Comments for Sample AD26535 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-16-2001 Time: 08:53:44

A scan of the extract prepared from the 508 analysis, from 35 to 550 amu using a mass spectrometer, reveals the presence of 1,1,2,2-Tetrachloroethane at an estimated level of 3.4 ug/L, and with a fit of 95 out of 100.

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 12/16/2001 Time: 08:51:00

Sample: AD26535
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/16/01 08:49 Ended: 12/16/01 08:49 Analyst: DLV
Page: 1

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC0701\26535.D Vial: 10
Acq On : 8 Dec 2001 2:19 am Operator: GALOS
Sample : gc/ms unknown 11/6 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 10 9:28 2001 Quant Results File: NP112701.RES

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration
DataAcq Meth : NP112701

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.88	152	967556	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.49	136	3662800	20.00	ng/uL	-0.06
31) Acenaphthene-d10	20.76	164	1664967	20.00	ng/uL	-0.06
49) Phenanthrene-d10	25.17	188	2351682	20.00	ng/uL	-0.07
59) Chrysene-d12	33.18	240	1629717	20.00	ng/uL	-0.07
68) Perylene-d12	37.26	264	917961	20.00	ng/uL	-0.07
System Monitoring Compounds						
4) 2-Fluorophenol	0.00	112	0	0.00	ng/uL	
Spiked Amount 75.000	Range 18 - 42		Recovery =	0.00%#		
5) Phenol-d5	0.00	99	0	0.00	ng/uL	
Spiked Amount 75.000	Range 19 - 32		Recovery =	0.00%#		
6) 2-Chlorophenol-d4	11.88	132	395	0.01	ng/uL	0.46
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.13	152	1083	0.03	ng/uL	-0.33
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.06%#		
20) Nitrobenzene-d5	0.00	82	0	0.00	ng/uL	
Spiked Amount 50.000	Range 55 - 98		Recovery =	0.00%#		
32) 2-Fluorobiphenyl	18.92	172	1725	0.02	ng/uL	0.08
Spiked Amount 50.000	Range 42 - 78		Recovery =	0.04%#		
33) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng/uL	
Spiked Amount 75.000	Range 45 - 96		Recovery =	0.00%#		
62) Terphenyl-d14	0.00	244	0	0.00	ng/uL	
Spiked Amount 50.000	Range 58 - 98		Recovery =	0.00%#		
Target Compounds						Qvalue
2) Pyridine	0.00	79	0	N.D.		
3) N-nitrosodimethylamine	0.00	74	0	N.D.		
8) Phenol	0.00	94	0	N.D.		
9) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
10) 2-Chlorophenol	0.00	128	0	N.D.		
11) 1,3-Dichlorobenzene	0.00	146	0	N.D.		
12) 1,4-Dichlorobenzene	0.00	146	0	N.D.		
13) 1,2-Dichlorobenzene	0.00	146	0	N.D.		
14) 2-Methylphenol	0.00	108	0	N.D.		
15) 2,2'-oxybis(1-Chloropropan	12.88	45	2738	N.D.		
16) 3&4-Methylphenol	0.00	108	0	N.D.		

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

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	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
17)	N-Nitrosodi-n-propylamine	0.00	70	0	N.D.	
18)	Hexachloroethane	0.00	117	0	N.D.	
21)	Nitrobenzene	0.00	77	0	N.D.	
22)	Isophorone	0.00	82	0	N.D.	
23)	2-Nitrophenol	0.00	139	0	N.D.	
24)	2,4-Dimethylphenol	0.00	107	0	N.D.	
25)	bis(2-Chloroethoxy)methane	0.00	93	0	N.D.	
26)	2,4-Dichlorophenol	0.00	162	0	N.D.	
27)	1,2,4-Trichlorobenzene	0.00	180	0	N.D.	
28)	Naphthalene	15.55	128	869	N.D.	
29)	Hexachlorobutadiene	0.00	225	0	N.D.	
30)	4-Chloro-3-methylphenol	0.00	107	0	N.D.	
34)	Hexachlorocyclopentadiene	0.00	237	0	N.D.	
35)	2,4,6-Trichlorophenol	0.00	196	0	N.D.	
36)	2,4,5-Trichlorophenol	0.00	196	0	N.D.	
37)	2-Chloronaphthalene	0.00	162	0	N.D.	
38)	Dimethylphthalate	0.00	163	0	N.D.	
39)	2,6-Dinitrotoluene	0.00	165	0	N.D.	d
40)	Acenaphthylene	0.00	152	0	N.D.	
41)	Acenaphthene	0.00	153	0	N.D.	
42)	2,4-Dinitrophenol	0.00	184	0	N.D.	
43)	4-Nitrophenol	0.00	65	0	N.D.	d
44)	2,4-Dinitrotoluene	0.00	165	0	N.D.	
45)	Diethylphthalate	0.00	149	0	N.D.	
46)	4-Chlorophenylphenylether	0.00	204	0	N.D.	
47)	Fluorene	0.00	166	0	N.D.	
48)	Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.	
50)	4,6-Dinitro-2-methylphenol	0.00	198	0	N.D.	
51)	N-Nitrosodiphenylamine	0.00	169	0	N.D.	
52)	4-Bromophenyl-phenylether	0.00	248	0	N.D.	
53)	Hexachlorobenzene	0.00	284	0	N.D.	
54)	Pentachlorophenol	0.00	266	0	N.D.	
55)	Phenanthrene	0.00	178	0	N.D.	
56)	Anthracene	0.00	178	0	N.D.	
57)	Di-n-butylphthalate	27.17	149	3991	N.D.	
58)	Fluoranthene	0.00	202	0	N.D.	
60)	3,3'-Dichlorobenzidine	0.00	252	0	N.D.	

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Response via : Initial Calibration
DataAcq Meth : NP112701

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
61)	Benzidine	0.00	184	0	N.D.	
63)	Pyrene	0.00	202	0	N.D.	
64)	Butylbenzylphthalate	0.00	149	0	N.D.	
65)	Benzo(a)anthracene	33.18	228	3038	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.44	149	5596	N.D.	
67)	Chrysene	33.18	228	3038	N.D.	
69)	Di-n-Octylphthalate	0.00	149	0	N.D.	
70)	Benzo(b)fluoranthene	0.00	252	0	N.D.	
71)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
72)	Benzo(a)pyrene	37.26	252	2672	N.D.	
73)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
74)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
75)	Benzo(g,h,i)perylene	0.00	276	0	N.D.	

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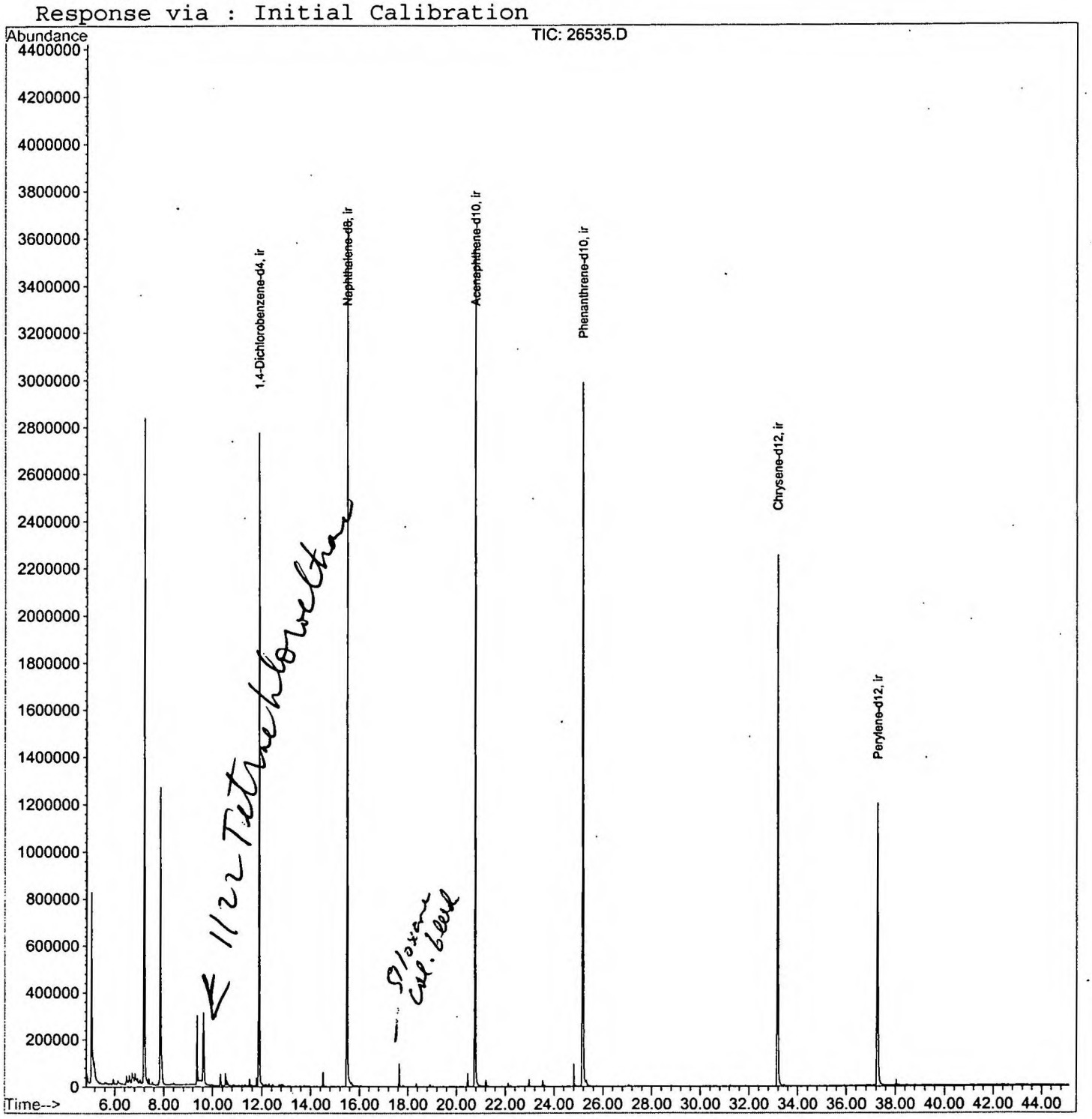
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 10 9:28 2001

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Results File: NP112701.RES

Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC0701\26535.D Vial: 10
Acq On : 8 Dec 2001 2:19 am Operator: GALOS
Sample : gc/ms unknown 11/6 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Signal : TIC

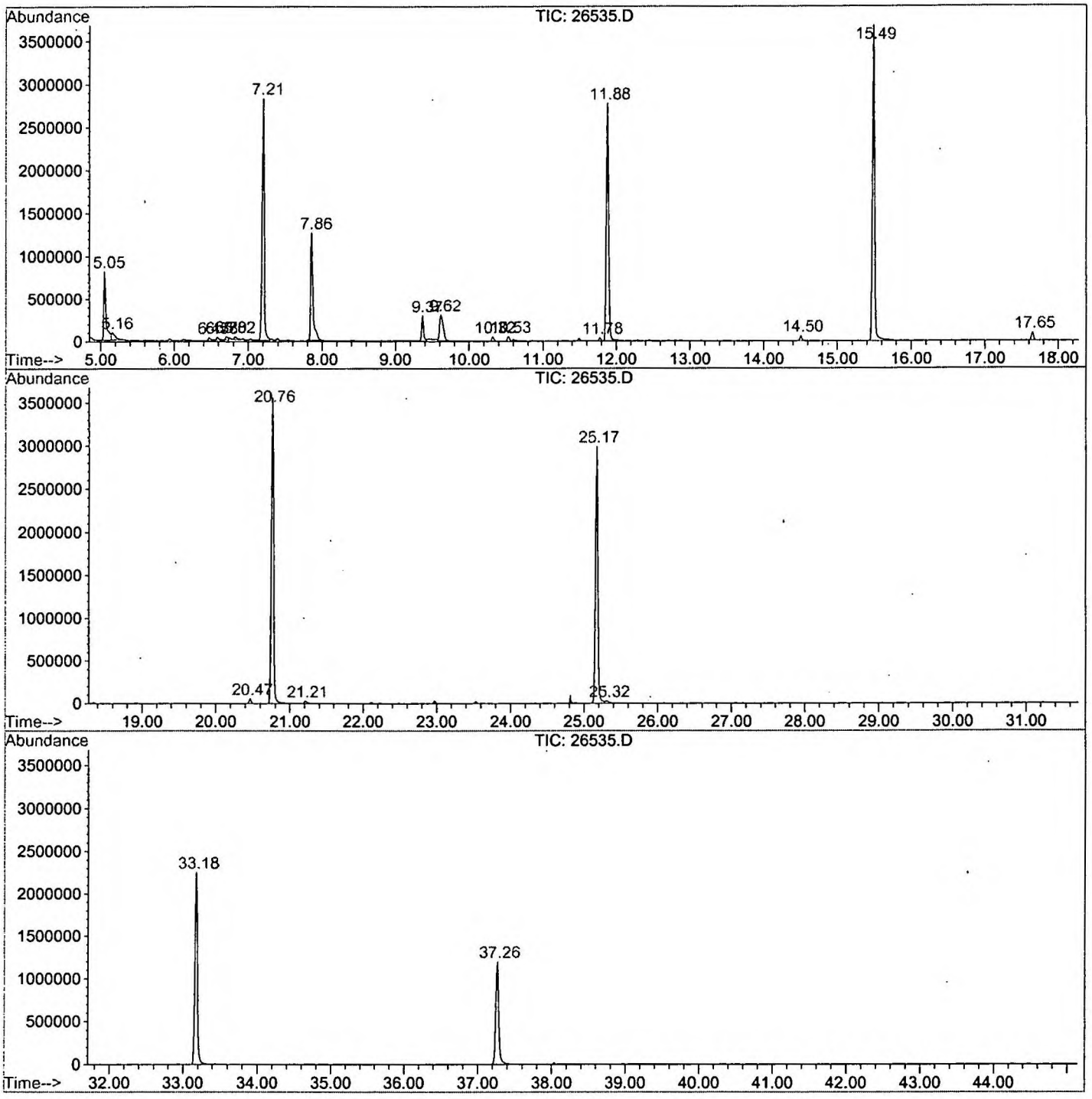
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.050	20	23	33	rBV	809686	1723746	23.22%	3.523%
2	5.160	33	35	48	rVB3	84024	267594	3.61%	0.547%
3	6.472	175	178	187	rBV	38228	115937	1.56%	0.237%
4	6.582	187	190	200	rBV	33126	80160	1.08%	0.164%
5	6.701	200	203	207	rBV	42064	104771	1.41%	0.214%
6	6.820	213	216	225	rBV	34372	100761	1.36%	0.206%
7	7.205	251	258	268	rBV	2823027	5457311	73.52%	11.153%
8	7.856	325	329	346	rVB	1259770	2734894	36.84%	5.589%
9	9.369	489	494	500	rBV	294161	546320	7.36%	1.117%
10	9.617	516	521	534	rVB	304204	1031008	13.89%	2.107%
11	10.323	593	598	606	rBV	51284	107802	1.45%	0.220%
12	10.534	616	621	627	rBV	54768	120030	1.62%	0.245%
13	11.781	751	757	762	rVB2	37687	76957	1.04%	0.157%
14	11.882	762	768	778	rBV	2774415	5975870	80.51%	12.213%
15	14.505	1049	1054	1058	rBV	62022	111314	1.50%	0.227%
16	15.486	1155	1161	1179	rBV	3682446	7422706	100.00%	15.170%
17	17.650	1390	1397	1405	rVB	98187	183034	2.47%	0.374%
18	20.465	1698	1704	1708	rBV	57430	101174	1.36%	0.207%
19	20.759	1730	1736	1757	rBV	3560576	7372863	99.33%	15.068%
20	21.208	1781	1785	1794	rBV2	28342	77280	1.04%	0.158%
21	25.170	2208	2217	2229	rBV2	2987326	6657057	89.69%	13.605%
22	25.316	2229	2233	2241	rVB2	23748	76544	1.03%	0.156%
23	33.175	3084	3090	3110	rBV2	2254181	5179964	69.79%	10.586%
24	37.265	3528	3536	3552	rBV	1197518	3305875	44.54%	6.756%

Sum of corrected areas: 48930972

26535.D NP112701.M Thu Dec 13 11:24:49 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0701\26535.D
Operator : GALOS
Acquired : 8 Dec 2001 2:19 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 11/6
Misc Info :
Vial Number: 10
Quant File :NP112701.RES (RTE Integrator)



26535.D NP112701.M Thu Dec 13 11:24:53 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0701\26535.D
Acq On : 8 Dec 2001 2:19 am
Sample : gc/ms unknown 11/6
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

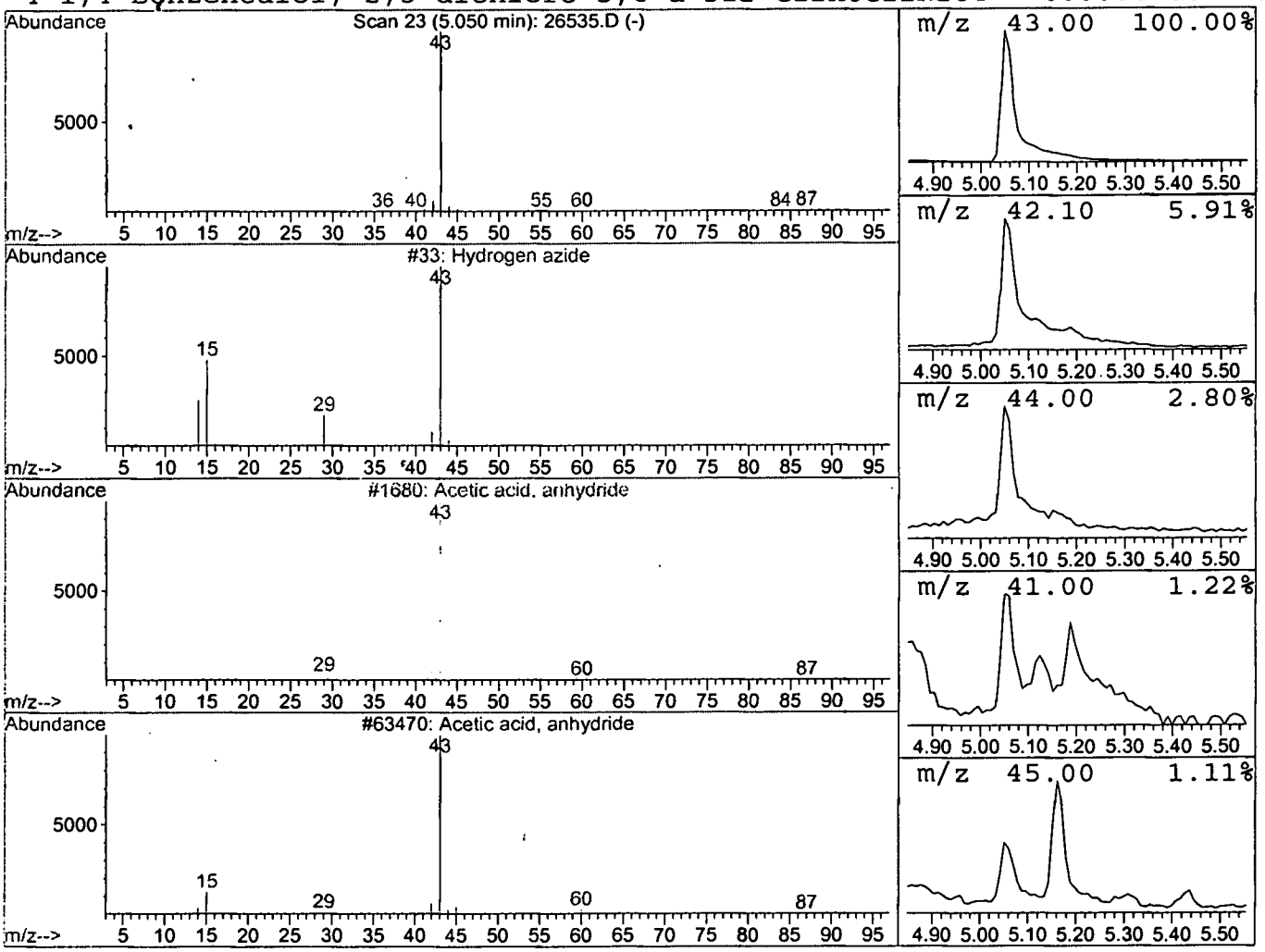
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Hydrogen azide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	5.77 ng/uL	1723750	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrogen azide	43	HN3	007782-79-8	4
2	Acetic acid, anhydride	102	C4H6O3	000108-24-7	2
3	Acetic acid, anhydride	102	C4H6O3	000108-24-7	2
4	1,4-Benzenediol, 2,3-dichloro-5,6-d	312	C12H6Cl2N2O4	000000-00-0	1



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Vial: 10
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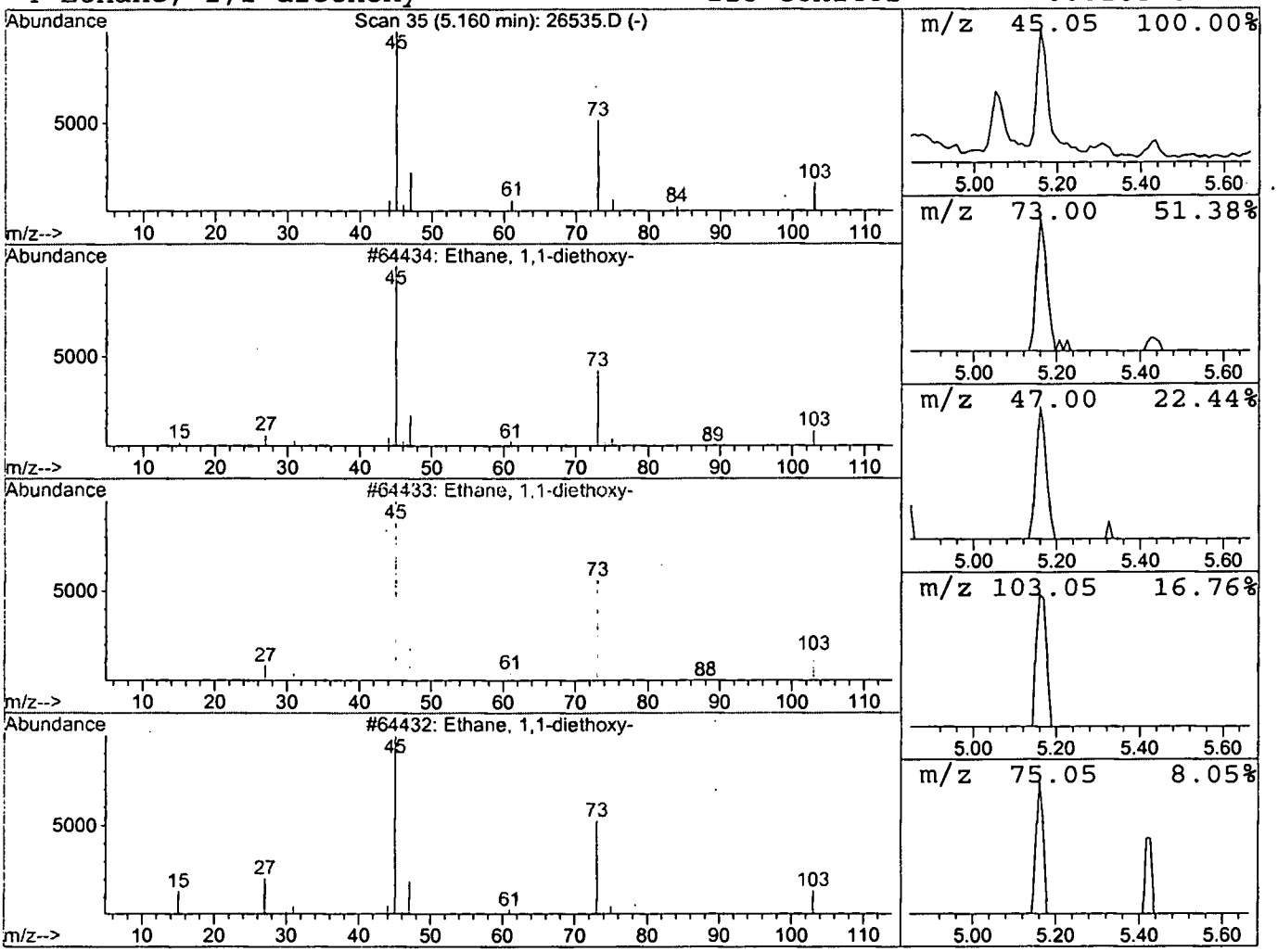
****

Peak Number 2 Ethane, 1,1-diethoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	0.90 ng/uL	267594	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethane, 1,1-diethoxy-	118	C6H14O2	000105-57-7	83
2	Ethane, 1,1-diethoxy-	118	C6H14O2	000105-57-7	83
3	Ethane, 1,1-diethoxy-	118	C6H14O2	000105-57-7	74
4	Ethane, 1,1-diethoxy-	118	C6H14O2	000105-57-7	74



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Misc :
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Vial: 10
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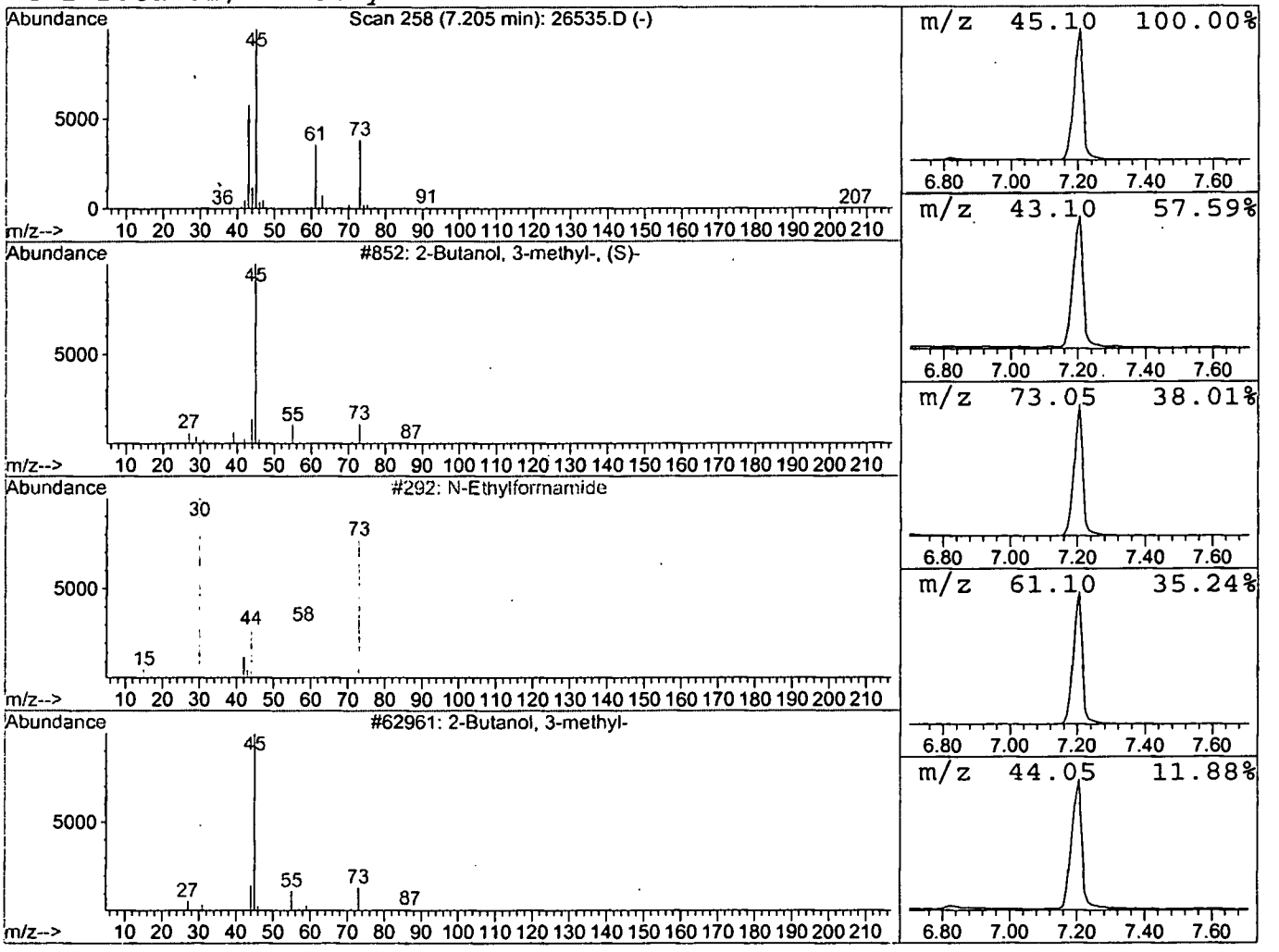
****

Peak Number 3 2-Butanol, 3-methyl-, (S)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.21	18.26 ng/uL	5457310	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 3-methyl-, (S)-	88	C5H12O	001517-66-4	40
2	N-Ethylformamide	73	C3H7NO	000627-45-2	9
3	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
4	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9



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Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
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Library : C:\DATABASE\NBS75K.L

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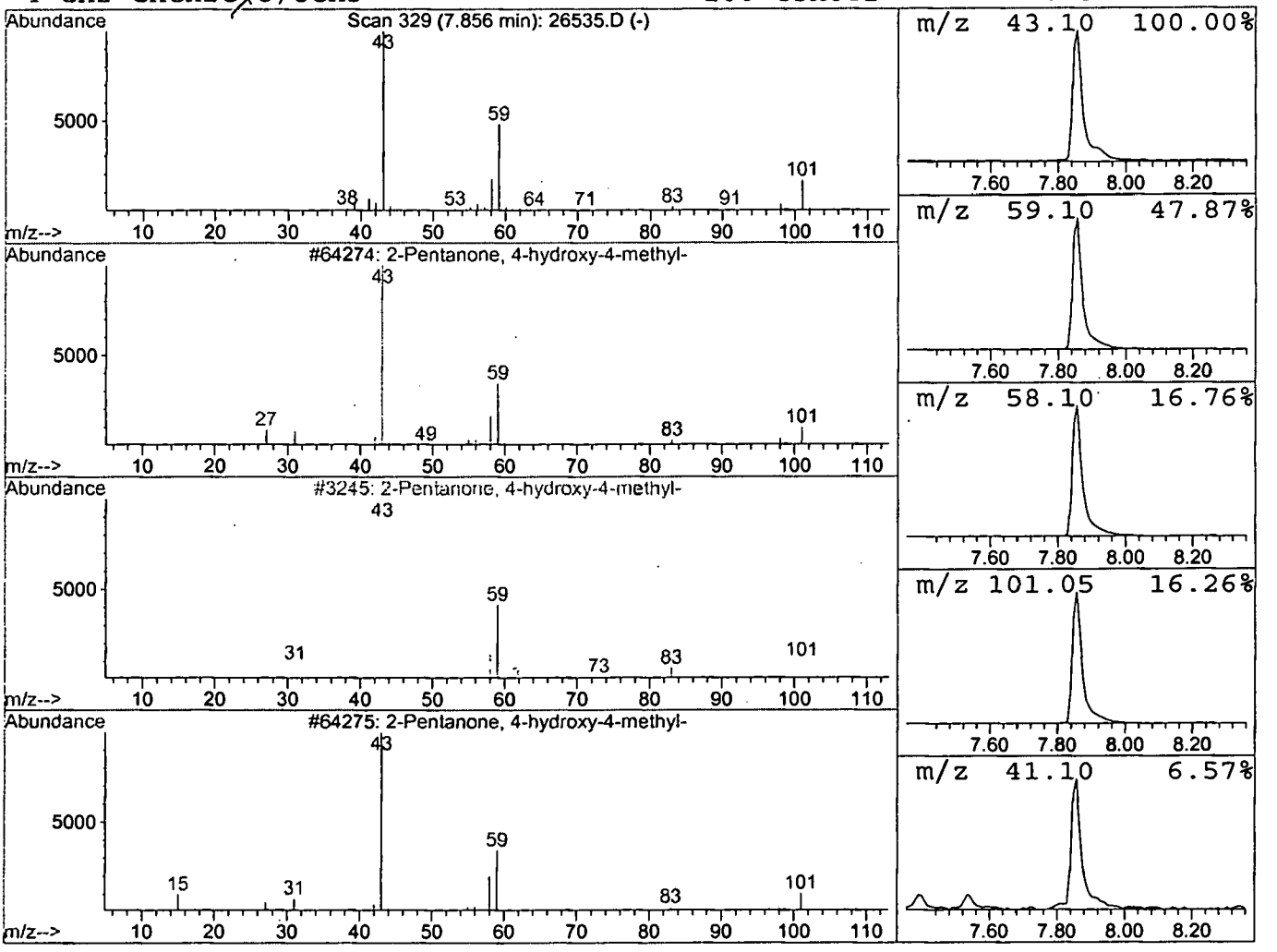
Peak Number 4 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	9.15 ng/uL	2734890	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	43
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	23
4	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	10



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Misc :
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Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
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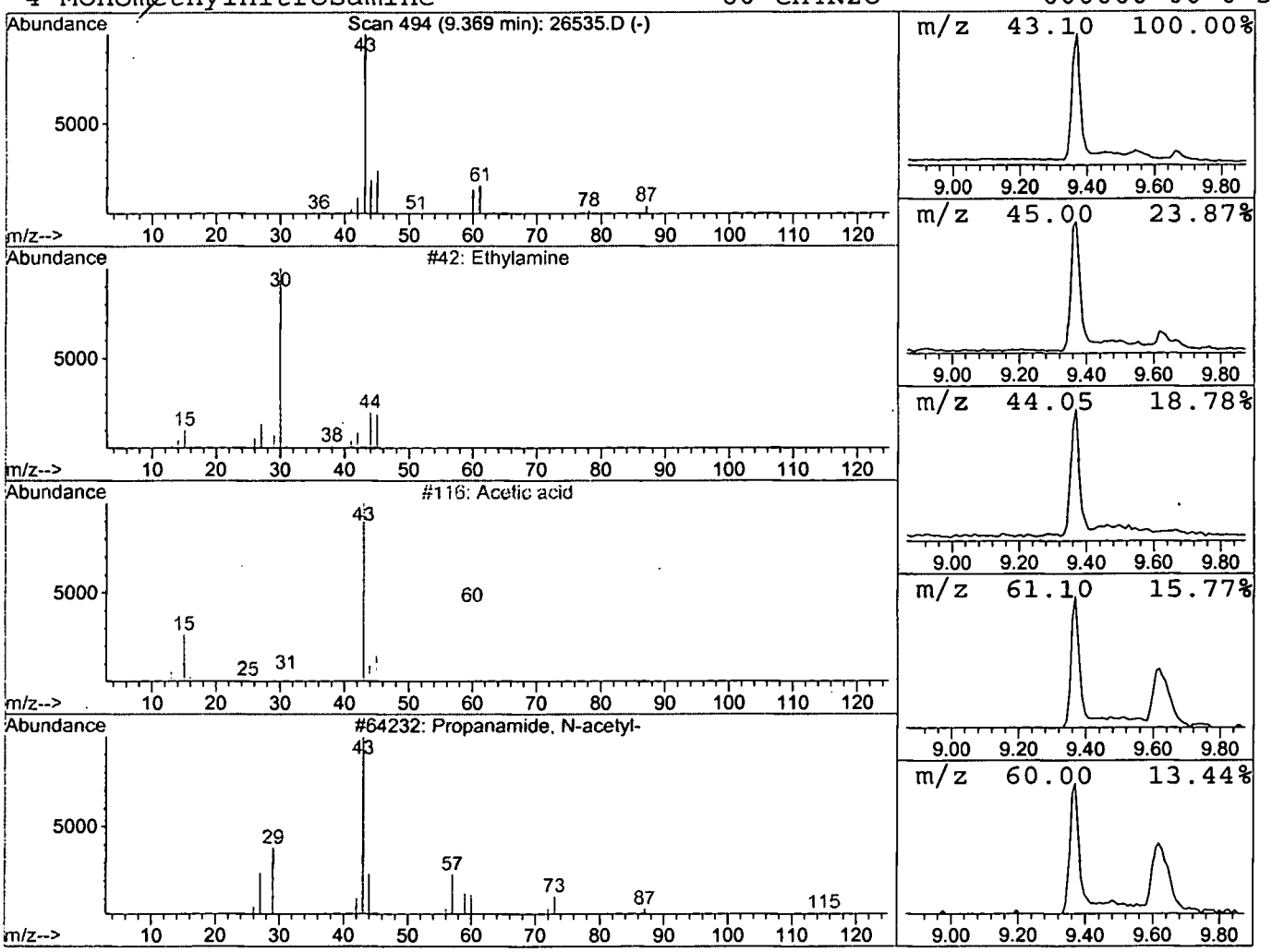
****

Peak Number 5 Ethylamine Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	1.83 ng/uL	546320	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylamine	45	C2H7N	000075-04-7	9
2	Acetic acid	60	C2H4O2	000064-19-7	9
3	Propanamide, N-acetyl-	115	C5H9NO2	019264-34-7	9
4	Monomethylnitrosamine	60	CH4N2O	000000-00-0	5



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Misc :
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Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
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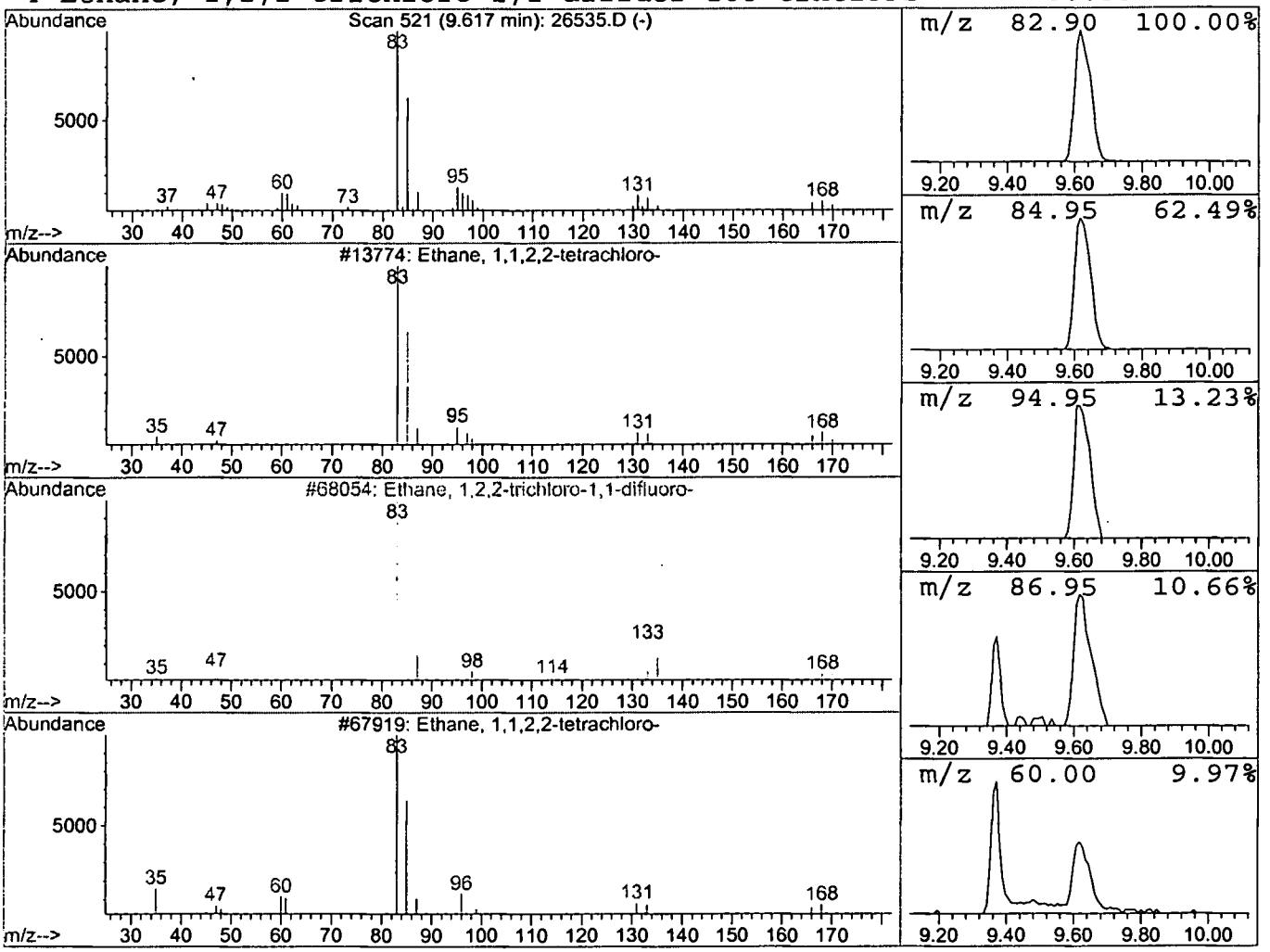
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Ethane, 1,1,2,2-tetrachloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	3.45 ng/uL	1031010	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethane, 1,1,2,2-tetrachloro-	166	C2H2Cl4	000079-34-5	95
2	Ethane, 1,2,2-trichloro-1,1-difluor	168	C2HCl3F2	000354-21-2	76
3	Ethane, 1,1,2,2-tetrachloro-	166	C2H2Cl4	000079-34-5	70
4	Ethane, 1,2,2-trichloro-1,1-difluor	168	C2HCl3F2	000354-21-2	53



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Acq On : 8 Dec 2001 2:19 am
Sample : gc/ms unknown 11/6
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

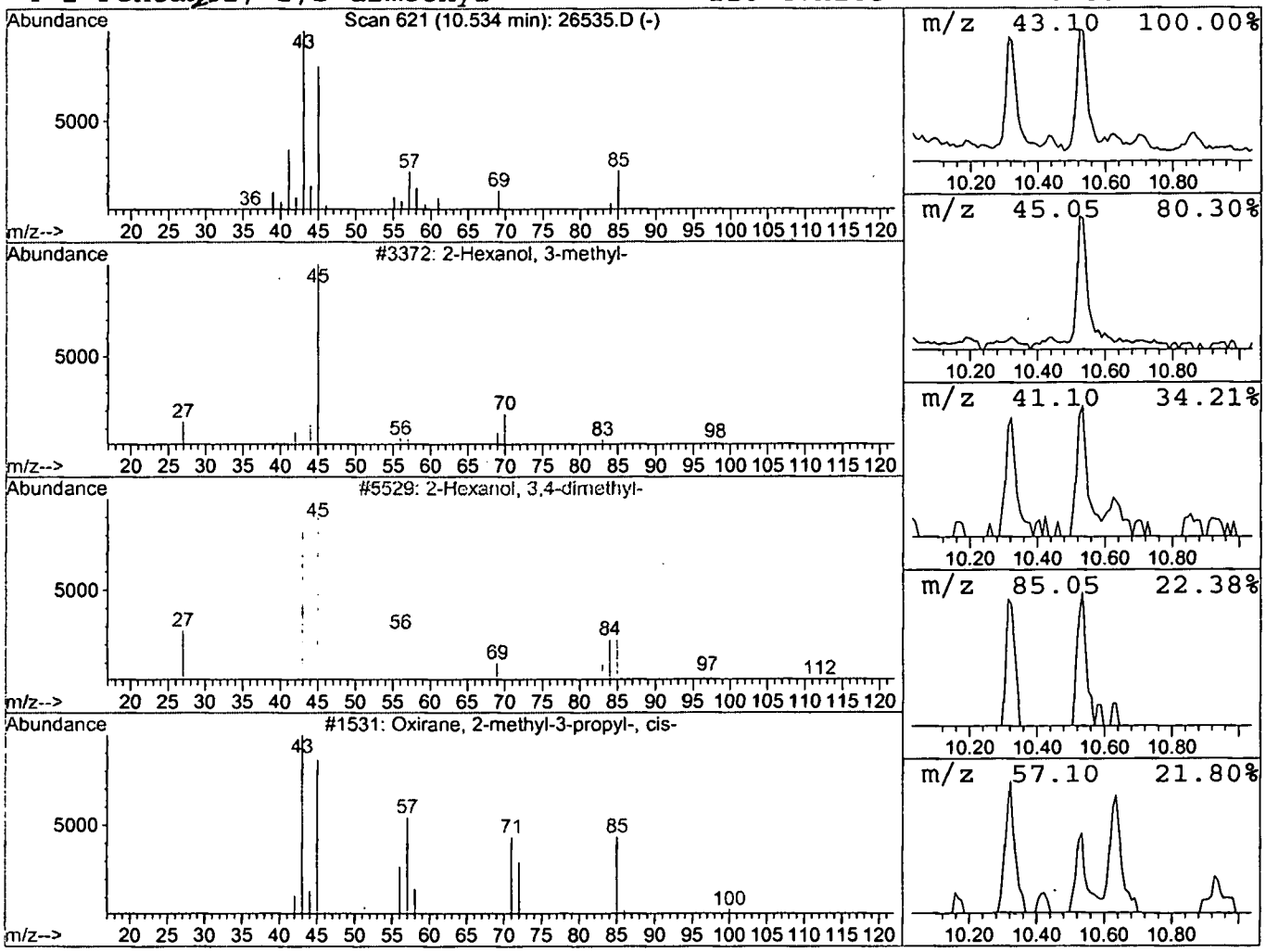
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 2-Hexanol, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	0.40 ng/uL	120030	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 3-methyl-	116	C7H16O	002313-65-7	40
2	2-Hexanol, 3,4-dimethyl-	130	C8H18O	019550-05-1	38
3	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	36
4	1-Pentanol, 2,2-dimethyl-	116	C7H16O	002370-12-9	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0701\26535.D
Acq On : 8 Dec 2001 2:19 am
Sample : gc/ms unknown 11/6
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

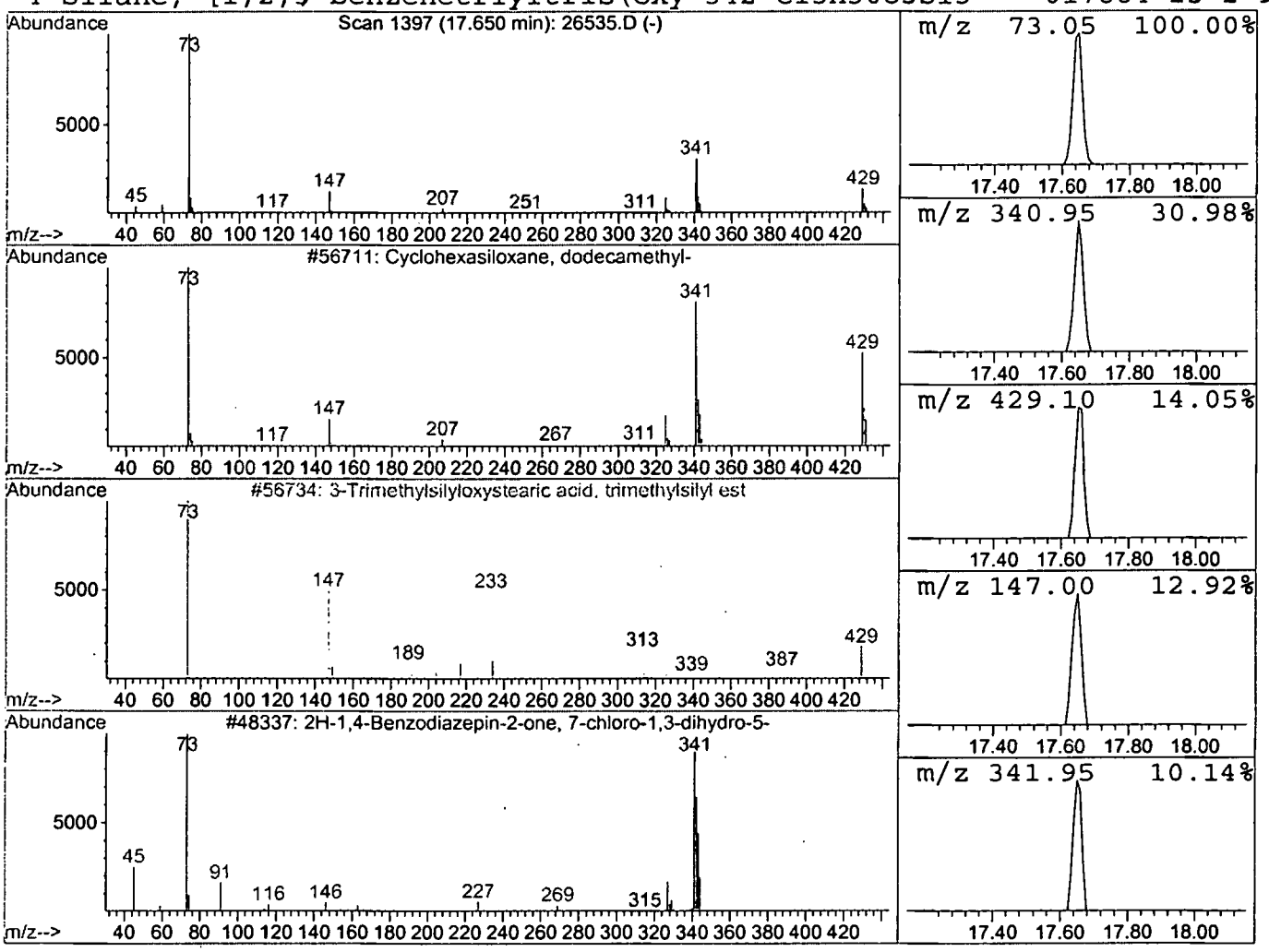
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Cyclohexasiloxane, dodecamethy Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.65	0.49 ng/uL	183034	Naphthalene-d8	15.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	78
2	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	12
3	2H-1,4-Benzodiazepin-2-one, 7-chlor	342	C18H19ClN2OSi	055299-24-6	9
4	Silane, [1,2,3-benzenetriyltris(oxy	342	C15H30O3Si3	017864-23-2	9



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 8 Dec 2001 2:19 am
Data File: C:\HPCHEM\1\DATA\DEC0701\26535.D
Name: gc/ms unknown 11/6
Misc:
Method: C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title: 208 625/TCLP calibration table
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Hydrogen azide	5.05	5.8	ng/uL	1723750	ISTD01	11.88	5975870	20.0
Ethane, 1,1-diethoxy	5.16	0.9	ng/uL	267594	ISTD01	11.88	5975870	20.0
2-Butanol, 3-methyl-	7.21	18.3	ng/uL	5457310	ISTD01	11.88	5975870	20.0
2-Pentanone, 4-hydro	7.86	9.2	ng/uL	2734890	ISTD01	11.88	5975870	20.0
Ethylamine	9.37	1.8	ng/uL	546320	ISTD01	11.88	5975870	20.0
Ethane, 1,1,2,2-tetr	9.62	3.5	ng/uL	1031010	ISTD01	11.88	5975870	20.0
2-Hexanol, 3-methyl-	10.53	0.4	ng/uL	120030	ISTD01	11.88	5975870	20.0
Cyclohexasiloxane, d	17.65	0.5	ng/uL	183034	ISTD02	15.49	7422710	20.0

26535.D NP112701.M Thu Dec 13 11:25:12 2001