User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/10/2001 Time: 14:54:53

Sample: AD26829
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/10/01 14:54 Ended: 12/10/01 14:54 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26829 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-10-2001 Time: 14:54:58

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\26829.D Vial: 6
Acq On : 6 Dec 2001 2:44 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 7 10:25 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	938937	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	3462269	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.78	164	1572174	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.18	188	2295741	20.00	ng/uL	-0.06
59) Chrysene-d12	33.18	240	1446339	20.00	ng/uL	-0.06
68) Perylene-d12	37.27	264	1001546	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	720	0.01	ng/uL	0.47
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.10	152	1730	0.04	ng/uL	-0.36
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.08%#		
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	1741	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.		
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

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Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
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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
71) Benzo(k)fluoranthene	0.00	252	0	N.D.	
72) Benzo(a)pyrene	37.27	252	3046	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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Quantitation Report

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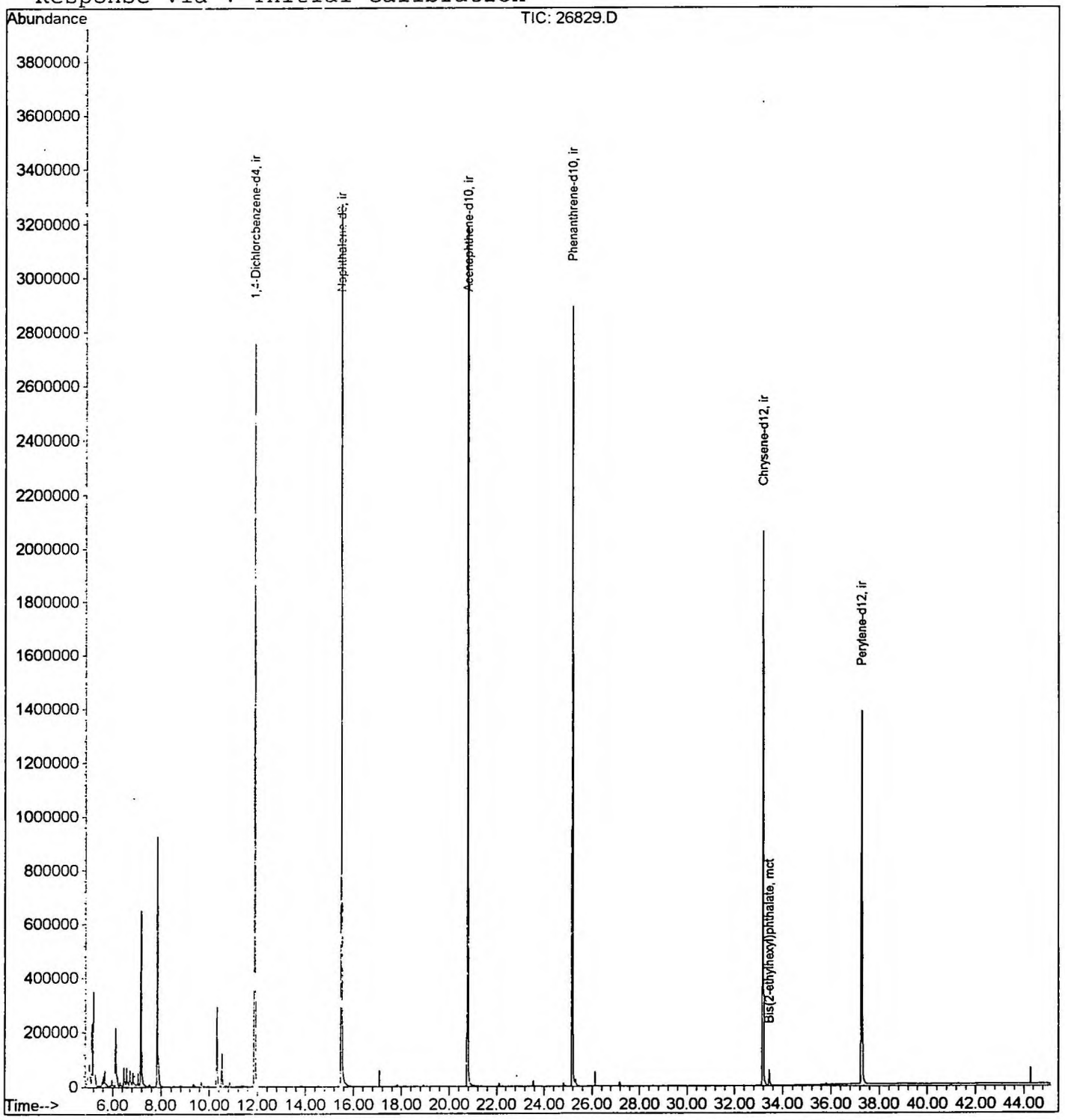
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 7 10:25 2001

Vial: 6
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 : Fri Dec 07 09:59:59 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D Vial: 6
Acq On : 6 Dec 2001 2:44 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

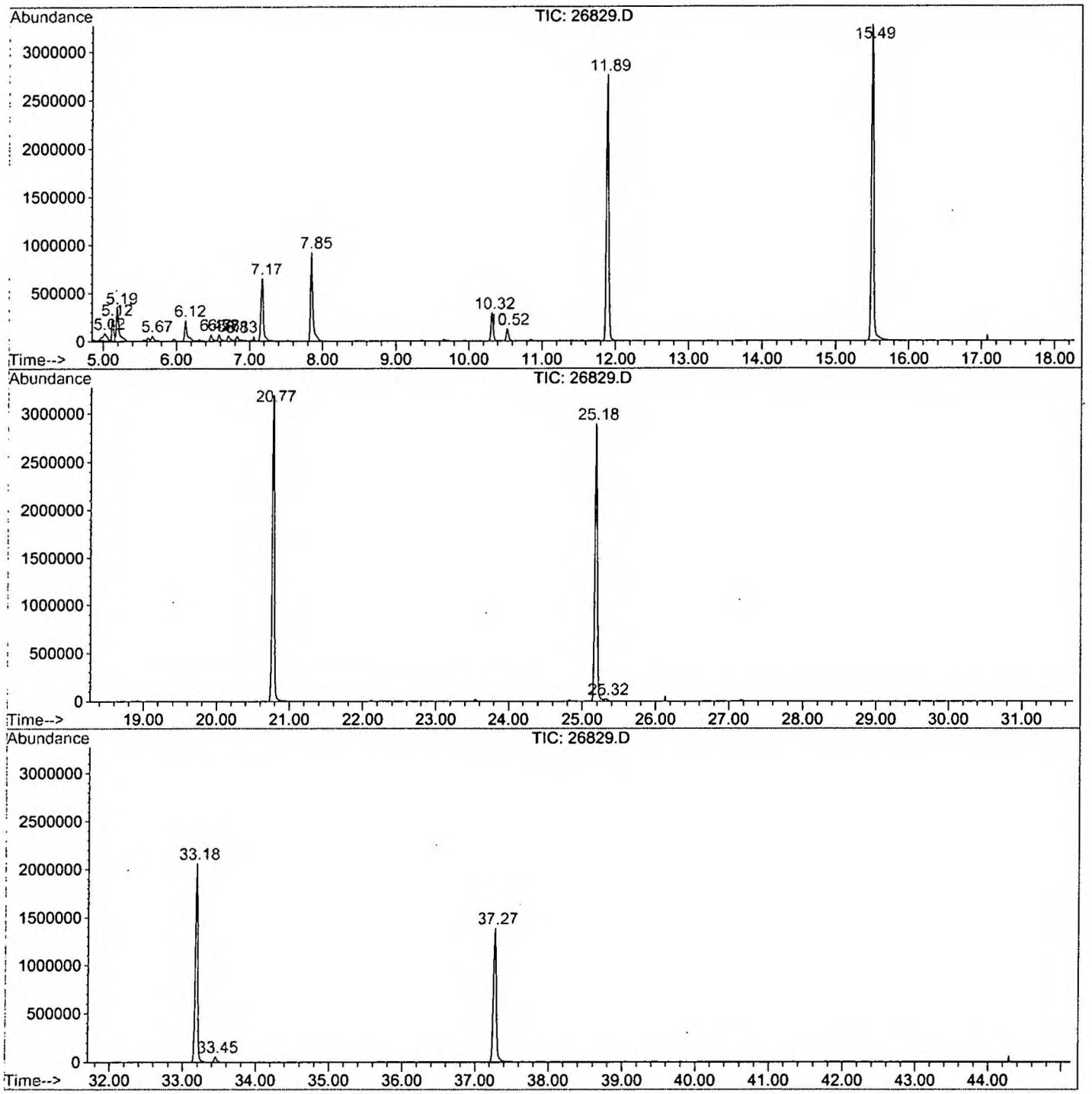
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.022	10	20	28	rBV2	76492	319428	4.56%	0.766%
2	5.123	28	31	35	rVV	223693	414360	5.92%	0.994%
3	5.187	35	38	55	rVB	345542	863314	12.32%	2.071%
4	5.673	87	91	103	rVB	58013	151713	2.17%	0.364%
5	6.123	136	140	153	rBV	215301	516307	7.37%	1.238%
6	6.471	174	178	186	rBV	69883	148870	2.13%	0.357%
7	6.581	186	190	200	rVB	69235	159430	2.28%	0.382%
8	6.710	200	204	213	rBV	58493	148836	2.12%	0.357%
9	6.829	213	217	221	rBV2	45660	93675	1.34%	0.225%
10	7.168	248	254	267	rBV	646250	1471639	21.01%	3.530%
11	7.847	324	328	346	rBV	924129	1899870	27.12%	4.557%
12	10.323	593	598	605	rBV	290999	590060	8.42%	1.415%
13	10.524	614	620	631	rBV	126416	282377	4.03%	0.677%
14	11.891	763	769	780	rBV	2754587	5753366	82.13%	13.799%
15	15.495	1157	1162	1180	rBV	3270908	7004779	100.00%	16.801%
16	20.767	1731	1737	1749	rBV	3189763	6930517	98.94%	16.623%
17	25.178	2212	2218	2230	rBV2	2892282	6435934	91.88%	15.436%
18	25.325	2230	2234	2244	rVB	25011	75178	1.07%	0.180%
19	33.184	3085	3091	3110	rBV2	2060600	4627610	66.06%	11.099%
20	33.450	3115	3120	3130	rVB	59693	145455	2.08%	0.349%
21	37.274	3529	3537	3555	rBV2	1381150	3660373	52.26%	8.779%

Sum of corrected areas: 41693091

26829.D NP112701.M Fri Dec 07 14:54:17 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0601\26829.D
Operator : GALOS
Acquired : 6 Dec 2001 2:44 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 6
Quant File :NP112701.RES (RTE Integrator)



26829.D NP112701.M Fri Dec 07 14:54:19 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D
Acq On : 6 Dec 2001 2:44 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

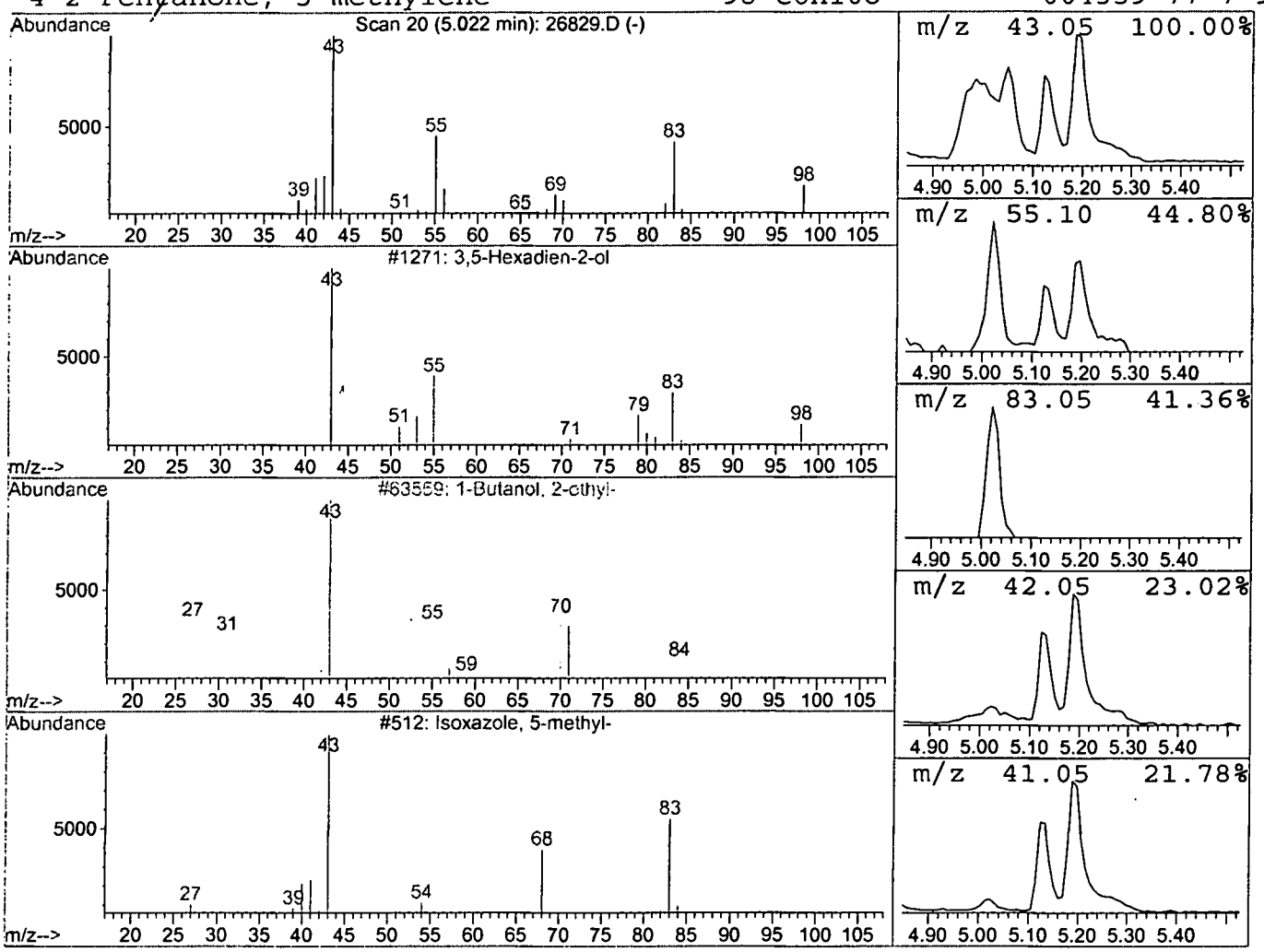
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 3,5-Hexadien-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	1.11 ng/uL	319428	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Hexadien-2-ol	98	C6H10O	003280-51-1	40
2	1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	23
3	Isoxazole, 5-methyl-	83	C4H5NO	005765-44-6	9
4	2-Pentanone, 3-methylene-	98	C6H10O	004359-77-7	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D
Acq On : 6 Dec 2001 2:44 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

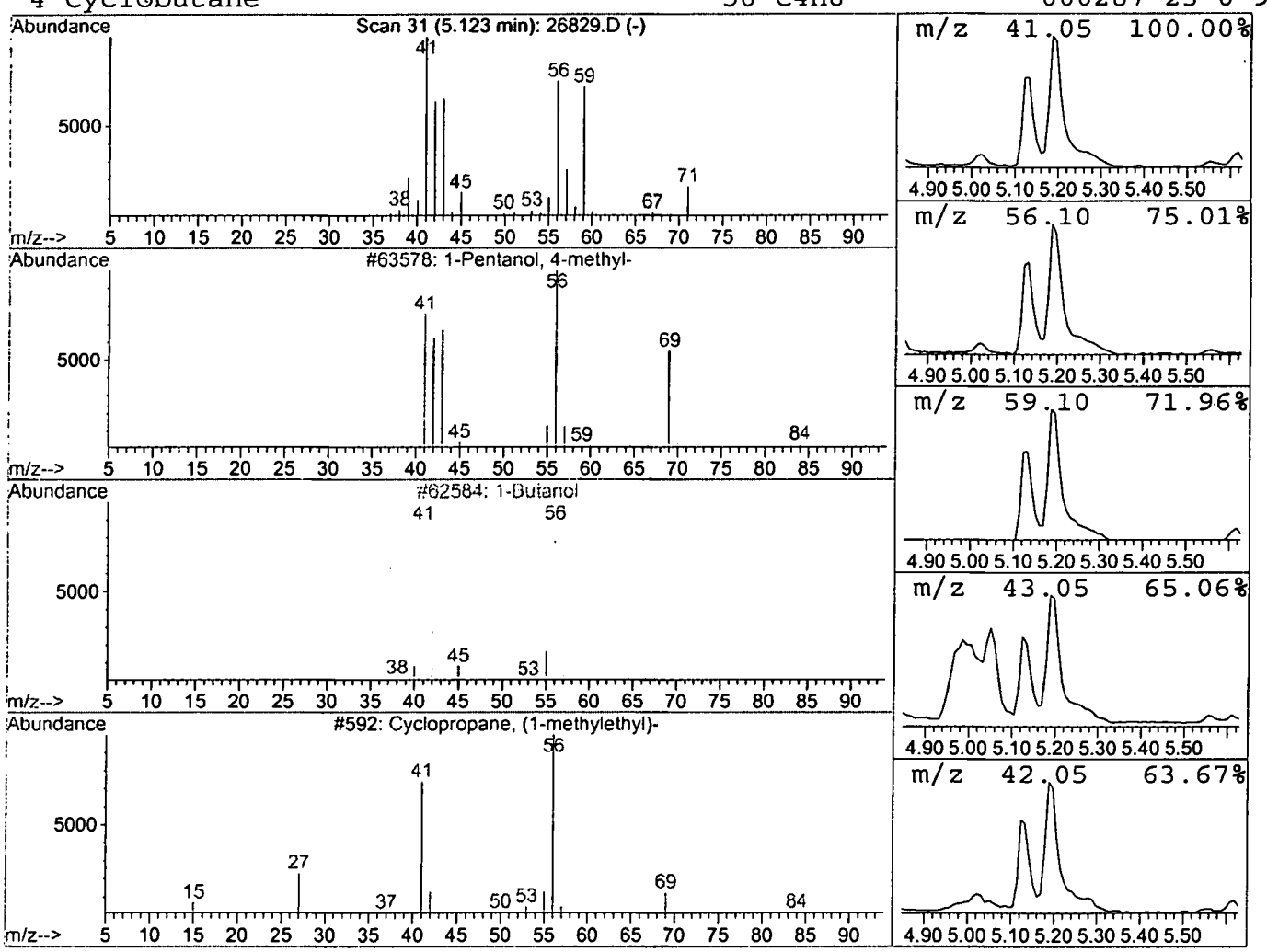
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	1.44 ng/uL	414360	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	36
2	1-Butanol	74	C4H10O	000071-36-3	33
3	Cyclopropane, (1-methylethyl)-	84	C6H12	003638-35-5	12
4	Cyclobutane	56	C4H8	000287-23-0	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D
Acq On : 6 Dec 2001 2:44 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

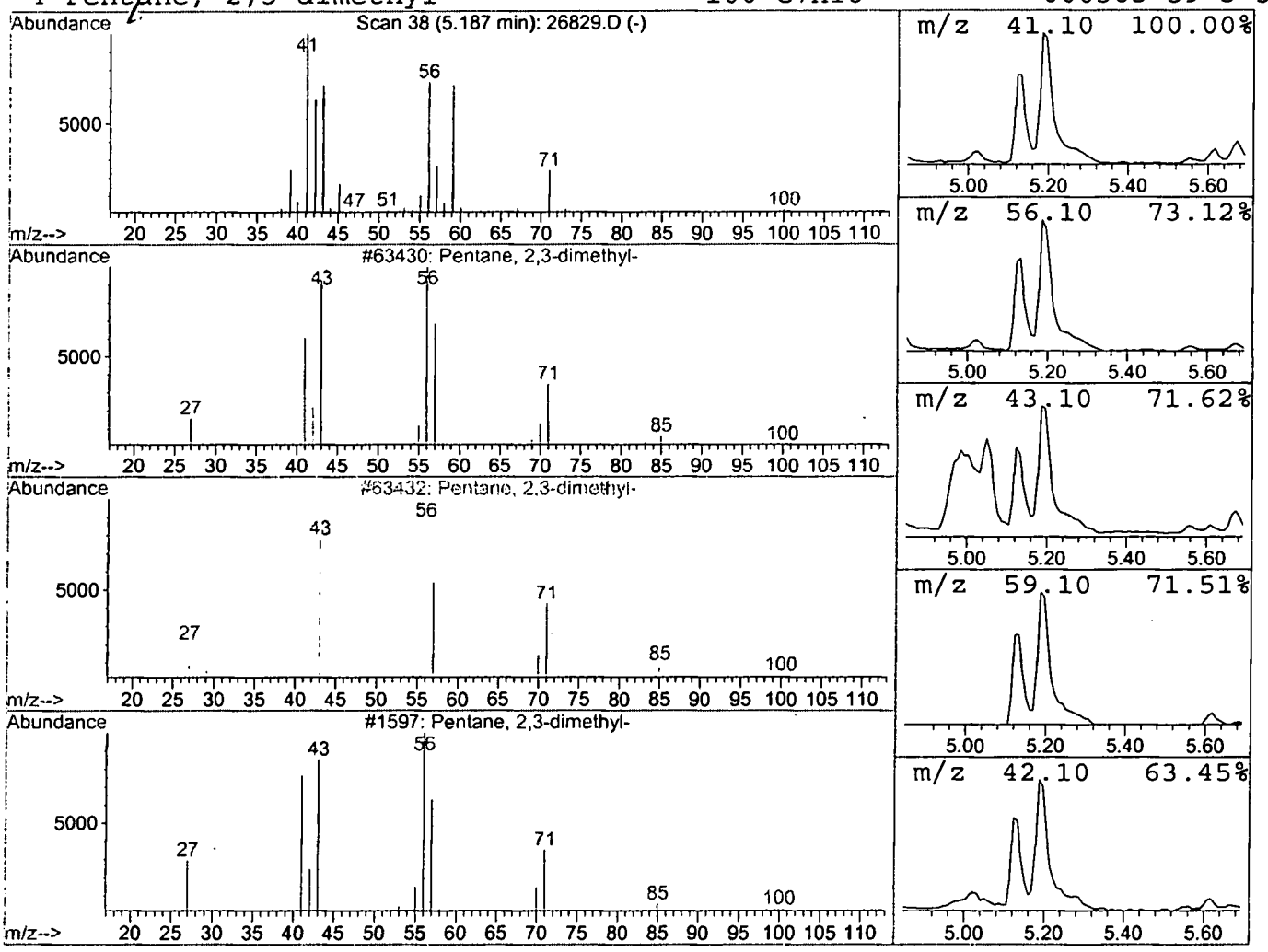
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.00 ng/uL	863314	1,4-Dichlorobenzene-d4	11.89

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	38
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	37
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	37



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D
Acq On : 6 Dec 2001 2:44 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

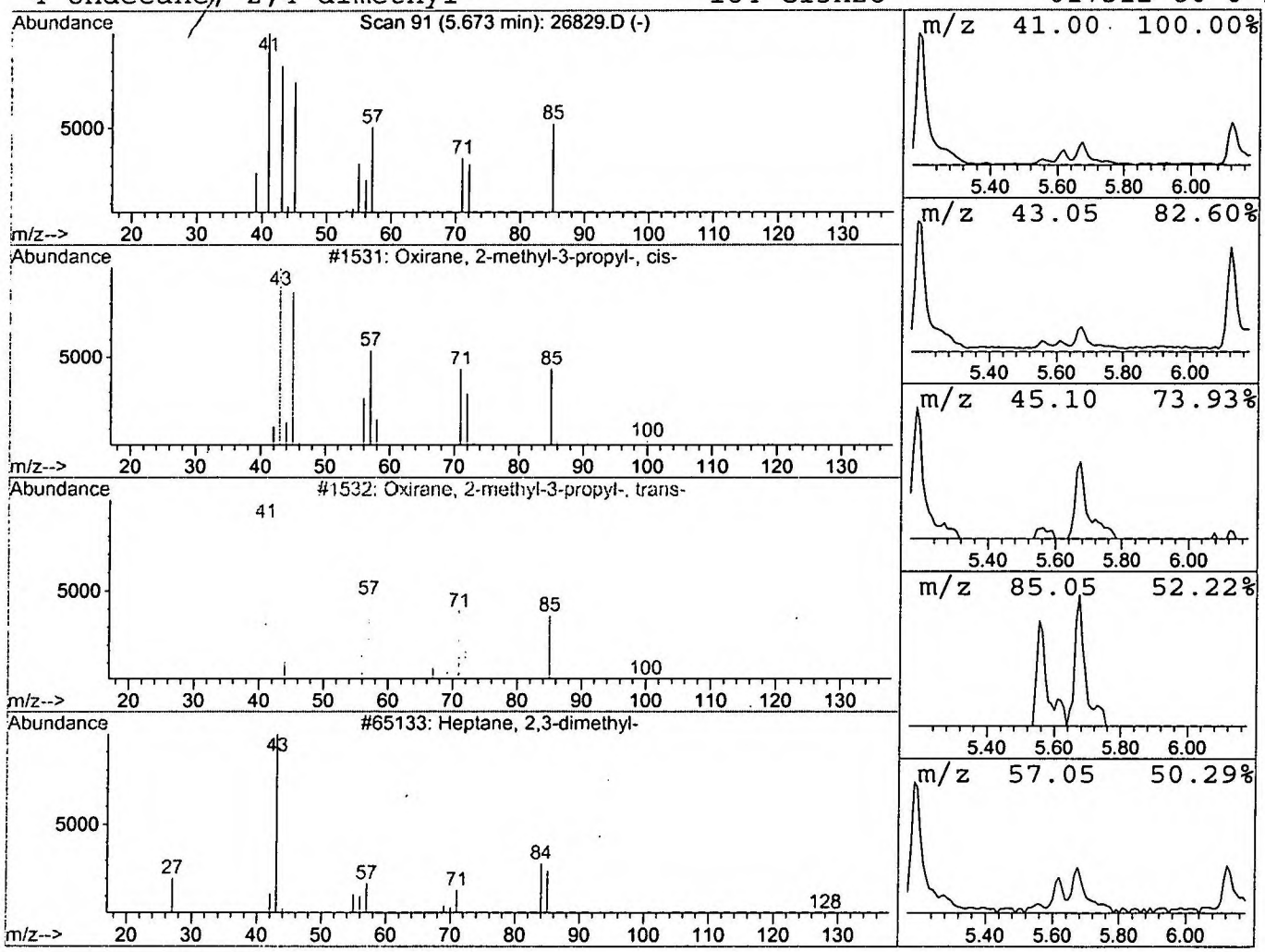
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Oxirane, 2-methyl-3-propyl-, c Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.67	0.53 ng/uL	151713	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	56
2	Oxirane, 2-methyl-3-propyl-, trans-	100	C6H12O	006124-91-0	17
3	Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	12
4	Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	12



Library Search Compound Report

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Acq On : 6 Dec 2001 2:44 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

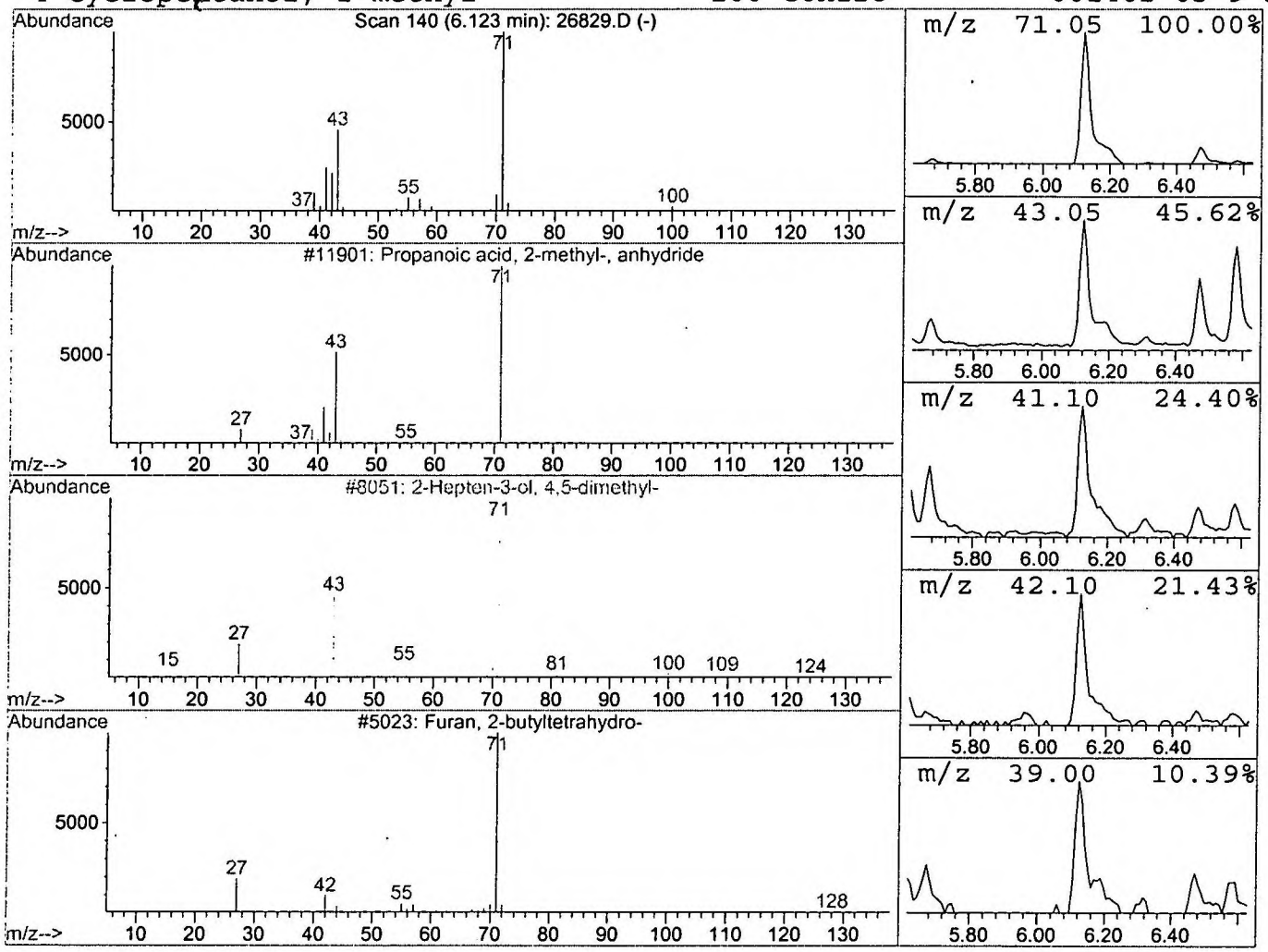
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Propanoic acid, 2-methyl-, anh Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	1.79 ng/uL	516307	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	53
2	2-Hepten-3-ol, 4,5-dimethyl-	142	C9H18O	055956-37-1	39
3	Furan, 2-butyltetrahydro-	128	C8H16O	001004-29-1	38
4	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	33



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Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

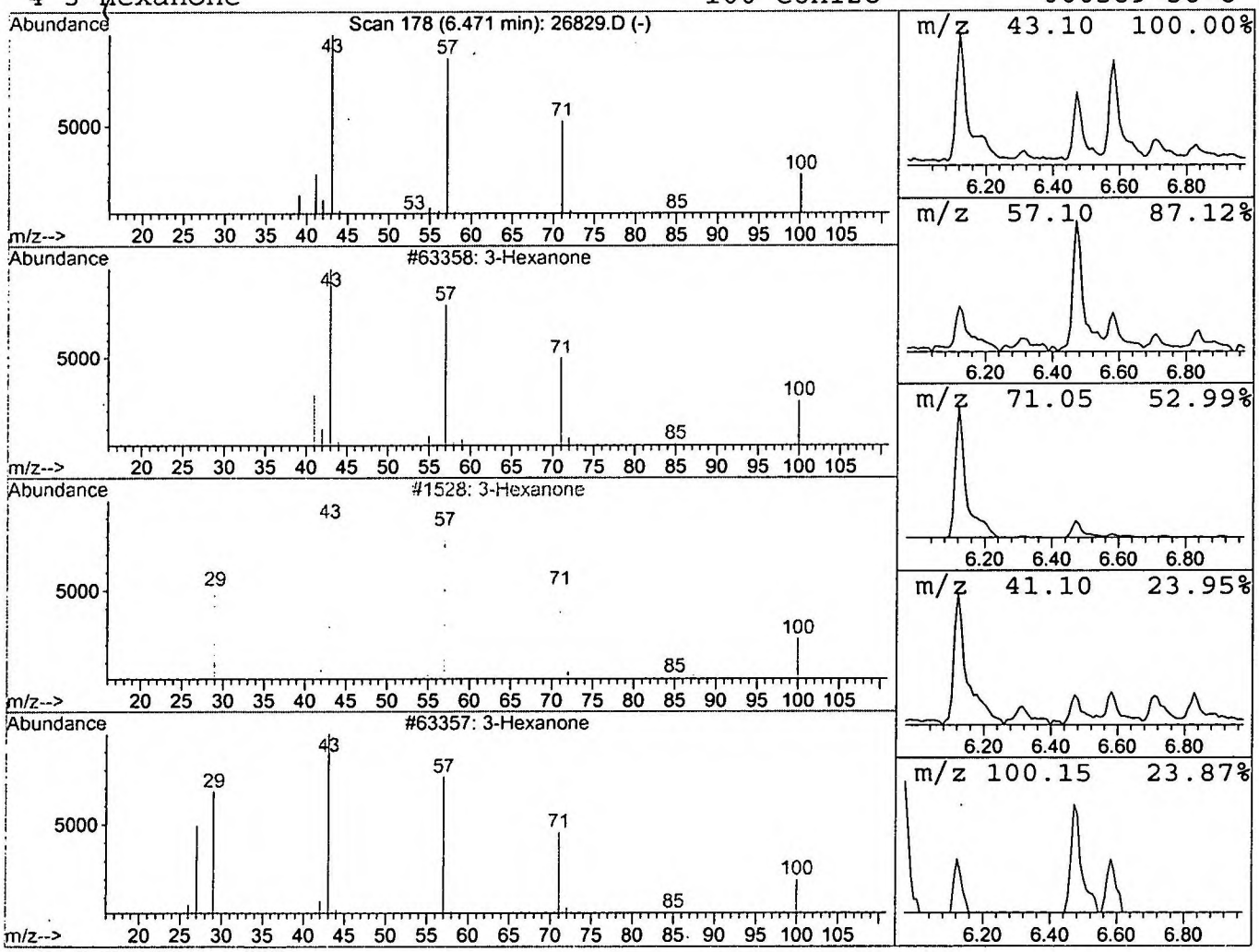
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 3-Hexanone — Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	0.52 ng/uL	148870	1,4-Dichlorobenzene-d4	11.89

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	90
3	3-Hexanone	100	C6H12O	000589-38-8	90
4	3-Hexanone	100	C6H12O	000589-38-8	72



Library Search Compound Report

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Acq On : 6 Dec 2001 2:44 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

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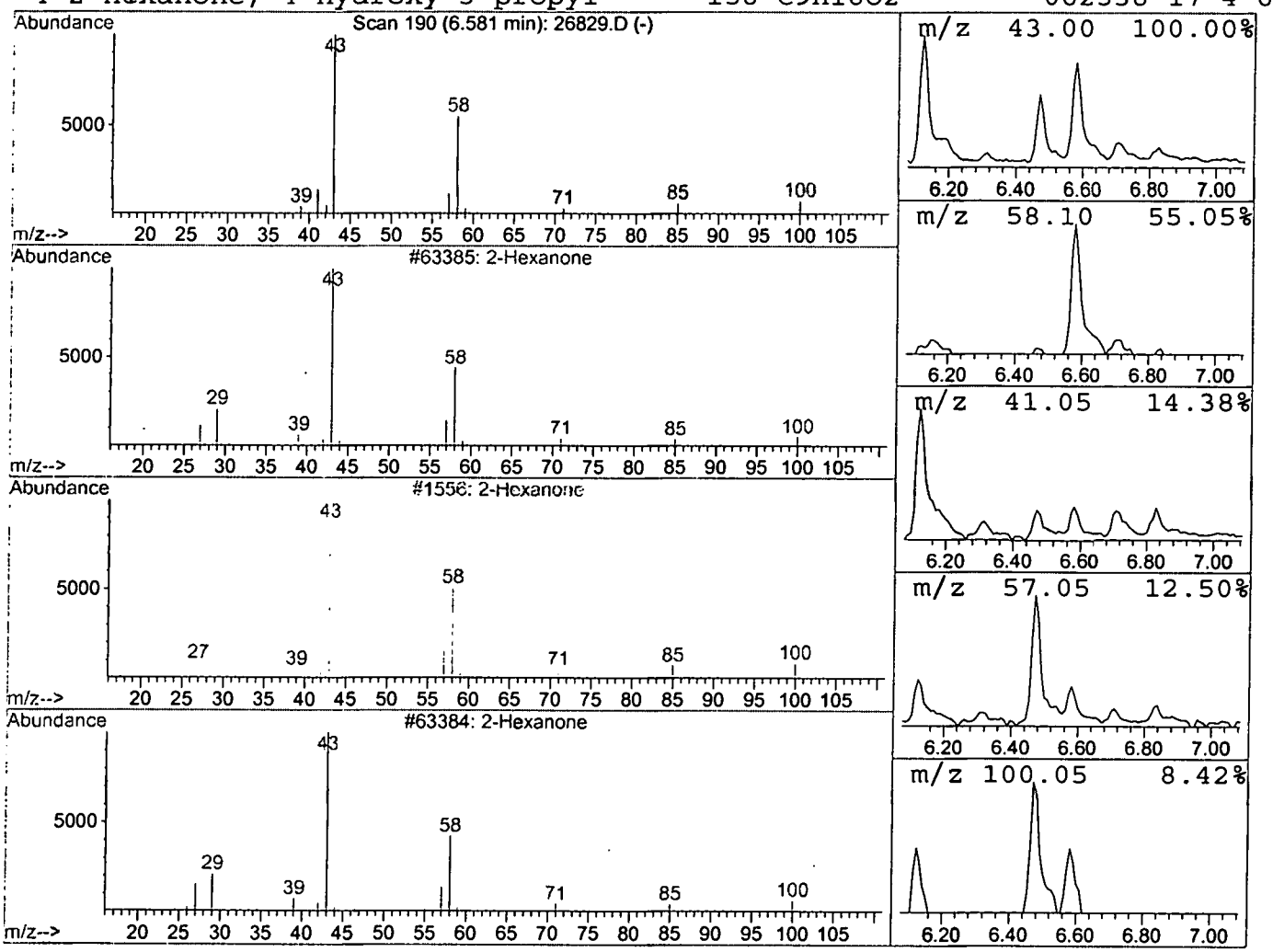
Peak Number 7 2-Hexanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	0.55 ng/uL	159430	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	90
2	2-Hexanone	100	C6H12O	000591-78-6	90
3	2-Hexanone	100	C6H12O	000591-78-6	90
4	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	64



Library Search Compound Report

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Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

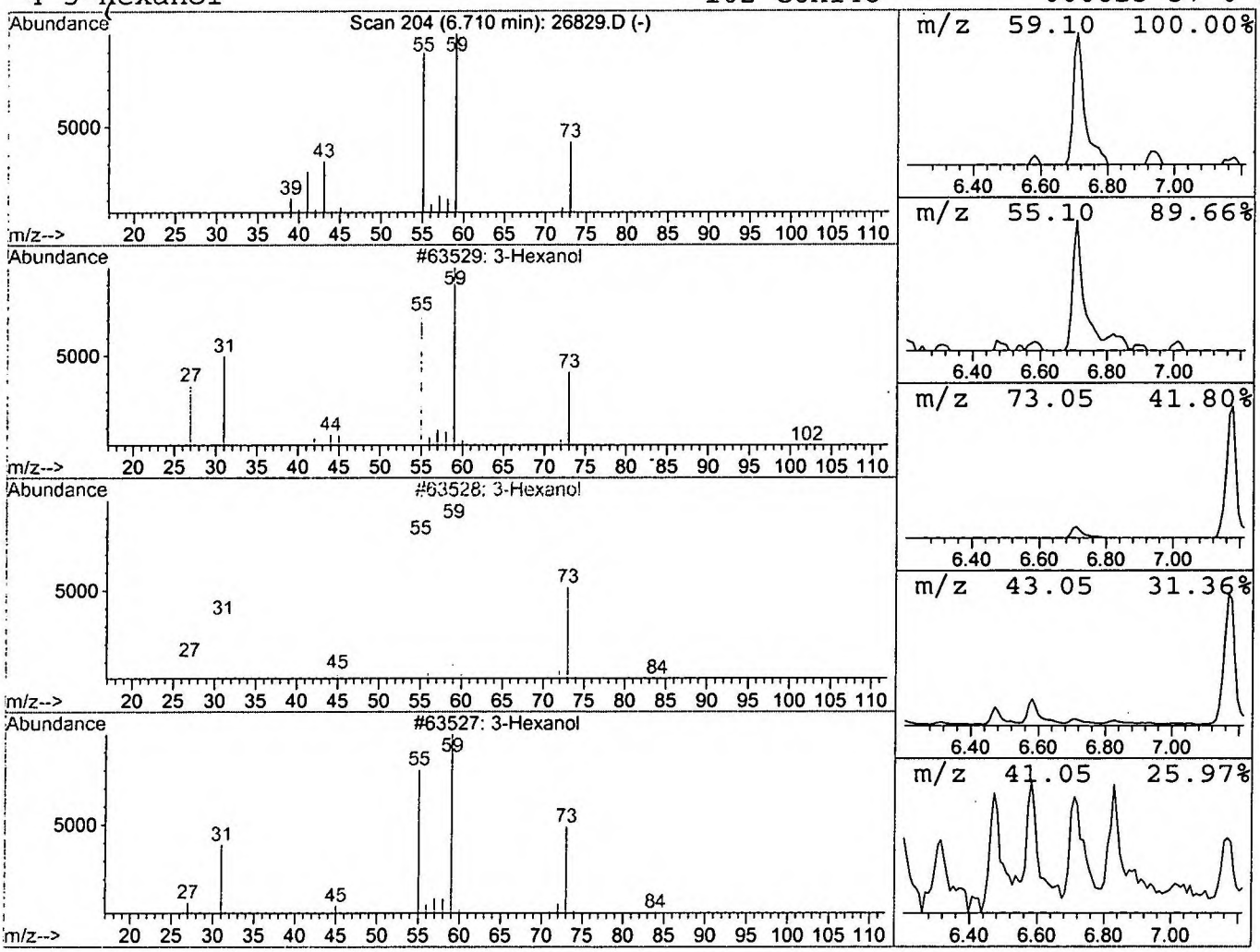
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 3-Hexanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	0.52 ng/uL	148836	1,4-Dichlorobenzene-d4	11.89

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

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D Vial: 6
Acq On : 6 Dec 2001 2:44 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

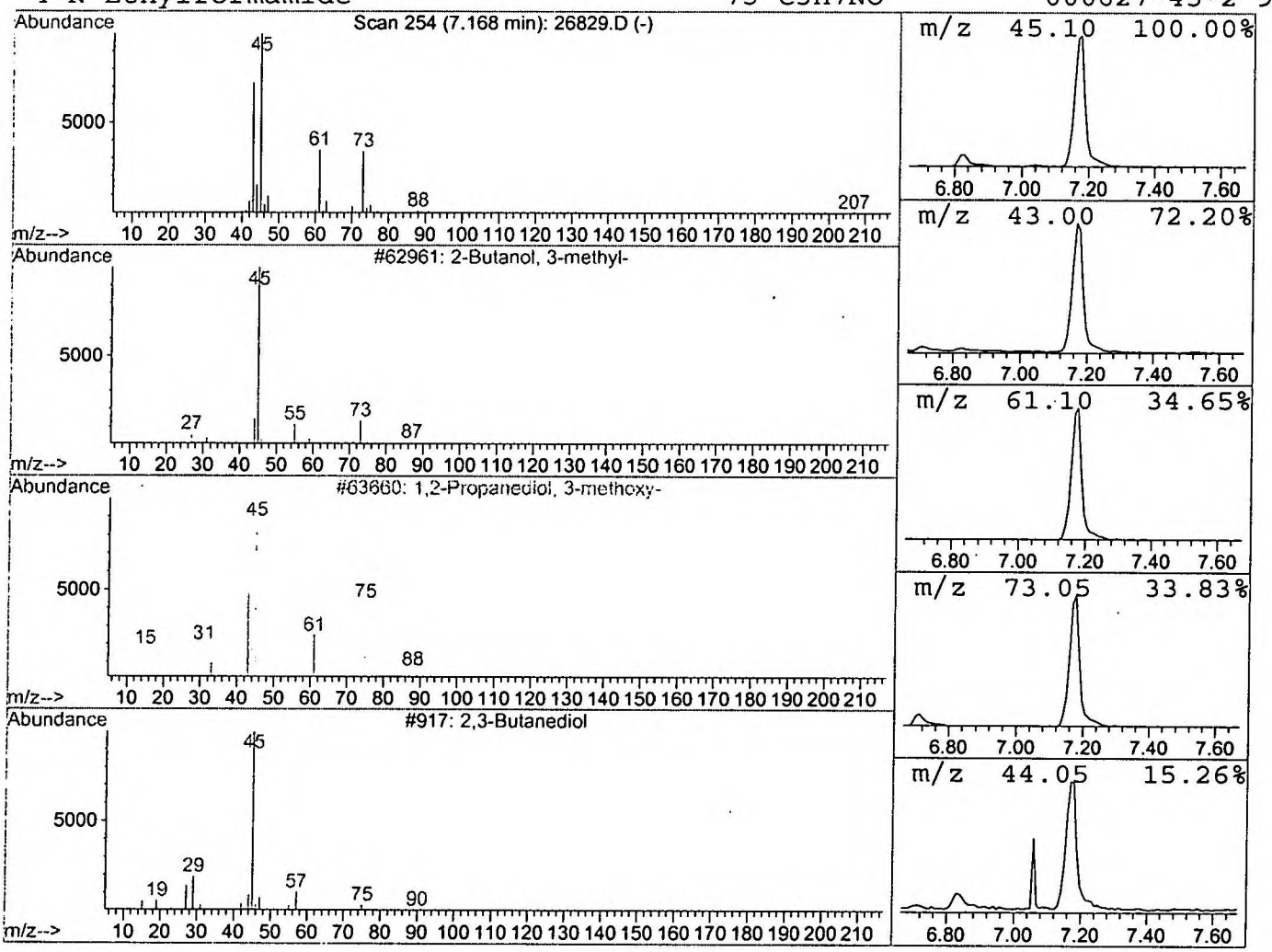
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 2-Butanol, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	5.12 ng/uL	1471640	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	40
2	1,2-Propanediol, 3-methoxy-	106	C4H10O3	000623-39-2	36
3	2,3-Butanediol	90	C4H10O2	000513-85-9	33
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D Vial: 6
Acq On : 6 Dec 2001 2:44 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

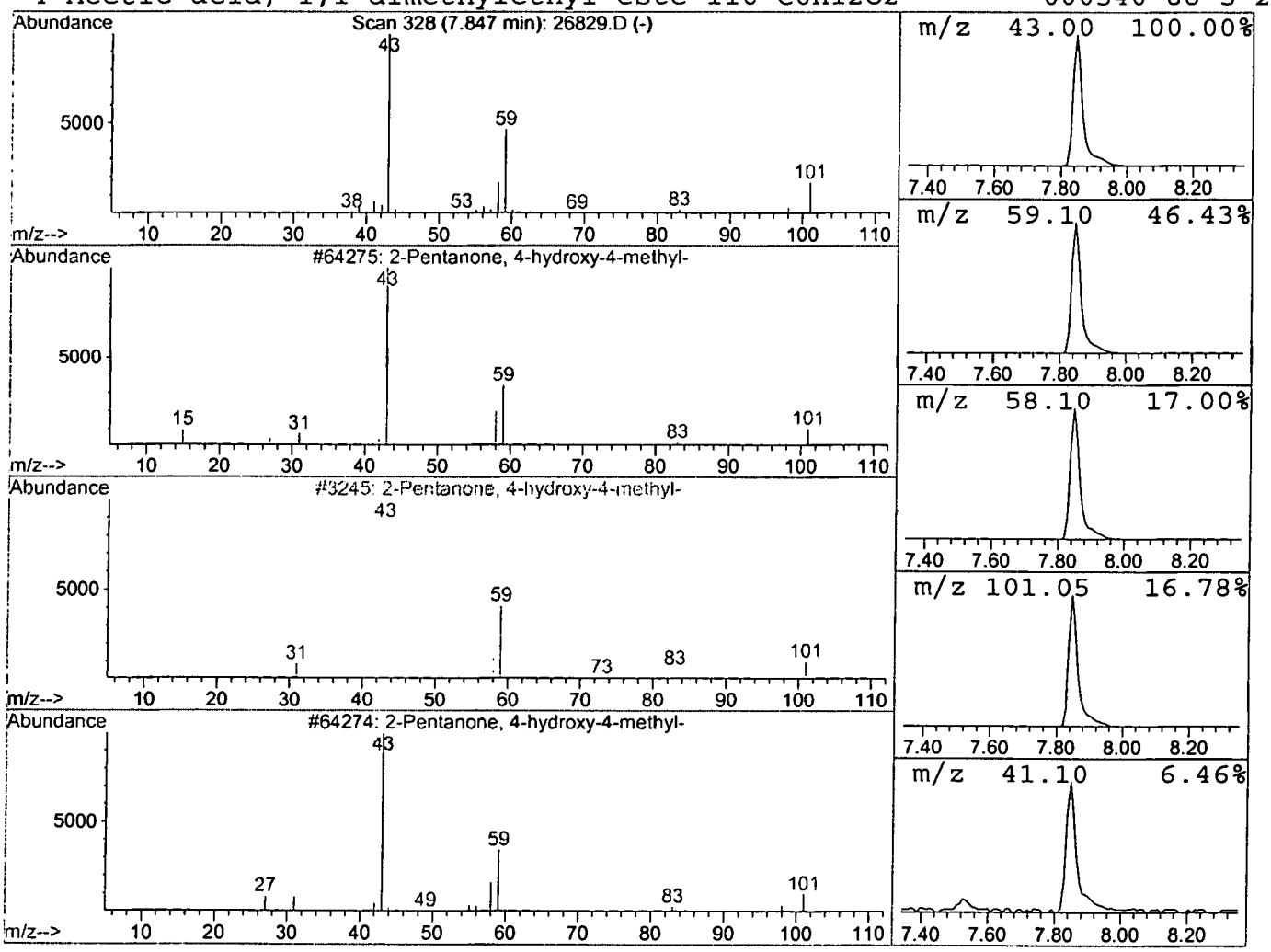
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 10 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	6.60 ng/uL	1899870	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	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	40
4	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28



Library Search Compound Report

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Acq On : 6 Dec 2001 2:44 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

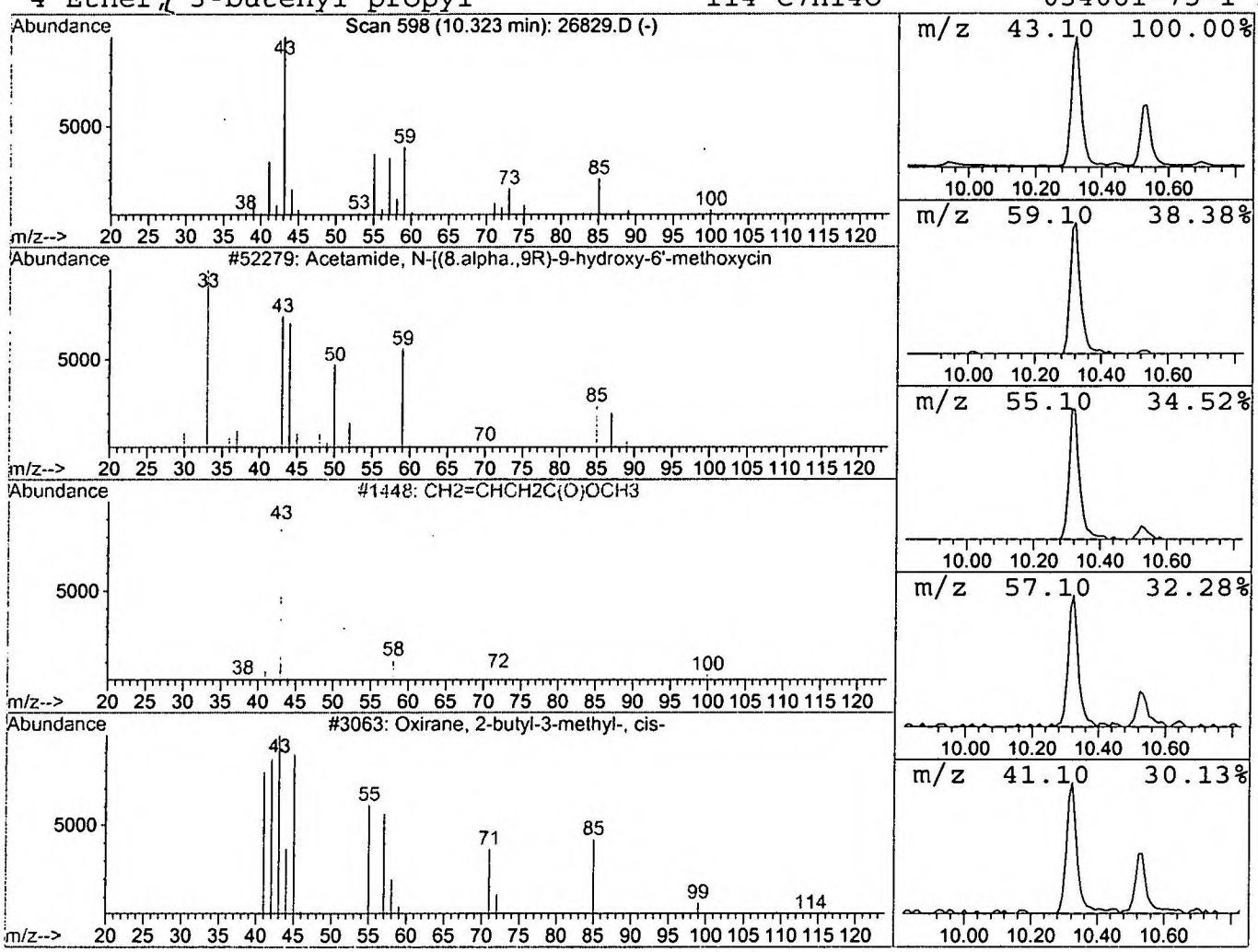
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 11 Acetamide, N-[(8.alpha.,9R)-9- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	2.05 ng/uL	590060	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-[(8.alpha.,9R)-9-hydro	381	C22H27N3O3	056847-12-2	28
2	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	14
3	Oxirane, 2-butyl-3-methyl-, cis-	114	C7H14O	056052-93-8	12
4	Ether, 3-butenyl propyl	114	C7H14O	034061-75-1	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26829.D
Acq On : 6 Dec 2001 2:44 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

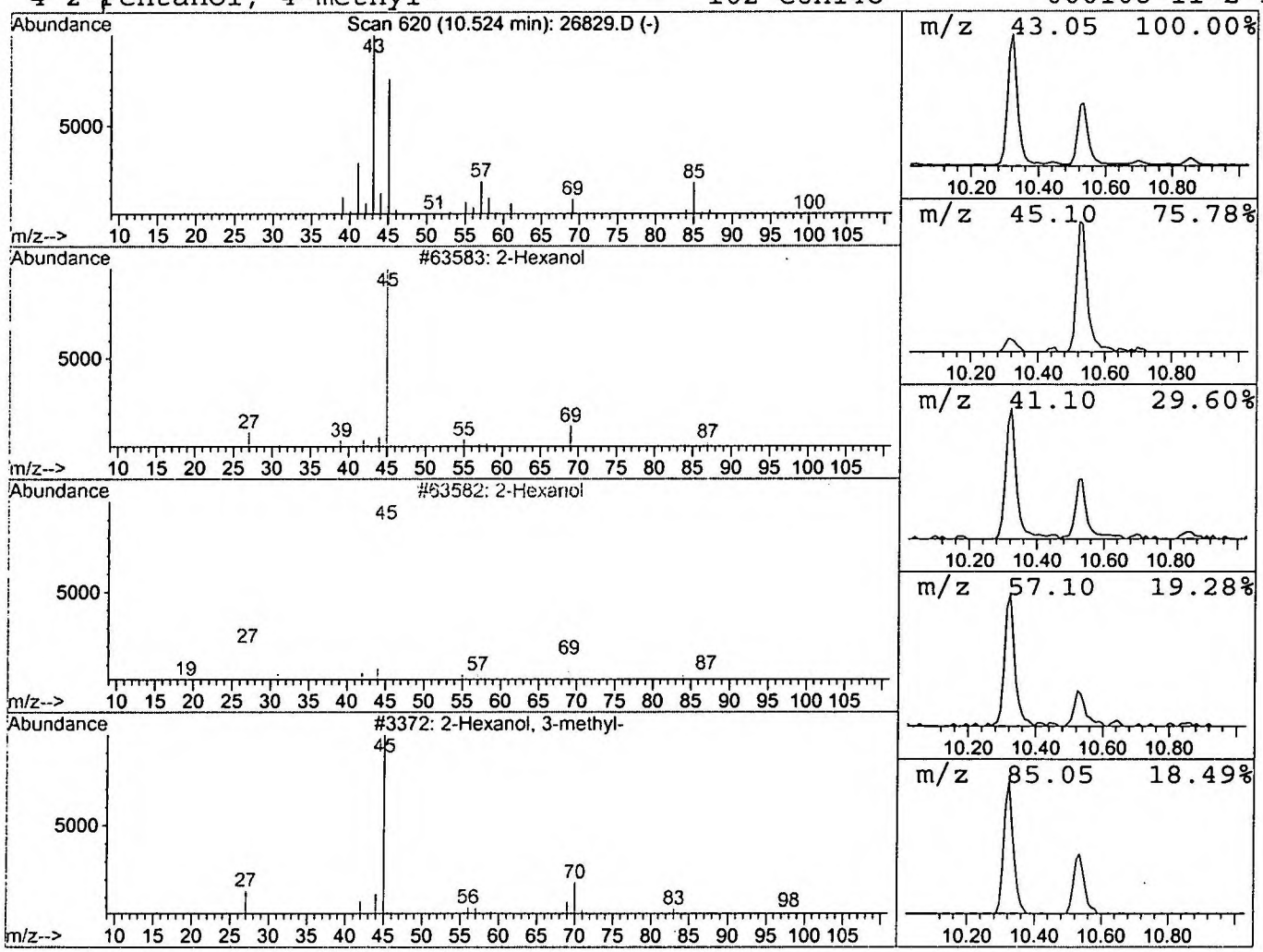
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 12 2-Hexanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.52	0.98 ng/uL	282377	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	50
2	2-Hexanol	102	C6H14O	000626-93-7	40
3	2-Hexanol, 3-methyl-	116	C7H16O	002313-65-7	40
4	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	36



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 6 Dec 2001 2:44 pm
Data File: C:\HPCHEM\1\DATA\DEC0601\26829.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
3,5-Hexadien-2-ol	5.02	1.1	ng/uL	319428	ISTD01	11.89	5753370	20.0
1-Pentanol, 4-methyl	5.12	1.4	ng/uL	414360	ISTD01	11.89	5753370	20.0
Pentane, 2,3-dimethy	5.19	3.0	ng/uL	863314	ISTD01	11.89	5753370	20.0
Oxirane, 2-methyl-3-	5.67	0.5	ng/uL	151713	ISTD01	11.89	5753370	20.0
Propanoic acid, 2-me	6.12	1.8	ng/uL	516307	ISTD01	11.89	5753370	20.0
3-Hexanone	6.47	0.5	ng/uL	148870	ISTD01	11.89	5753370	20.0
2-Hexanone	6.58	0.6	ng/uL	159430	ISTD01	11.89	5753370	20.0
3-Hexanol	6.71	0.5	ng/uL	148836	ISTD01	11.89	5753370	20.0
2-Butanol, 3-methyl-	7.17	5.1	ng/uL	1471640	ISTD01	11.89	5753370	20.0
2-Pentanone, 4-hydro	7.85	6.6	ng/uL	1899870	ISTD01	11.89	5753370	20.0
Acetamide, N-[(8.alp	10.32	2.1	ng/uL	590060	ISTD01	11.89	5753370	20.0
2-Hexanol	10.52	1.0	ng/uL	282377	ISTD01	11.89	5753370	20.0

26829.D NP112701.M Fri Dec 07 14:54:45 2001

Comments for Sample AD26829 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-16-2001 Time: 11:51:03

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 0.50 ug/L.