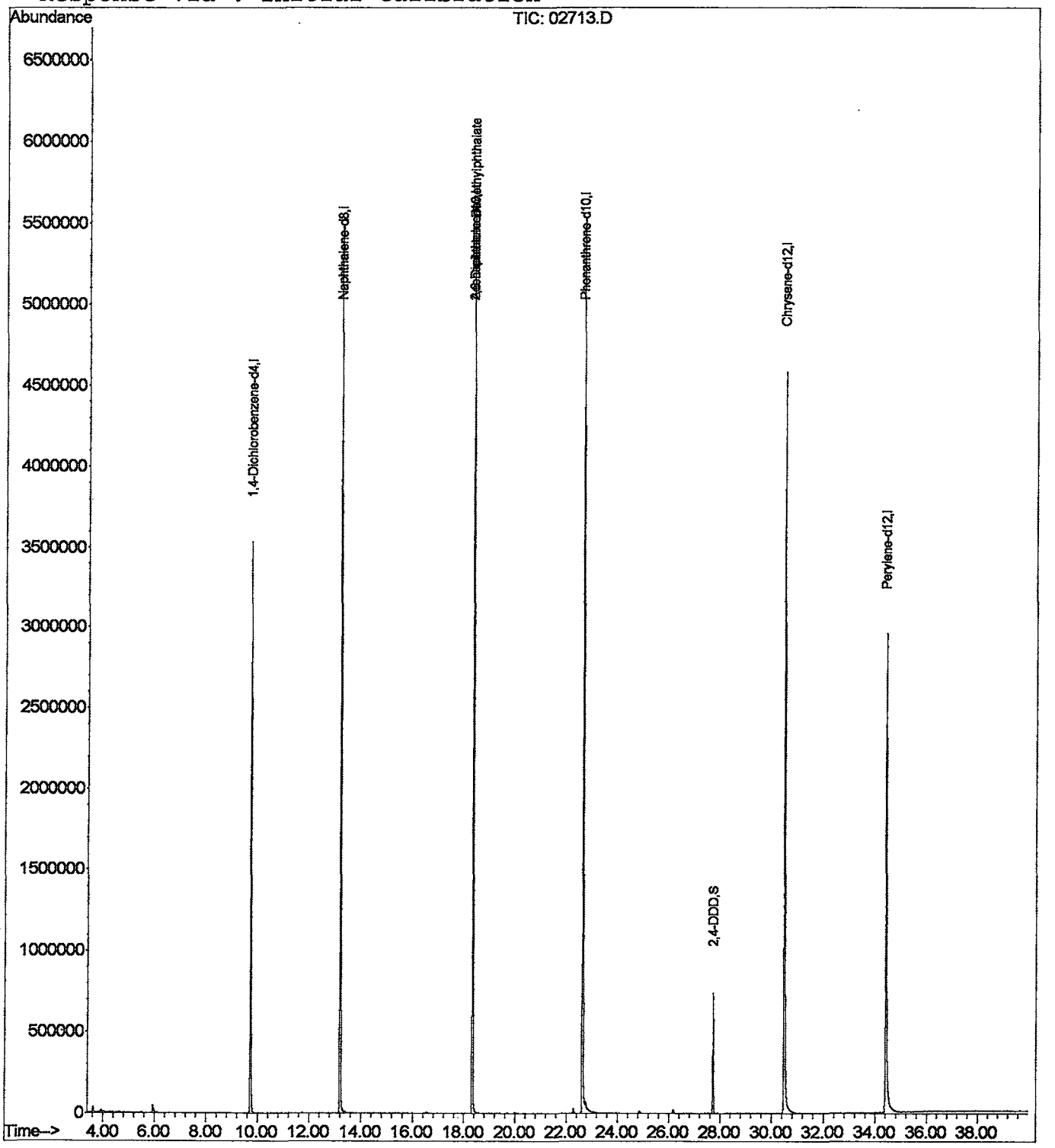
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.72	152	1322826	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	5634841	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	2832598	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	4886132	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	3721363	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2408190	20.00	ug/L	-0.04
System Monitoring Compounds						
37) 2,4-DDD	27.72	235	260749	4.00	ug/L	0.00
Spiked Amount 5.000			Recovery = 80.00%			
Target Compounds						Qvalue
20) Dimethylphthalate	18.34	163	468468	2.73	ug/L	# 57
21) 2,6-Dinitrotoluene	18.33	165	360762	9.09	ug/L	# 27

(#) = qualifier out of range (m) = manual integration
02713.D HP022102.M Fri Mar 01 09:02:49 2002 Page 1

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02713.D Vial: 3
Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Mar 1 9:02 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\02FEB28\02713.D Vial: 3
Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 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

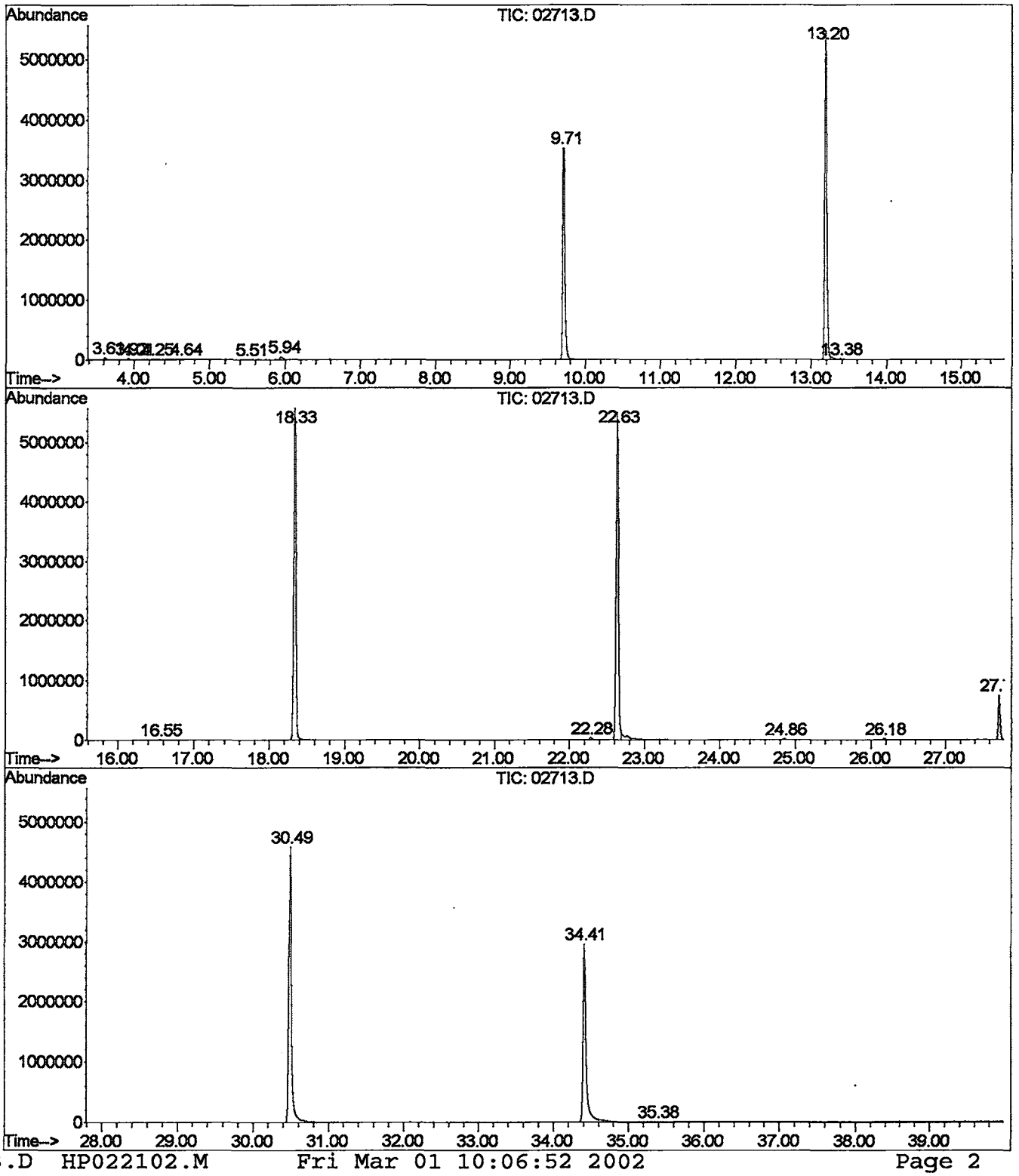
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.611	19	21	29	rBV	34102	58932	0.49%	0.094%
2	3.916	45	48	53	rVB	15104	26319	0.22%	0.042%
3	4.006	53	56	67	rVB2	5200	18448	0.15%	0.029%
4	4.254	73	78	85	rVB6	5168	23619	0.20%	0.038%
5	4.638	107	112	119	rVB3	7574	18997	0.16%	0.030%
6	5.507	185	189	199	rVB3	3307	11950	0.10%	0.019%
7	5.936	225	227	237	rBV	48659	130387	1.08%	0.208%
8	9.707	557	561	575	rVV	3537868	7416377	61.55%	11.858%
9	13.195	865	870	875	rBV	5478156	10627684	88.20%	16.993%
10	13.376	883	886	901	rVB3	10815	33222	0.28%	0.053%
11	16.548	1163	1167	1175	rBV2	6536	15158	0.13%	0.024%
12	18.332	1319	1325	1331	rBV	5580907	11530866	95.69%	18.437%
13	22.283	1671	1675	1681	rBV	30662	59076	0.49%	0.094%
14	22.633	1701	1706	1711	rBV2	5506347	12050116	100.00%	19.267%
15	24.857	1897	1903	1909	rBV	13592	33312	0.28%	0.053%
16	26.177	2017	2020	2027	rBV2	22454	44003	0.37%	0.070%
17	27.724	2153	2157	2163	rVV	744460	1263874	10.49%	2.021%
18	30.490	2397	2402	2423	rBV	4582027	10911516	90.55%	17.447%
19	34.407	2743	2749	2787	rBV	2951046	8248864	68.45%	13.189%
20	35.378	2831	2835	2841	rBV4	5514	19382	0.16%	0.031%

Sum of corrected areas: 62542102

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02FEB28\02713.D
Operator :
Acquired : 28 Feb 2002 12:10 pm using AcqMethod HP020702
Instrument : Instrumen
Sample Name: 508 scan
Misc Info : February 28, 2002
Vial Number: 3
Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02713.D Vial: 3
Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

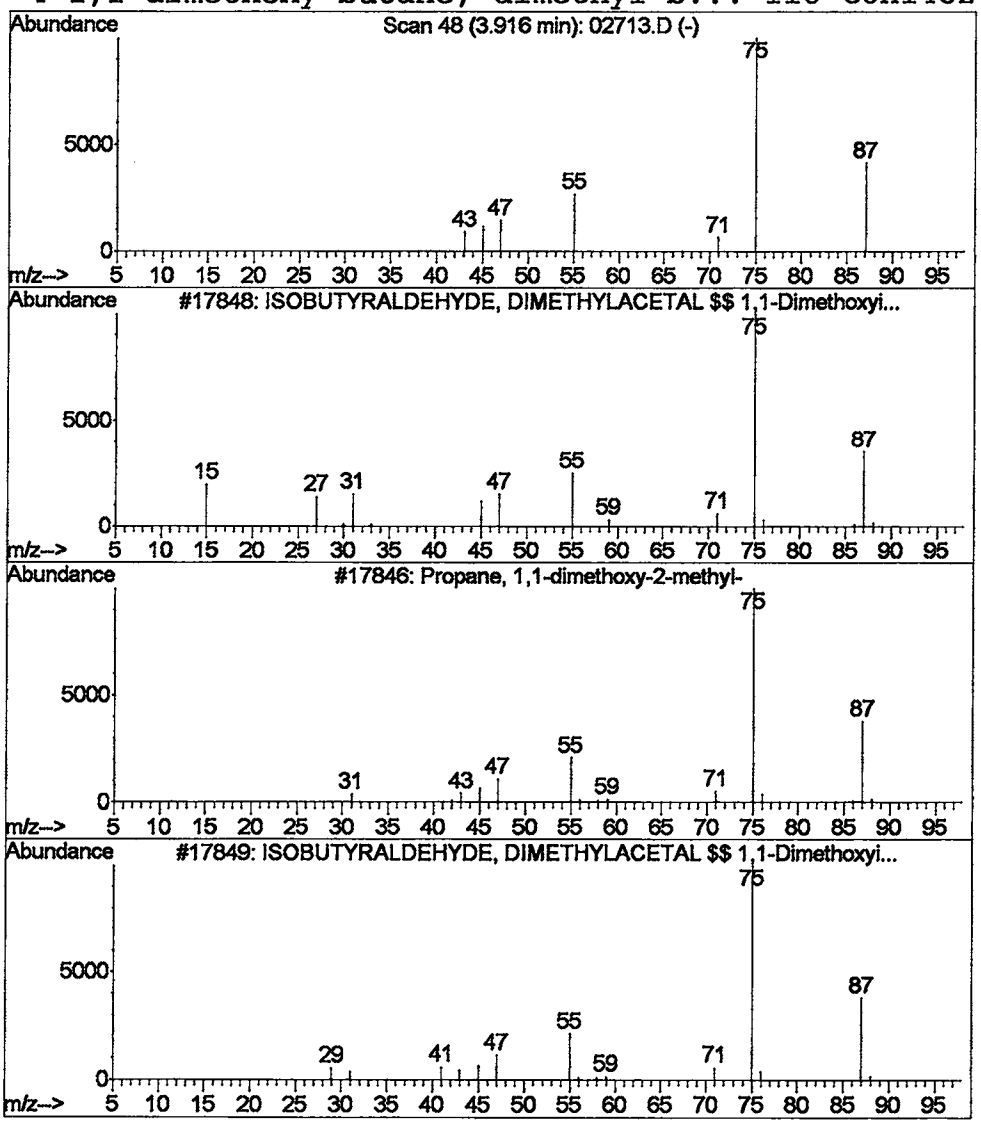
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

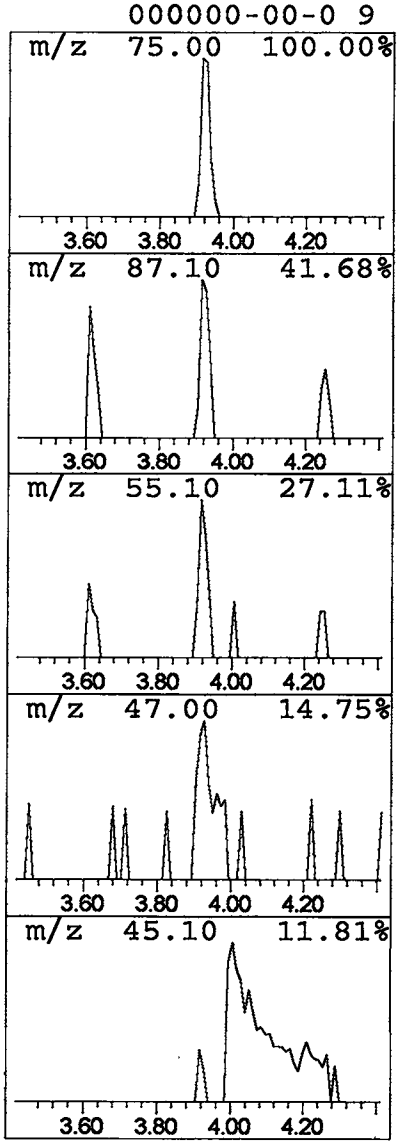
****

Peak Number 2 ISOBUTYRALDEHYDE, DIMETHYLA... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.92	0.07 ug/L	26319	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	83
2	Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	83
3	ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	83
4	1,1-dimethoxy butane, dimethyl b...	118	C6H14O2	000000-00-0	9





Library Search Compound Report

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Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
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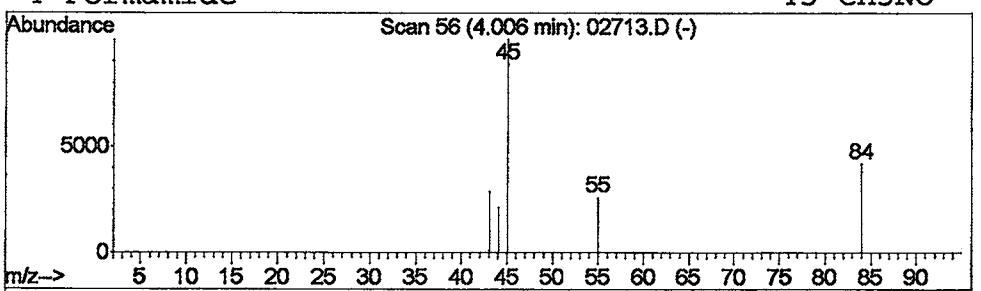
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

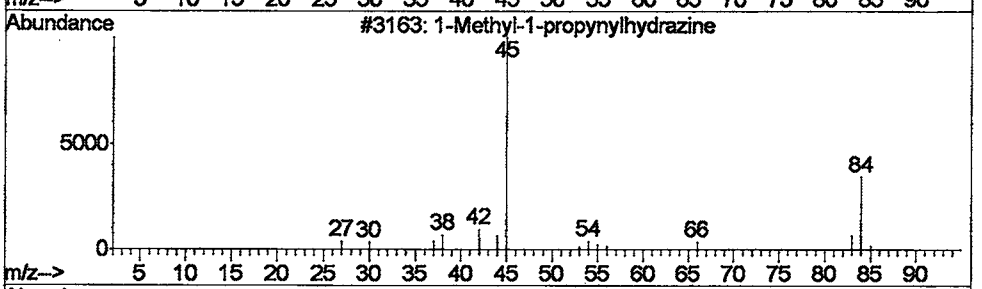
****

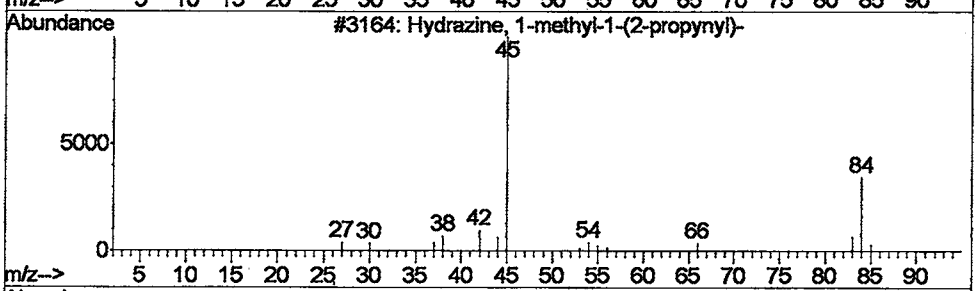
Peak Number 3 1-Methyl-1-propynylhydrazine Concentration Rank 10

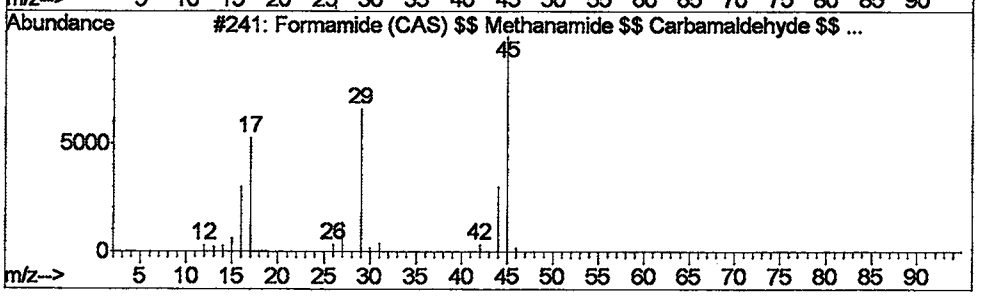
R.T.	EstConc	Area	Relative to ISTD	R.T.
4.01	0.05 ug/L	18448	1,4-Dichlorobenzene-d4	9.72

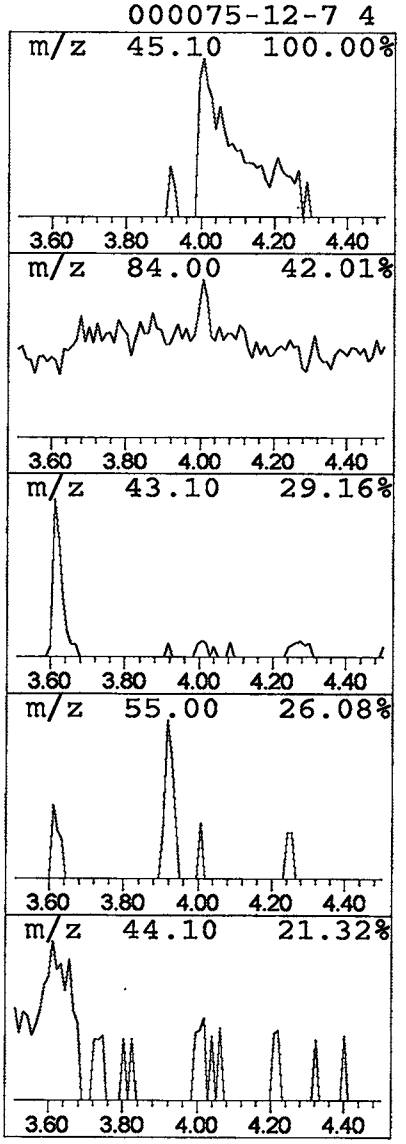
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-1-propynylhydrazine	84	C4H8N2	000000-00-0	4
2	Hydrazine, 1-methyl-1-(2-propynyl)-	84	C4H8N2	007422-82-4	4
3	Formamide (CAS) $$ Methanamide $...	45	CH3NO	000075-12-7	4
4	Formamide	45	CH3NO	000075-12-7	4











Library Search Compound Report

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Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

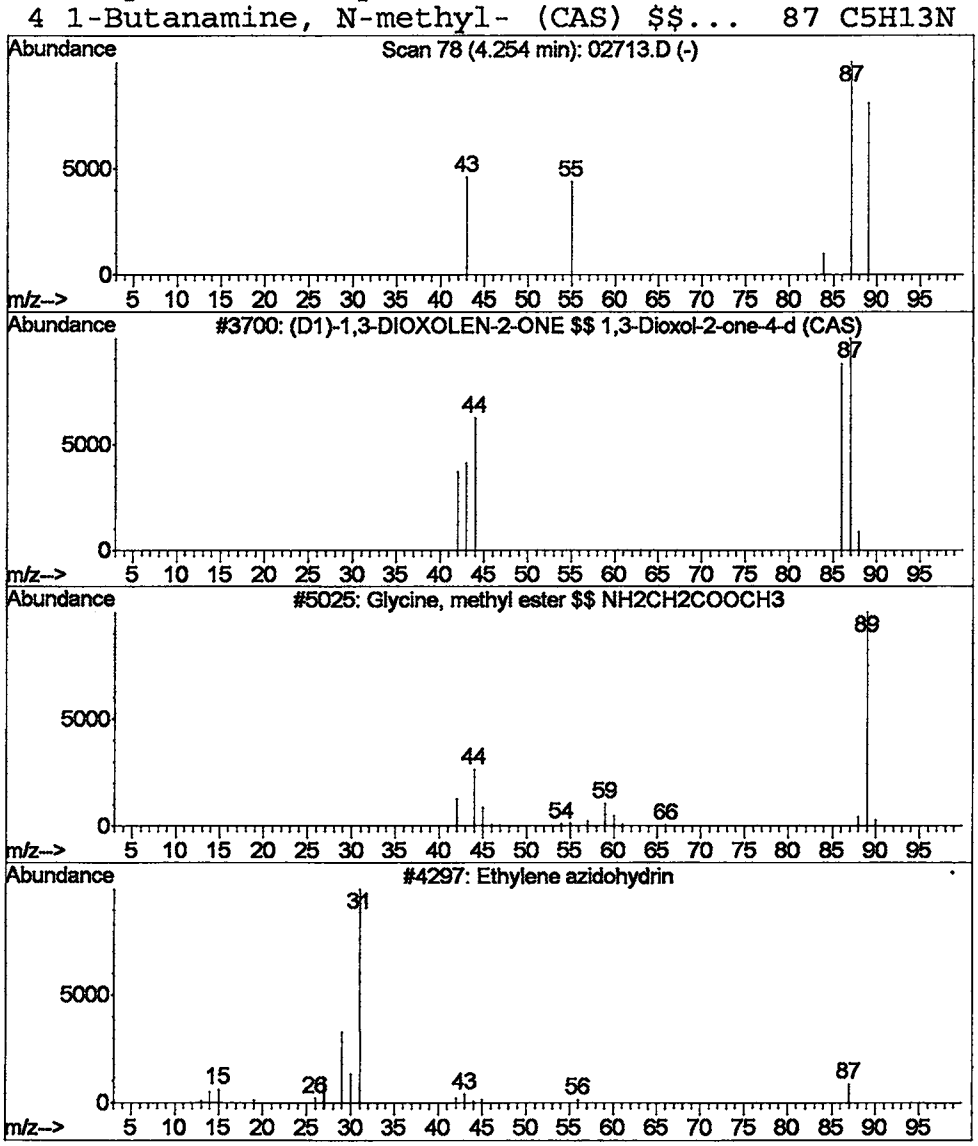
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

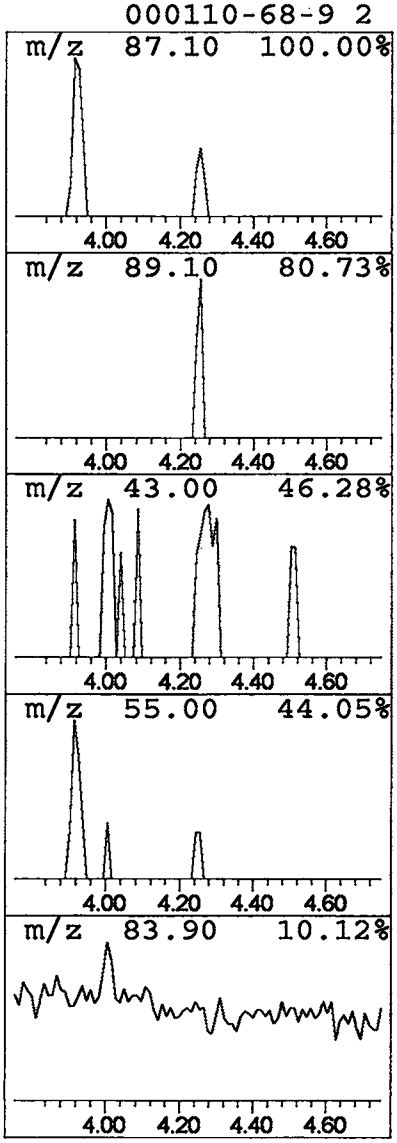
****

Peak Number 4 (D1)-1,3-DIOXOLEN-2-ONE $$ ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.25	0.06 ug/L	23619	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(D1)-1,3-DIOXOLEN-2-ONE $$ 1,3-D...	87	C3HDO3	067790-46-9	4
2	Glycine, methyl ester $$ NH2CH2C...	89	C3H7NO2	000616-34-2	3
3	Ethylene azidohydrin	87	C2H5N3O	000000-00-0	3
4	1-Butanamine, N-methyl- (CAS) $$...	87	C5H13N	000110-68-9	2





Library Search Compound Report

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Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

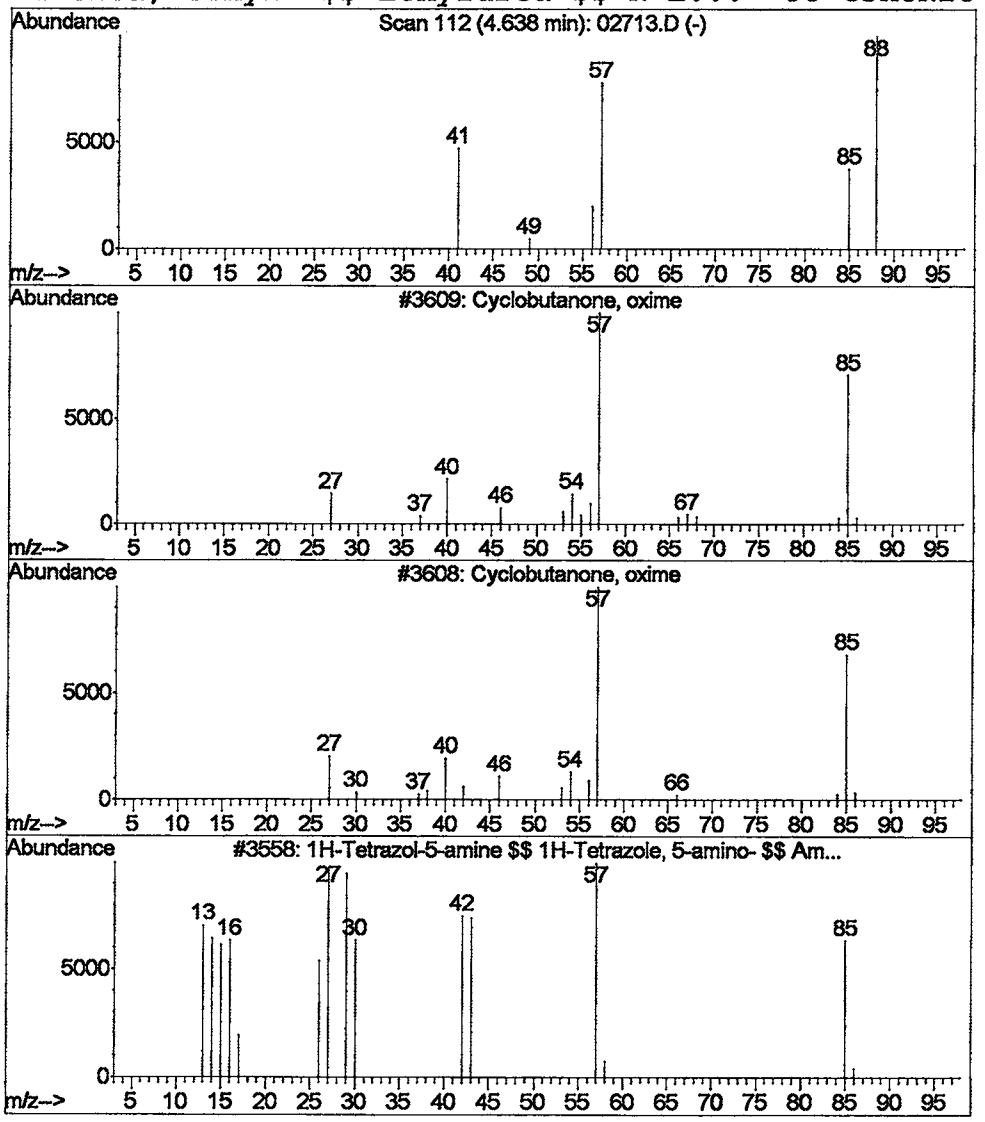
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

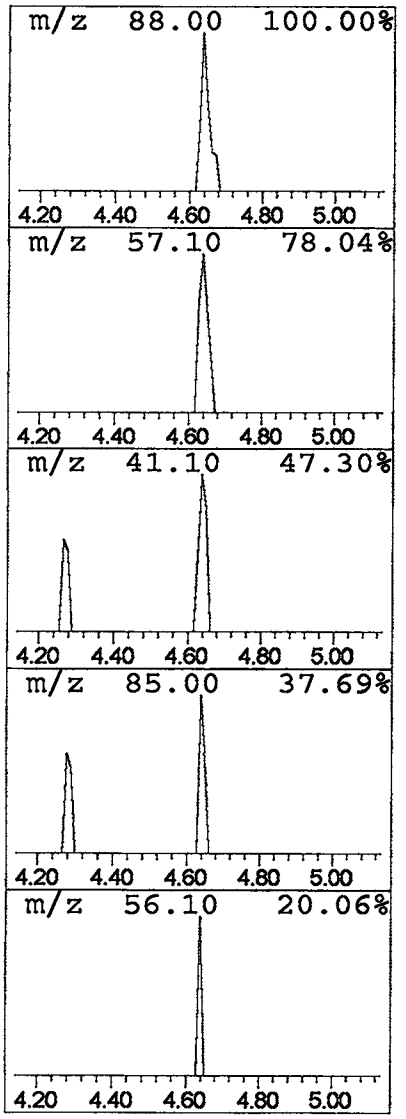
****

Peak Number 5 Cyclobutanone, oxime Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.64	0.05 ug/L	18997	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutanone, oxime	85	C4H7NO	002972-05-6	4
2	Cyclobutanone, oxime	85	C4H7NO	002972-05-6	4
3	1H-Tetrazol-5-amine $$ 1H-Tetraz...	85	CH3N5	004418-61-5	4
4	Urea, ethyl- $$ Ethylurea $$ N-E...	88	C3H8N2O	000625-52-5	4





Library Search Compound Report

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Acq On : 28 Feb 2002 12:10 pm Operator:
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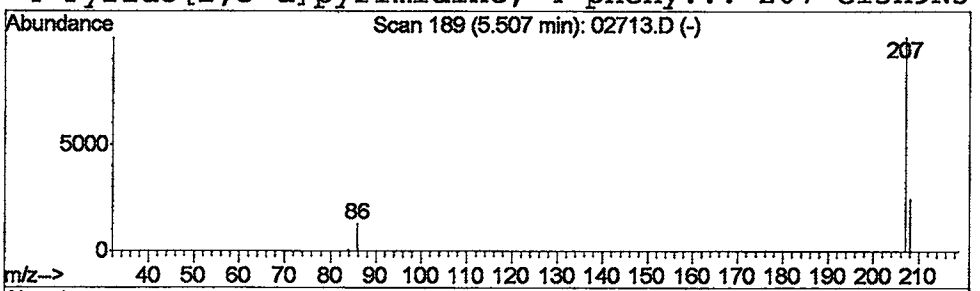
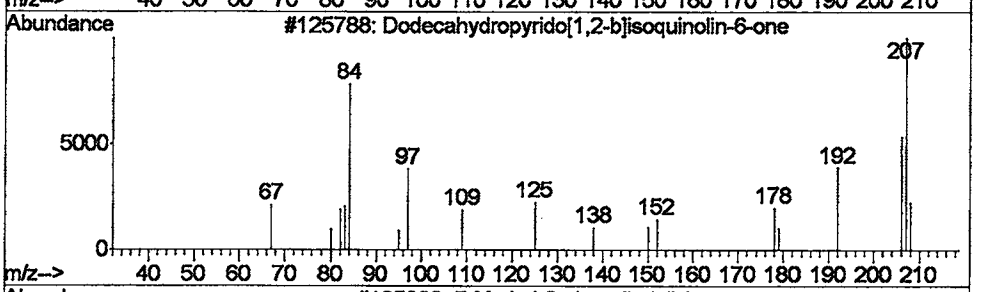
****

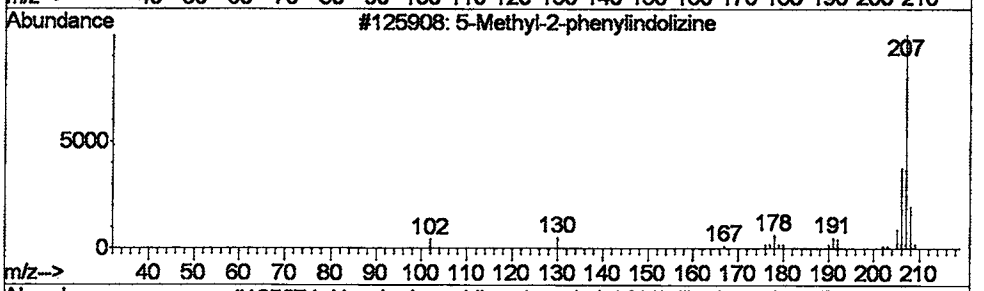
Peak Number 6 Dodecahydropyrido[1,2-b]iso... Concentration Rank 12

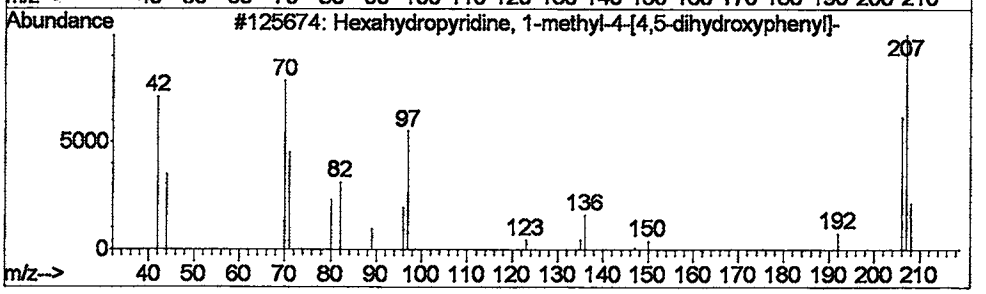
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.51	0.03 ug/L	11950	1,4-Dichlorobenzene-d4	9.72

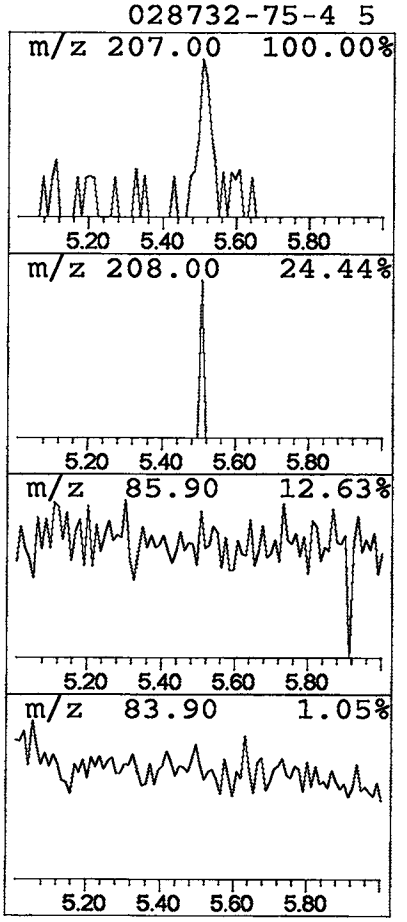
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecahydropyrido[1,2-b]isoquino...	207	C13H21NO	000000-00-0	5
2	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	5
3	Hexahydropyridine, 1-methyl-4-[4...	207	C12H17NO2	094427-47-1	5
4	Pyrido[2,3-d]pyrimidine, 4-pheny...	207	C13H9N3	028732-75-4	5











Library Search Compound Report

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Acq On : 28 Feb 2002 12:10 pm Operator:
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Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

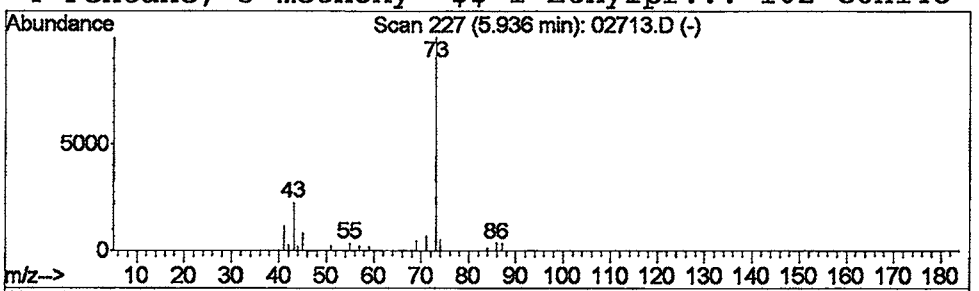
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

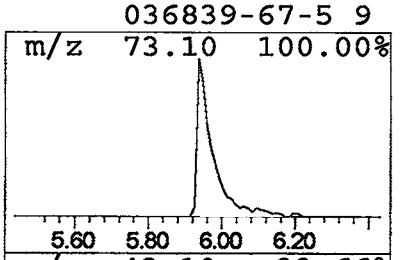
****

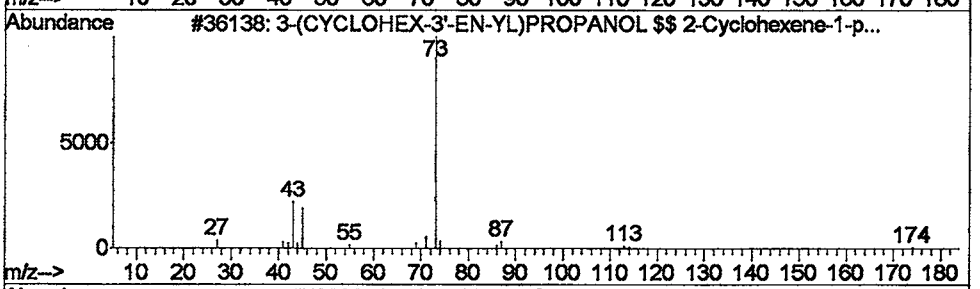
Peak Number 7 3-(CYCLOHEX-3'-EN-YL)PROPAN... Concentration Rank 1

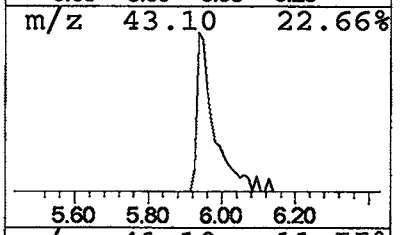
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.94	0.35 ug/L	130387	1,4-Dichlorobenzene-d4	9.72

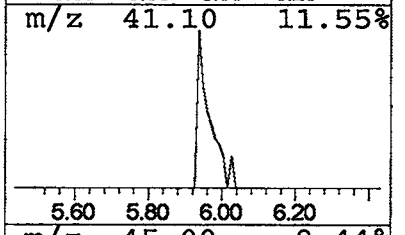
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(CYCLOHEX-3'-EN-YL)PROPANOL $$...	174	C8H14O4	015745-87-6	50
2	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	50
3	N-Ethylformamide $$ N-Formylethy...	73	C3H7NO	000627-45-2	9
4	Pentane, 3-methoxy- $$ 1-Ethylpr...	102	C6H14O	036839-67-5	9

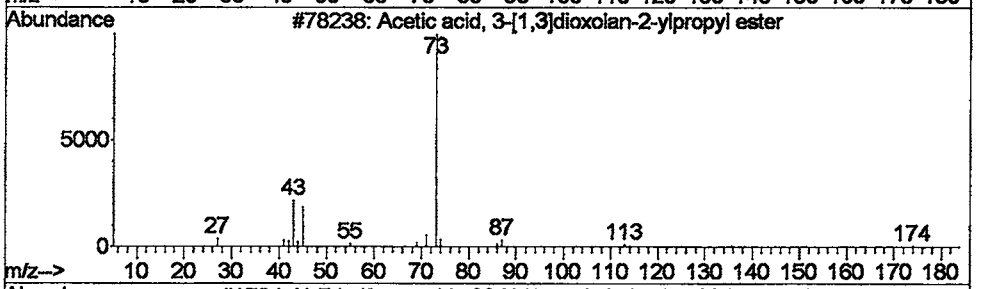


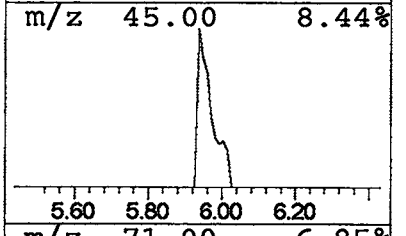


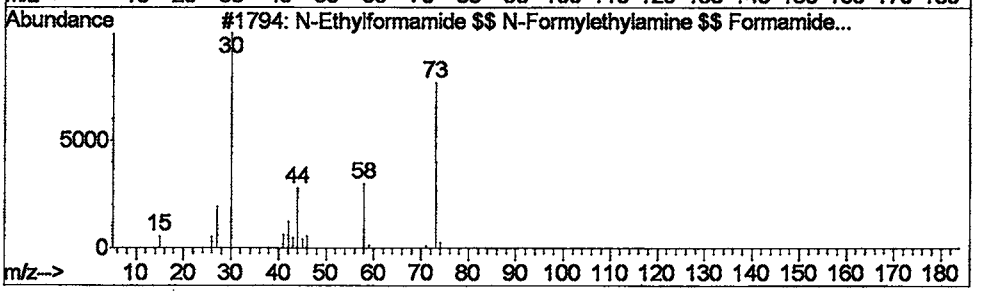


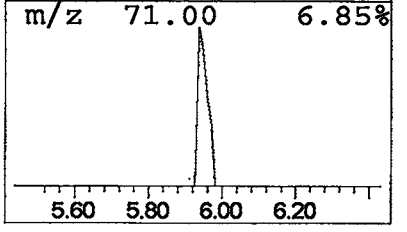












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02713.D Vial: 3
Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

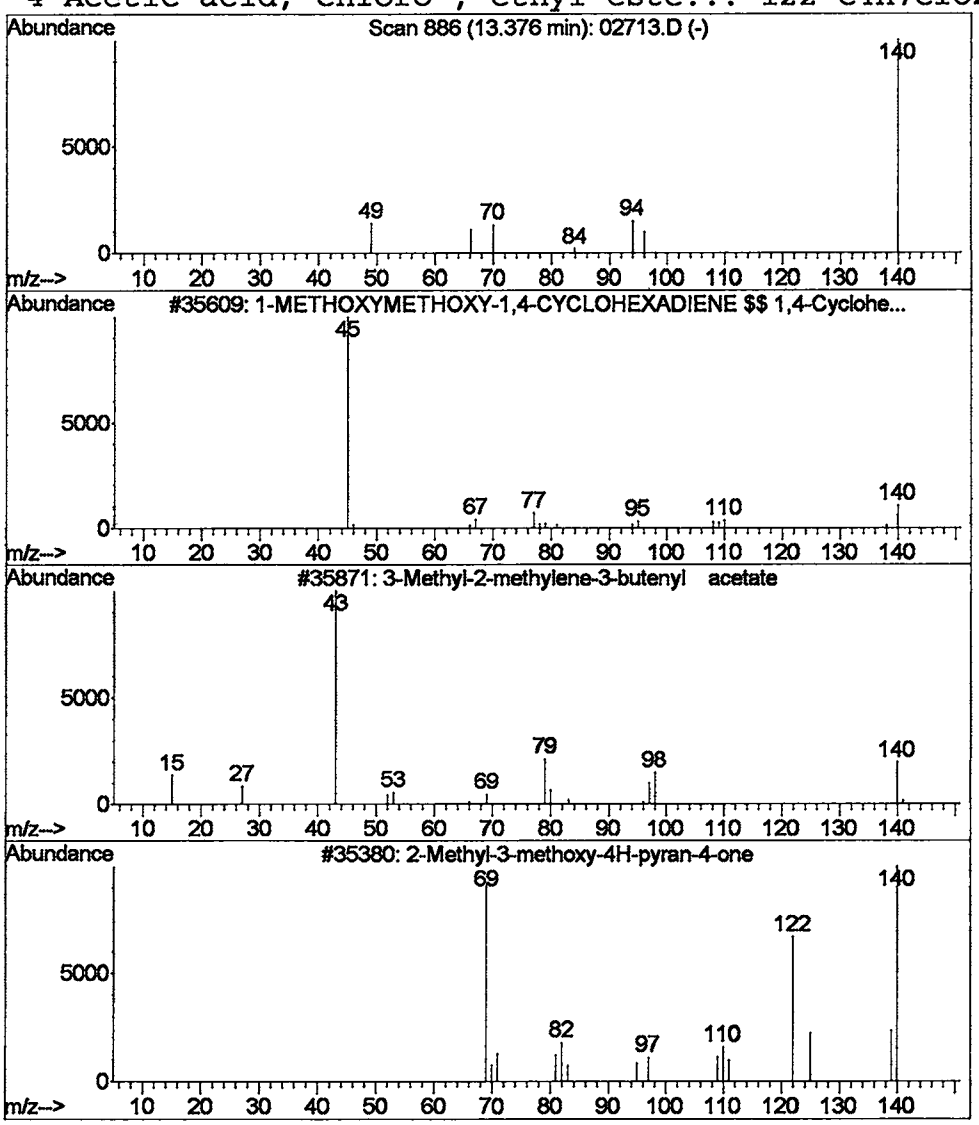
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

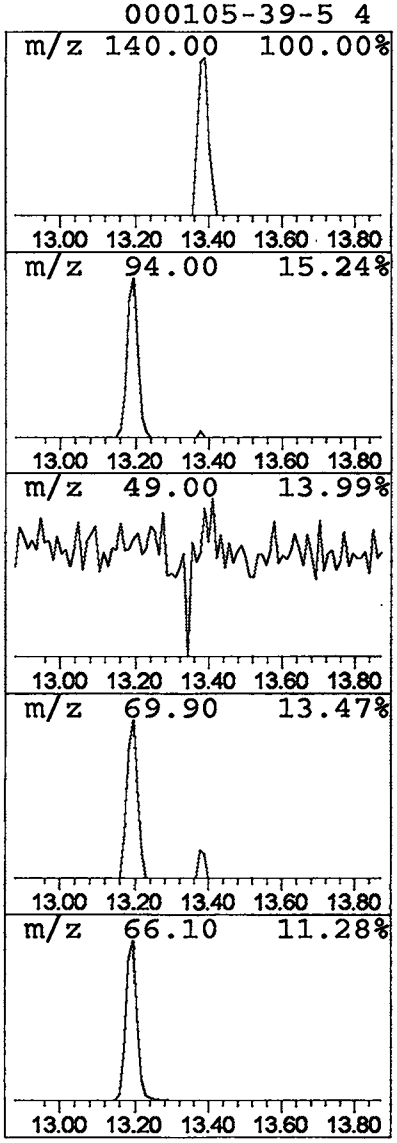
****

Peak Number 8 1-METHOXYMETHOXY-1,4-CYCLOH... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	0.06 ug/L	33222	Naphthalene-d8	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-METHOXYMETHOXY-1,4-CYCLOHEXADI...	140	C8H12O2	057234-30-7	5
2	3-Methyl-2-methylene-3-butenyl ...	140	C8H12O2	000000-00-0	4
3	2-Methyl-3-methoxy-4H-pyran-4-one	140	C7H8O3	000000-00-0	4
4	Acetic acid, chloro-, ethyl este...	122	C4H7ClO2	000105-39-5	4





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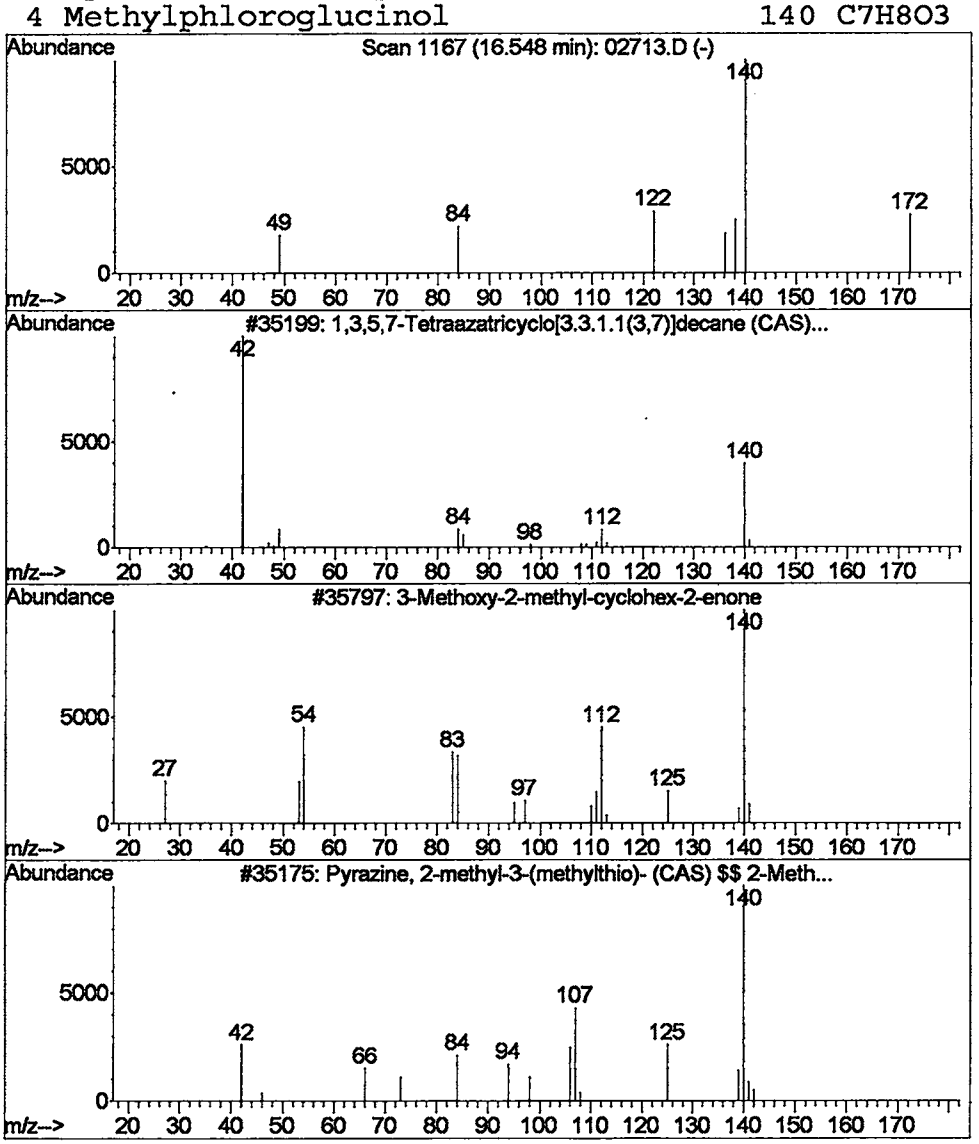
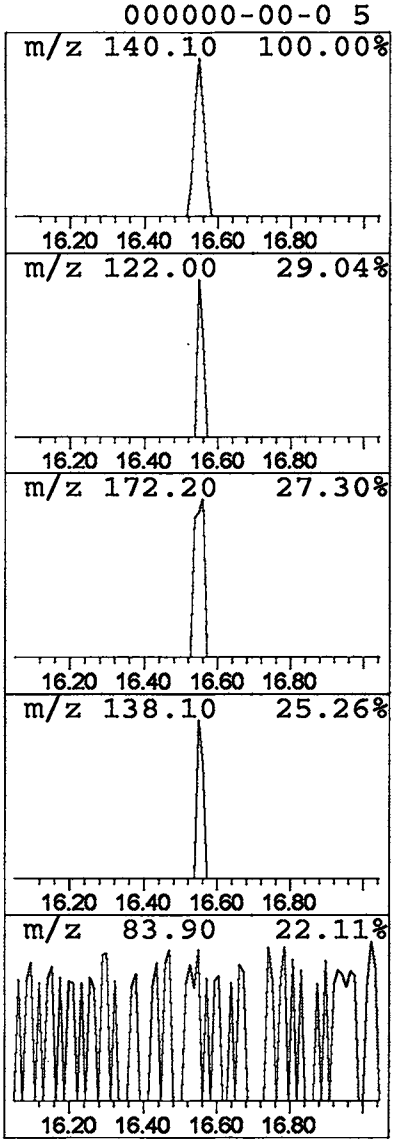
Peak Number 9 1,3,5,7-Tetraazatricyclo[3.... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	0.03 ug/L	15158	Acenaphthene-d10	18.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5,7-Tetraazatricyclo[3.3.1.1...	140	C6H12N4	000100-97-0	7
2	3-Methoxy-2-methyl-cyclohex-2-enone	140	C8H12O2	025112-91-8	5
3	Pyrazine, 2-methyl-3-(methylthio...	140	C6H8N2S	002882-20-4	5
4	Methylphloroglucinol	140	C7H8O3	000000-00-0	5





Library Search Compound Report

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Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 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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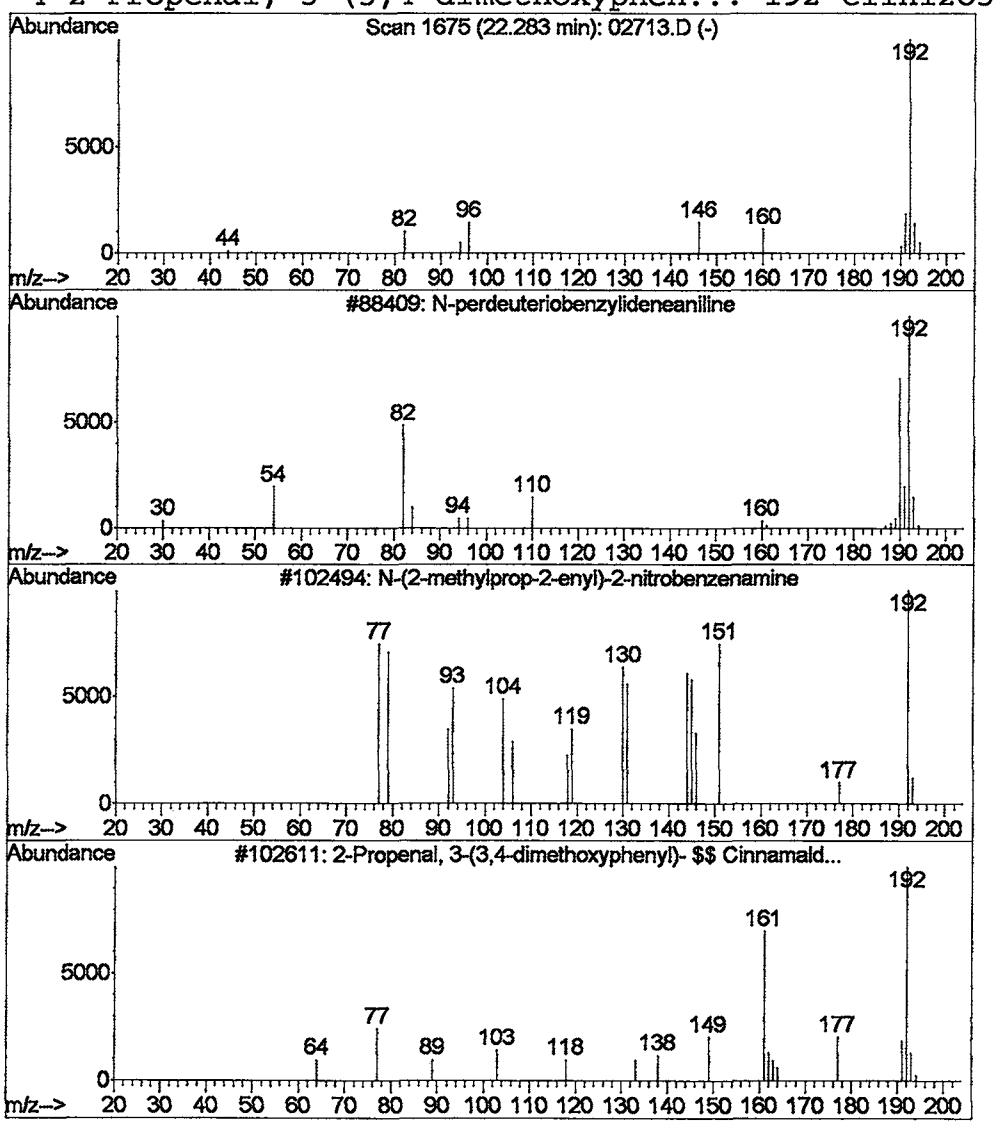
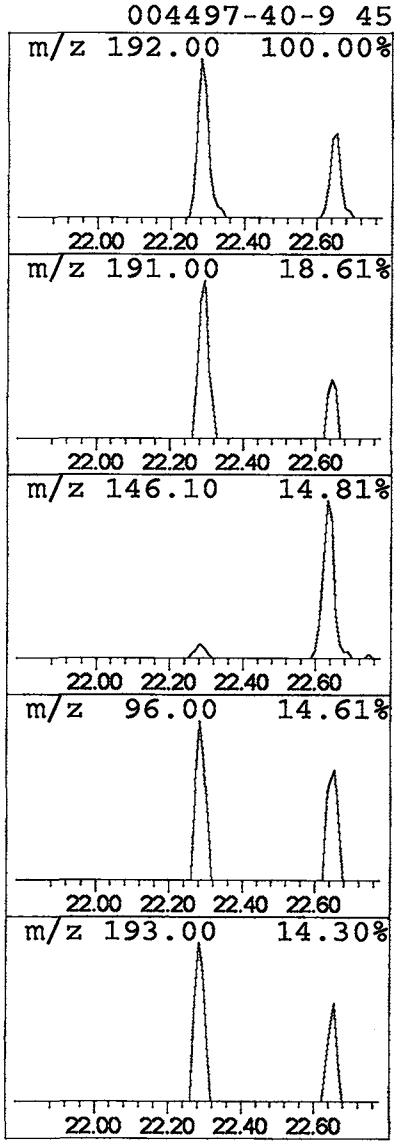
Peak Number 10 N-perdeuteriobenzylideneani... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.10 ug/L	59076	Phenanthrene-d10	22.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	64
2	N-(2-methylprop-2-enyl)-2-nitrob...	192	C10H12N2O2	000000-00-0	52
3	2-Propenal, 3-(3,4-dimethoxyphen...	192	C11H12O3	004497-40-9	45
4	2-Propenal, 3-(3,4-dimethoxyphen...	192	C11H12O3	004497-40-9	45





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Title : 508 scan method
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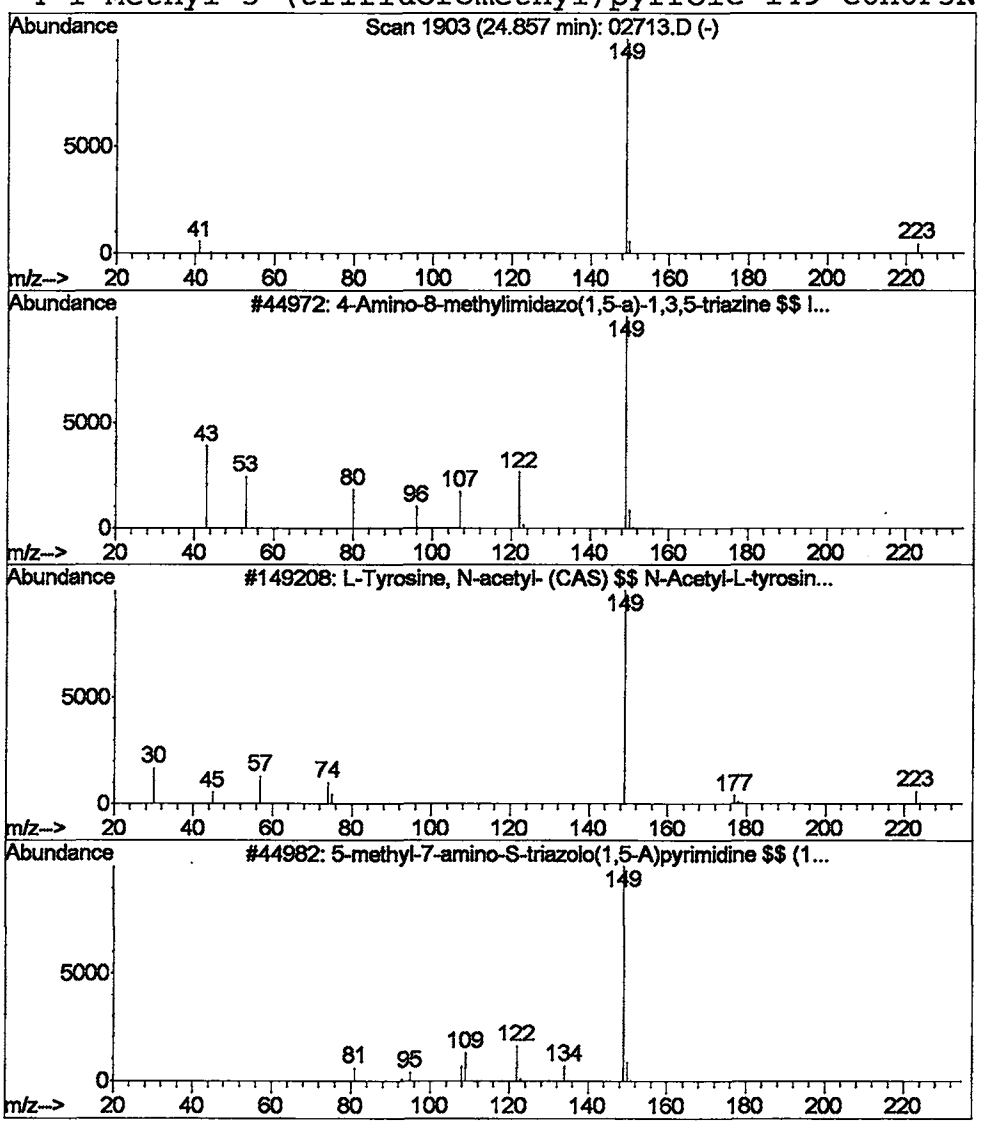
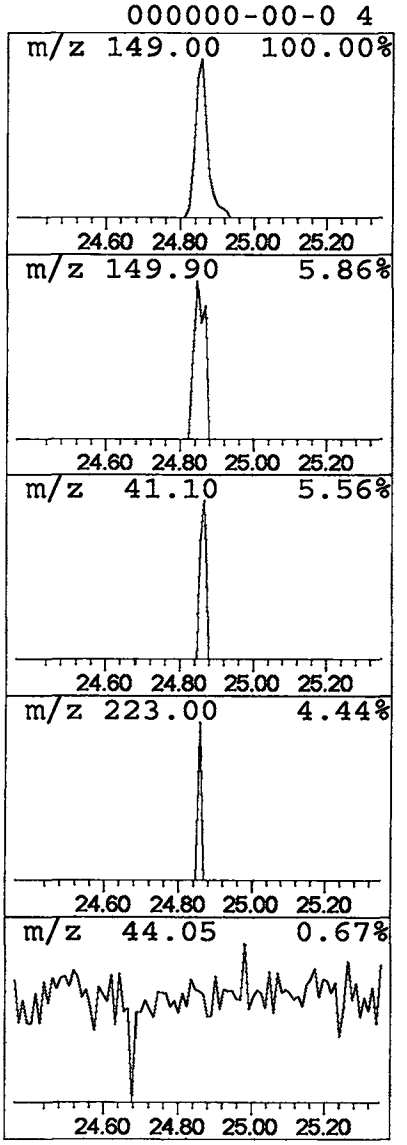
Peak Number 11 4-Amino-8-methylimidazo(1,5... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.86	0.06 ug/L	33312	Phenanthrene-d10	22.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Amino-8-methylimidazo(1,5-a)-1...	149	C6H7N5	079681-01-9	4
2	L-Tyrosine, N-acetyl- (CAS) $$ N...	223	C11H13NO4	000537-55-3	4
3	5-methyl-7-amino-S-triazolo(1,5-...	149	C6H7N5	033376-96-4	4
4	1-Methyl-5-(trifluoromethyl)pyrrole	149	C6H6F3N	000000-00-0	4





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Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
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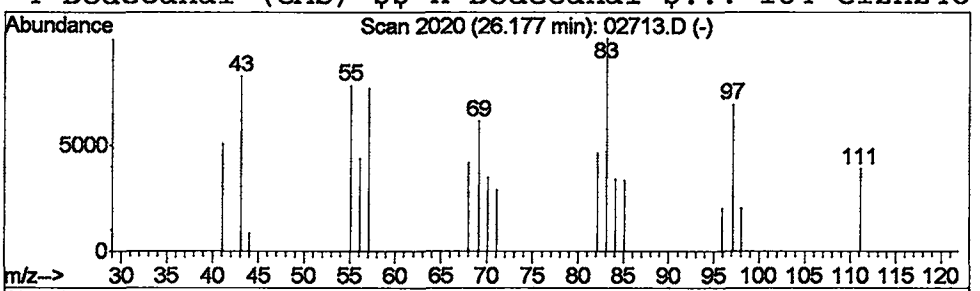
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

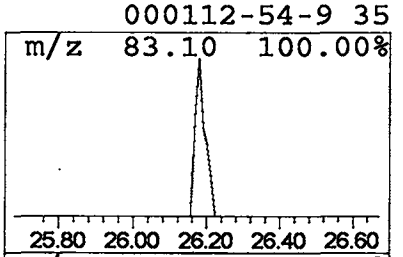
****

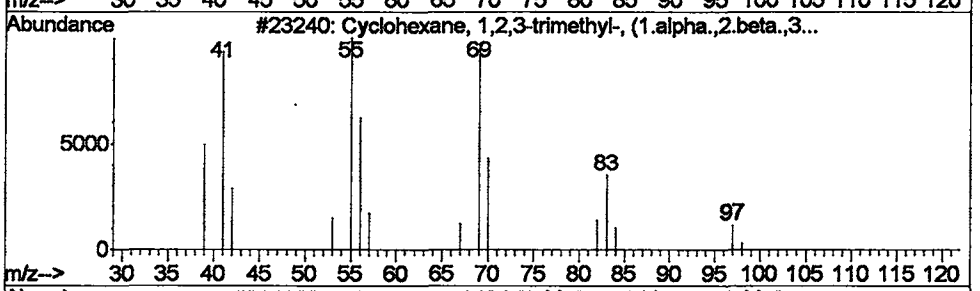
Peak Number 12 Cyclohexane, 1,2,3-trimethy... Concentration Rank 4

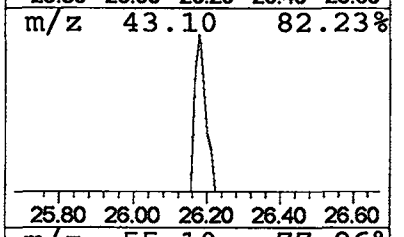
R.T.	EstConc	Area	Relative to ISTD	R.T.
26.18	0.07 ug/L	44003	Phenanthrene-d10	22.63

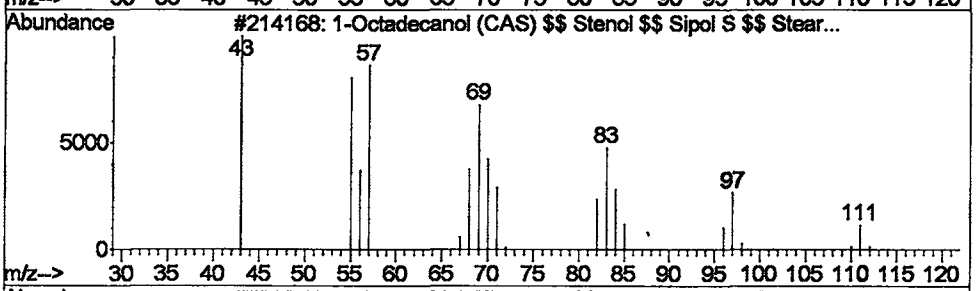
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,3-trimethyl-, (...	126	C9H18	001678-81-5	50
2	1-Octadecanol (CAS) $$ Stenol $$...	270	C18H38O	000112-92-5	43
3	7-Oxabicyclo[4.1.0]heptane $$ Cy...	98	C6H10O	000286-20-4	38
4	Dodecanal (CAS) $$ n-Dodecanal $...	184	C12H24O	000112-54-9	35

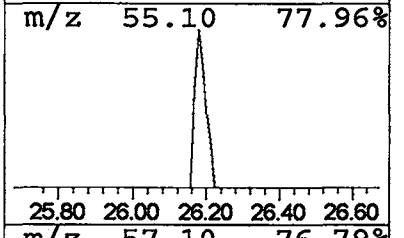


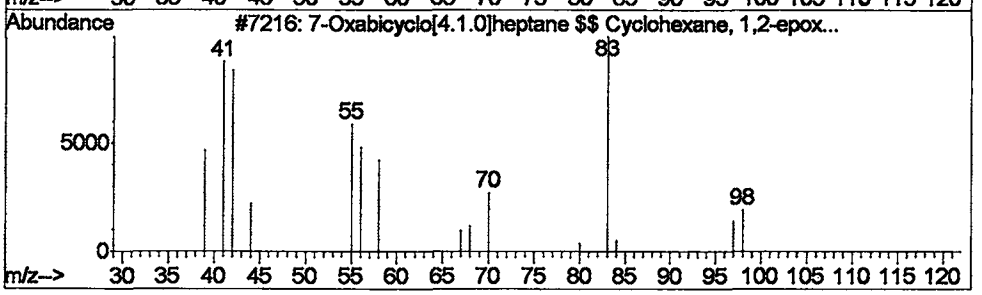


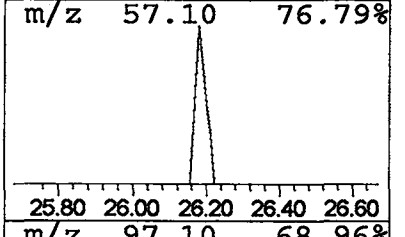


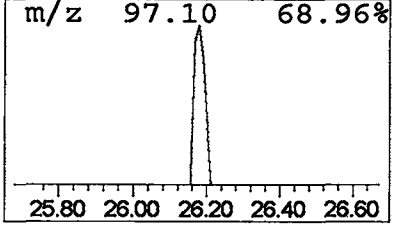












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Acq On : 28 Feb 2002 12:10 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : February 28, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

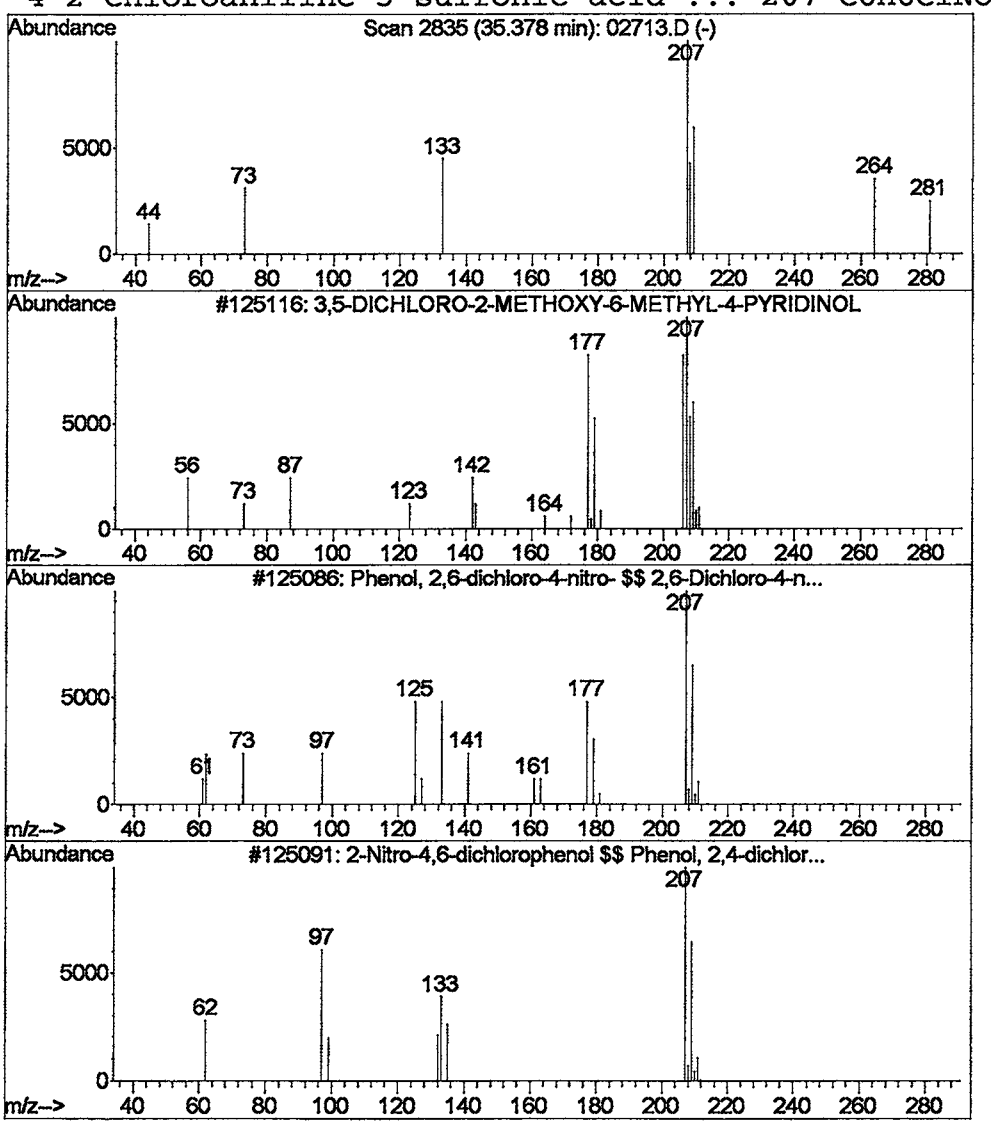
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

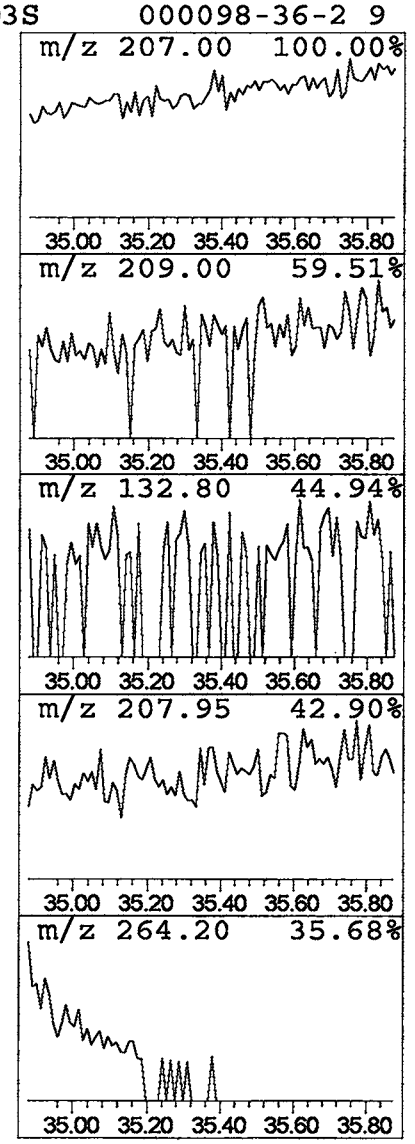
****

Peak Number 13 3,5-DICHLORO-2-METHOXY-6-ME... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.38	0.05 ug/L	19382	Perylene-d12	34.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-DICHLORO-2-METHOXY-6-METHYL-...	207	C7H7Cl2NO2	000000-00-0	50
2	Phenol, 2,6-dichloro-4-nitro- $$...	207	C6H3Cl2NO3	000618-80-4	47
3	2-Nitro-4,6-dichlorophenol $$ Ph...	207	C6H3Cl2NO3	000609-89-2	38
4	2-Chloroaniline-5-sulfonic acid ...	207	C6H6ClNO3S	000098-36-2	9





Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 28 Feb 2002 12:10 pm
Data File: C:\MSDCHEM\1\5973N\02FEB28\02713.D
Name: 508 scan
Misc: February 28, 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...	3.61	0.2	ug/L	58932	1	9.72	7416380	20.0
ISOBUTYRALDEHYDE,...	3.92	0.1	ug/L	26319	1	9.72	7416380	20.0
1-Methyl-1-propyn...	4.01	0.0	ug/L	18448	1	9.72	7416380	20.0
(D1)-1,3-DIOXOLEN...	4.25	0.1	ug/L	23619	1	9.72	7416380	20.0
Cyclobutanone, oxime	4.64	0.1	ug/L	18997	1	9.72	7416380	20.0
Dodecahydropyrido...	5.51	0.0	ug/L	11950	1	9.72	7416380	20.0
3-(CYCLOHEX-3'-EN...	5.94	0.4	ug/L	130387	1	9.72	7416380	20.0
1-METHOXYMETHOXY-...	13.38	0.1	ug/L	33222	2	13.20	10627700	20.0
1,3,5,7-Tetraazat...	16.55	0.0	ug/L	15158	3	18.33	11530900	20.0
N-perdeuteriobenz...	22.28	0.1	ug/L	59076	4	22.63	12050100	20.0
4-Amino-8-methyli...	24.86	0.1	ug/L	33312	4	22.63	12050100	20.0
Cyclohexane, 1,2,...	26.18	0.1	ug/L	44003	4	22.63	12050100	20.0
3,5-DICHLORO-2-ME...	35.38	0.0	ug/L	19382	6	34.41	8248860	20.0