Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D Vial: 3
Acq On : 5 Dec 2001 10:35 am Operator: GALOS
Sample : gc/ms unknown 11/7 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 7 9:55 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	11.86	152	1131948	20.00	ng/uL	-0.08
19) Naphthalene-d8	15.47	136	4229441	20.00	ng/uL	-0.08
31) Acenaphthene-d10	20.75	164	1885261	20.00	ng/uL	-0.08
49) Phenanthrene-d10	25.17	188	2548107	20.00	ng/uL	-0.08
59) Chrysene-d12	33.16	240	1408776	20.00	ng/uL	-0.09
68) Perylene-d12	37.25	264	906689	20.00	ng/uL	-0.09
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	11.19	99	303	0.00	ng/uL	0.06
Spiked Amount 75.000	Range 19 - 32		Recovery =	0.00%#		
6) 2-Chlorophenol-d4	11.87	132	1023	0.01	ng/uL	0.45
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.16	152	557	0.01	ng/uL	-0.30
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.02%#		
20) Nitrobenzene-d5	0.00	82	0	0.00	ng/uL	
Spiked Amount 50.000	Range 55 - 98		Recovery =	0.00%#		
32) 2-Fluorobiphenyl	18.89	172	1602	0.01	ng/uL	0.06
Spiked Amount 50.000	Range 42 - 78		Recovery =	0.02%#		
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	12.86	45	1603	N.D.		
16) 3&4-Methylphenol	0.00	108	0	N.D.		
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.		

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

Quantitation Report (QT Reviewed)

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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

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
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	0.00	128	0	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	0.00	178	0	N.D.	
56) Anthracene	0.00	178	0	N.D.	
57) Di-n-butylphthalate	27.16	149	23620	N.D.	
58) Fluoranthene	0.00	202	0	N.D.	
60) 3,3'-Dichlorobenzidine	0.00	252	0	N.D.	
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	33.16	228	2755	N.D.	
66) Bis(2-ethylhexyl)phthalate	33.43	149	27822	0.34 ng/uL	98
67) Chrysene	33.16	228	2755	N.D.	
69) Di-n-Octylphthalate	0.00	149	0	N.D.	
70) Benzo(b)fluoranthene	0.00	252	0	N.D.	

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

Quantitation Report (QT Reviewed)

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Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 7 9:55 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

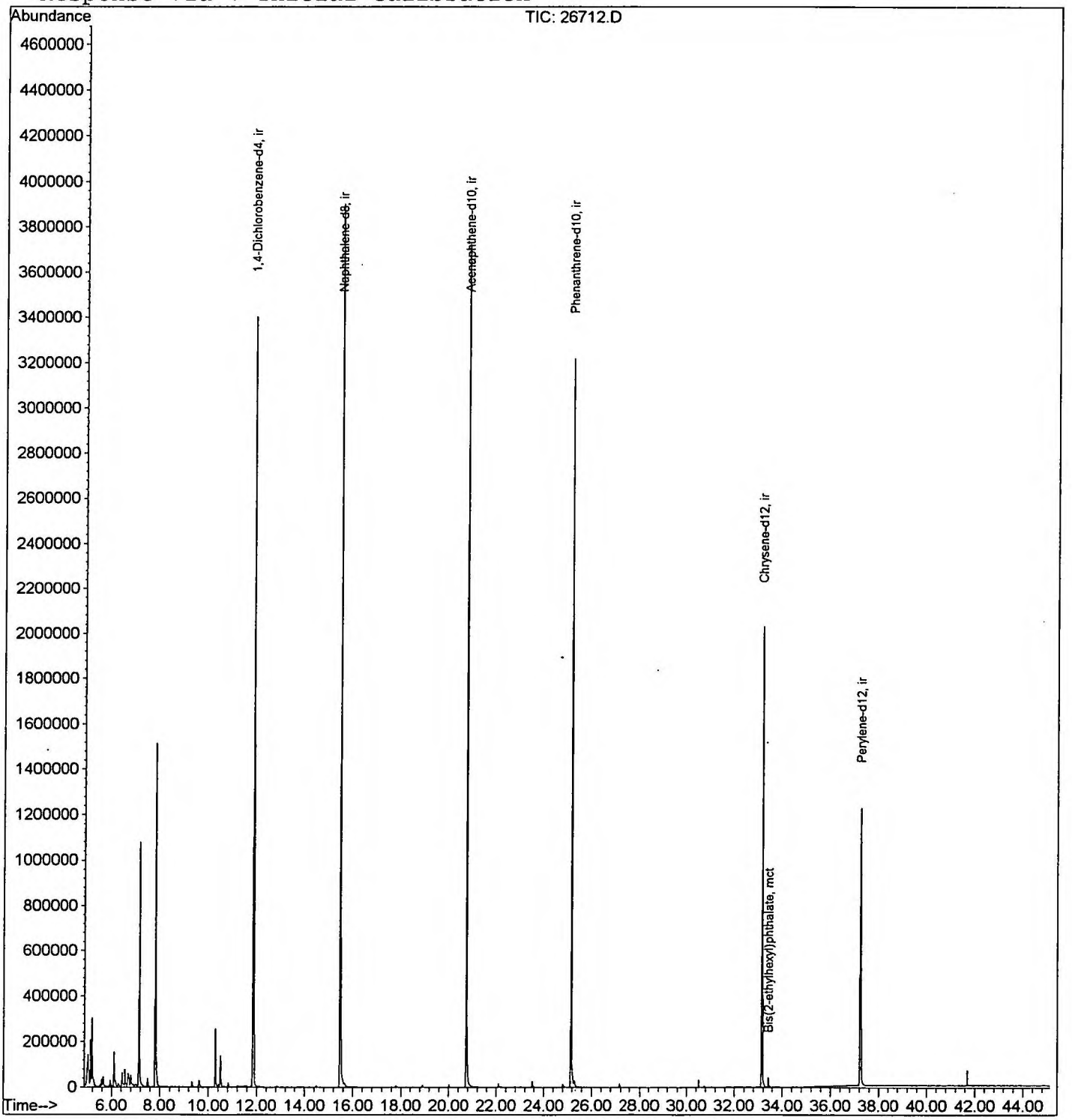
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71) Benzo(k)fluoranthene	0.00	252	0	N.D.	
72) Benzo(a)pyrene	37.25	252	2993	N.D.	
73) Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
74) Dibenz(a,h)anthracene	0.00	278	0	N.D.	
75) Benzo(g,h,i)perylene	0.00	276	0	N.D.	

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

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D Vial: 3
Acq On : 5 Dec 2001 10:35 am Operator: GALOS
Sample : gc/ms unknown 11/7 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 7 9:55 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\DEC0501\26712.D Vial: 3
Acq On : 5 Dec 2001 10:35 am Operator: GALOS
Sample : gc/ms unknown 11/7 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

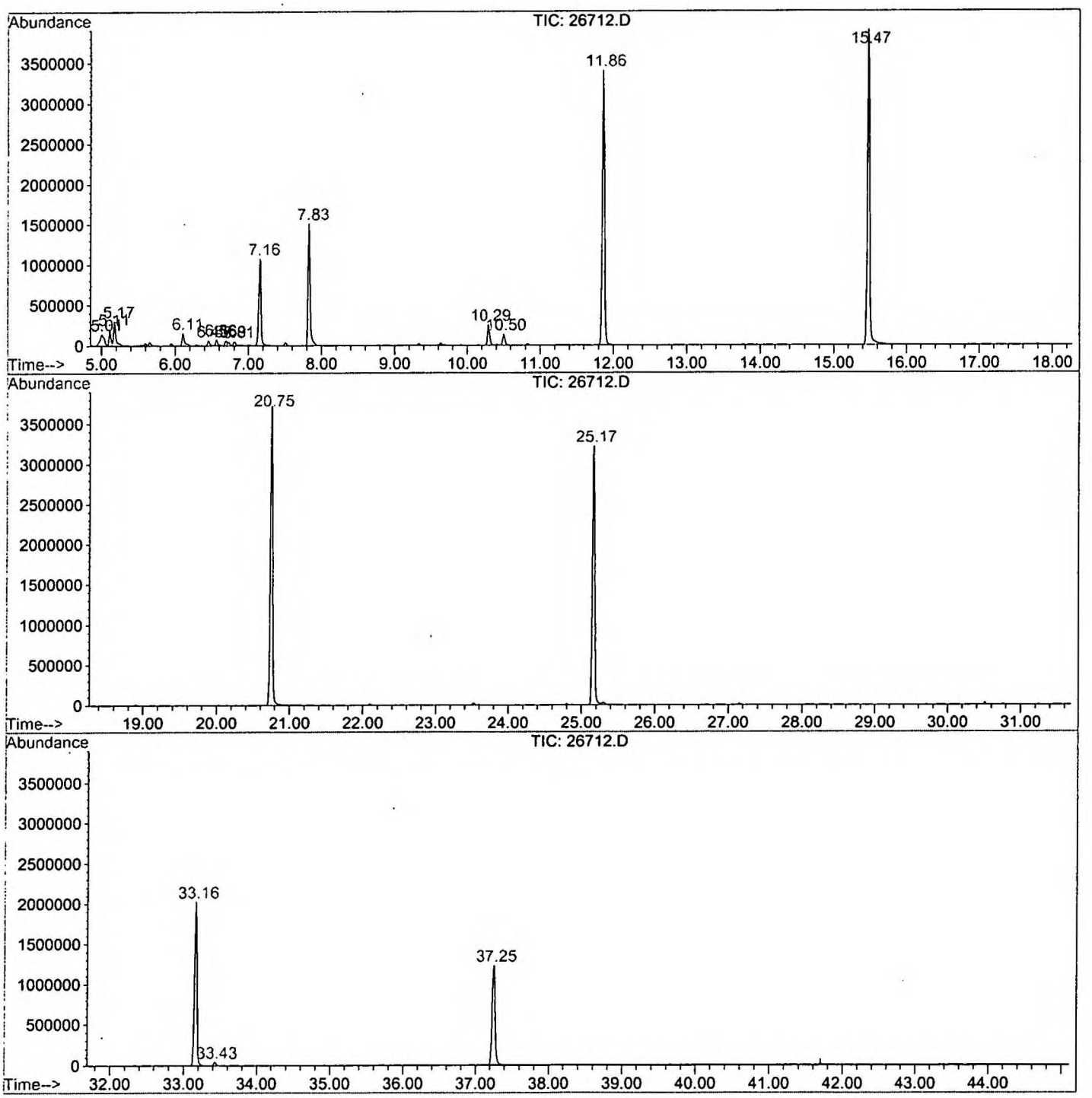
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.010	7	18	26	rBV2	135619	552784	6.47%	1.170%
2	5.111	26	29	33	rVV	203057	354684	4.15%	0.751%
3	5.175	33	36	53	rVB	300425	656290	7.68%	1.389%
4	6.110	134	138	150	rVB	153333	373184	4.37%	0.790%
5	6.450	171	175	184	rBV	60407	136893	1.60%	0.290%
6	6.560	184	187	191	rBV	71682	123644	1.45%	0.262%
7	6.688	197	201	203	rBV	58186	109413	1.28%	0.232%
8	6.807	210	214	219	rBV2	45450	96182	1.13%	0.204%
9	7.156	244	252	263	rBV	1072003	2288021	26.77%	4.842%
10	7.825	321	325	338	rBV	1511482	2861959	33.49%	6.057%
11	10.292	590	594	602	rBV	253543	513104	6.00%	1.086%
12	10.503	613	617	627	rBV2	136242	283922	3.32%	0.601%
13	11.860	759	765	782	rVB	3400770	6992465	81.81%	14.798%
14	15.473	1153	1159	1177	rBV	3900040	8546900	100.00%	18.088%
15	20.746	1728	1734	1752	rBV	3722602	8315576	97.29%	17.598%
16	25.166	2209	2216	2227	rBV2	3219215	7160404	83.78%	15.153%
17	33.162	3082	3088	3101	rBV2	2029317	4483585	52.46%	9.489%
18	33.428	3112	3117	3125	rVB	41403	95649	1.12%	0.202%
19	37.252	3526	3534	3547	rBV2	1222164	3308072	38.70%	7.001%

Sum of corrected areas: 47252731

26712.D NP112701.M Fri Dec 07 13:15:01 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0501\26712.D
Operator : GALOS
Acquired : 5 Dec 2001 10:35 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 11/7
Misc Info :
Vial Number: 3
Quant File :NP112701.RES (RTE Integrator)



26712.D NP112701.M Fri Dec 07 13:15:03 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D
Acq On : 5 Dec 2001 10:35 am
Sample : gc/ms unknown 11/7
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

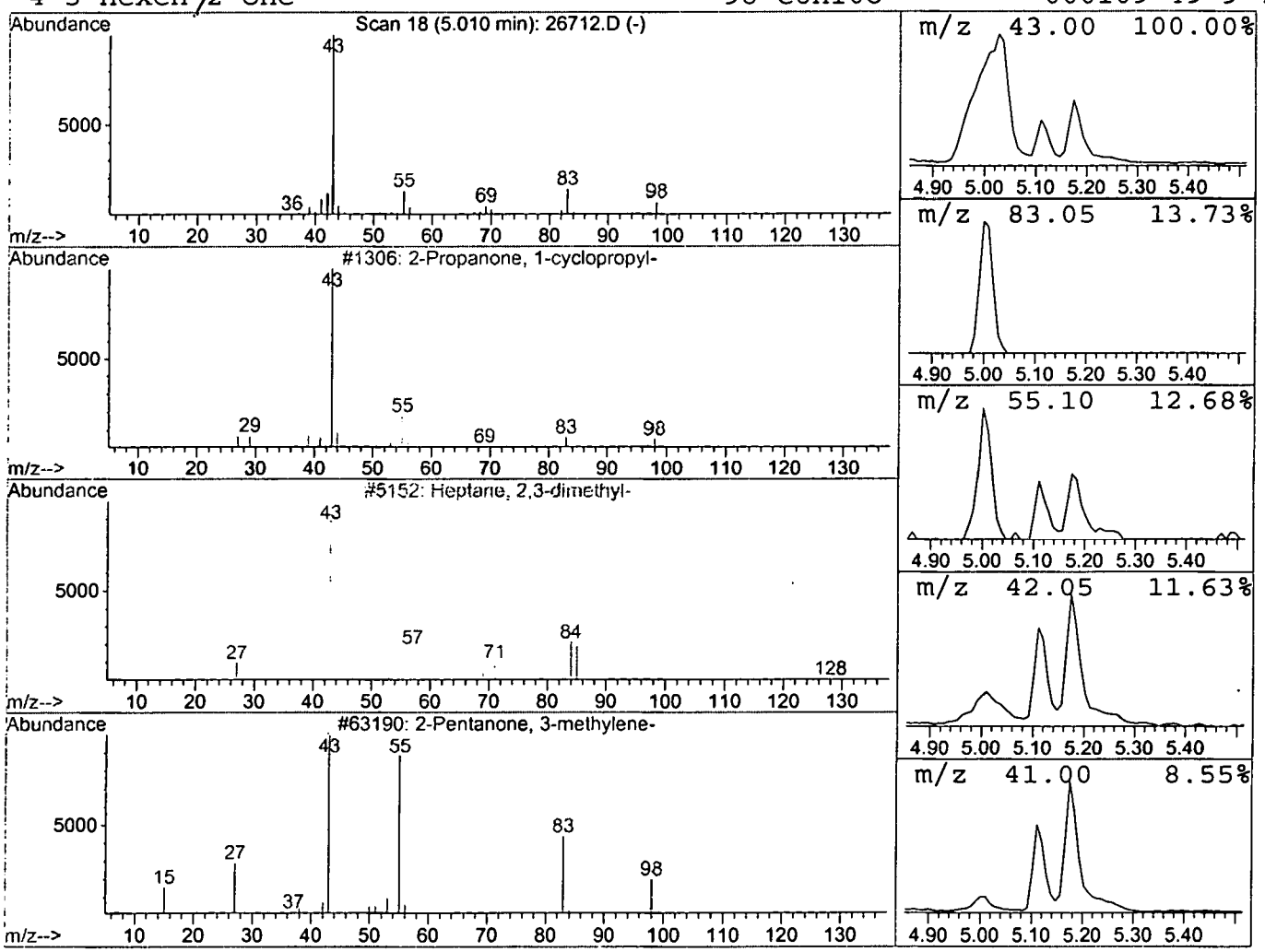
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 2-Propanone, 1-cyclopropyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	1.58 ng/uL	552784	1,4-Dichlorobenzene-d4	11.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	9
2	Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	9
3	2-Pentanone, 3-methylene-	98	C6H10O	004359-77-7	7
4	5-Hexen-2-one	98	C6H10O	000109-49-9	7



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D
Acq On : 5 Dec 2001 10:35 am
Sample : gc/ms unknown 11/7
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

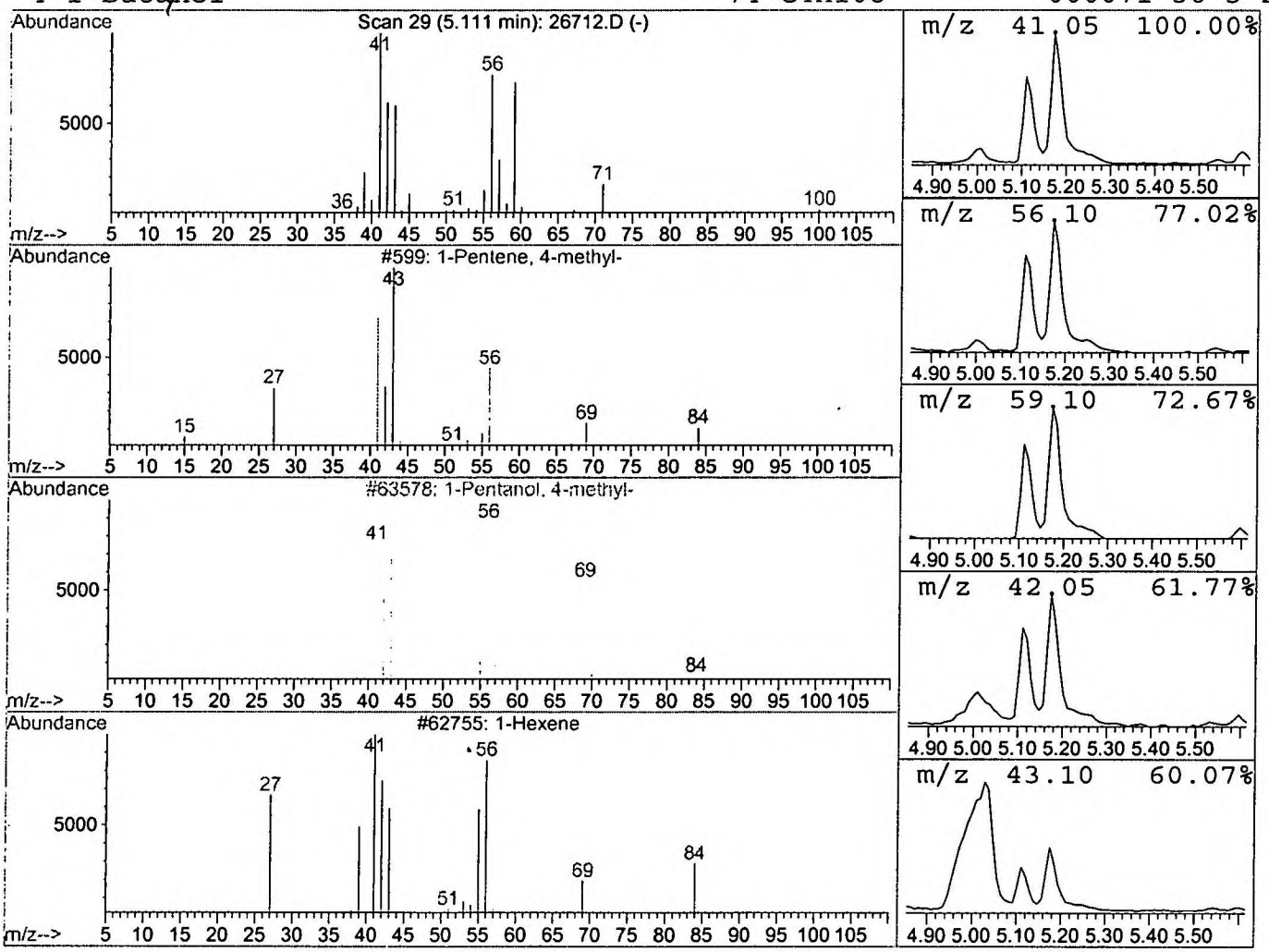
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 1-Pentene, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	1.01 ng/uL	354684	1,4-Dichlorobenzene-d4	11.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	33
2	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	33
3	1-Hexene	84	C6H12	000592-41-6	33
4	1-Butanol	74	C4H10O	000071-36-3	28



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D
Acq On : 5 Dec 2001 10:35 am
Sample : gc/ms unknown 11/7
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

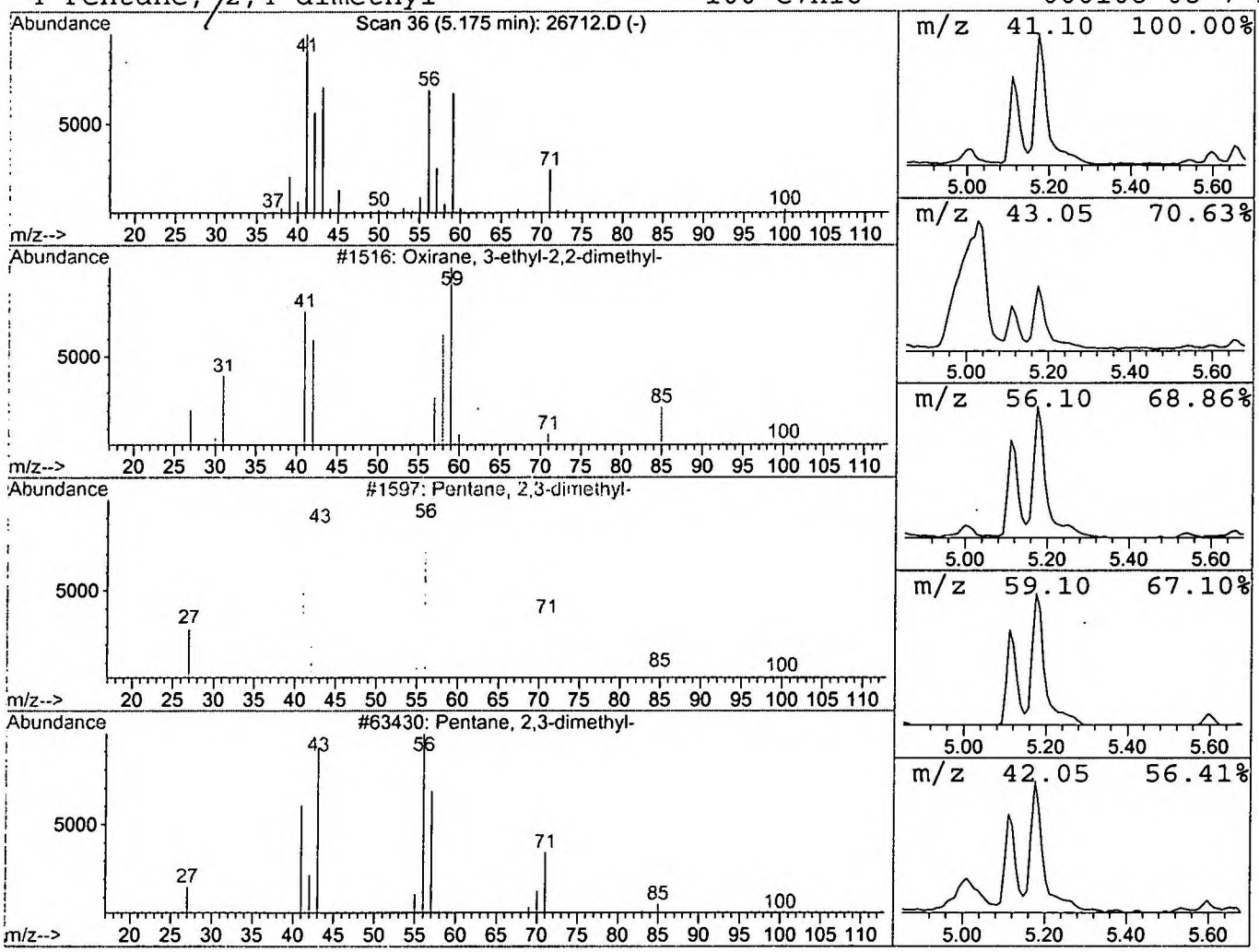
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	1.88 ng/uL	656290	1,4-Dichlorobenzene-d4	11.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	43
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35
4	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	32



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D Vial: 3
Acq On : 5 Dec 2001 10:35 am Operator: GALOS
Sample : gc/ms unknown 11/7 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

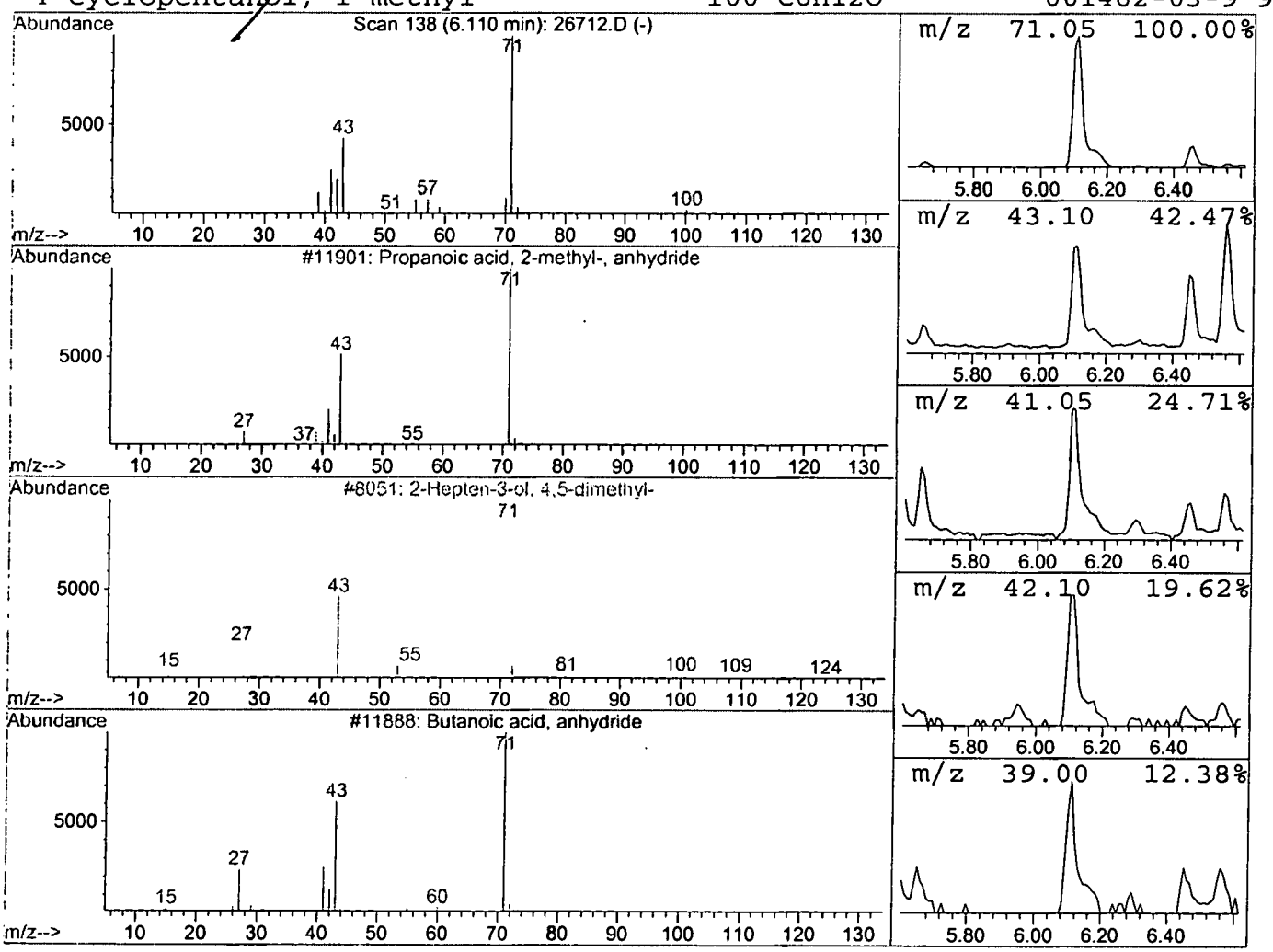
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Propanoic acid, 2-methyl-, anh Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.11	1.07 ng/uL	373184	1,4-Dichlorobenzene-d4	11.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	50
2	2-Hepten-3-ol, 4,5-dimethyl-	142	C9H18O	055956-37-1	38
3	Butanoic acid, anhydride	158	C8H14O3	000106-31-0	36
4	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	9



Library Search Compound Report

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Acq On : 5 Dec 2001 10:35 am Operator: GALOS
Sample : gc/ms unknown 11/7 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

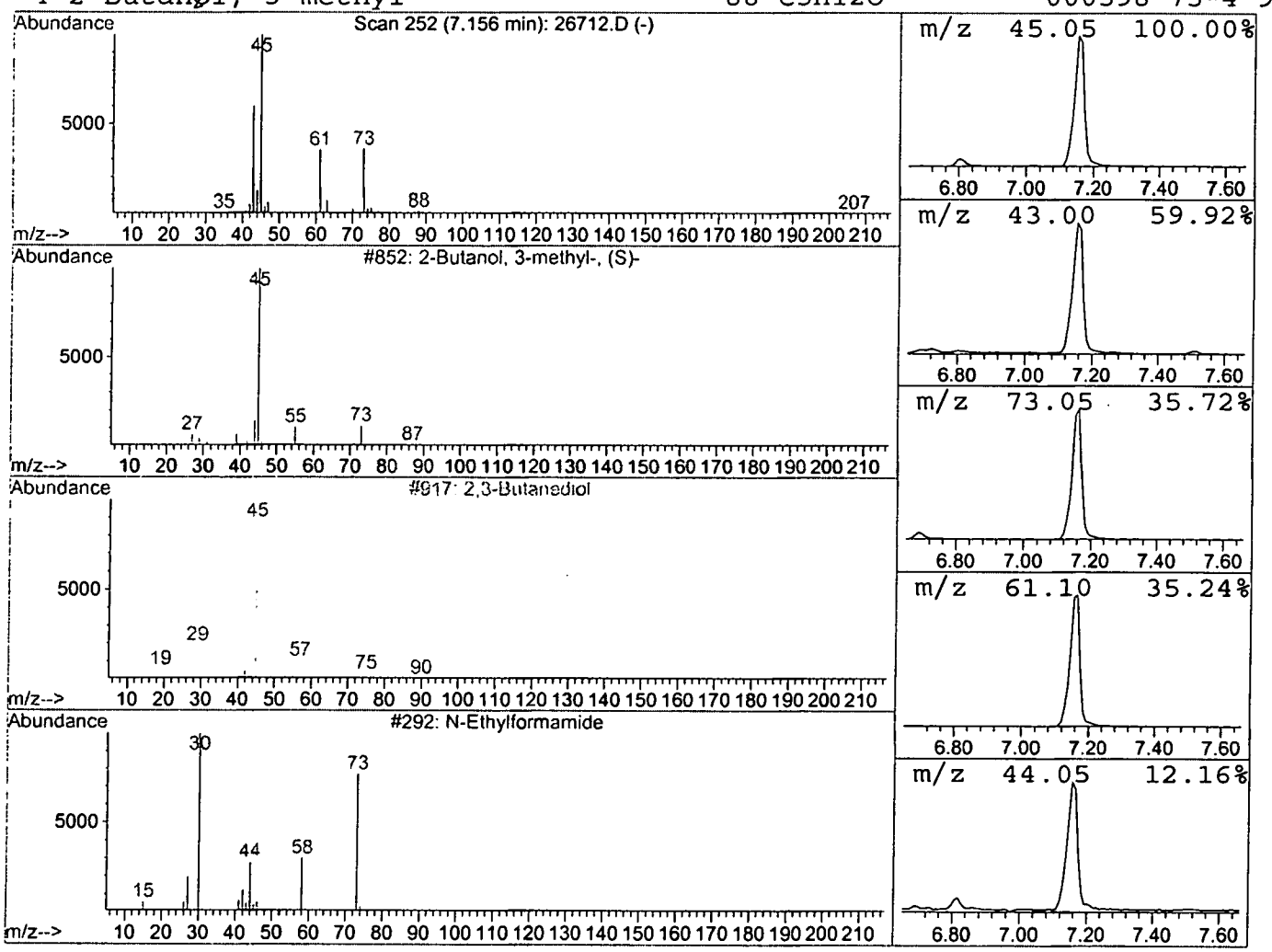
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 2-Butanol, 3-methyl-, (S)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	6.54 ng/uL	2288020	1,4-Dichlorobenzene-d4	11.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 3-methyl-, (S)-	88	C5H12O	001517-66-4	42
2	2,3-Butanediol	90	C4H10O2	000513-85-9	33
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9



Library Search Compound Report

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Acq On : 5 Dec 2001 10:35 am Operator: GALOS
Sample : gc/ms unknown 11/7 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

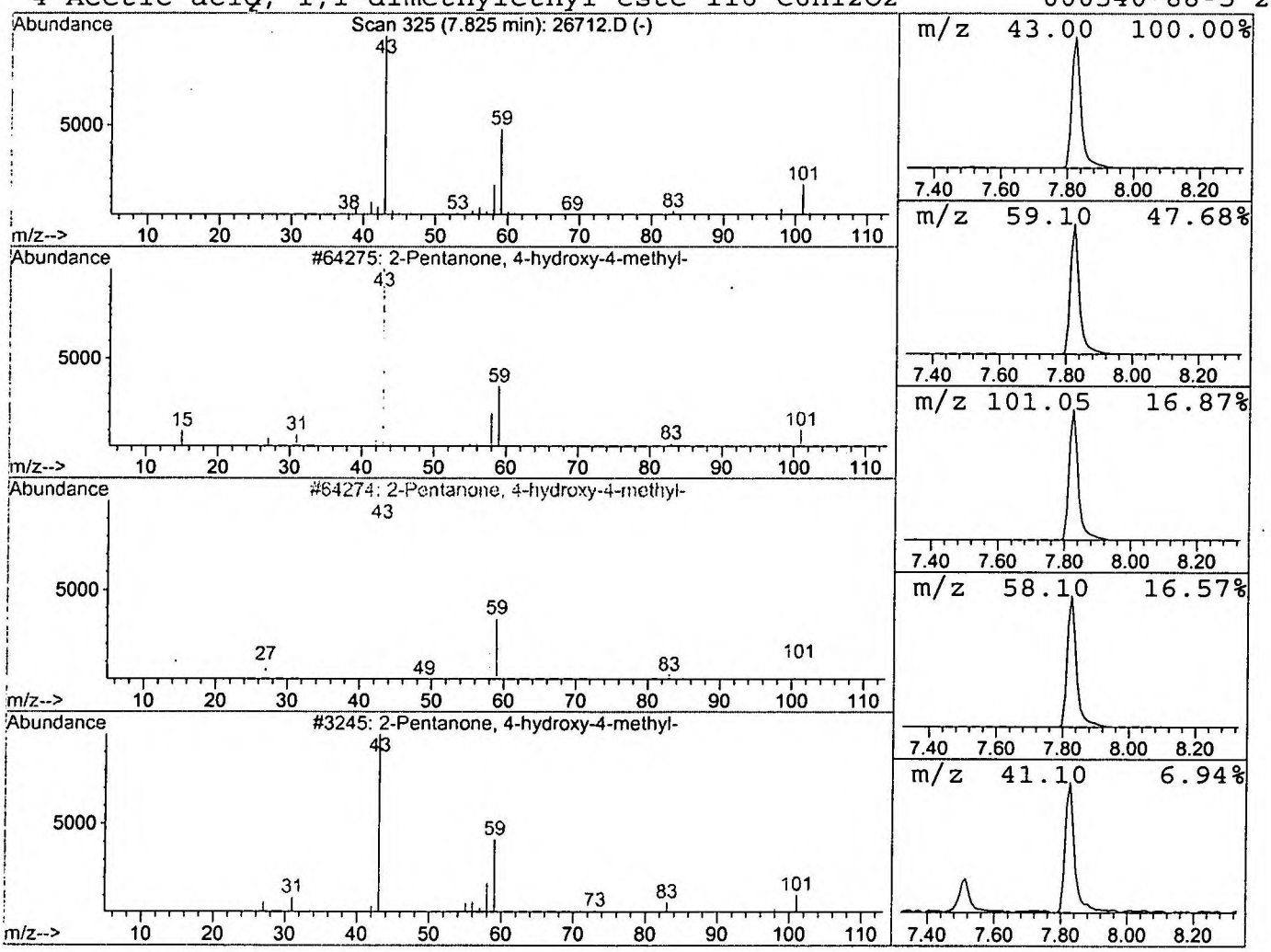
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	8.19 ng/uL	2861960	1,4-Dichlorobenzene-d4	11.86

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	43
4	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D
Acq On : 5 Dec 2001 10:35 am
Sample : gc/ms unknown 11/7
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

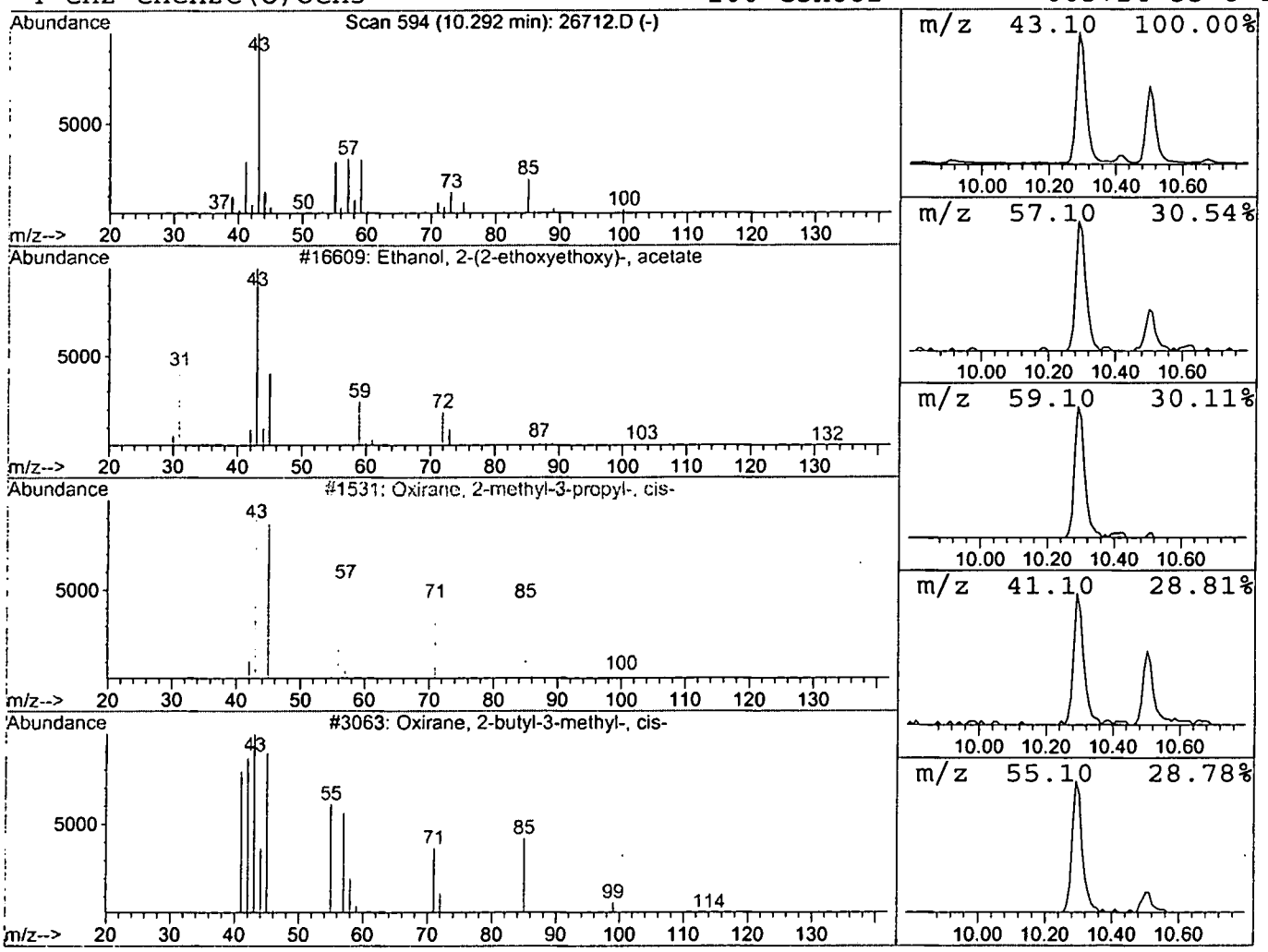
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Ethanol, 2-(2-ethoxyethoxy)-, Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.29	1.47 ng/uL	513104	1,4-Dichlorobenzene-d4	11.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-, aceta	176	C8H16O4	000112-15-2	33
2	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	17
3	Oxirane, 2-butyl-3-methyl-, cis-	114	C7H14O	056052-93-8	12
4	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0501\26712.D
Acq On : 5 Dec 2001 10:35 am
Sample : gc/ms unknown 11/7
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

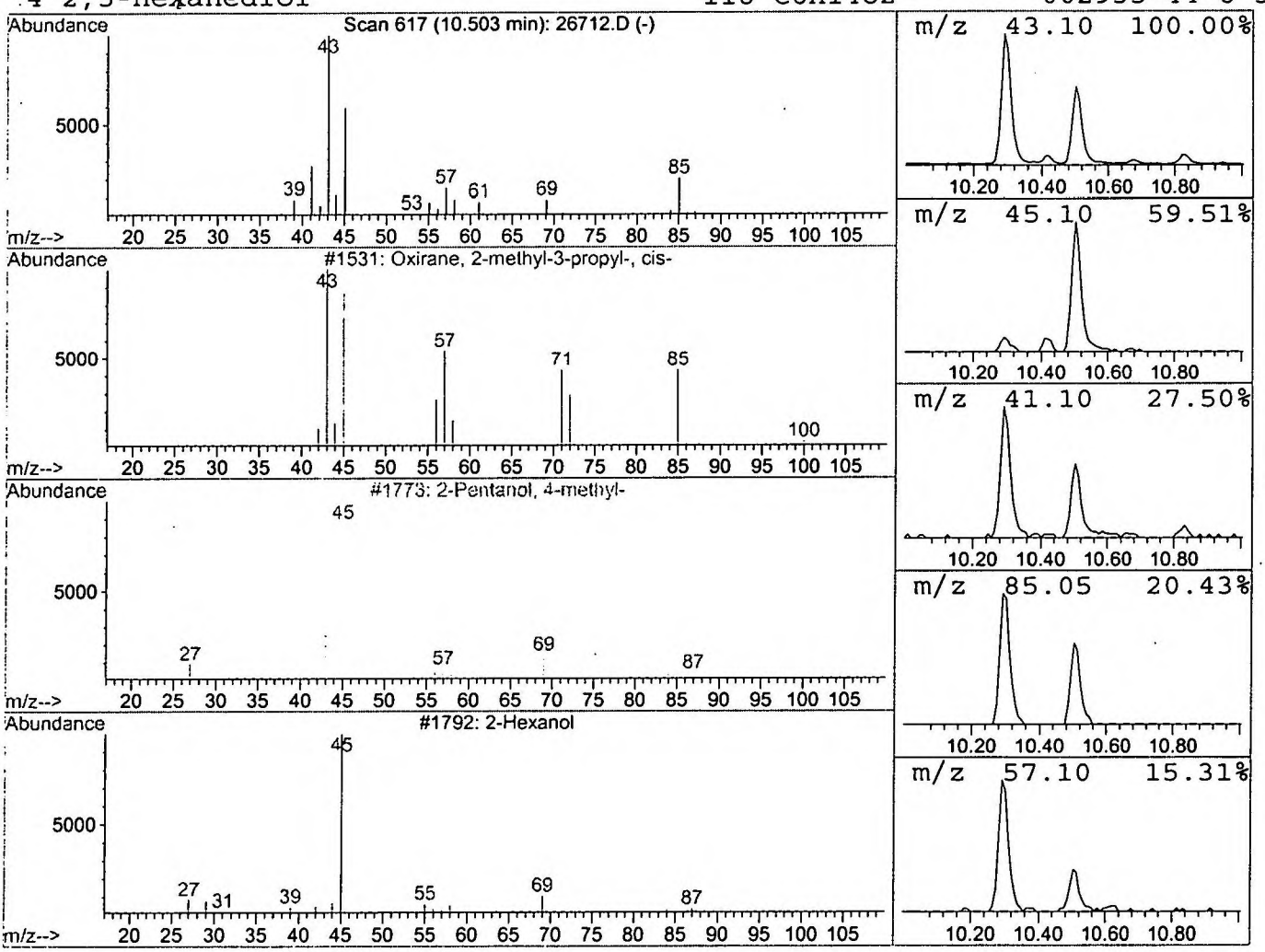
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Oxirane, 2-methyl-3-propyl-, c Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.50	0.81 ng/uL	283922	1,4-Dichlorobenzene-d4	11.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	38
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	38
3	2-Hexanol	102	C6H14O	000626-93-7	35
4	2,5-Hexanediol	118	C6H14O2	002935-44-6	33



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 5 Dec 2001 10:35 am
Data File: C:\HPCHEM\1\DATA\DEC0501\26712.D
Name: gc/ms unknown 11/7
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
2-Propanone, 1-cyclo	5.01	1.6	ng/uL	552784	ISTD01	11.86	6992470	20.0
1-Pentene, 4-methyl-	5.11	1.0	ng/uL	354684	ISTD01	11.86	6992470	20.0
Oxirane, 3-ethyl-2,2	5.17	1.9	ng/uL	656290	ISTD01	11.86	6992470	20.0
Propanoic acid, 2-me	6.11	1.1	ng/uL	373184	ISTD01	11.86	6992470	20.0
2-Butanol, 3-methyl-	7.16	6.5	ng/uL	2288020	ISTD01	11.86	6992470	20.0
2-Pentanone, 4-hydro	7.83	8.2	ng/uL	2861960	ISTD01	11.86	6992470	20.0
Ethanol, 2-(2-ethoxy	10.29	1.5	ng/uL	513104	ISTD01	11.86	6992470	20.0
Oxirane, 2-methyl-3-	10.50	0.8	ng/uL	283922	ISTD01	11.86	6992470	20.0

26712.D NP112701.M Fri Dec 07 13:15:21 2001

Comments for Sample AD26712 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-16-2001 Time: 11:42:09

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, except for the presence of bis(2-Ethylhexyl)phthalate at an estimated level of 0.34 ug/L.

User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/10/2001 Time: 14:53:35

Sample: AD26712
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/10/01 14:53 Ended: 12/10/01 14:53 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26712 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-10-2001 Time: 14:53:40

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.