User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/11/2001 Time: 16:56:40

Sample: AD26982
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/11/01 16:55 Ended: 12/11/01 16:55 Analyst: MG
Page: 1

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

Comments for Sample AD26982 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-11-2001 Time: 16:56:45

A scan of the extract prepared for the 508 analysis, from 35-550 amu using a mass spectrometer, does not reveal the presence of non-target compounds.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC0601\26982.D Vial: 24
Acq On : 7 Dec 2001 7:05 am Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:43 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	11.89	152	1347997	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	5164583	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.77	164	2352969	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.18	188	3486506	20.00	ng/uL	-0.06
59) Chrysene-d12	33.19	240	2434067	20.00	ng/uL	-0.06
68) Perylene-d12	37.28	264	1749158	20.00	ng/uL	-0.06
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	11.89	132	959	0.01	ng/uL	0.46
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.17	152	931	0.02	ng/uL	-0.29
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.04%#		
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	2554	0.02	ng/uL	0.09
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	0.00	45	0	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	15.56	128	290	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.	
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	25.18	178	348	N.D.	
56)	Anthracene	25.18	178	348	N.D.	
57)	Di-n-butylphthalate	27.18	149	6386	N.D.	
58)	Fluoranthene	0.00	202	0	N.D.	
60)	3,3'-Dichlorobenzidine	0.00	252	0	N.D.	

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

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Last Update : Wed Nov 28 15:33:44 2001
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	33.19	228	5312	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.45	149	9413	N.D.	
67)	Chrysene	33.19	228	5312	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	37.27	252	5980	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.	

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

Quantitation Report

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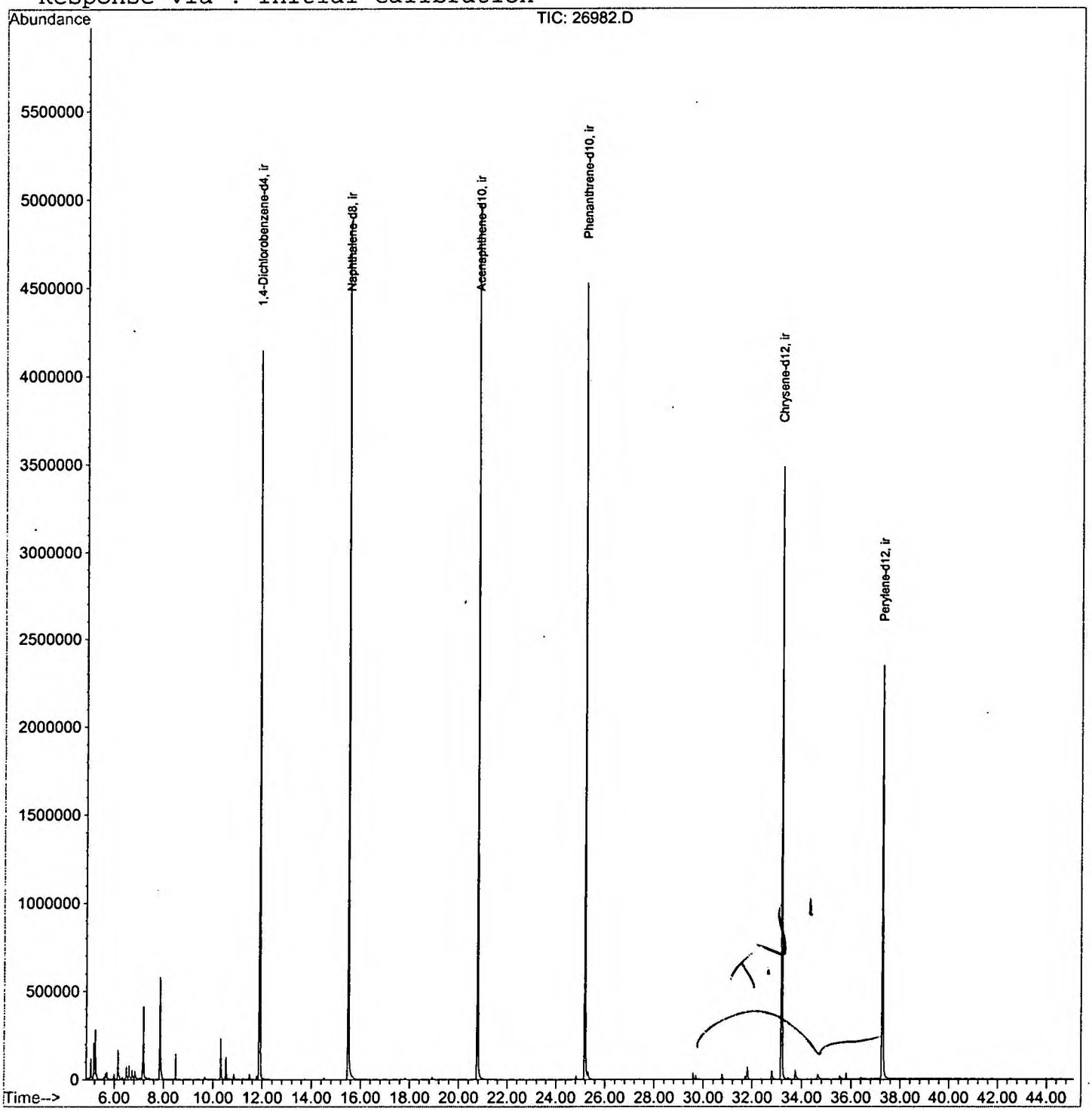
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:43 2001 Quant Results File: NP112701.RES

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\DEC0601\26982.D Vial: 24
Acq On : 7 Dec 2001 7:05 am Operator: GALOS
Sample : gc/ms unknown 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

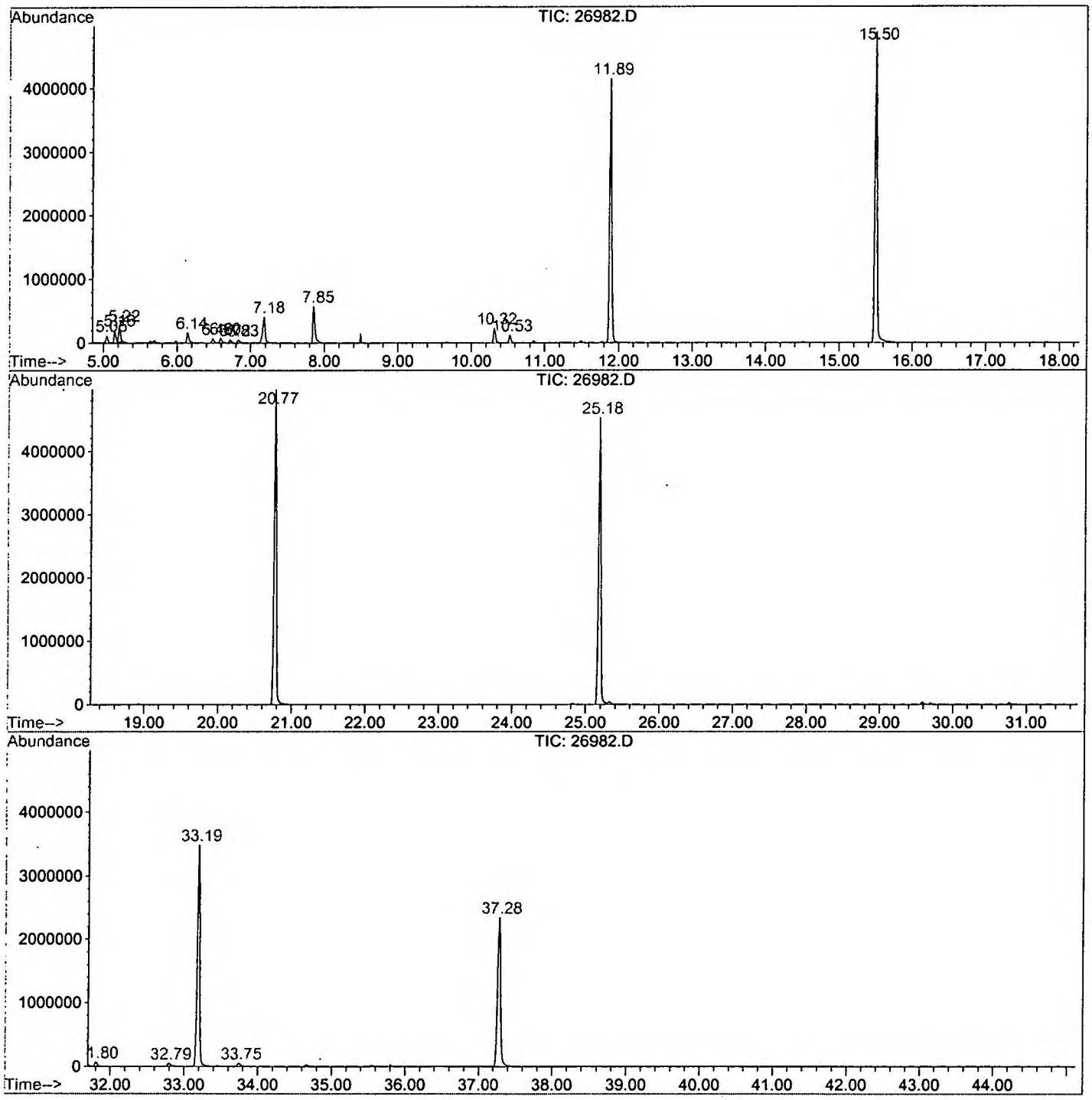
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.046	15	22	31	rVB	117155	220909	2.11%	0.378%
2	5.156	31	34	38	rBV	204743	344128	3.29%	0.588%
3	5.220	38	41	47	rVB	260280	450827	4.31%	0.771%
4	6.137	138	141	154	rVB	167009	368145	3.52%	0.629%
5	6.486	175	179	187	rBV	72300	146172	1.40%	0.250%
6	6.596	187	191	201	rVB	78112	162912	1.56%	0.278%
7	6.724	201	205	213	rBV	52298	107162	1.02%	0.183%
8	6.834	213	217	223	rBV2	45661	108042	1.03%	0.185%
9	7.183	247	255	263	rBV	410038	932184	8.91%	1.593%
10	7.852	324	328	346	rVB	578623	1244482	11.90%	2.127%
11	10.319	592	597	604	rBV	233098	449783	4.30%	0.769%
12	10.530	613	620	629	rBV	127814	248918	2.38%	0.425%
13	11.887	762	768	781	rBV	4145098	8300527	79.35%	14.188%
14	15.500	1155	1162	1179	rBV	4874156	10460441	100.00%	17.879%
15	20.773	1730	1737	1756	rBV	4976717	10412500	99.54%	17.797%
16	25.184	2211	2218	2230	rBV2	4526263	9757888	93.28%	16.679%
17	31.805	2934	2940	2950	rVB	69216	171797	1.64%	0.294%
18	32.795	3043	3048	3056	rBV2	50939	145703	1.39%	0.249%
19	33.189	3084	3091	3114	rBV2	3482387	7788973	74.46%	13.313%
20	33.749	3146	3152	3162	rBV3	51699	162627	1.55%	0.278%
21	37.279	3529	3537	3556	rBV2	2340530	6521345	62.34%	11.147%

Sum of corrected areas: 58505465

26982.D NP112701.M Tue Dec 11 14:48:37 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0601\26982.D
Operator : GALOS
Acquired : 7 Dec 2001 7:05 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 24
Quant File :NP112701.RES (RTE Integrator)



26982.D NP112701.M Tue Dec 11 14:48:39 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26982.D
Acq On : 7 Dec 2001 7:05 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 24
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

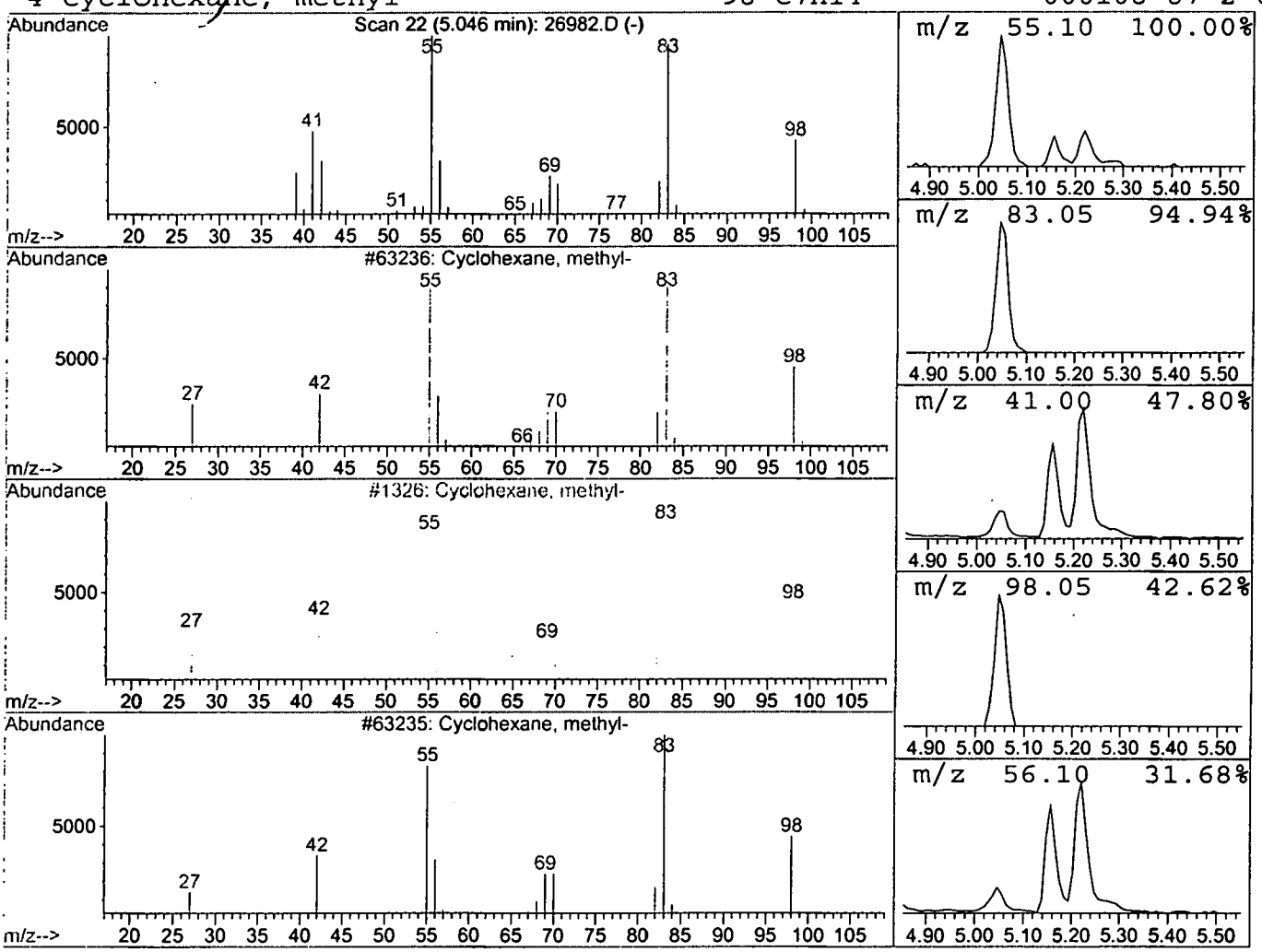
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Cyclohexane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	0.53 ng/uL	220909	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, methyl-	98	C7H14	000108-87-2	97
2	Cyclohexane, methyl-	98	C7H14	000108-87-2	95
3	Cyclohexane, methyl-	98	C7H14	000108-87-2	91
4	Cyclohexane, methyl-	98	C7H14	000108-87-2	60



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26982.D
Acq On : 7 Dec 2001 7:05 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 24
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

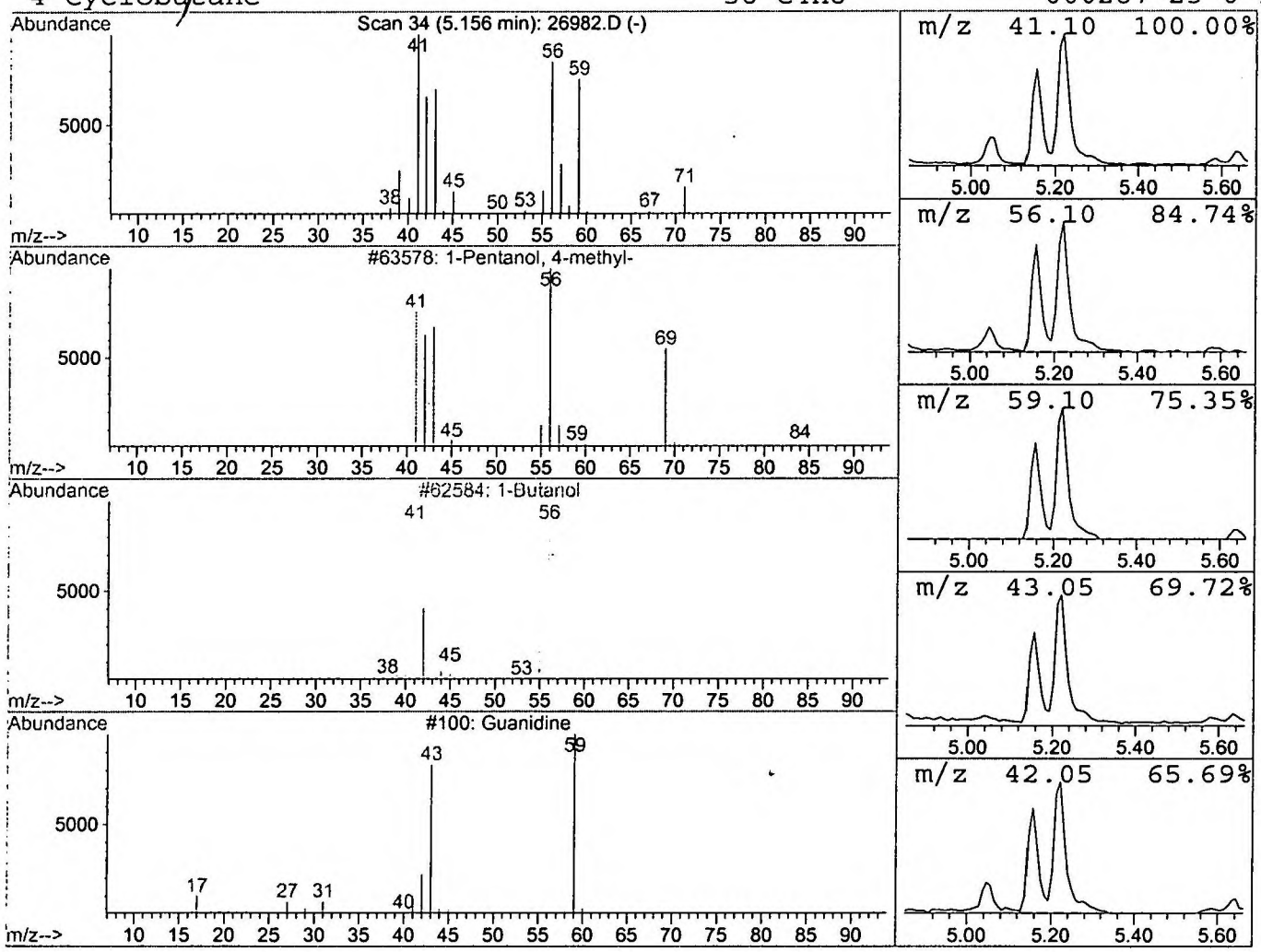
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 1-Pentanol, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	0.83 ng/uL	344128	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	28
2	1-Butanol	74	C4H10O	000071-36-3	25
3	Guanidine	59	CH5N3	000113-00-8	9
4	Cyclobutane	56	C4H8	000287-23-0	9



Library Search Compound Report

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Acq On : 7 Dec 2001 7:05 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 24
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
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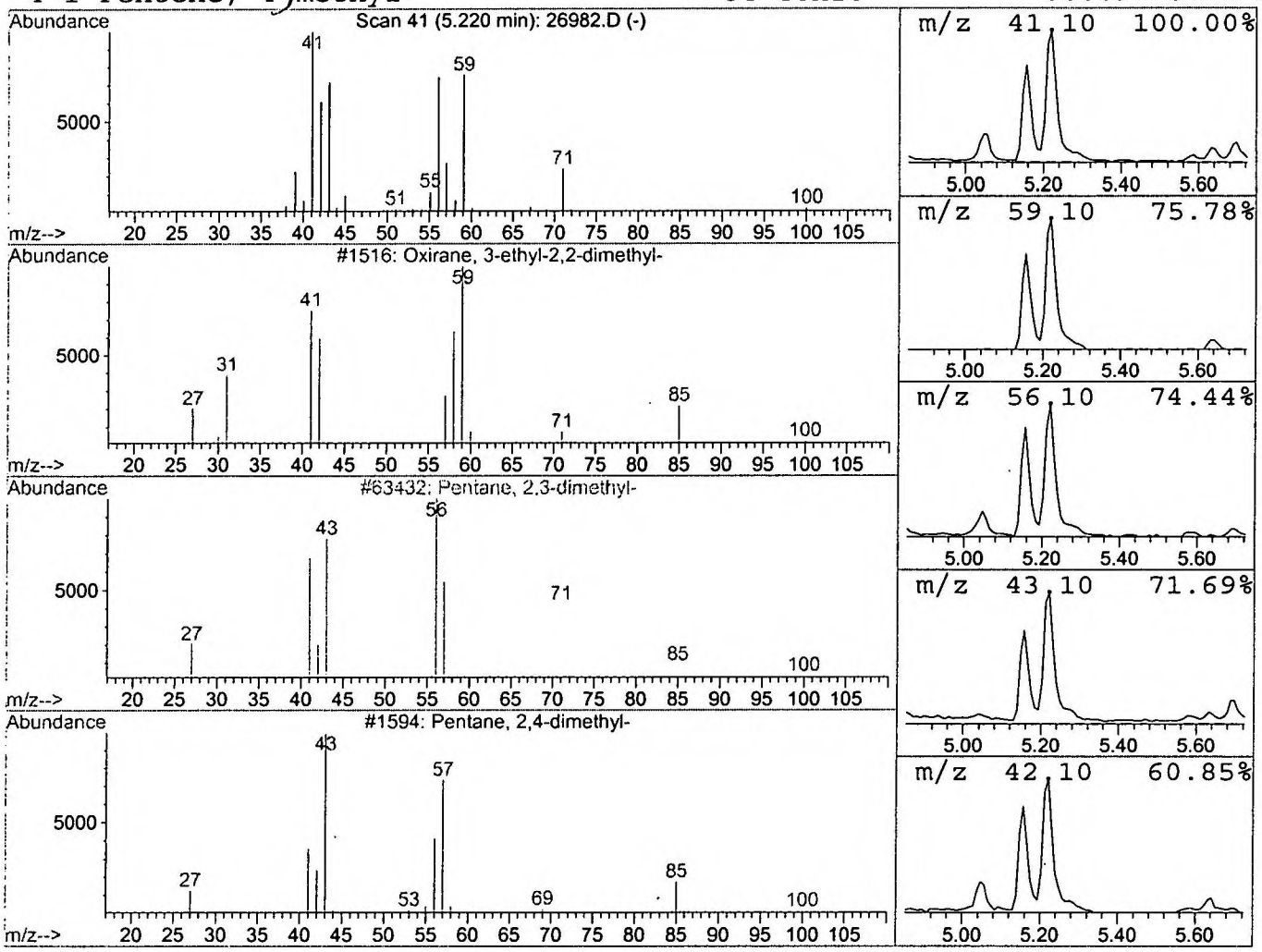
****

Peak Number 3 Oxirane, 3-ethyl-2,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	1.09 ng/uL	450827	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	38
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	38
3	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	32
4	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	32



Library Search Compound Report

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Acq On : 7 Dec 2001 7:05 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 24
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

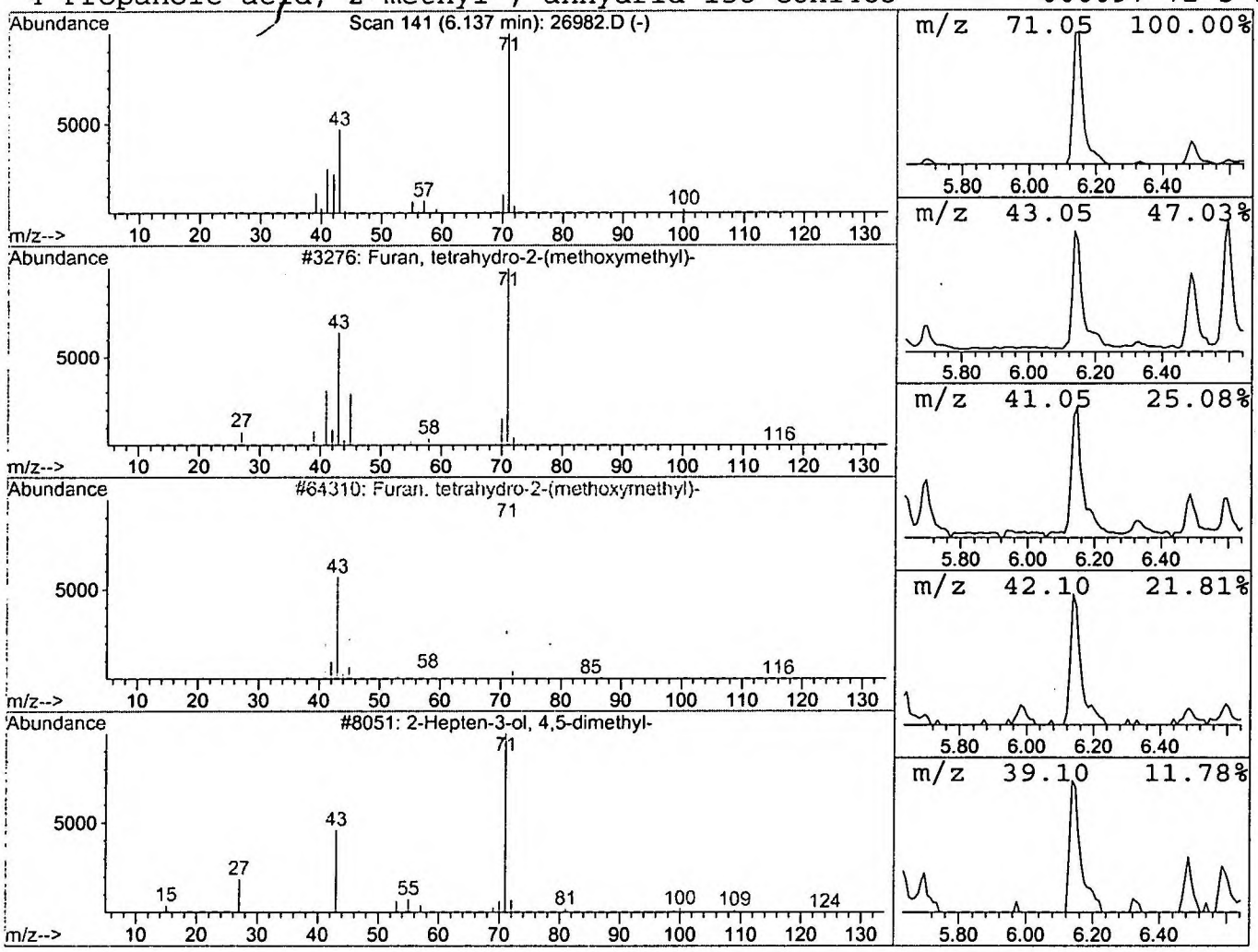
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Furan, tetrahydro-2-(methoxyme Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	0.89 ng/uL	368145	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, tetrahydro-2-(methoxymethyl)	116	C6H12O2	019354-27-9	59
2	Furan, tetrahydro-2-(methoxymethyl)	116	C6H12O2	019354-27-9	45
3	2-Hepten-3-ol, 4,5-dimethyl-	142	C9H18O	055956-37-1	38
4	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	36



Library Search Compound Report

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Acq On : 7 Dec 2001 7:05 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 24
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

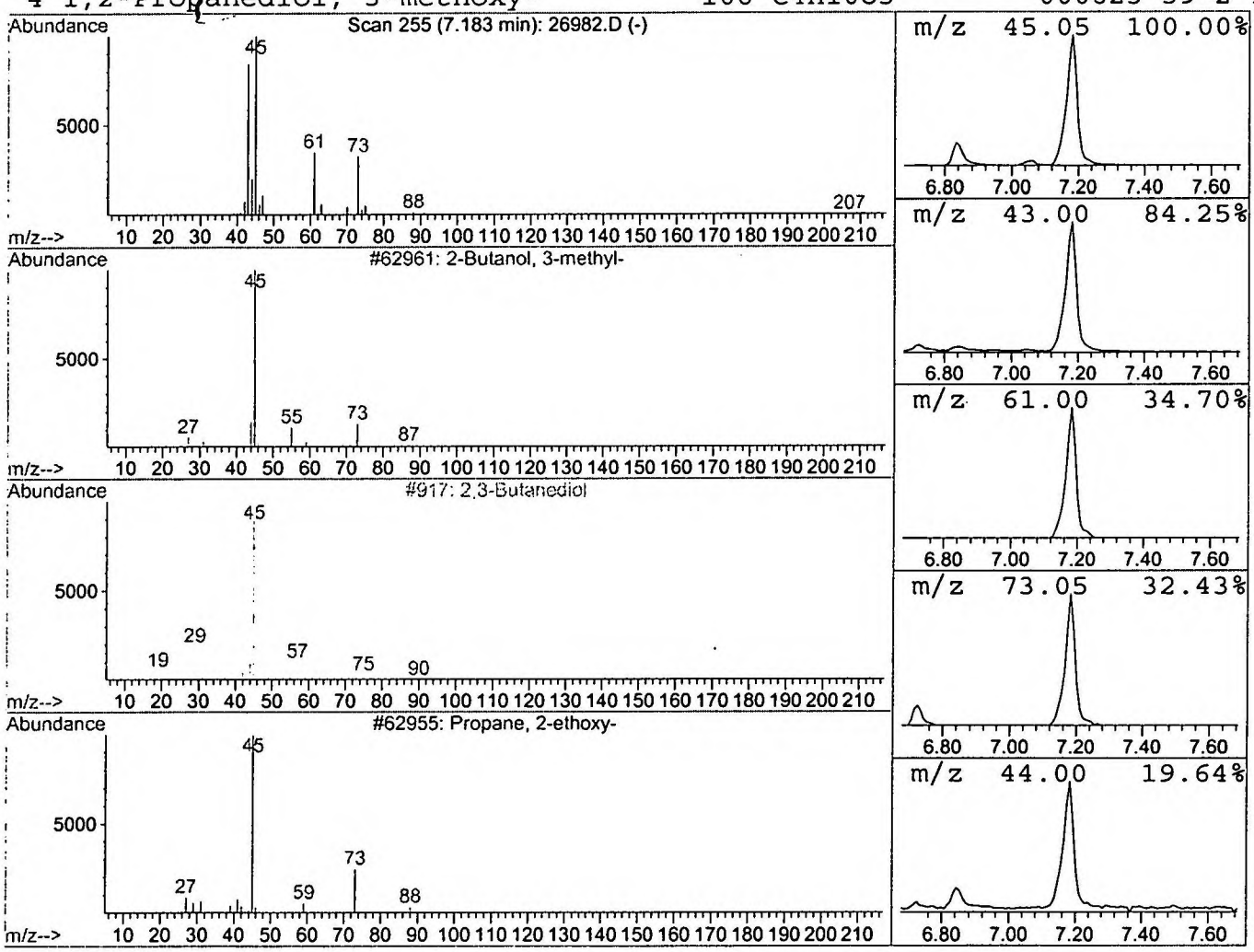
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 2-Butanol, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	2.25 ng/uL	932184	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
2	2,3-Butanediol	90	C4H10O2	000513-85-9	9
3	Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9
4	1,2-Propanediol, 3-methoxy-	106	C4H10O3	000623-39-2	9



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Acq On : 7 Dec 2001 7:05 am Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

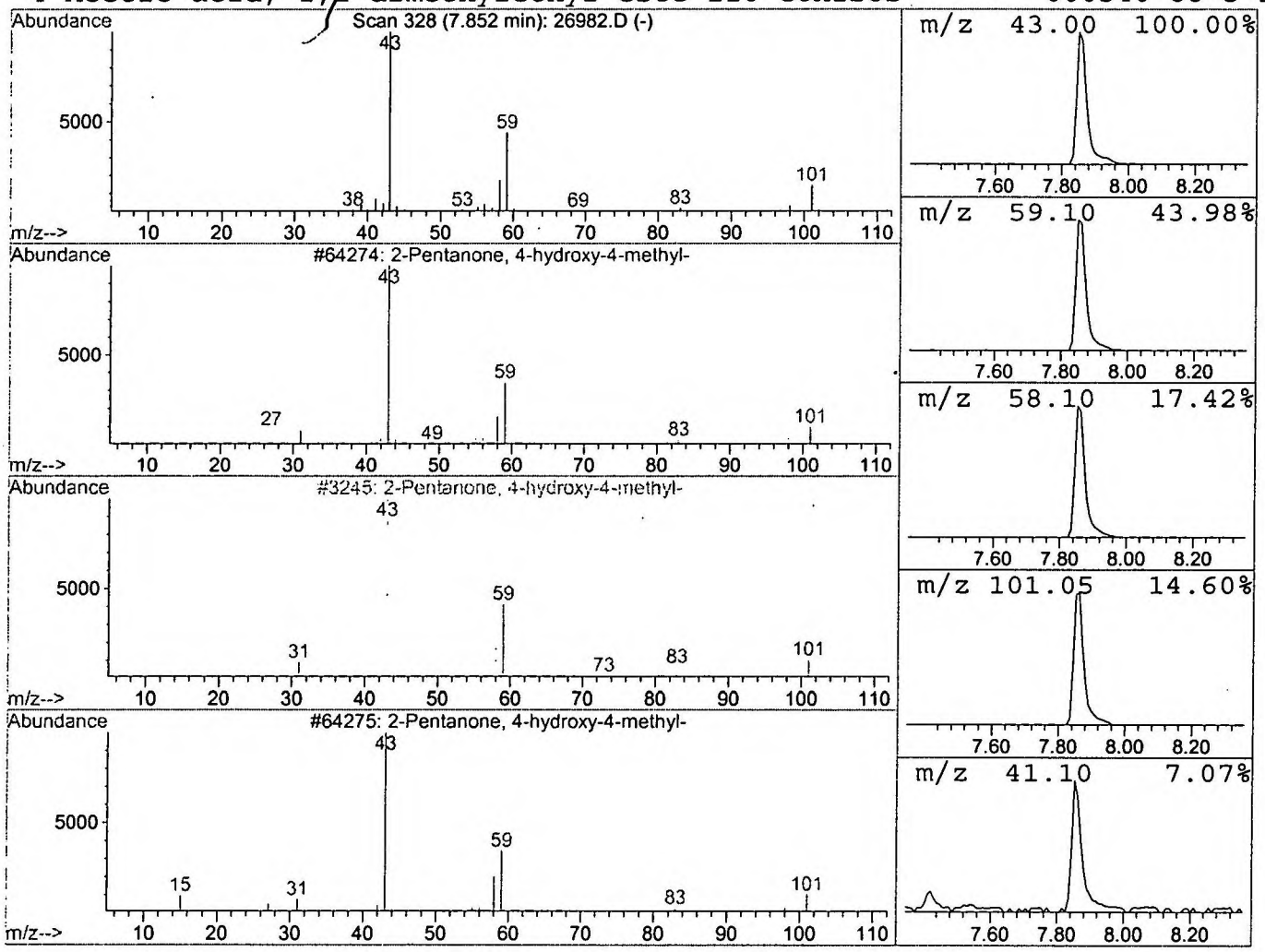
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	3.00 ng/uL	1244480	1,4-Dichlorobenzene-d4	11.89

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	64
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
4	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28



Library Search Compound Report

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Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

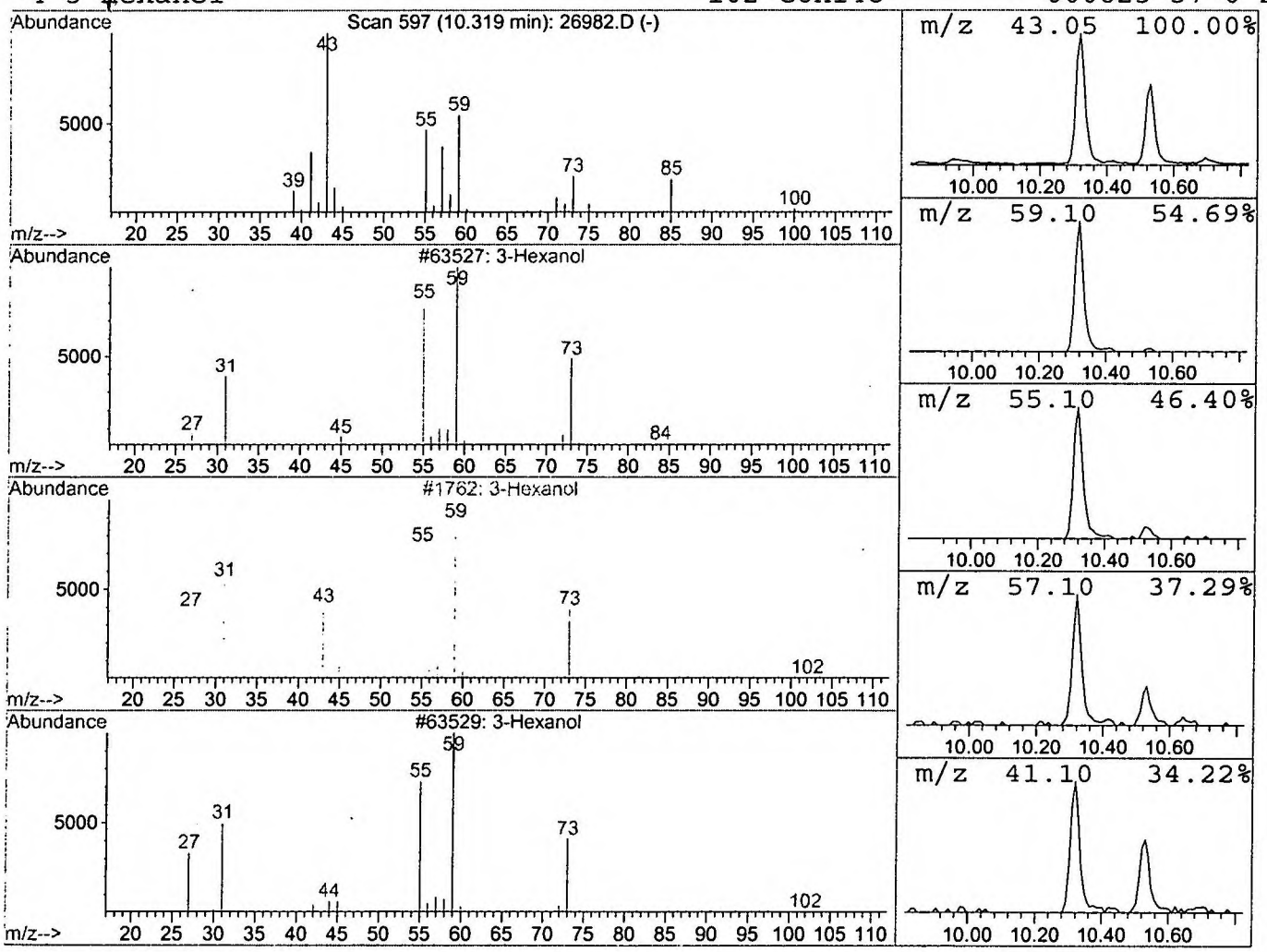
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 3-Hexanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	1.08 ng/uL	449783	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	43
2	3-Hexanol	102	C6H14O	000623-37-0	38
3	3-Hexanol	102	C6H14O	000623-37-0	37
4	3-Hexanol	102	C6H14O	000623-37-0	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26982.D
Acq On : 7 Dec 2001 7:05 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 24
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

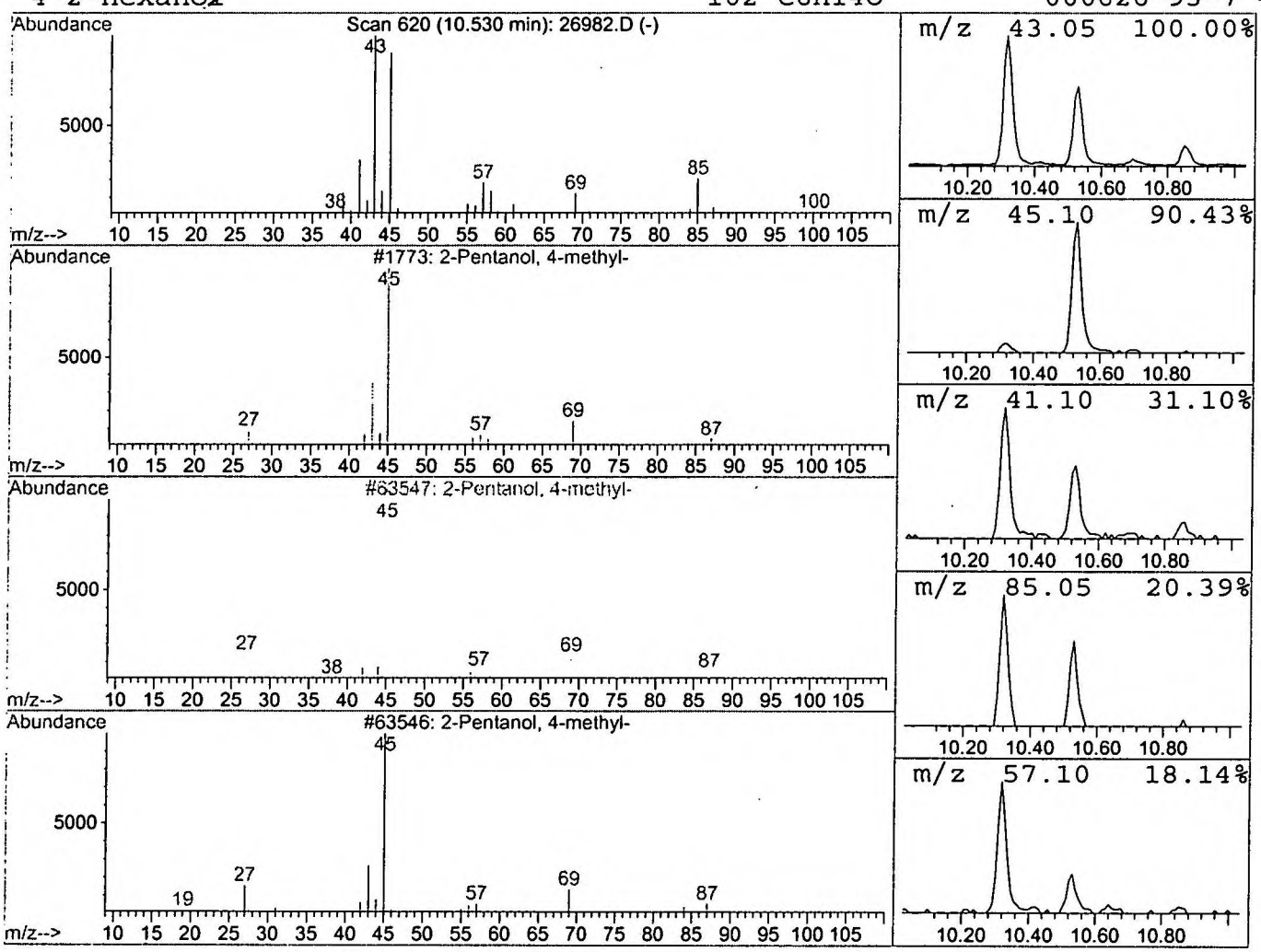
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 2-Pentanol, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	0.60 ng/uL	248918	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	53
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	53
3	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	50
4	2-Hexanol	102	C6H14O	000626-93-7	42



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26982.D
Acq On : 7 Dec 2001 7:05 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 24
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

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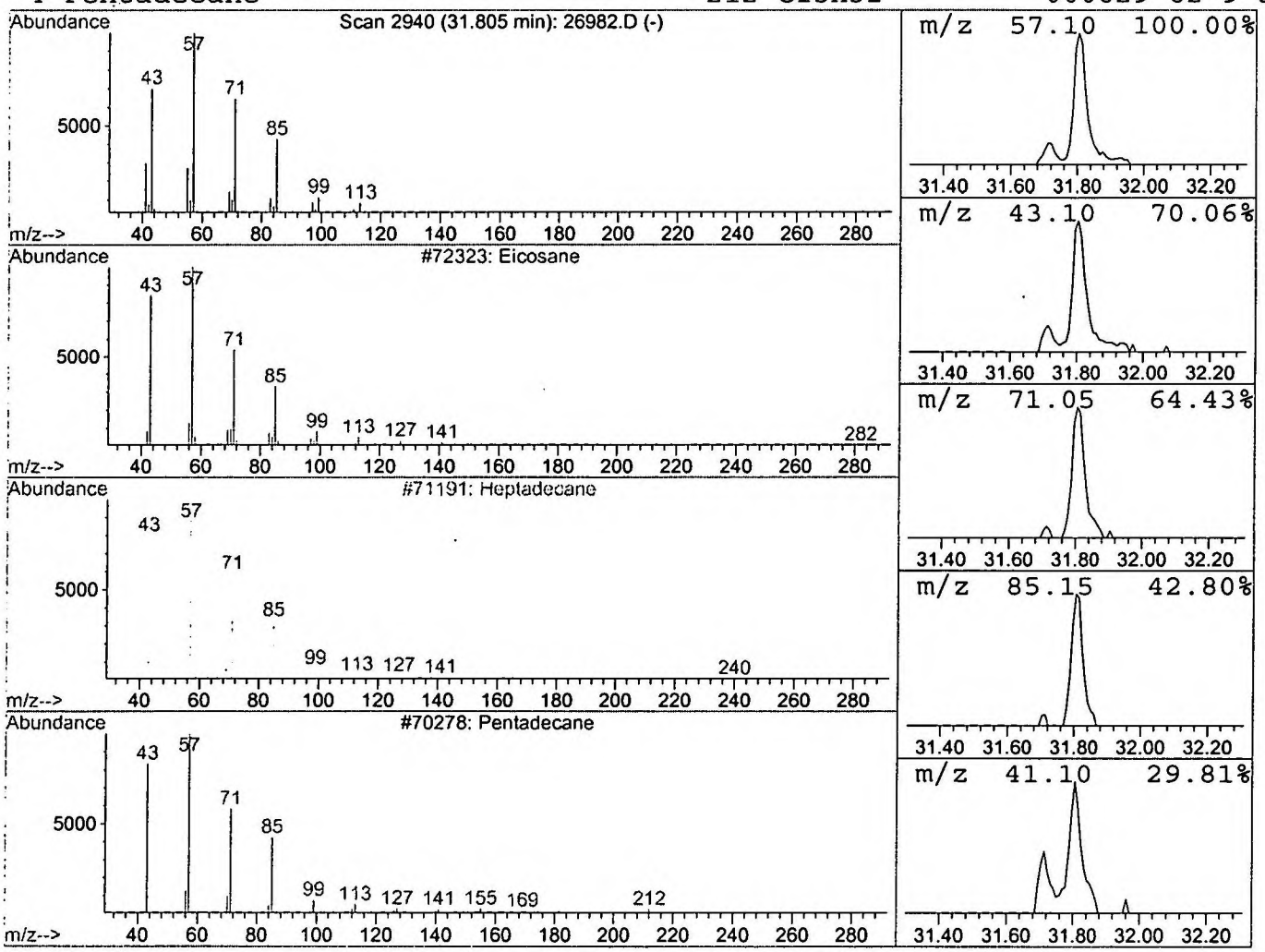
Peak Number 9 Eicosane — Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.80	0.44 ng/uL	171797	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	64
2	Heptadecane	240	C17H36	000629-78-7	64
3	Pentadecane	212	C15H32	000629-62-9	64
4	Pentadecane	212	C15H32	000629-62-9	56



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26982.D Vial: 24
Acq On : 7 Dec 2001 7:05 am Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

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

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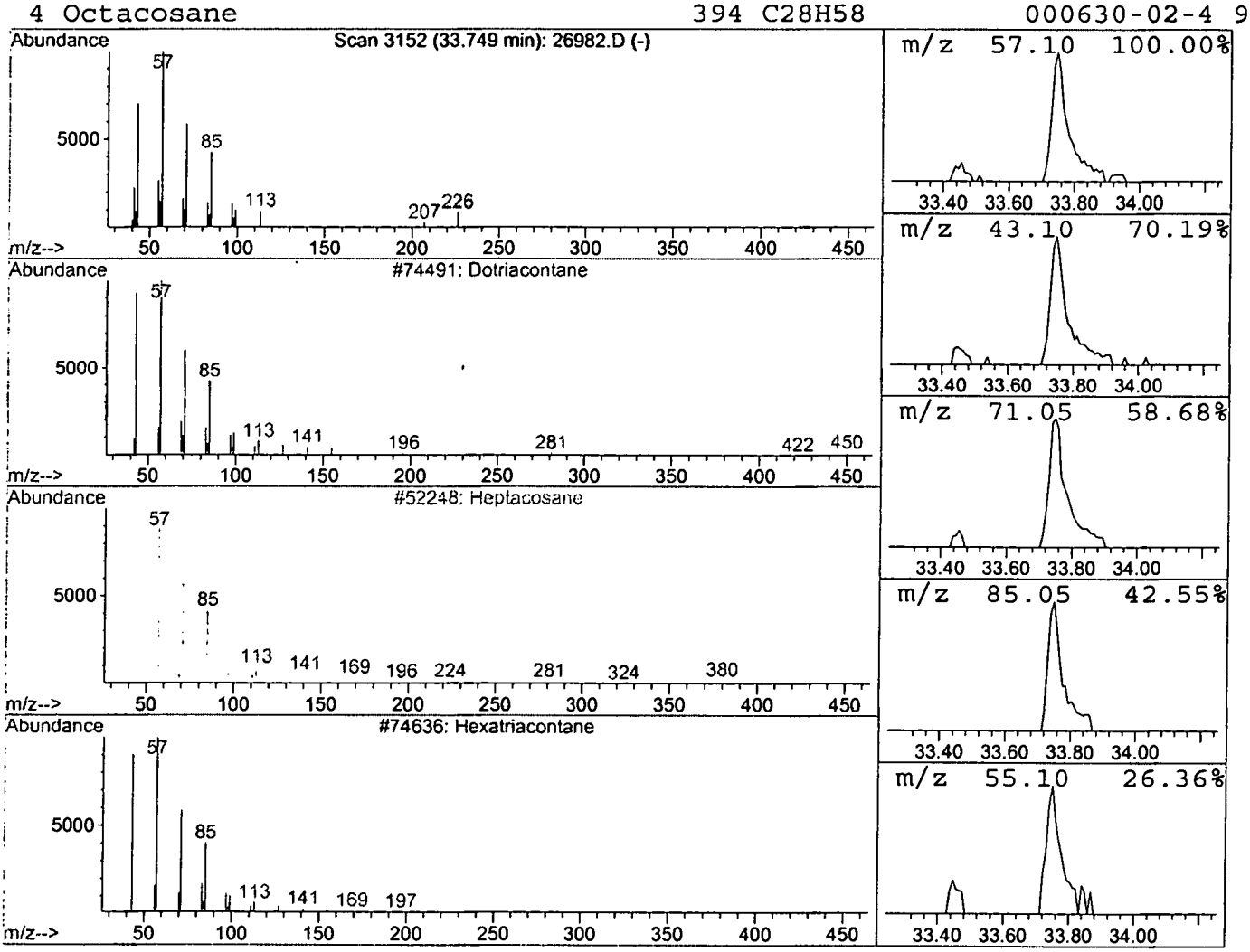
Peak Number 10 Dotriacontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.75	0.42 ng/uL	162627	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dotriacontane	451	C32H66	000544-85-4	91
2	Heptacosane	380	C27H56	000593-49-7	90
3	Hexatriacontane	507	C36H74	000630-06-8	90
4	Octacosane	394	C28H58	000630-02-4	90



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 7 Dec 2001 7:05 am
Data File: C:\HPCHEM\1\DATA\DEC0601\26982.D
Name: gc/ms unknown
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
Cyclohexane, methyl-	5.05	0.5	ng/uL	220909	ISTD01	11.89	8300530	20.0
1-Pentanol, 4-methyl	5.16	0.8	ng/uL	344128	ISTD01	11.89	8300530	20.0
Oxirane, 3-ethyl-2,2	5.22	1.1	ng/uL	450827	ISTD01	11.89	8300530	20.0
Furan, tetrahydro-2-	6.14	0.9	ng/uL	368145	ISTD01	11.89	8300530	20.0
2-Butanol, 3-methyl-	7.18	2.2	ng/uL	932184	ISTD01	11.89	8300530	20.0
2-Pentanone, 4-hydro	7.85	3.0	ng/uL	1244480	ISTD01	11.89	8300530	20.0
3-Hexanol	10.32	1.1	ng/uL	449783	ISTD01	11.89	8300530	20.0
2-Pentanol, 4-methyl	10.53	0.6	ng/uL	248918	ISTD01	11.89	8300530	20.0
Eicosane	31.80	0.4	ng/uL	171797	ISTD05	33.19	7788970	20.0
Dotriacontane	33.75	0.4	ng/uL	162627	ISTD05	33.19	7788970	20.0

26982.D NP112701.M Tue Dec 11 14:49:02 2001