User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/10/2001 Time: 14:58:30

Sample: AD26625
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/10/01 14:58 Ended: 12/10/01 14:58 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26625 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-10-2001 Time: 14:58:34

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\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 7 10:14 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	1051219	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	3946725	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.77	164	1831908	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.18	188	2588974	20.00	ng/uL	-0.06
59) Chrysene-d12	33.18	240	1586155	20.00	ng/uL	-0.06
68) Perylene-d12	37.27	264	1077730	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	674	0.01	ng/uL	0.47
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.15	152	337	0.01	ng/uL	-0.31
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	18.92	172	1845	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	12.88	45	706	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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Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration
DataAcq Meth : NP112701

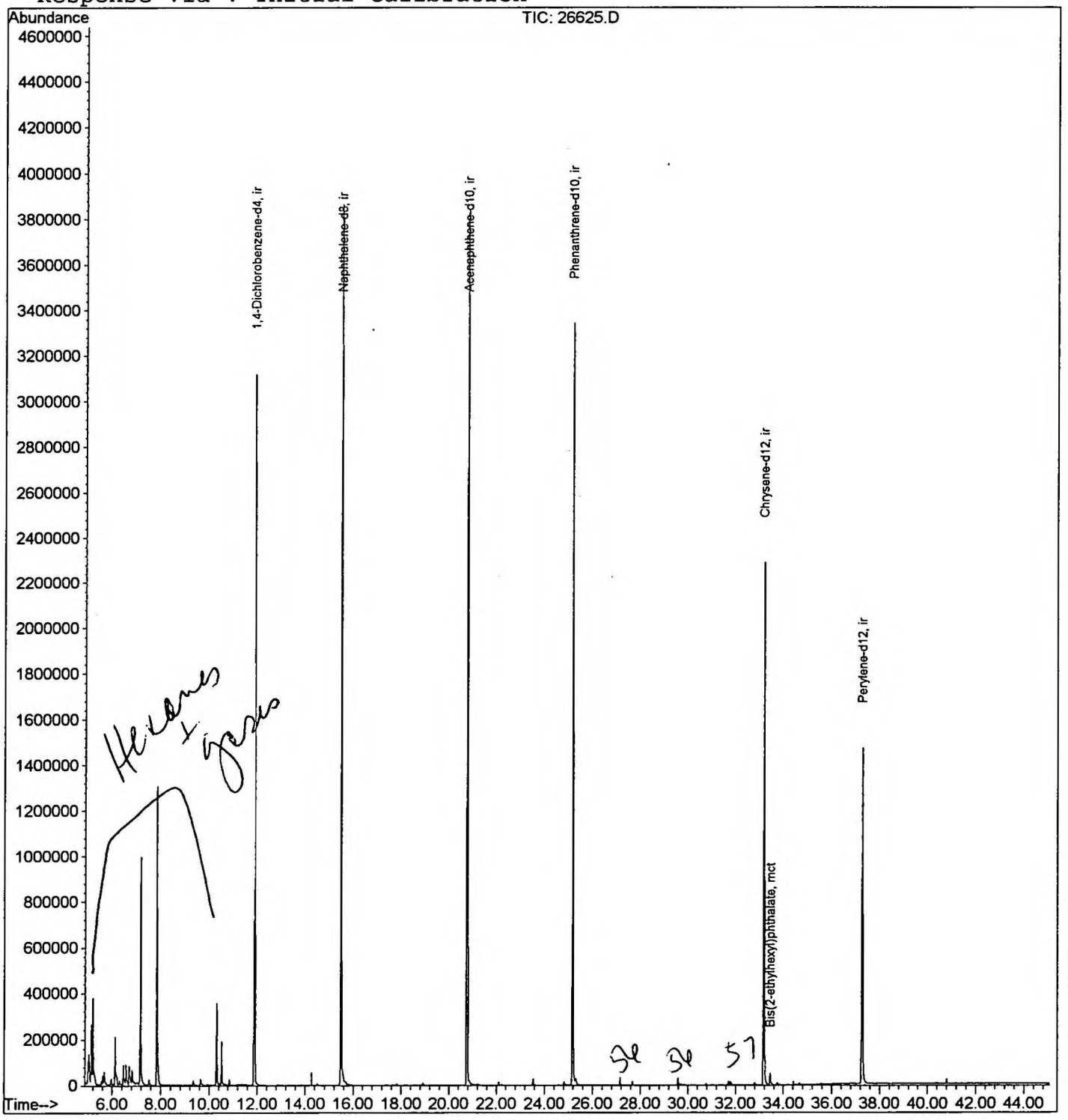
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
72)	Benzo(a)pyrene	37.27	252	3188	N.D.	
73)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
74)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
75)	Benzo(g,h,i)perylene	0.00	276	0	N.D.	

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

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 7 10:14 2001 Quant Results File: NP112701.RES

Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Last Update : Fri Dec 07 09:59:59 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm 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

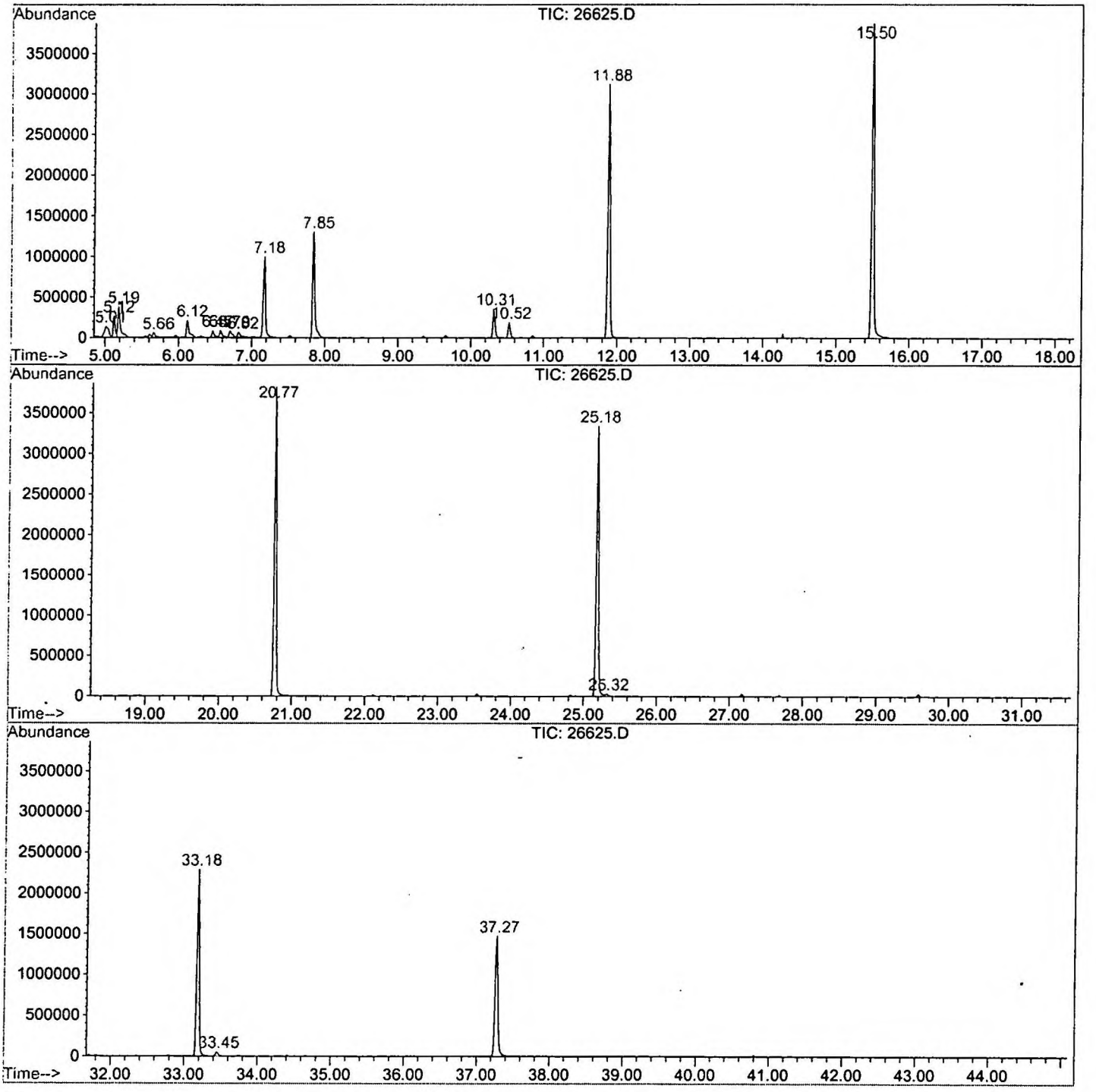
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.014	12	19	28	rBV	130861	515744	6.42%	1.069%
2	5.124	28	31	35	rVV	258408	471635	5.87%	0.978%
3	5.188	35	38	56	rVB	373996	907764	11.31%	1.882%
4	5.665	87	90	98	rVB	52391	110041	1.37%	0.228%
5	6.123	136	140	153	rVB	209157	534735	6.66%	1.109%
6	6.472	174	178	186	rBV	83912	188108	2.34%	0.390%
7	6.573	186	189	200	rVB	84880	208816	2.60%	0.433%
8	6.701	200	203	212	rBV2	81173	232354	2.89%	0.482%
9	6.820	212	216	220	rBV2	56833	109414	1.36%	0.227%
10	7.178	246	255	274	rVB	995399	2096232	26.11%	4.346%
11	7.847	324	328	343	rBV	1303136	2772176	34.53%	5.747%
12	10.314	593	597	605	rBV	355894	724345	9.02%	1.502%
13	10.525	615	620	631	rBV	191287	383907	4.78%	0.796%
14	11.882	763	768	782	rBV	3118260	6462817	80.49%	13.398%
15	15.495	1156	1162	1181	rBV	3865515	8029244	100.00%	16.645%
16	20.768	1731	1737	1757	rBV	3821247	8018774	99.87%	16.624%
17	25.179	2212	2218	2230	rBV2	3341158	7238355	90.15%	15.006%
18	25.316	2230	2233	2246	rVB2	28360	89483	1.11%	0.186%
19	33.184	3084	3091	3109	rBV2	2289624	5054174	62.95%	10.478%
20	33.450	3114	3120	3133	rVB2	52653	141240	1.76%	0.293%
21	37.274	3529	3537	3552	rBV2	1467655	3947945	49.17%	8.184%

Sum of corrected areas: 48237303

26625.D NP112701.M Fri Dec 07 14:48:25 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0601\26625.D
Operator : GALOS
Acquired : 6 Dec 2001 4:34 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 8
Quant File :NP112701.RES (RTE Integrator)



26625.D NP112701.M Fri Dec 07 14:48:27 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

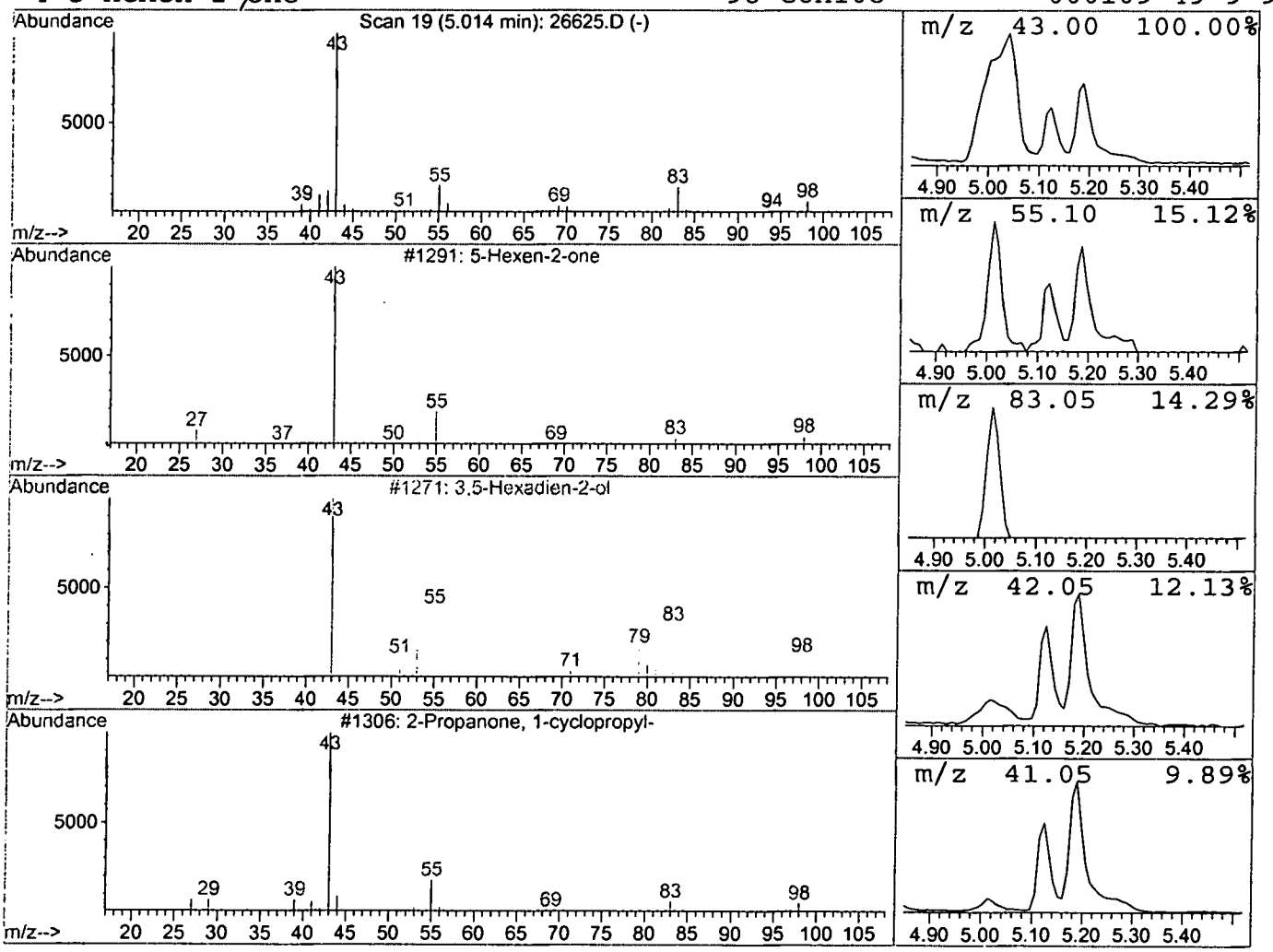
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 5-Hexen-2-one Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	1.60 ng/uL	515744	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Hexen-2-one	98	C6H10O	000109-49-9	9
2	3,5-Hexadien-2-ol	98	C6H10O	003280-51-1	9
3	2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

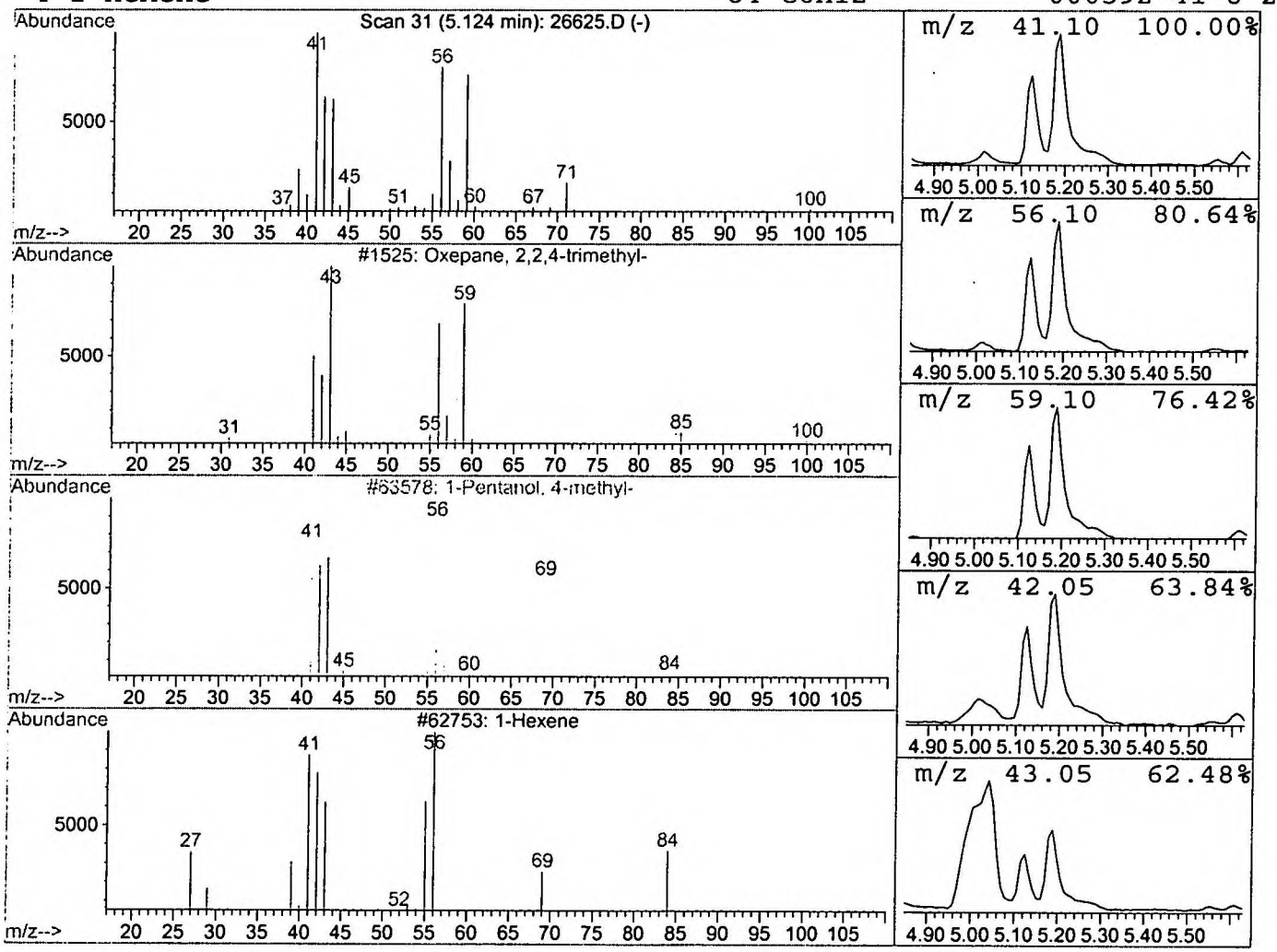
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Oxepane, 2,2,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	1.46 ng/uL	471635	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxepane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	25
2	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	25
3	1-Hexene	84	C6H12	000592-41-6	25
4	1-Hexene	84	C6H12	000592-41-6	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D
Acq On : 6 Dec 2001 4:34 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

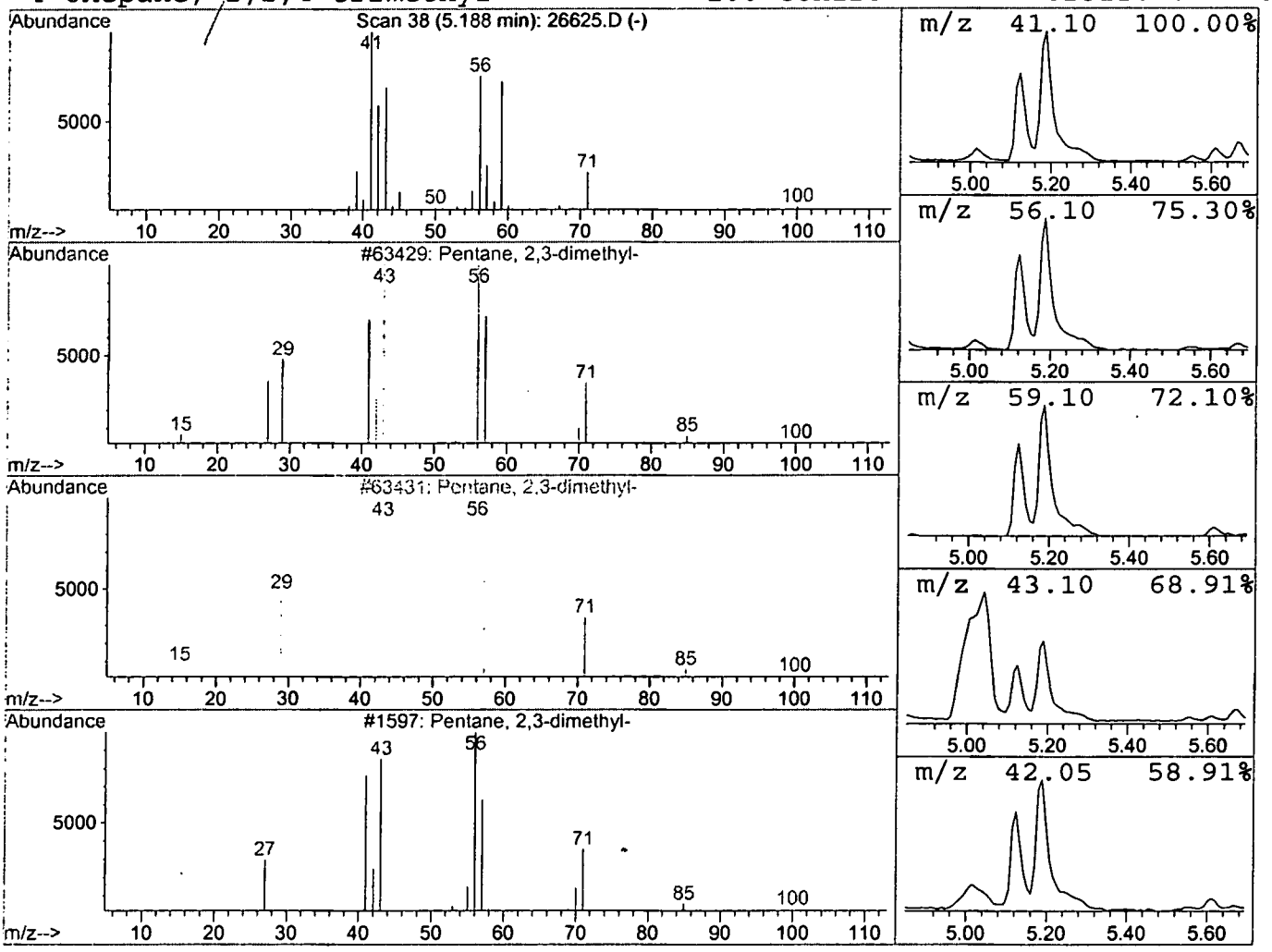
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	2.81 ng/uL	907764	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	43
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	43
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	37
4	Oxepane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	33



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

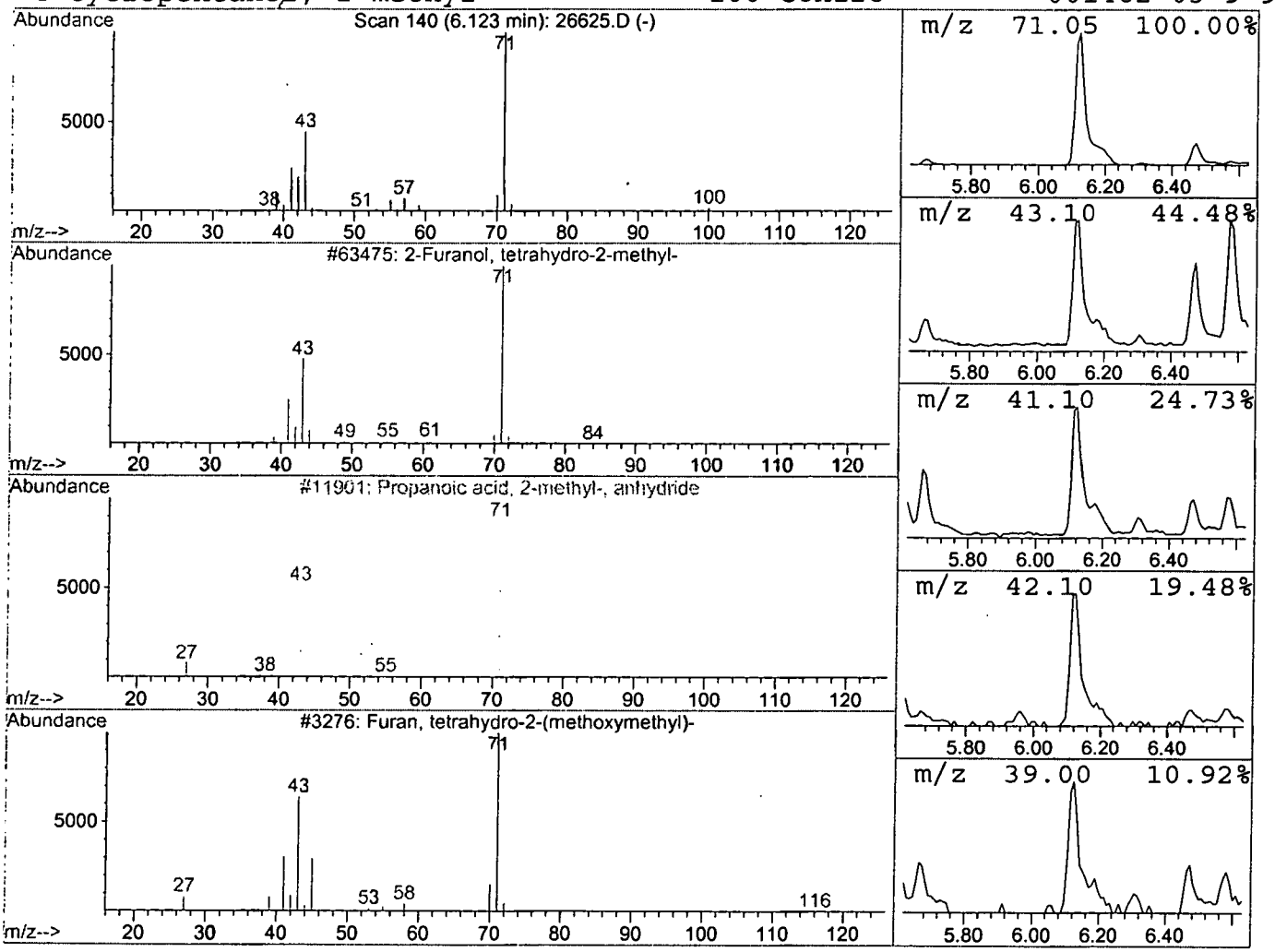
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 2-Furanol, tetrahydro-2-methyl Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	1.65 ng/uL	534735	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furanol, tetrahydro-2-methyl-	102	C5H10O2	007326-46-7	43
2	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	40
3	Furan, tetrahydro-2-(methoxymethyl)	116	C6H12O2	019354-27-9	23
4	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D
Acq On : 6 Dec 2001 4:34 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

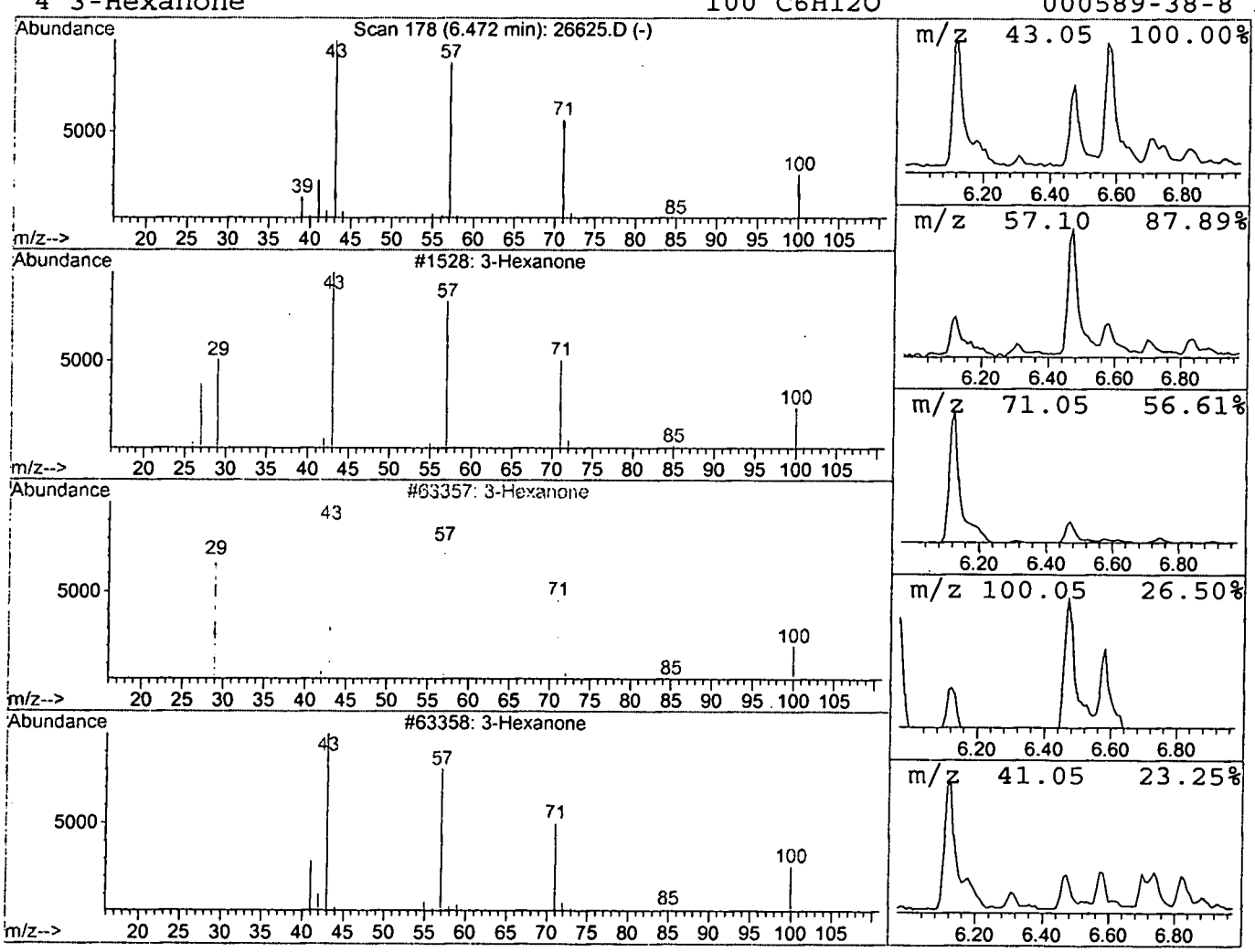
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 3-Hexanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	0.58 ng/uL	188108	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	91
2	3-Hexanone	100	C6H12O	000589-38-8	91
3	3-Hexanone	100	C6H12O	000589-38-8	86
4	3-Hexanone	100	C6H12O	000589-38-8	53



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D
Acq On : 6 Dec 2001 4:34 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

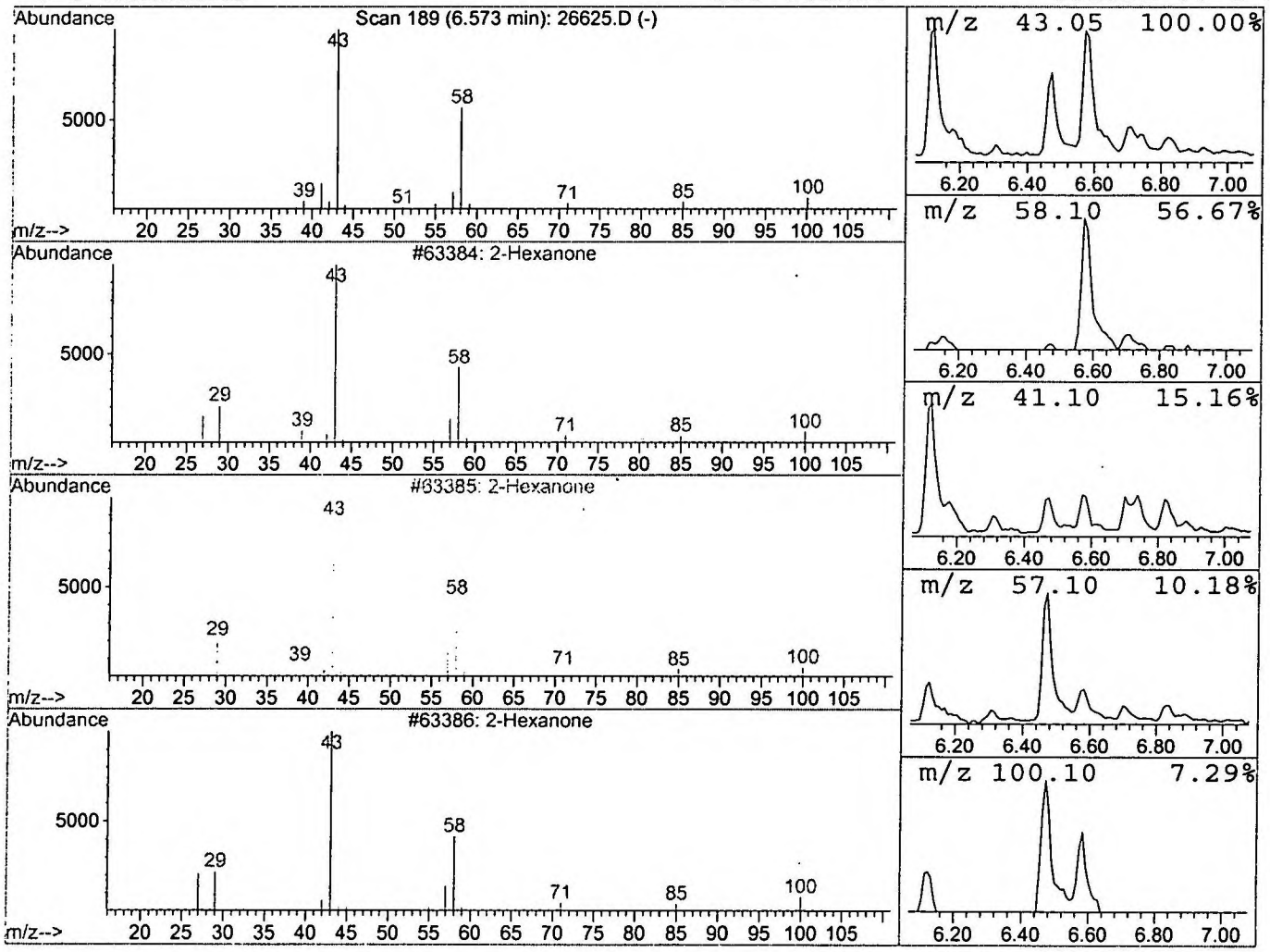
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 2-Hexanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	0.65 ng/uL	208816	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	64
2	2-Hexanone	100	C6H12O	000591-78-6	64
3	2-Hexanone	100	C6H12O	000591-78-6	53
4	2-Hexanone	100	C6H12O	000591-78-6	53



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

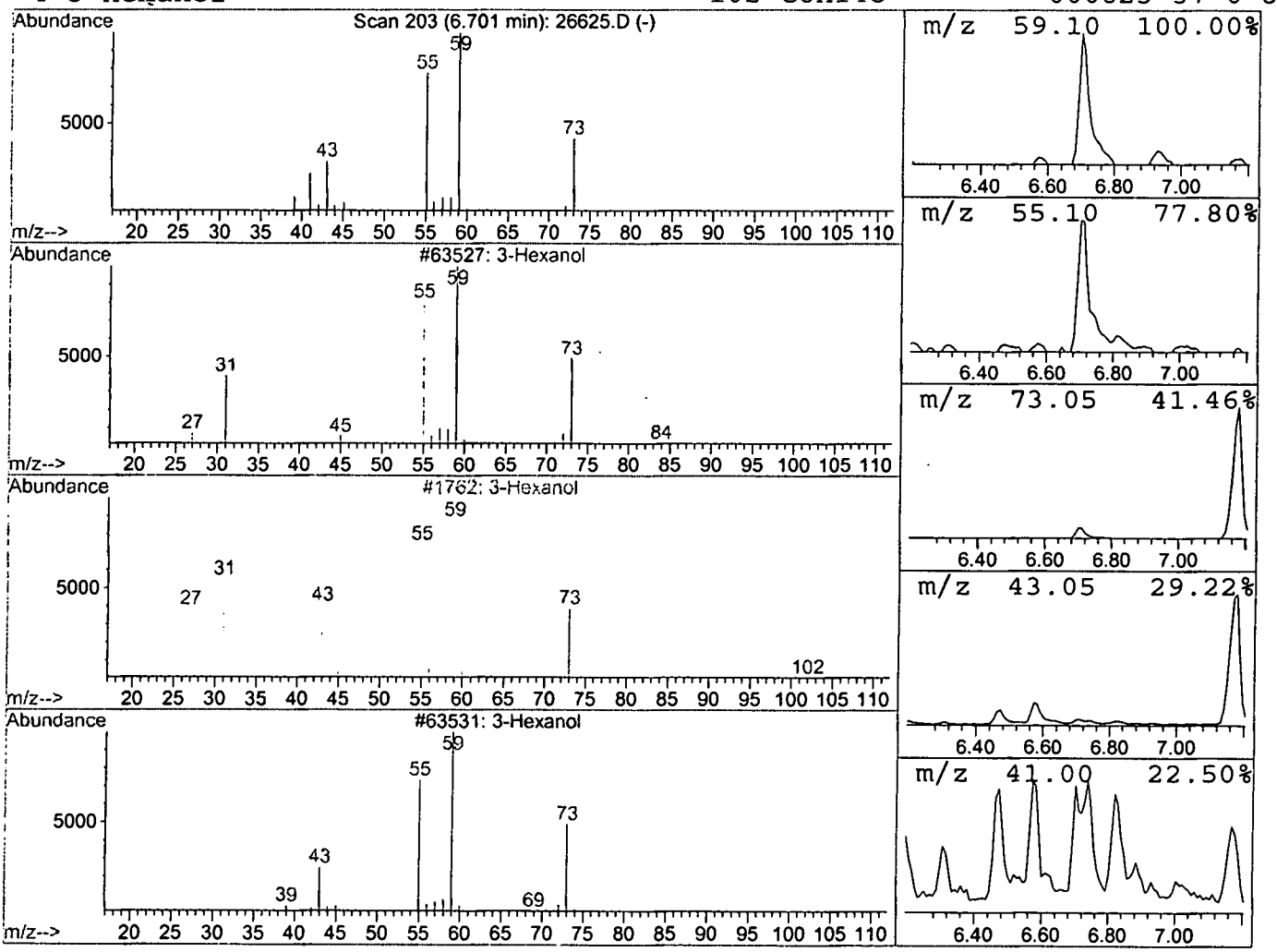
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 3-Hexanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	0.72 ng/uL	232354	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	83
2	3-Hexanol	102	C6H14O	000623-37-0	83
3	3-Hexanol	102	C6H14O	000623-37-0	83
4	3-Hexanol	102	C6H14O	000623-37-0	83



Library Search Compound Report

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Acq On : 6 Dec 2001 4:34 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

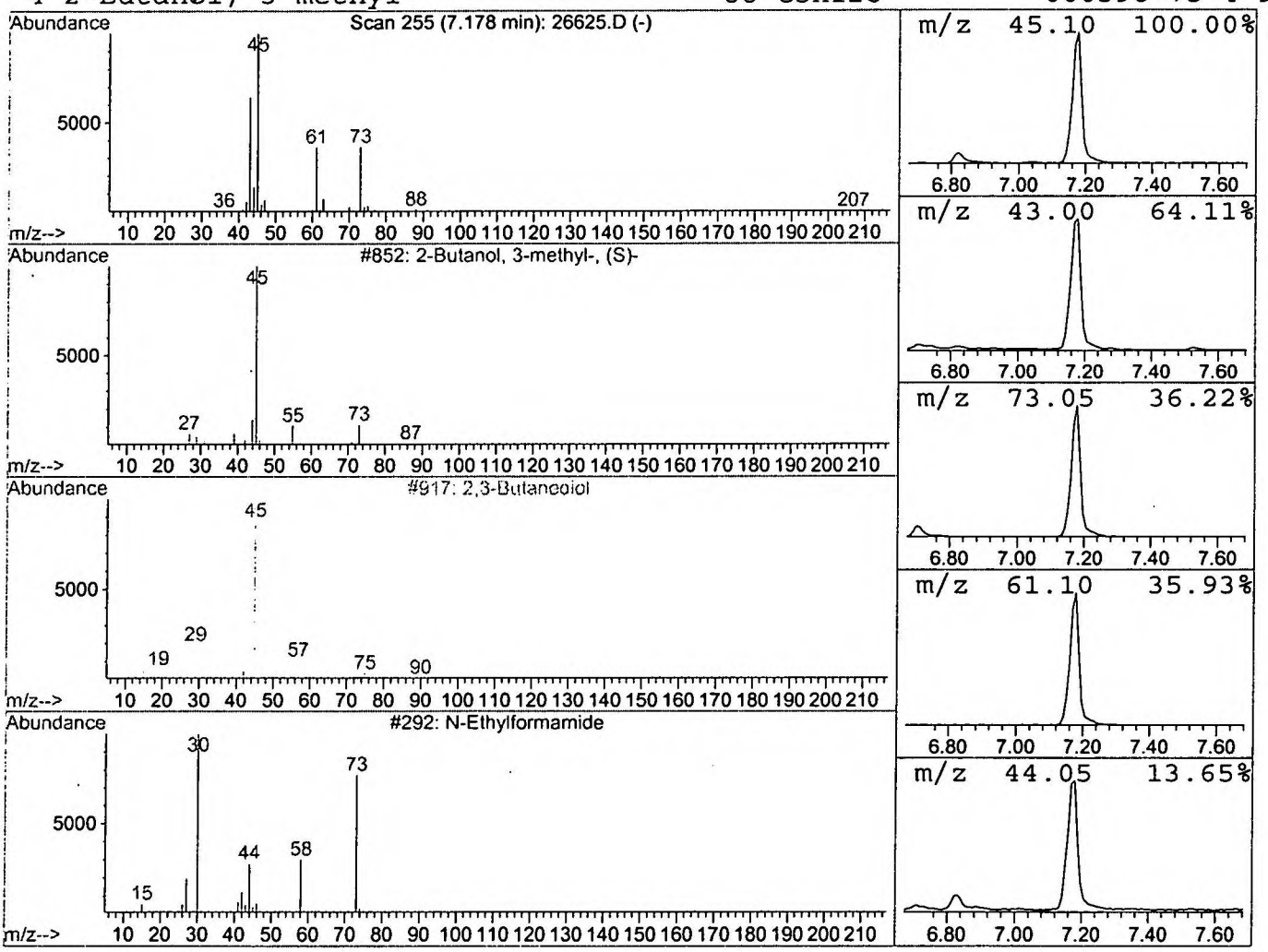
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 2-Butanol, 3-methyl-, (S)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	6.49 ng/uL	2096230	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	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

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D
Acq On : 6 Dec 2001 4:34 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

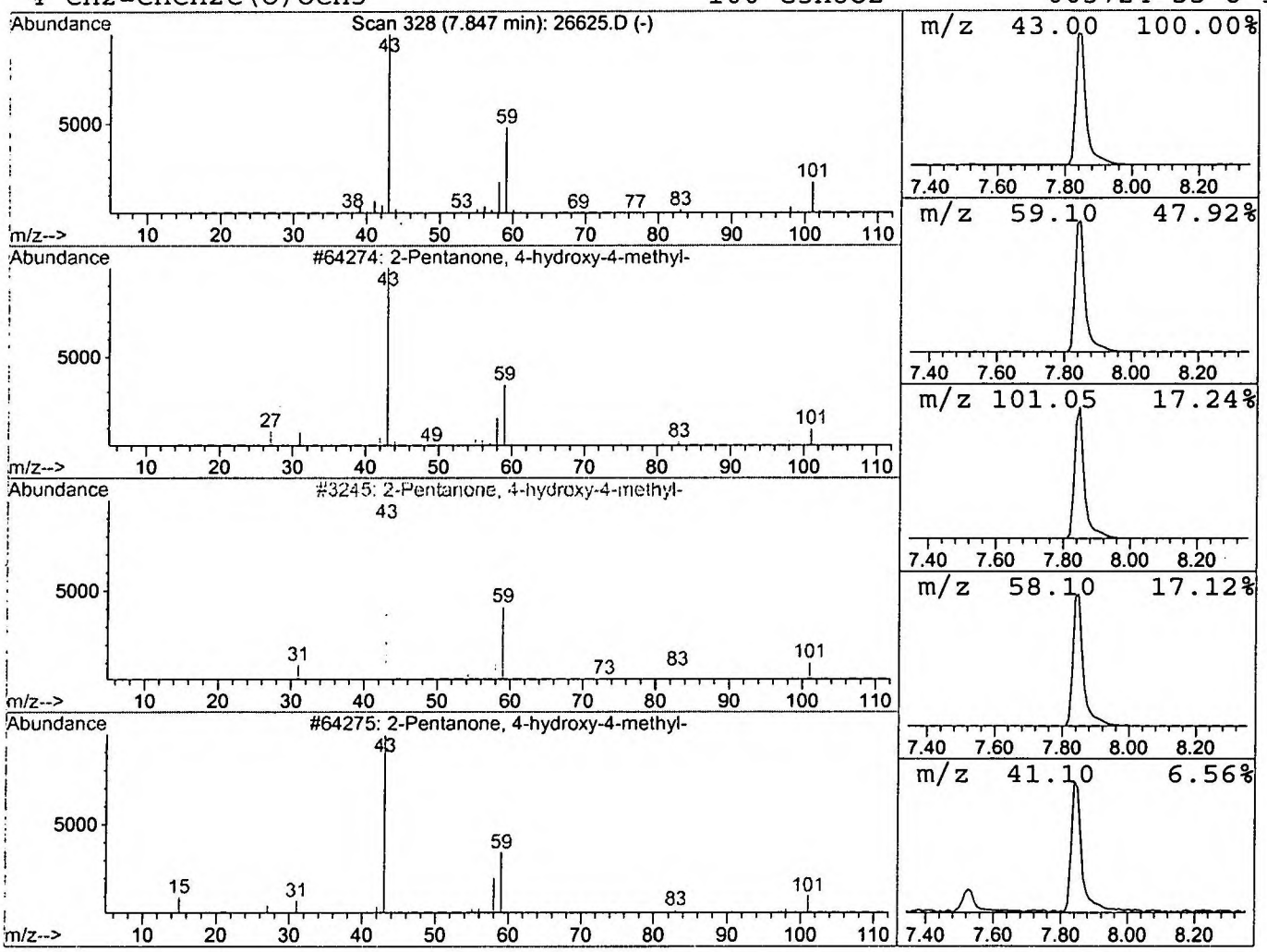
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	8.58 ng/uL	2772180	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	59
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	25
4	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D Vial: 8
Acq On : 6 Dec 2001 4:34 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

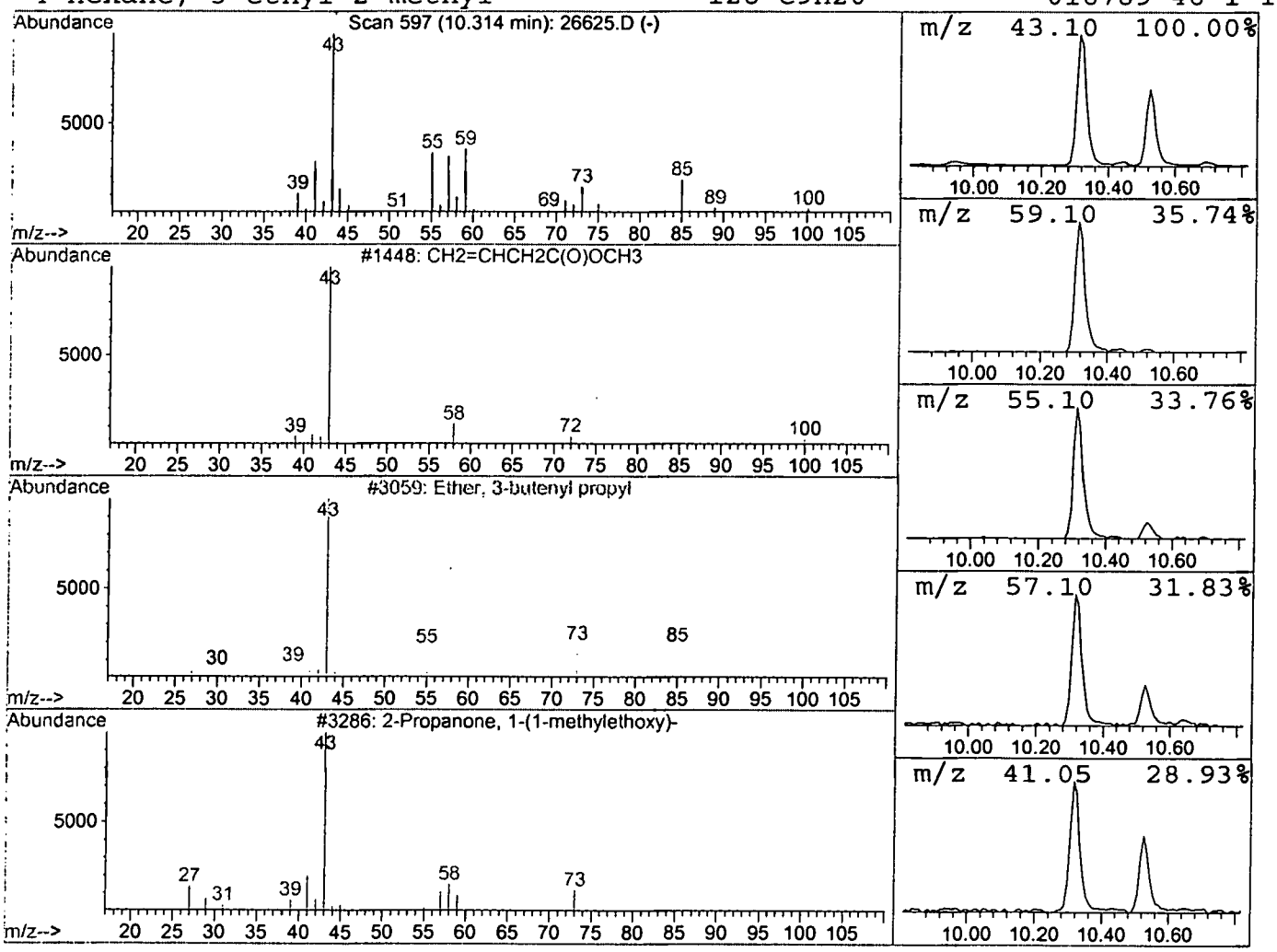
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 10 CH2=CHCH2C(O)OCH3 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.31	2.24 ng/uL	724345	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	22
2	Ether, 3-butenyl propyl	114	C7H14O	034061-75-1	17
3	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	12
4	Hexane, 3-ethyl-2-methyl-	128	C9H20	016789-46-1	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26625.D
Acq On : 6 Dec 2001 4:34 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

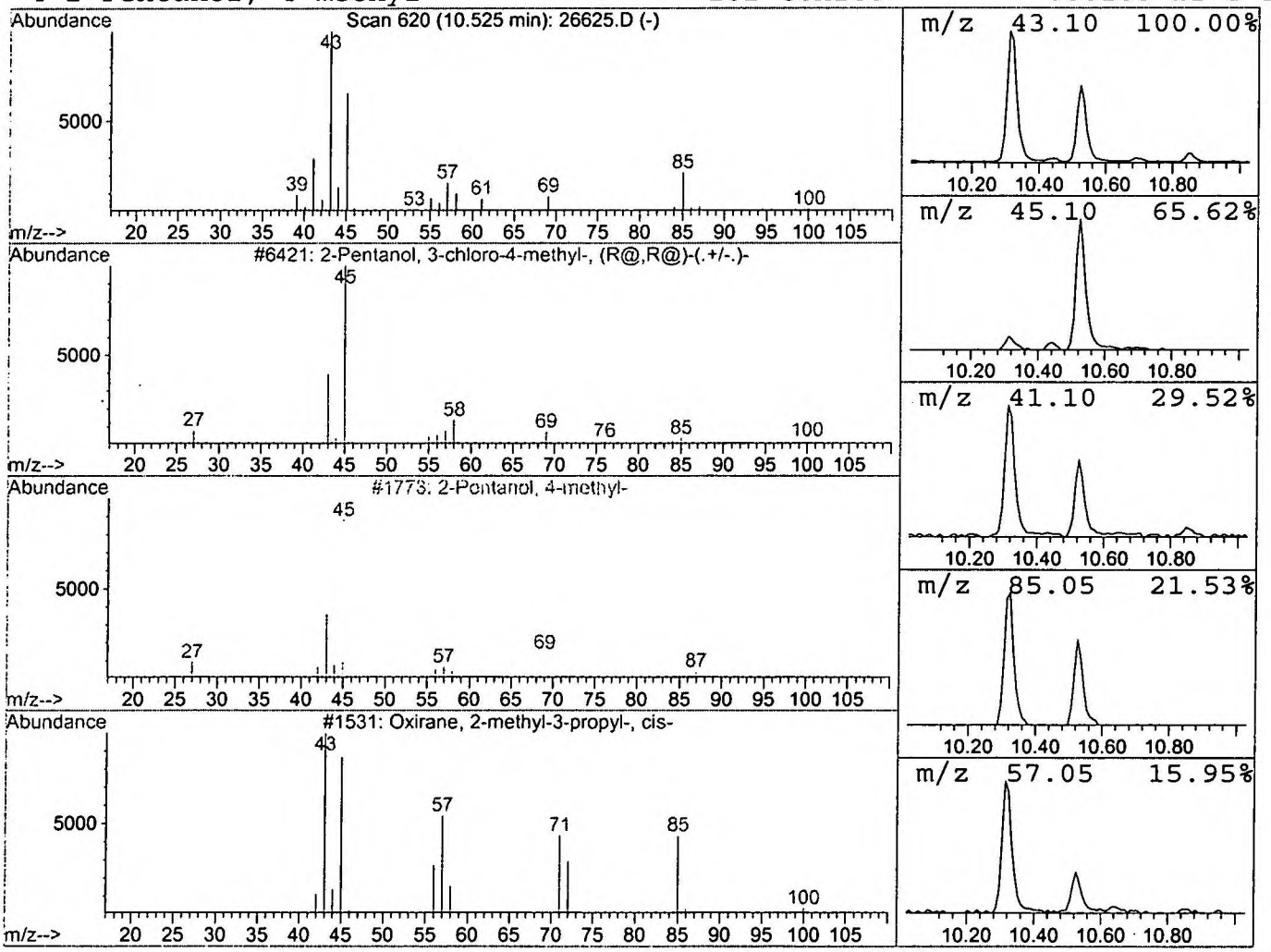
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 11 2-Pentanol, 3-chloro-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	1.19 ng/uL	383907	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 3-chloro-4-methyl-, (R@	136	C6H13ClO	074685-47-5	43
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	38
3	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	36
4	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	35



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 6 Dec 2001 4:34 pm
Data File: C:\HPCHEM\1\DATA\DEC0601\26625.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
5-Hexen-2-one	5.01	1.6	ng/uL	515744	ISTD01	11.88	6462820	20.0
Oxepane, 2,2,4-trime	5.12	1.5	ng/uL	471635	ISTD01	11.88	6462820	20.0
Pentane, 2,3-dimethy	5.19	2.8	ng/uL	907764	ISTD01	11.88	6462820	20.0
2-Furanol, tetrahydr	6.12	1.7	ng/uL	534735	ISTD01	11.88	6462820	20.0
3-Hexanone	6.47	0.6	ng/uL	188108	ISTD01	11.88	6462820	20.0
2-Hexanone	6.57	0.6	ng/uL	208816	ISTD01	11.88	6462820	20.0
3-Hexanol	6.70	0.7	ng/uL	232354	ISTD01	11.88	6462820	20.0
2-Butanol, 3-methyl-	7.18	6.5	ng/uL	2096230	ISTD01	11.88	6462820	20.0
2-Pentanone, 4-hydro	7.85	8.6	ng/uL	2772180	ISTD01	11.88	6462820	20.0
CH2=CHCH2C(O)OCH3	10.31	2.2	ng/uL	724345	ISTD01	11.88	6462820	20.0
2-Pentanol, 3-chloro	10.53	1.2	ng/uL	383907	ISTD01	11.88	6462820	20.0

26625.D NP112701.M Fri Dec 07 14:48:51 2001

Comments for Sample AD26625 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-16-2001 Time: 11:31:00

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 0.44 ug/L.

0.33 in Blank