User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/11/2001 Time: 16:46:41

Sample: AD27009
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/11/01 16:45 Ended: 12/11/01 16:45 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 AD27009 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-11-2001 Time: 16:46:45

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\27009.D Vial: 28
Acq On : 7 Dec 2001 11:38 am Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:49 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	1655286	20.00	ng/uL	-0.04
19) Naphthalene-d8	15.50	136	6353415	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.78	164	2853451	20.00	ng/uL	-0.04
49) Phenanthrene-d10	25.19	188	4162465	20.00	ng/uL	-0.05
59) Chrysene-d12	33.19	240	2768973	20.00	ng/uL	-0.05
68) Perylene-d12	37.28	264	2059989	20.00	ng/uL	-0.05

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	1928	0.02	ng/uL	0.47
Spiked Amount 75.000	Range 34 - 84		Recovery = 0.03%#			
7) 1,2-Dichlorobenzene-d4	12.36	152	275	0.00	ng/uL	-0.10
Spiked Amount 50.000	Range 26 - 73		Recovery = 0.00%#			
20) Nitrobenzene-d5	0.00	82	0	0.00	ng/uL	
Spiked Amount 50.000	Range 55 - 98		Recovery = 0.00%#			
32) 2-Fluorobiphenyl	18.93	172	3469	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

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
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	1388	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	22.26	149	584	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	1130	N.D.	
56)	Anthracene	25.18	178	1130	N.D.	
57)	Di-n-butylphthalate	27.17	149	10678	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

Quantitation Report

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Sample : gc/ms unknown
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:49 2001

Vial: 28
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Title : 208 625/TCLP calibration table
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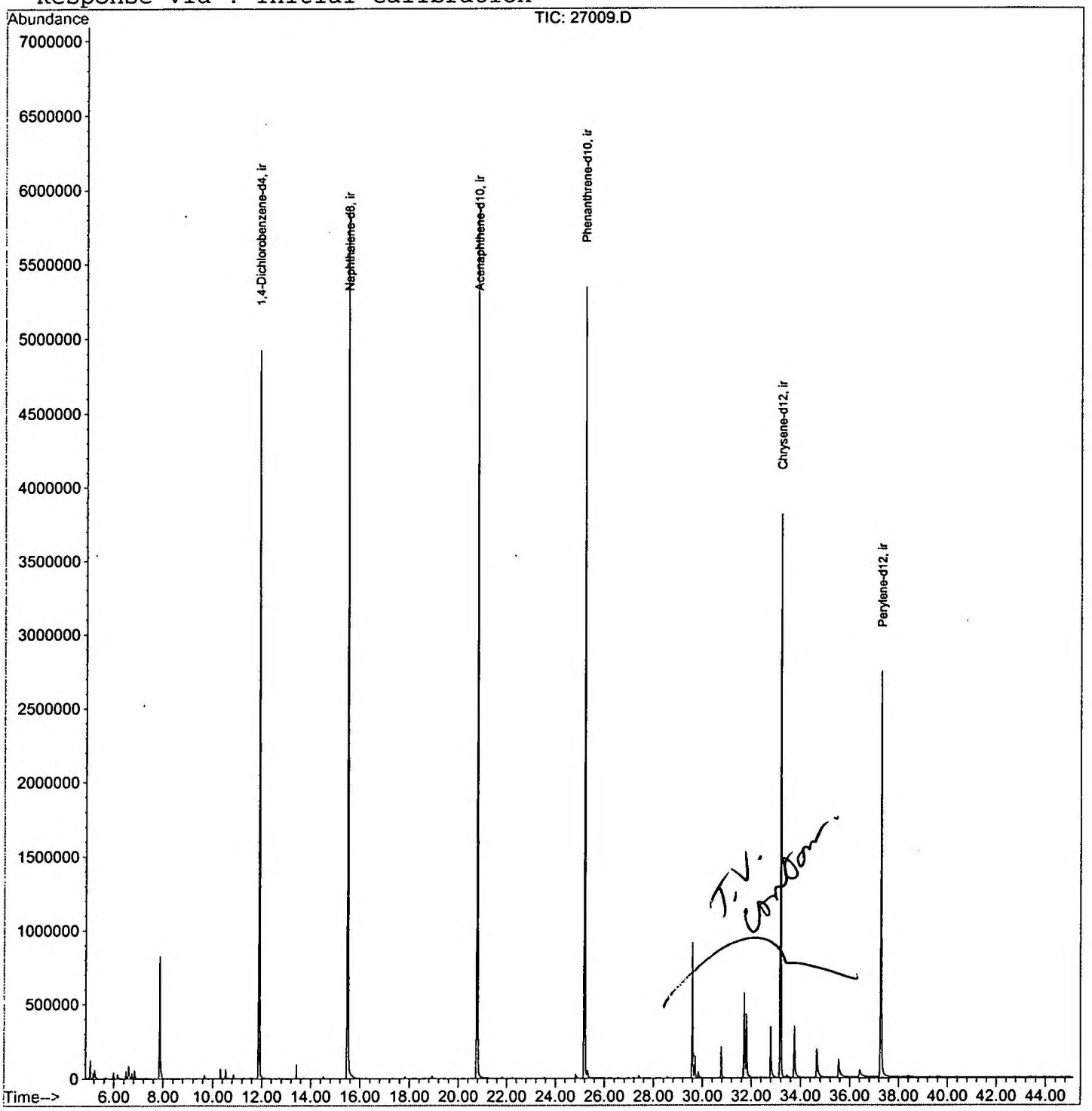


LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC0601\27009.D Vial: 28
Acq On : 7 Dec 2001 11:38 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

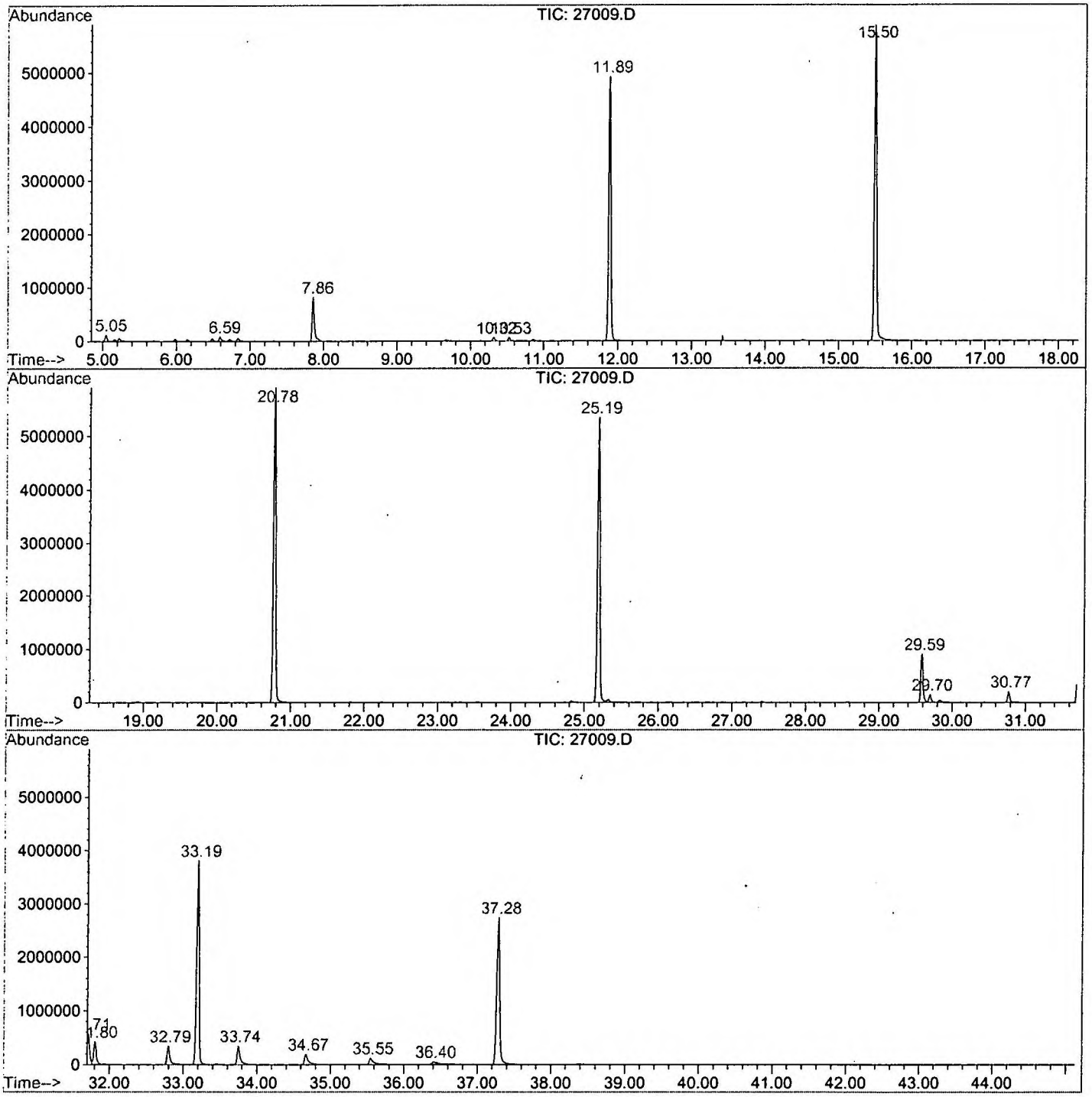
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.052	17	23	27	rBV	122523	227146	1.77%	0.305%
2	6.592	187	191	201	rVV	85674	187507	1.46%	0.251%
3	7.858	325	329	347	rVB	823116	1758532	13.68%	2.357%
4	10.324	594	598	606	rVB	66422	129658	1.01%	0.174%
5	10.526	614	620	629	rBV	65361	133336	1.04%	0.179%
6	11.893	763	769	780	rBV	4916146	10174139	79.15%	13.639%
7	15.496	1156	1162	1179	rBV	5875006	12854050	100.00%	17.232%
8	20.778	1731	1738	1755	rBV	5909769	12558666	97.70%	16.836%
9	25.189	2212	2219	2230	rBV2	5344975	11655117	90.67%	15.624%
10	29.591	2693	2699	2707	rBV	913171	1896884	14.76%	2.543%
11	29.701	2707	2711	2719	rVV	151008	323601	2.52%	0.434%
12	30.774	2823	2828	2839	rBV	207411	449521	3.50%	0.603%
13	31.709	2925	2930	2936	rBV	573672	1336175	10.39%	1.791%
14	31.801	2936	2940	2951	rVB	417868	1012677	7.88%	1.358%
15	32.791	3042	3048	3061	rBV	340468	907033	7.06%	1.216%
16	33.195	3085	3092	3106	rBV2	3810837	8876435	69.06%	11.899%
17	33.745	3146	3152	3169	rBV	343239	1016132	7.91%	1.362%
18	34.671	3247	3253	3275	rBV	190723	702774	5.47%	0.942%
19	35.551	3343	3349	3363	rBV	120567	480707	3.74%	0.644%
20	36.404	3437	3442	3446	rBV4	49123	146512	1.14%	0.196%
21	37.284	3529	3538	3566	rBV2	2734060	7768863	60.44%	10.415%

Sum of corrected areas: 74595465

27009.D NP112701.M Tue Dec 11 14:59:13 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0601\27009.D
Operator : GALOS
Acquired : 7 Dec 2001 11:38 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 28
Quant File :NP112701.RES (RTE Integrator)



27009.D NP112701.M Tue Dec 11 14:59:17 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\27009.D
Acq On : 7 Dec 2001 11:38 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

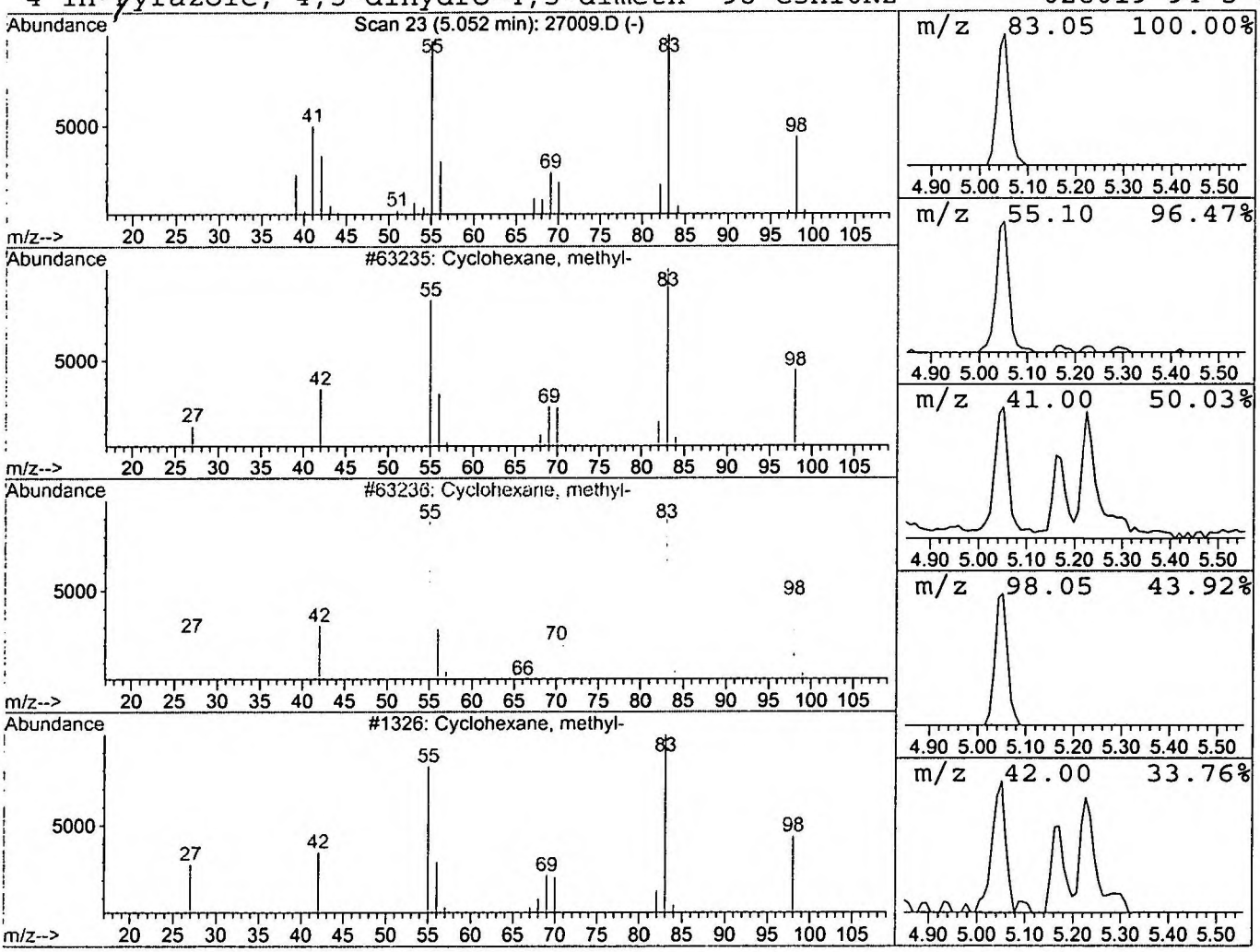
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Cyclohexane, methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	0.45 ng/uL	227146	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, methyl-	98	C7H14	000108-87-2	95
2	Cyclohexane, methyl-	98	C7H14	000108-87-2	95
3	Cyclohexane, methyl-	98	C7H14	000108-87-2	95
4	1H-Pyrazole, 4,5-dihydro-4,5-dimeth	98	C5H10N2	028019-94-5	64



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\27009.D
Acq On : 7 Dec 2001 11:38 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

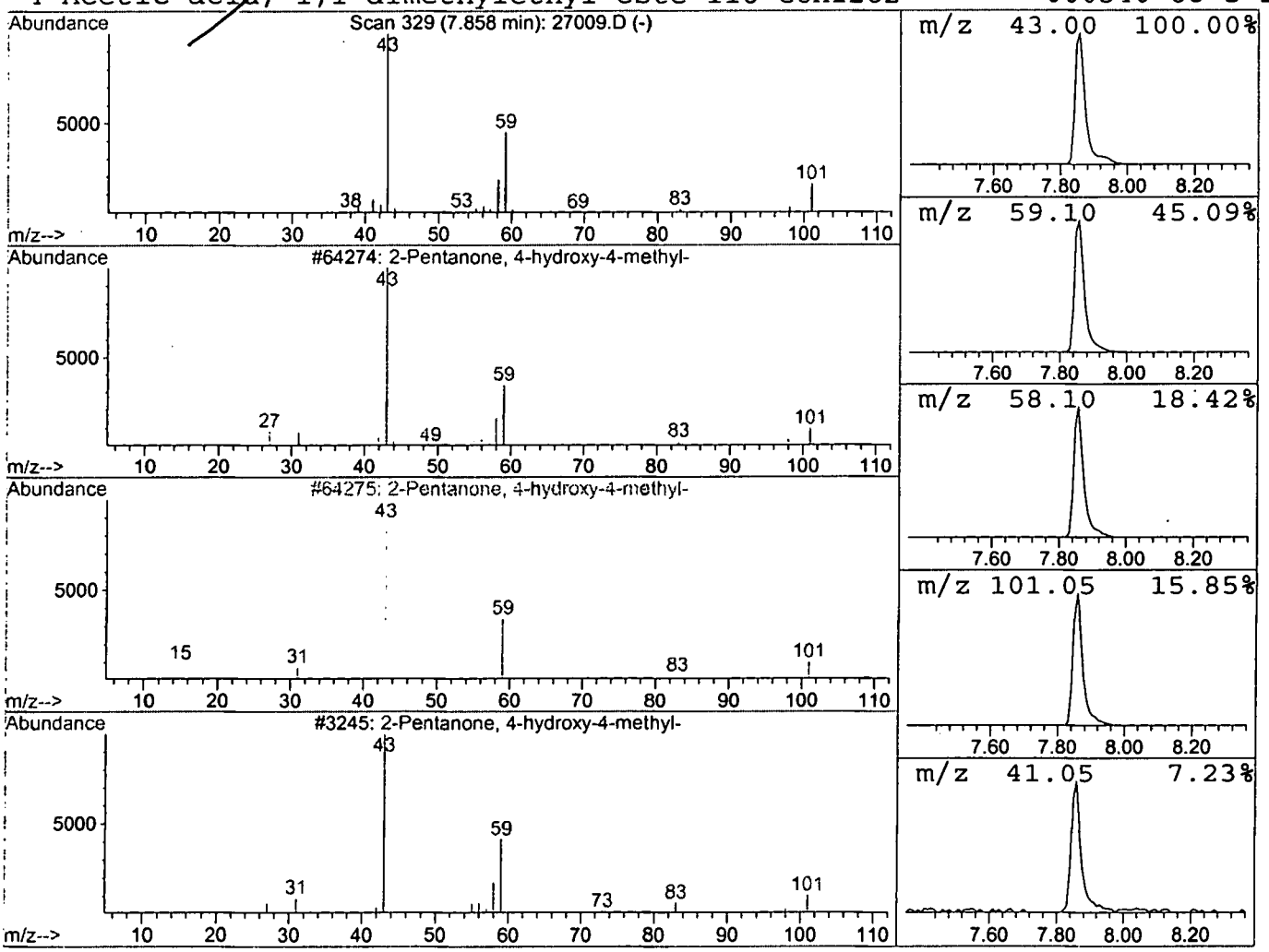
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	3.46 ng/uL	1758530	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	53
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

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Acq On : 7 Dec 2001 11:38 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

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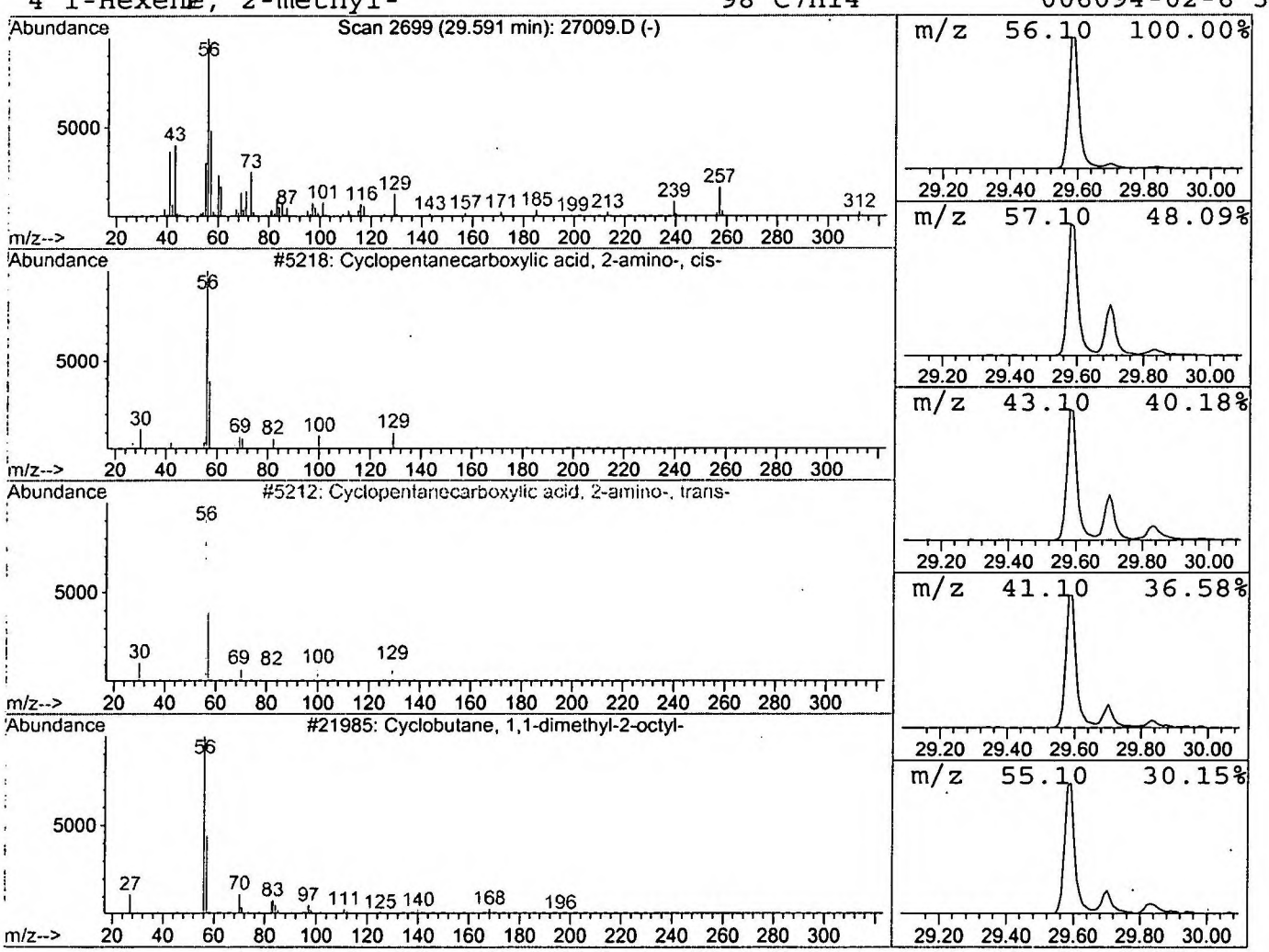
****

Peak Number 3 Cyclopentanecarboxylic acid, 2 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.59	4.27 ng/uL	1896880	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	037910-65-9	43
2	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	040482-05-1	43
3	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	38
4	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	38



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Acq On : 7 Dec 2001 11:38 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

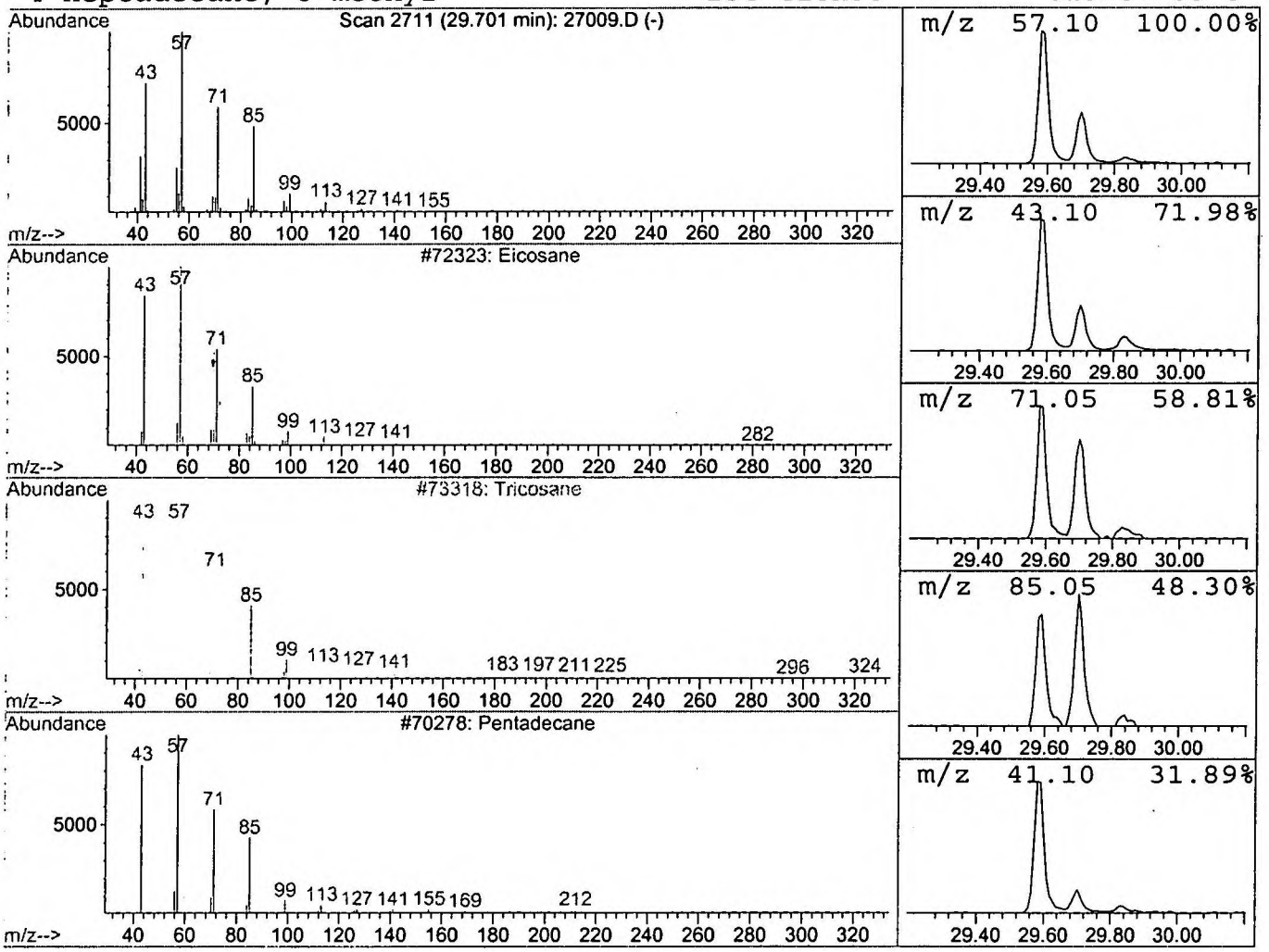
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Eicosane — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.70	0.73 ng/uL	323601	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	83
2	Tricosane	324	C23H48	000638-67-5	83
3	Pentadecane	212	C15H32	000629-62-9	80
4	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	72



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Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

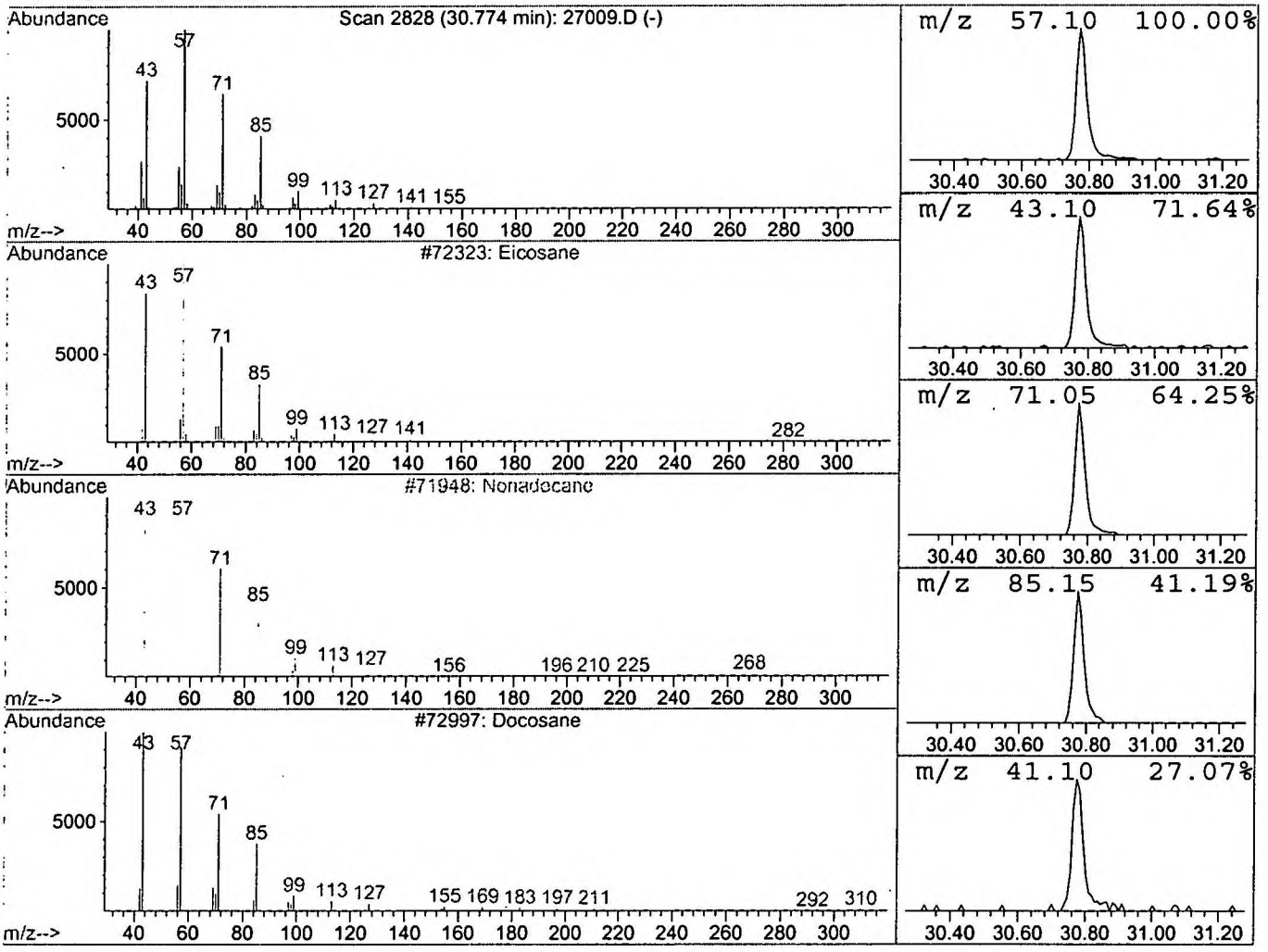
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Eicosane — Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.77	1.01 ng/uL	449521	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Nonadecane	268	C19H40	000629-92-5	90
3	Docosane	310	C22H46	000629-97-0	90
4	Tetracosane	338	C24H50	000646-31-1	87



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

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

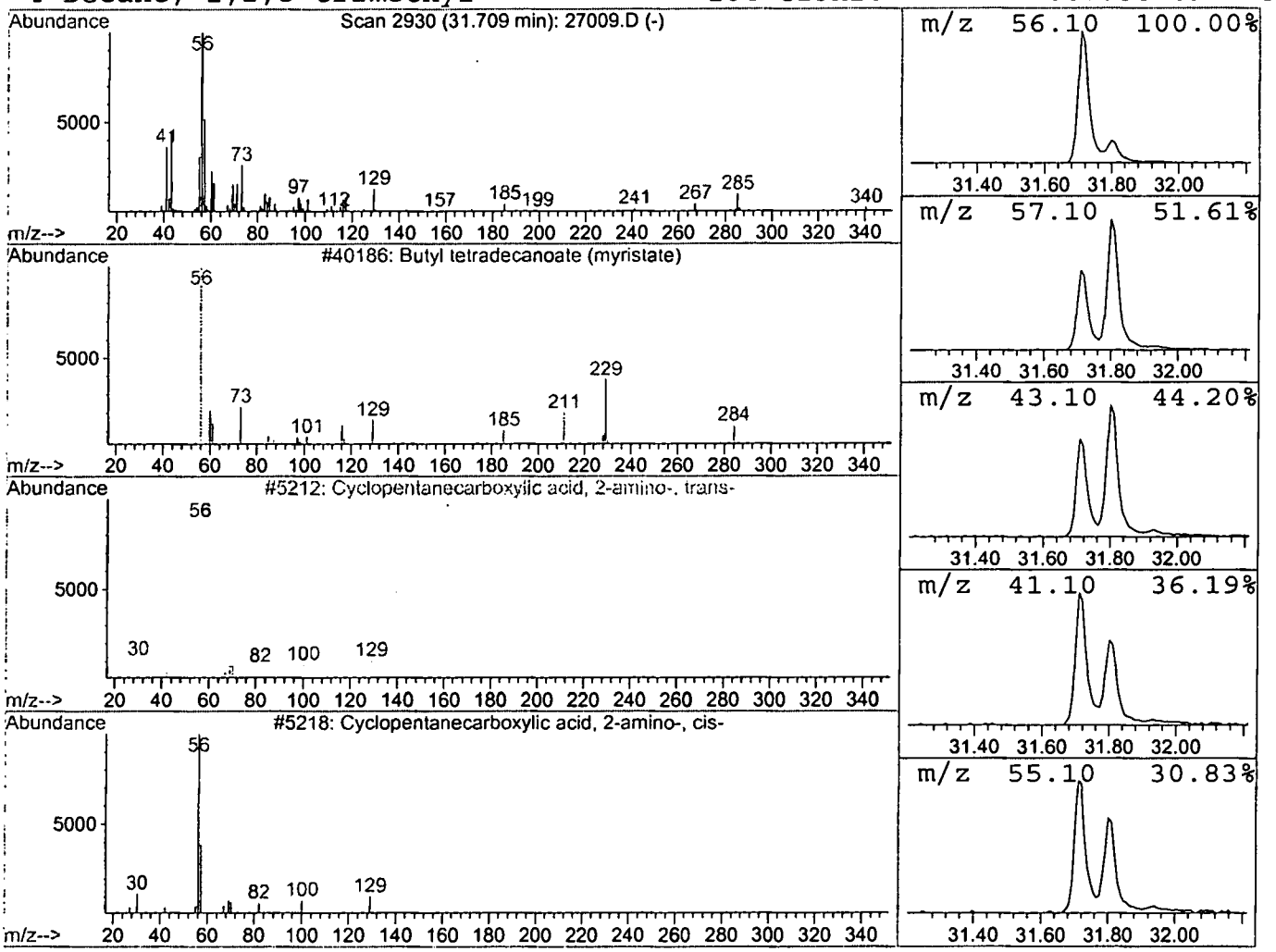
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Butyl tetradecanoate (myristat Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.71	3.01 ng/uL	1336180	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	80
2	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	040482-05-1	43
3	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	037910-65-9	43
4	Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	38



Library Search Compound Report

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

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

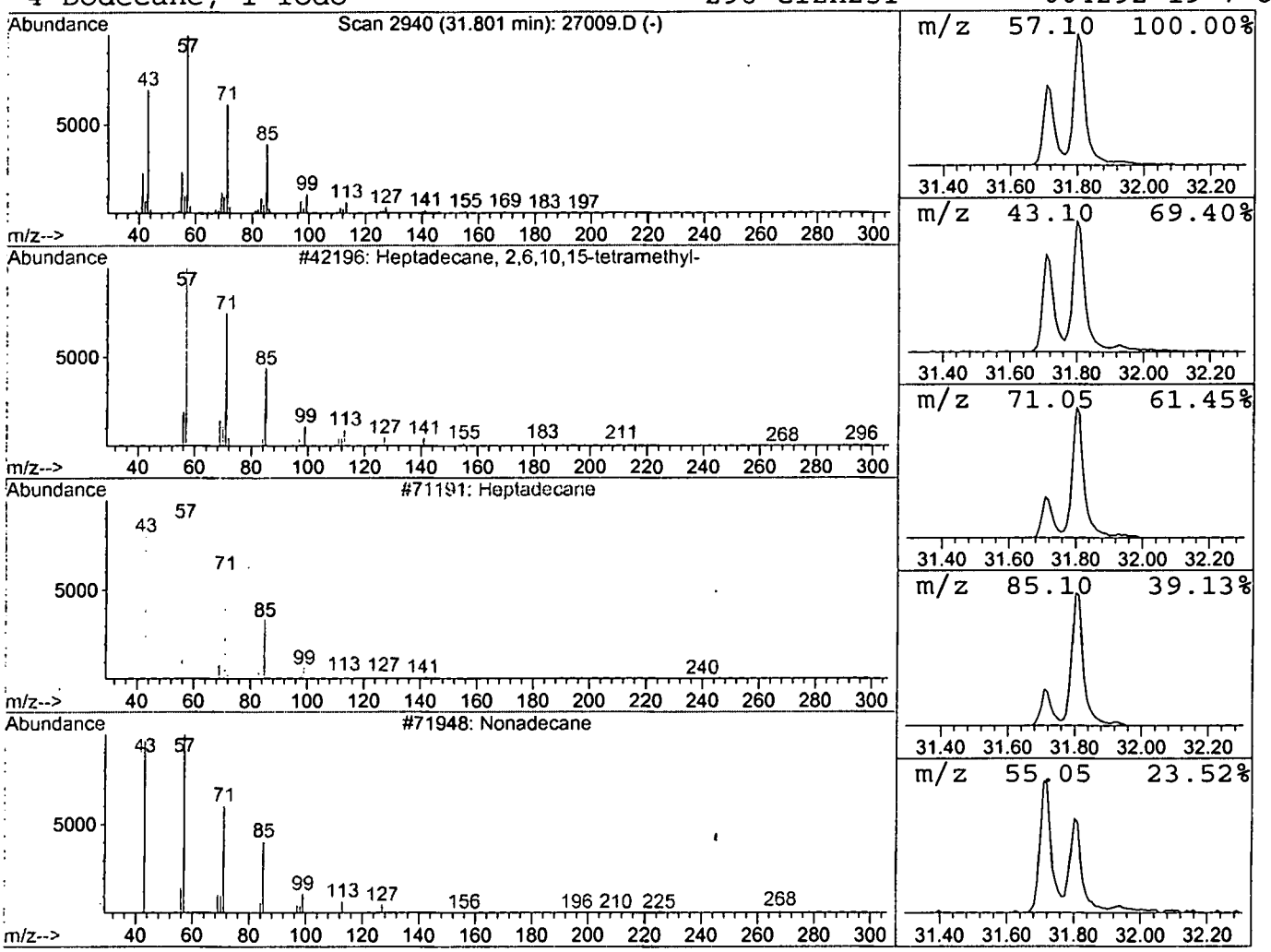
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Heptadecane, 2,6,10,15-tetrame Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.80	2.28 ng/uL	1012680	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
2	Heptadecane	240	C17H36	000629-78-7	86
3	Nonadecane	268	C19H40	000629-92-5	86
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86



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Vial: 28
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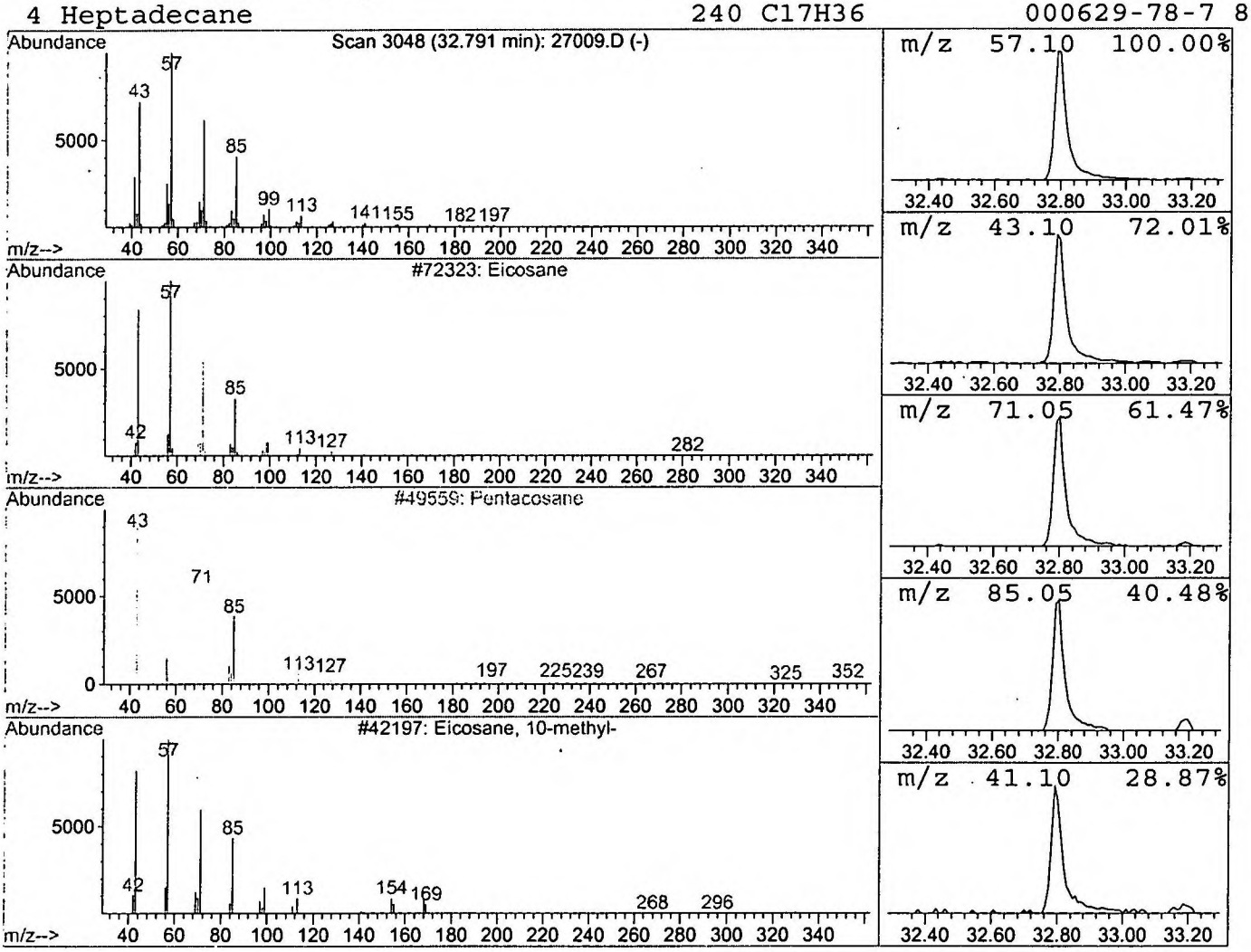
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Eicosane — Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.79	2.04 ng/uL	907033	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Pentacosane	352	C25H52	000629-99-2	90
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	86
4	Heptadecane	240	C17H36	000629-78-7	86



Library Search Compound Report

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Acq On : 7 Dec 2001 11:38 am
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Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

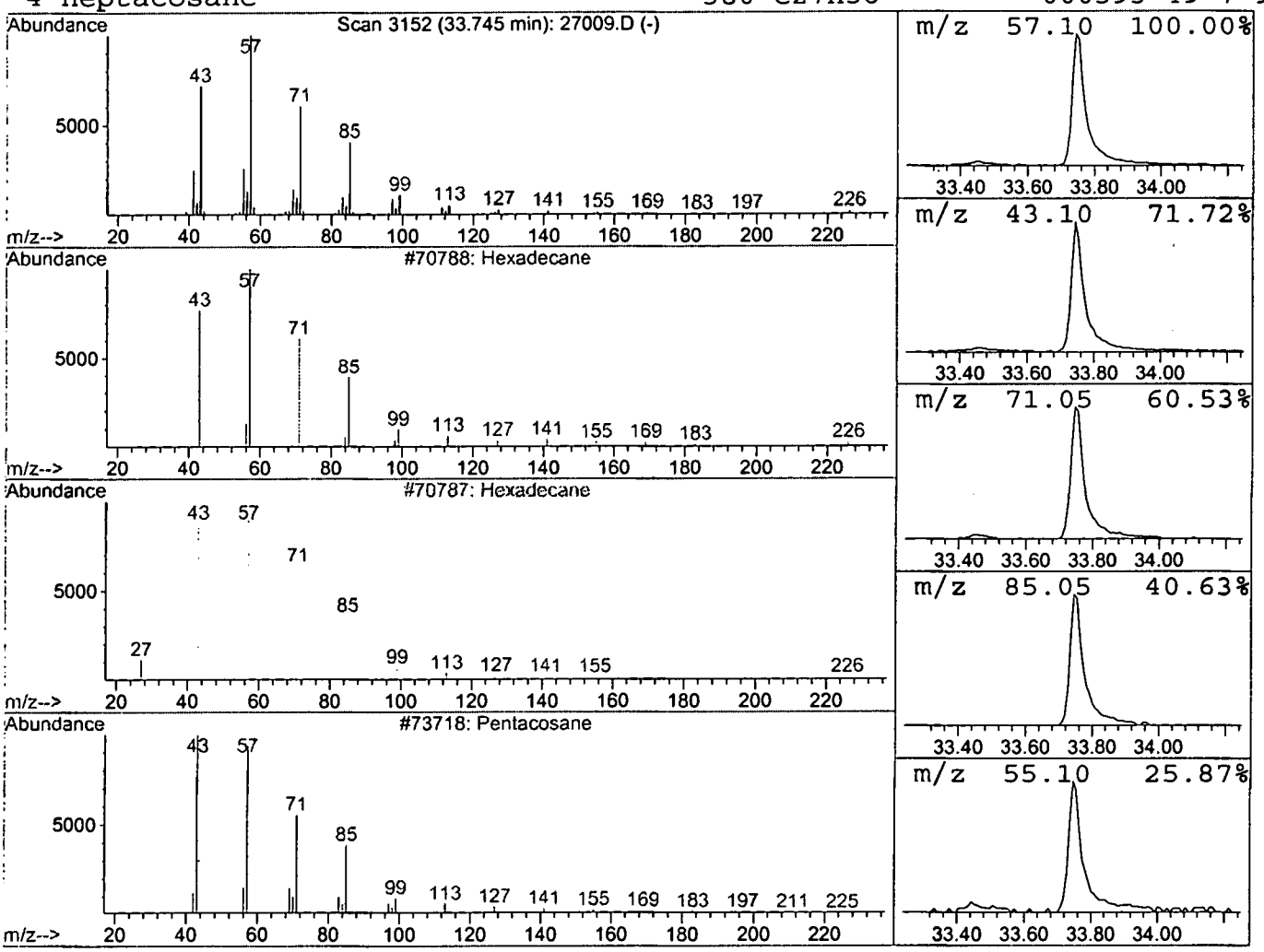
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 Hexadecane — Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.74	2.29 ng/uL	1016130	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Hexadecane	226	C16H34	000544-76-3	91
3	Pentacosane	352	C25H52	000629-99-2	91
4	Heptacosane	380	C27H56	000593-49-7	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\27009.D
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Misc :
MS Integration Params: LSCINT.P

Vial: 28
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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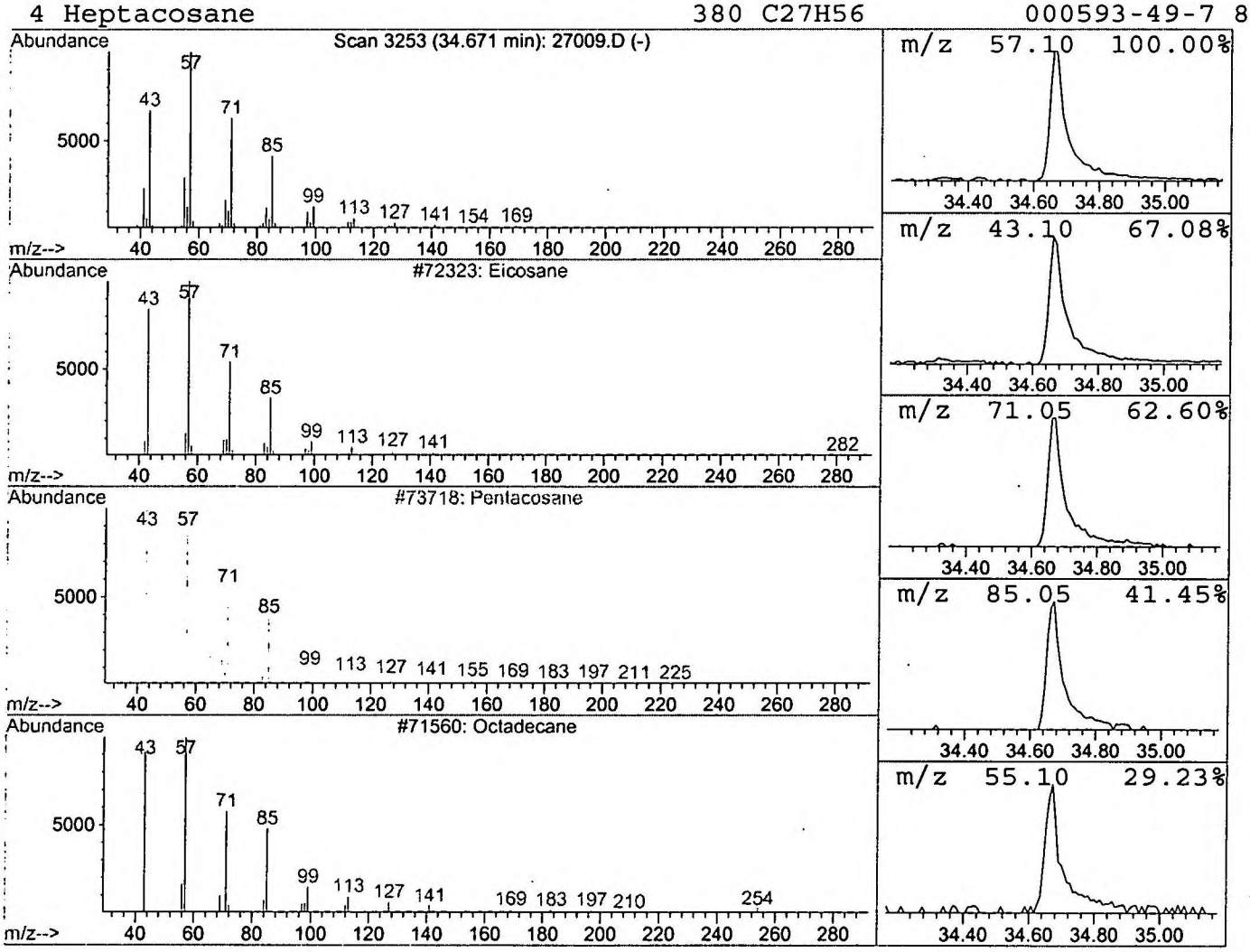
Peak Number 10 Eicosane — Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.67	1.58 ng/uL	702774	Chrysene-d12	33.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	87
2	Pentacosane	352	C25H52	000629-99-2	87
3	Octadecane	254	C18H38	000593-45-3	86
4	Heptacosane	380	C27H56	000593-49-7	86



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\27009.D Vial: 28
Acq On : 7 Dec 2001 11:38 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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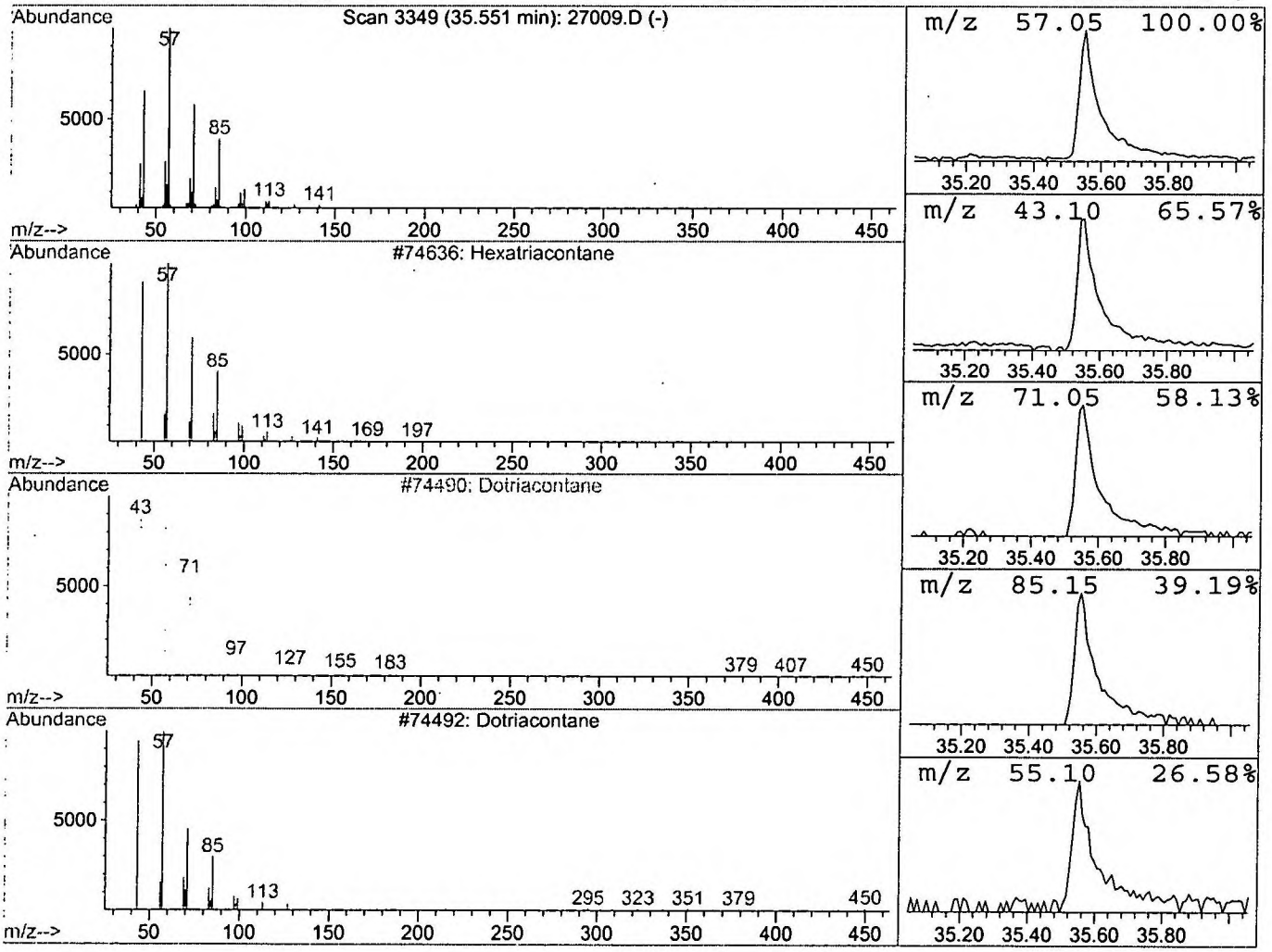
Peak Number 11 Hexatriacontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.55	1.24 ng/uL	480707	Perylene-d12	37.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	91
2	Dotriacontane	451	C32H66	000544-85-4	86
3	Dotriacontane	451	C32H66	000544-85-4	86
4	Hexatriacontane	507	C36H74	000630-06-8	78



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 7 Dec 2001 11:38 am
Data File: C:\HPCHEM\1\DATA\DEC0601\27009.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	0.4	ng/uL	227146	ISTD01	11.89	10174100	20.0
2-Pentanone, 4-hydro	7.86	3.5	ng/uL	1758530	ISTD01	11.89	10174100	20.0
Cyclopentanecarboxyl	29.59	4.3	ng/uL	1896880	ISTD05	33.19	8876440	20.0
Eicosane	29.70	0.7	ng/uL	323601	ISTD05	33.19	8876440	20.0
Eicosane	30.77	1.0	ng/uL	449521	ISTD05	33.19	8876440	20.0
Butyl tetradecanoate	31.71	3.0	ng/uL	1336180	ISTD05	33.19	8876440	20.0
Heptadecane, 2,6,10,	31.80	2.3	ng/uL	1012680	ISTD05	33.19	8876440	20.0
Eicosane	32.79	2.0	ng/uL	907033	ISTD05	33.19	8876440	20.0
Hexadecane	33.74	2.3	ng/uL	1016130	ISTD05	33.19	8876440	20.0
Eicosane	34.67	1.6	ng/uL	702774	ISTD05	33.19	8876440	20.0
Hexatriacontane	35.55	1.2	ng/uL	480707	ISTD06	37.28	7768860	20.0

27009.D NP112701.M Tue Dec 11 14:59:44 2001