User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/11/2001 Time: 16:51:22

Sample: AD26991
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/11/01 16:50 Ended: 12/11/01 16:50 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 AD26991 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-11-2001 Time: 16:51:27

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\26991.D Vial: 17
Acq On : 7 Dec 2001 12:43 am Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:36 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	1053883	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	4014269	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.78	164	1826801	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.19	188	2678688	20.00	ng/uL	-0.05
59) Chrysene-d12	33.18	240	1799078	20.00	ng/uL	-0.06
68) Perylene-d12	37.27	264	1301675	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	725	0.01	ng/uL	0.47
Spiked Amount 75.000	Range 34 - 84		Recovery = 0.01%#			
7) 1,2-Dichlorobenzene-d4	12.22	152	277	0.01	ng/uL	-0.24
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	1529	0.01	ng/uL	0.09
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	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	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	22.27	149	472	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	27.18	149	5865	N.D.	
58)	Fluoranthene	0.00	202	0	N.D.	
60)	3,3'-Dichlorobenzidine	0.00	252	0	N.D.	

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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
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	4250	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.45	149	7918	N.D.	
67)	Chrysene	33.19	228	4250	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	37.28	252	4246	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 : Multiplr: 1.00
MS Integration Params: RTEINT.P
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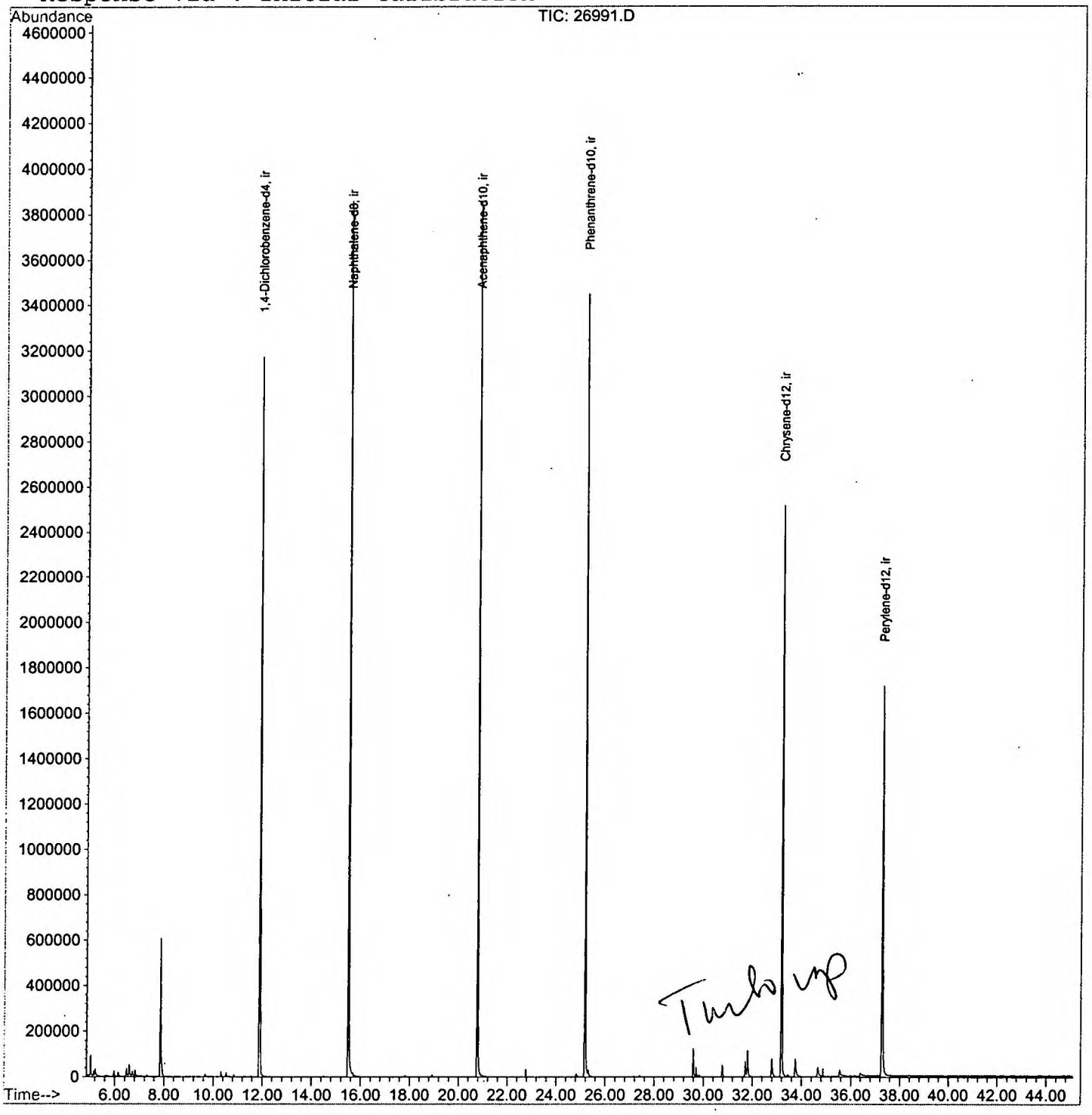
Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26991.D Vial: 17
Acq On : 7 Dec 2001 12:43 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

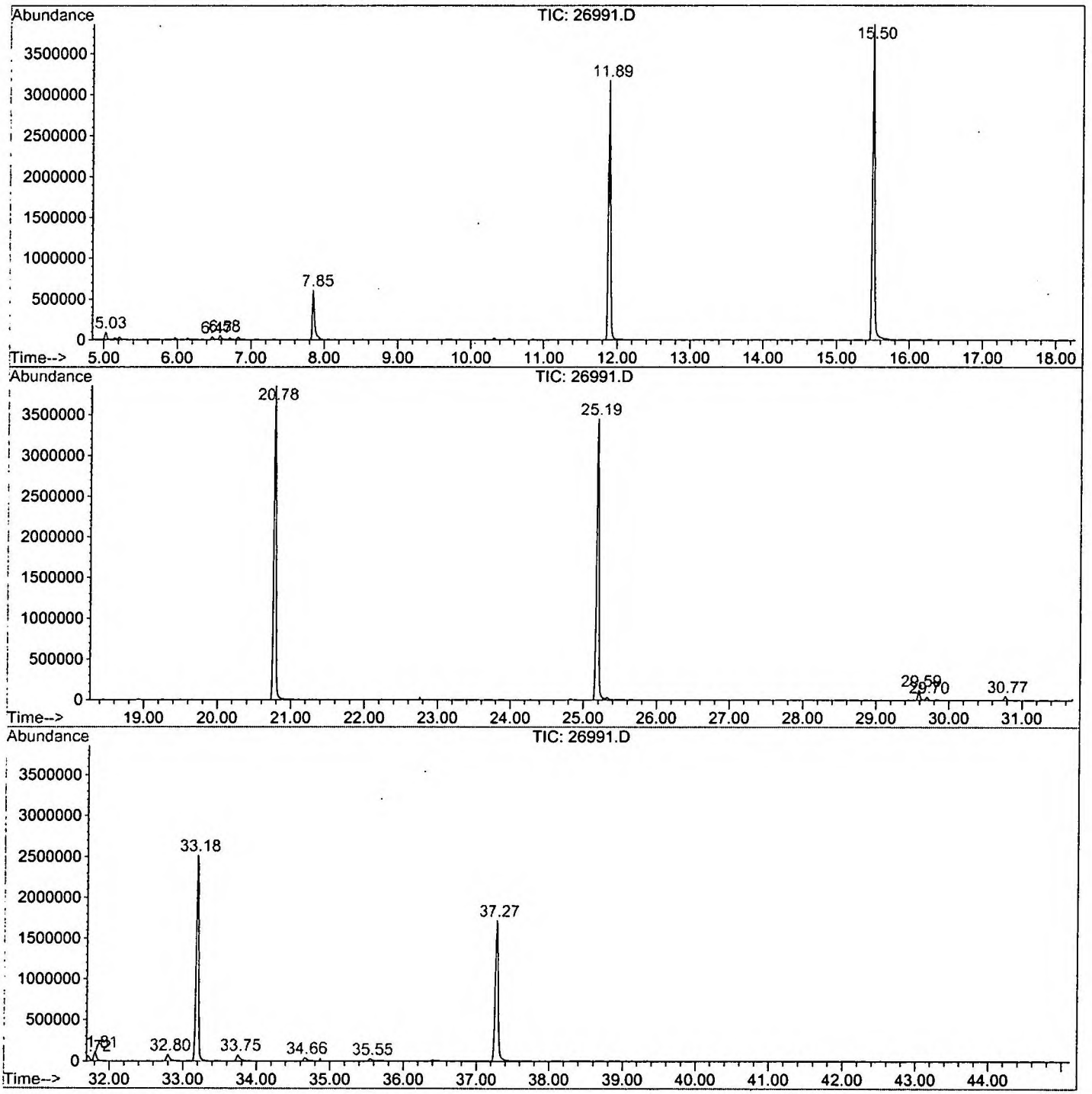
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.032	15	21	26	rBV	89639	168172	2.07%	0.382%
2	6.472	174	178	186	rBV	36416	85105	1.05%	0.193%
3	6.582	186	190	193	rBV	51454	89775	1.11%	0.204%
4	7.847	324	328	343	rVB	606133	1287409	15.88%	2.922%
5	11.891	763	769	787	rBV	3172499	6497666	80.16%	14.747%
6	15.504	1157	1163	1180	rBV	3858779	8105741	100.00%	18.397%
7	20.777	1732	1738	1755	rBV	3856752	8003250	98.74%	18.164%
8	25.188	2212	2219	2231	rBV2	3449650	7519091	92.76%	17.065%
9	29.589	2694	2699	2707	rBV	125635	262827	3.24%	0.597%
10	29.699	2707	2711	2717	rVB	41058	86088	1.06%	0.195%
11	30.772	2823	2828	2838	rVB	51820	120174	1.48%	0.273%
12	31.717	2926	2931	2936	rBV	67674	170499	2.10%	0.387%
13	31.809	2936	2941	2953	rVB	114915	295236	3.64%	0.670%
14	32.799	3043	3049	3060	rBV	81093	241724	2.98%	0.549%
15	33.184	3085	3091	3109	rBV2	2517399	5774343	71.24%	13.105%
16	33.753	3147	3153	3166	rBV	77696	252486	3.11%	0.573%
17	34.661	3247	3252	3264	rBV2	42604	154643	1.91%	0.351%
18	35.550	3342	3349	3361	rBV	30226	114035	1.41%	0.259%
19	37.274	3529	3537	3555	rBV2	1716914	4832745	59.62%	10.968%

Sum of corrected areas: 44061009

26991.D NP112701.M Tue Dec 11 14:54:13 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0601\26991.D
Operator : GALOS
Acquired : 7 Dec 2001 12:43 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 17
Quant File :NP112701.RES (RTE Integrator)



26991.D NP112701.M Tue Dec 11 14:54:16 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26991.D Vial: 17
Acq On : 7 Dec 2001 12:43 am Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

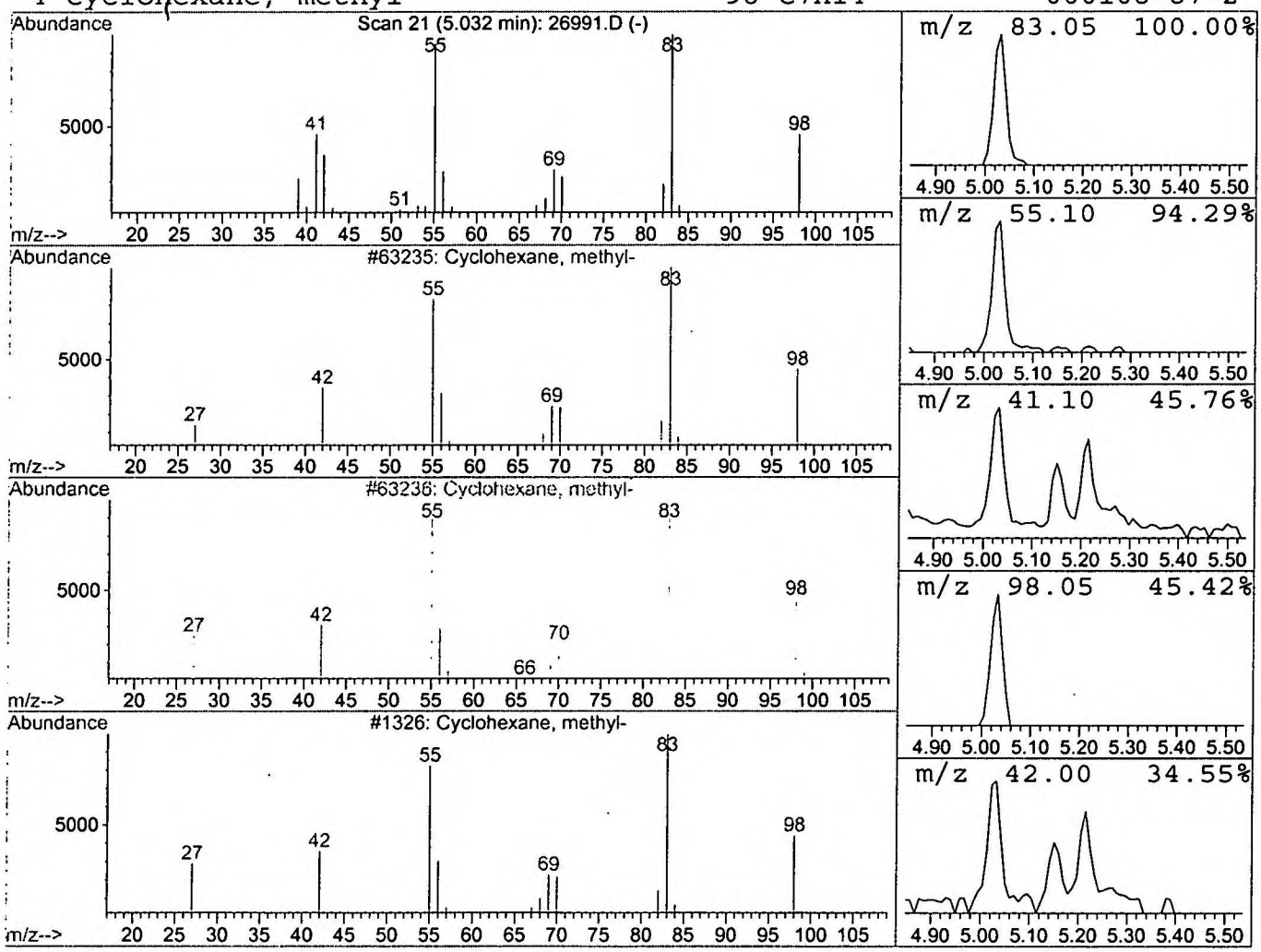
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Cyclohexane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	0.52 ng/uL	168172	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	94
4	Cyclohexane, methyl-	98	C7H14	000108-87-2	78



Library Search Compound Report

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Acq On : 7 Dec 2001 12:43 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 17
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

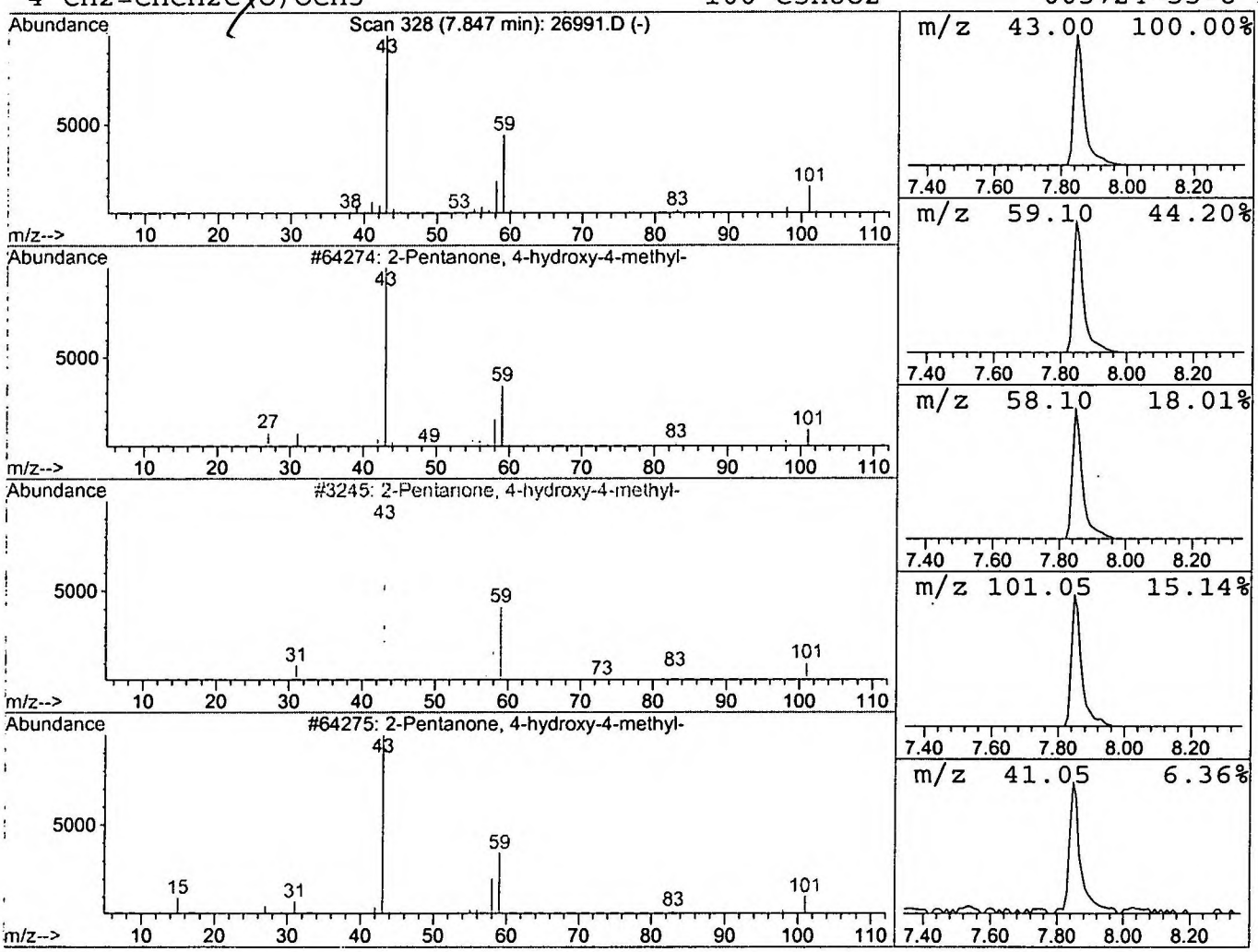
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	3.96 ng/uL	1287410	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	59
4	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	10



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Misc :
MS Integration Params: LSCINT.P

Vial: 17
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

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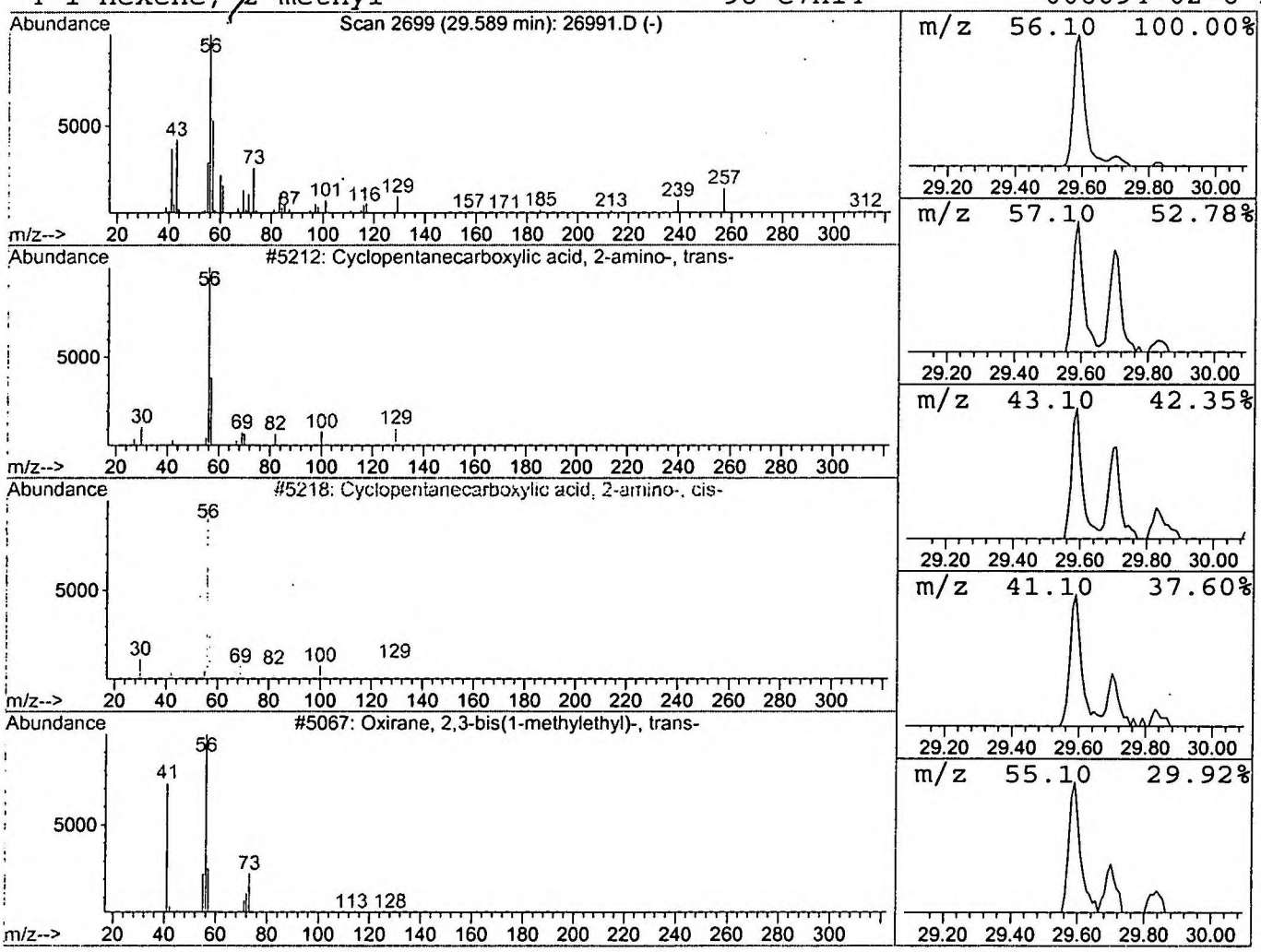
****

Peak Number 3 Cyclopentanecarboxylic acid, 2 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.59	0.91 ng/uL	262827	Chrysene-d12	33.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	040482-05-1	43
2	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	037910-65-9	43
3	Oxirane, 2,3-bis(1-methylethyl)-, t	128	C8H16O	054644-32-5	37
4	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	35



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Misc :
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Vial: 17
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

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Library : C:\DATABASE\NBS75K.L

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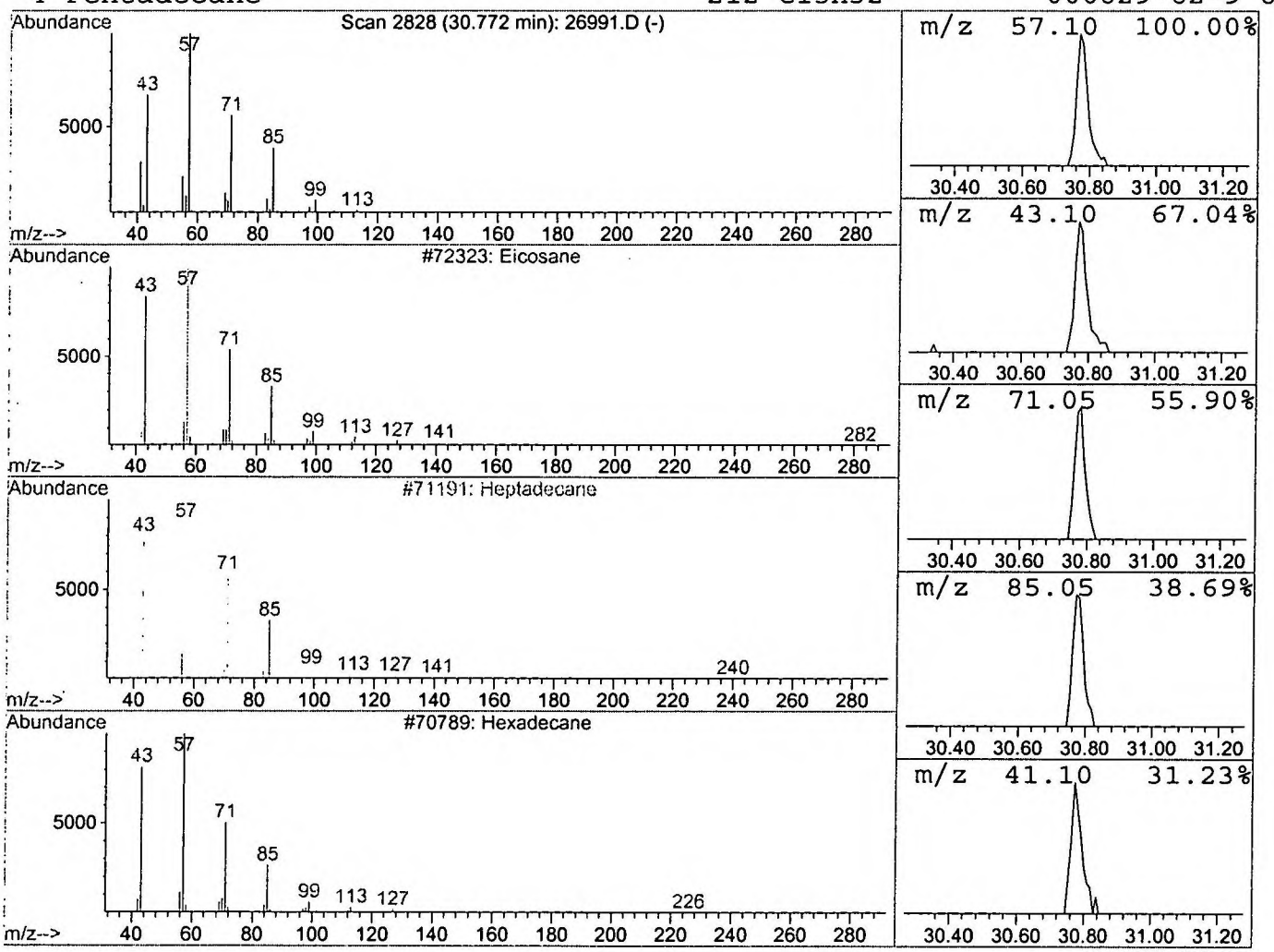
Peak Number 4 Eicosane — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.77	0.42 ng/uL	120174	Chrysene-d12	33.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	90
2	Heptadecane	240	C17H36	000629-78-7	78
3	Hexadecane	226	C16H34	000544-76-3	72
4	Pentadecane	212	C15H32	000629-62-9	64



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

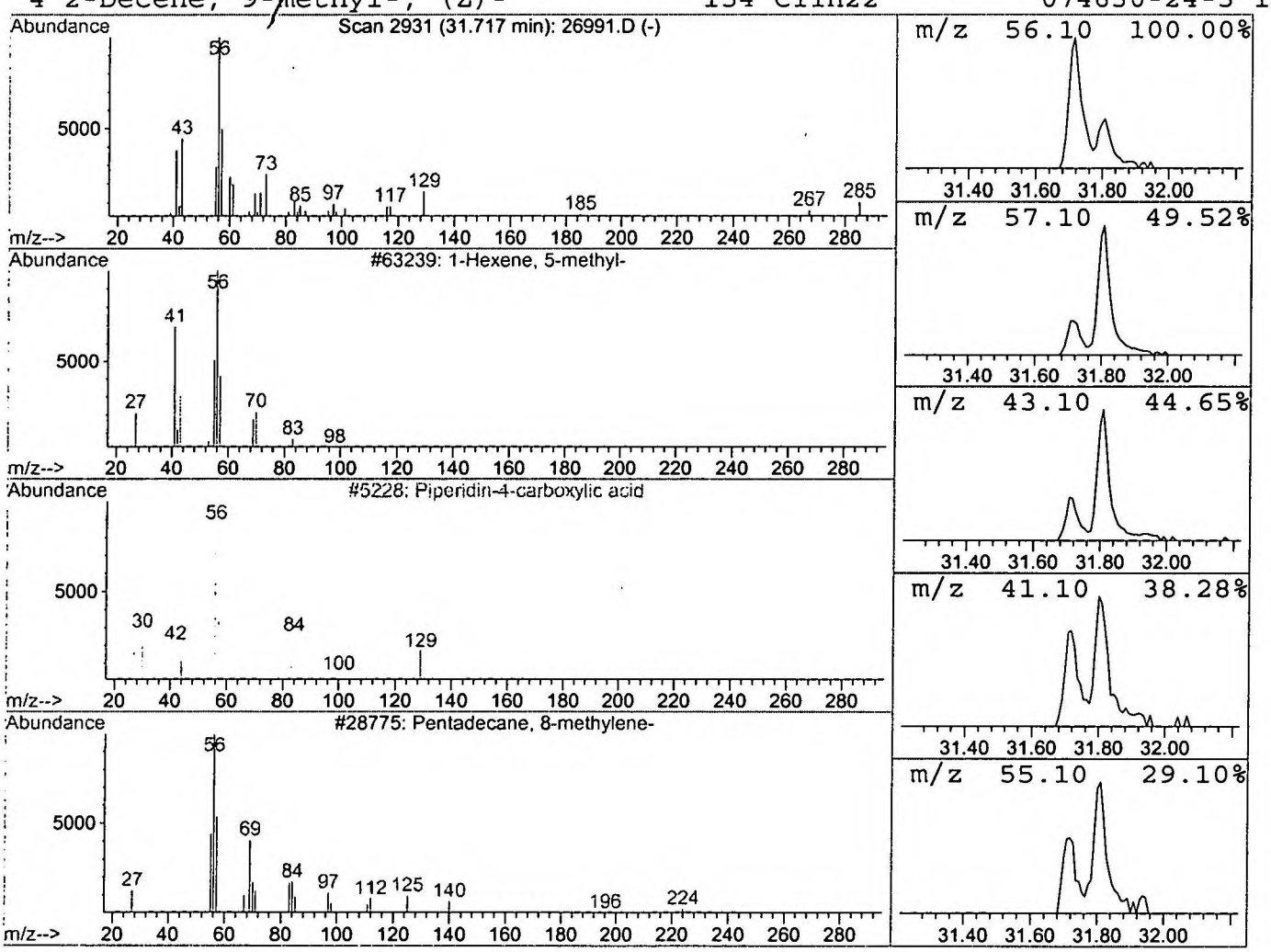
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 1-Hexene, 5-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.72	0.59 ng/uL	170499	Chrysene-d12	33.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene, 5-methyl-	98	C7H14	003524-73-0	27
2	Piperidin-4-carboxylic acid	129	C6H11NO2	000498-94-2	23
3	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	17
4	2-Decene, 9-methyl-, (Z)-	154	C11H22	074630-24-3	17



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Misc :
MS Integration Params: LSCINT.P

Vial: 17
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

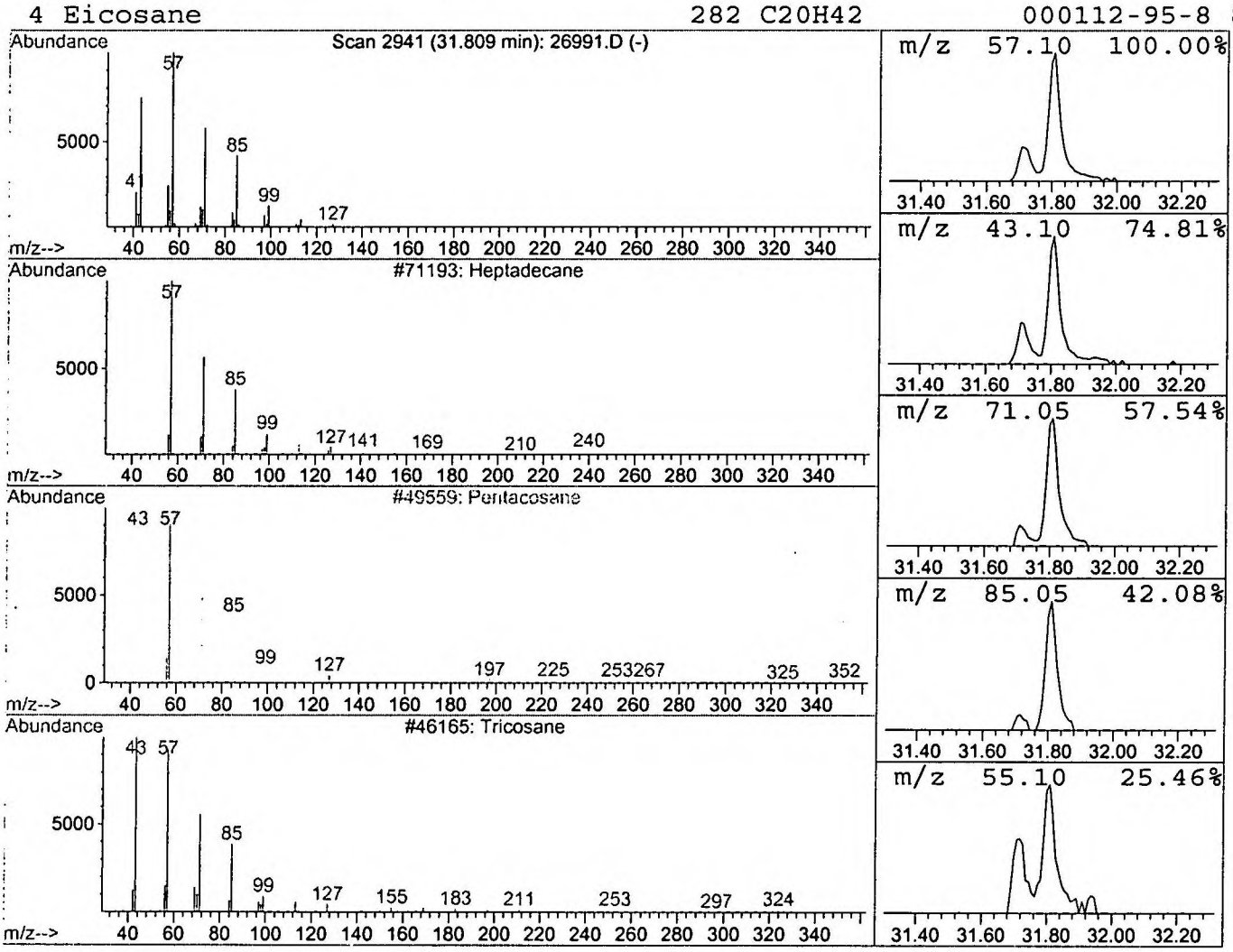
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Heptadecane — Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.81	1.02 ng/uL	295236	Chrysene-d12	33.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	90
2	Pentacosane	352	C25H52	000629-99-2	87
3	Tricosane	324	C23H48	000638-67-5	87
4	Eicosane	282	C20H42	000112-95-8	87



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MS Integration Params: LSCINT.P

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Inst : GC/MS Ins
Multiplr: 1.00

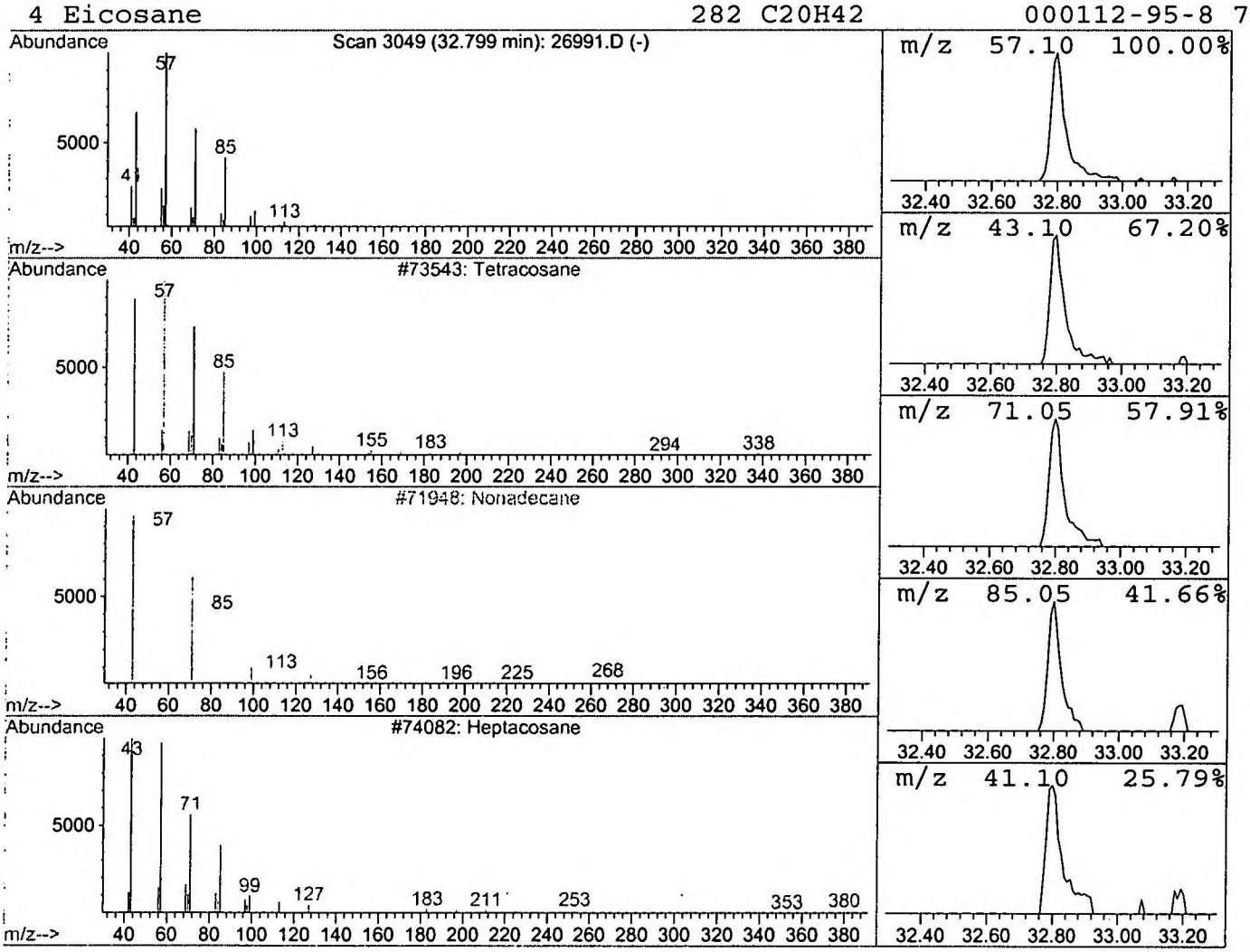
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Tetracosane — Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.80	0.84 ng/uL	241724	Chrysene-d12	33.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Nonadecane	268	C19H40	000629-92-5	90
3	Heptacosane	380	C27H56	000593-49-7	90
4	Eicosane	282	C20H42	000112-95-8	72



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26991.D
Acq On : 7 Dec 2001 12:43 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 17
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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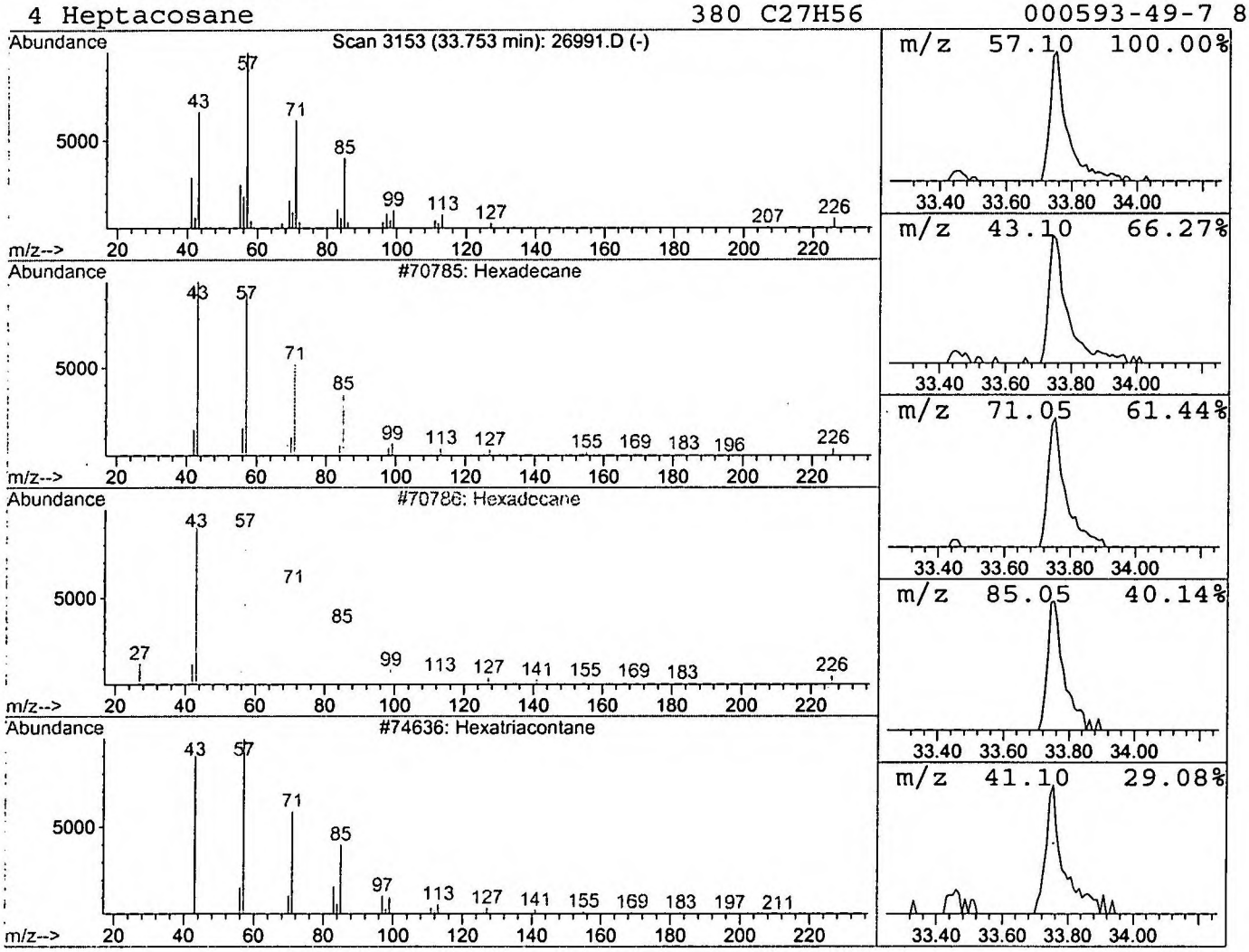
Peak Number 8 Hexadecane — Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.75	0.87 ng/uL	252486	Chrysene-d12	33.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Hexadecane	226	C16H34	000544-76-3	87
3	Hexatriacontane	507	C36H74	000630-06-8	86
4	Heptacosane	380	C27H56	000593-49-7	86



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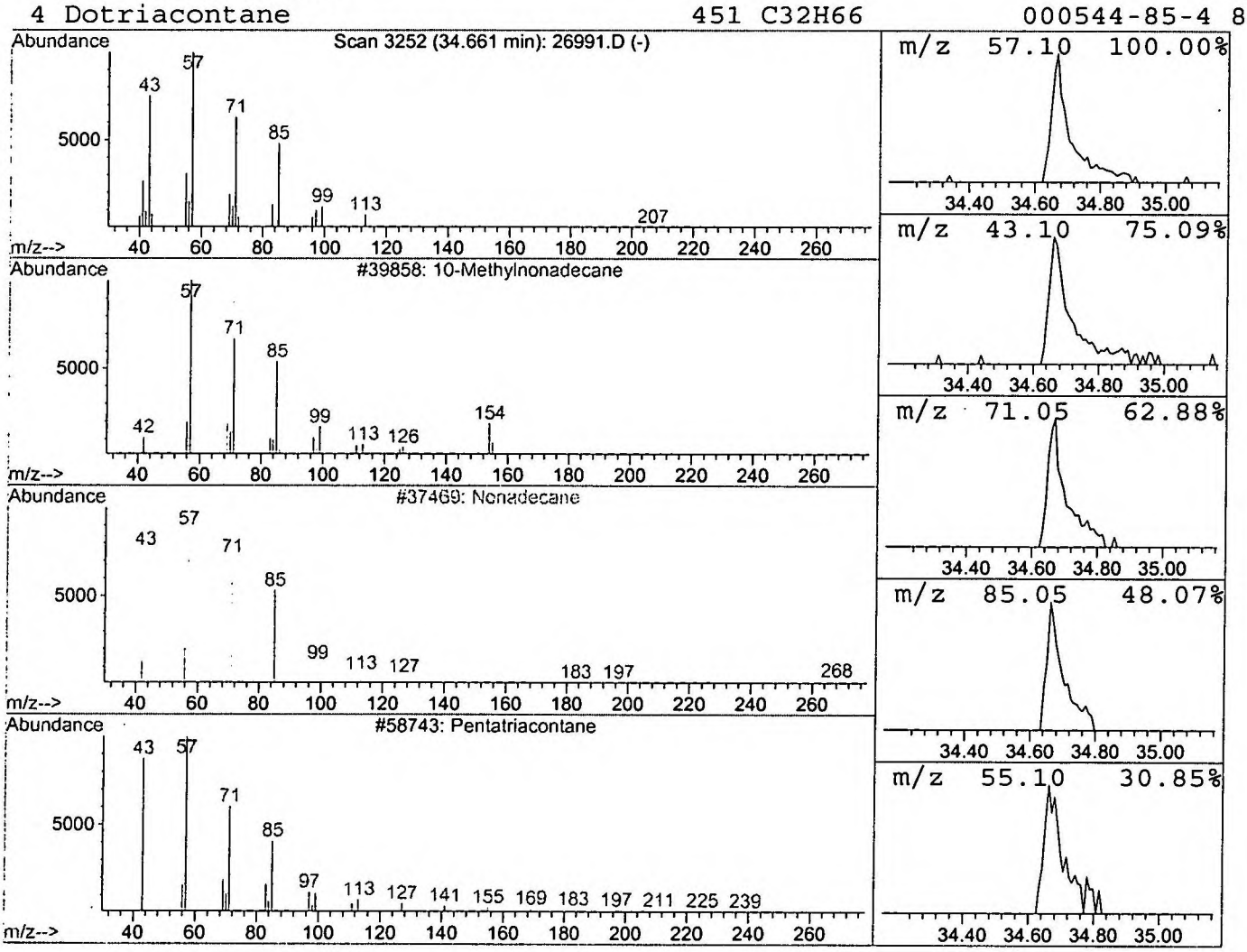
****

Peak Number 9 10-Methylnonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.66	0.54 ng/uL	154643	Chrysene-d12	33.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methylnonadecane	282	C20H42	000000-00-0	90
2	Nonadecane	268	C19H40	000629-92-5	90
3	Pentatriacontane	493	C35H72	000630-07-9	86
4	Dotriacontane	451	C32H66	000544-85-4	86



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Acq On : 7 Dec 2001 12:43 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)
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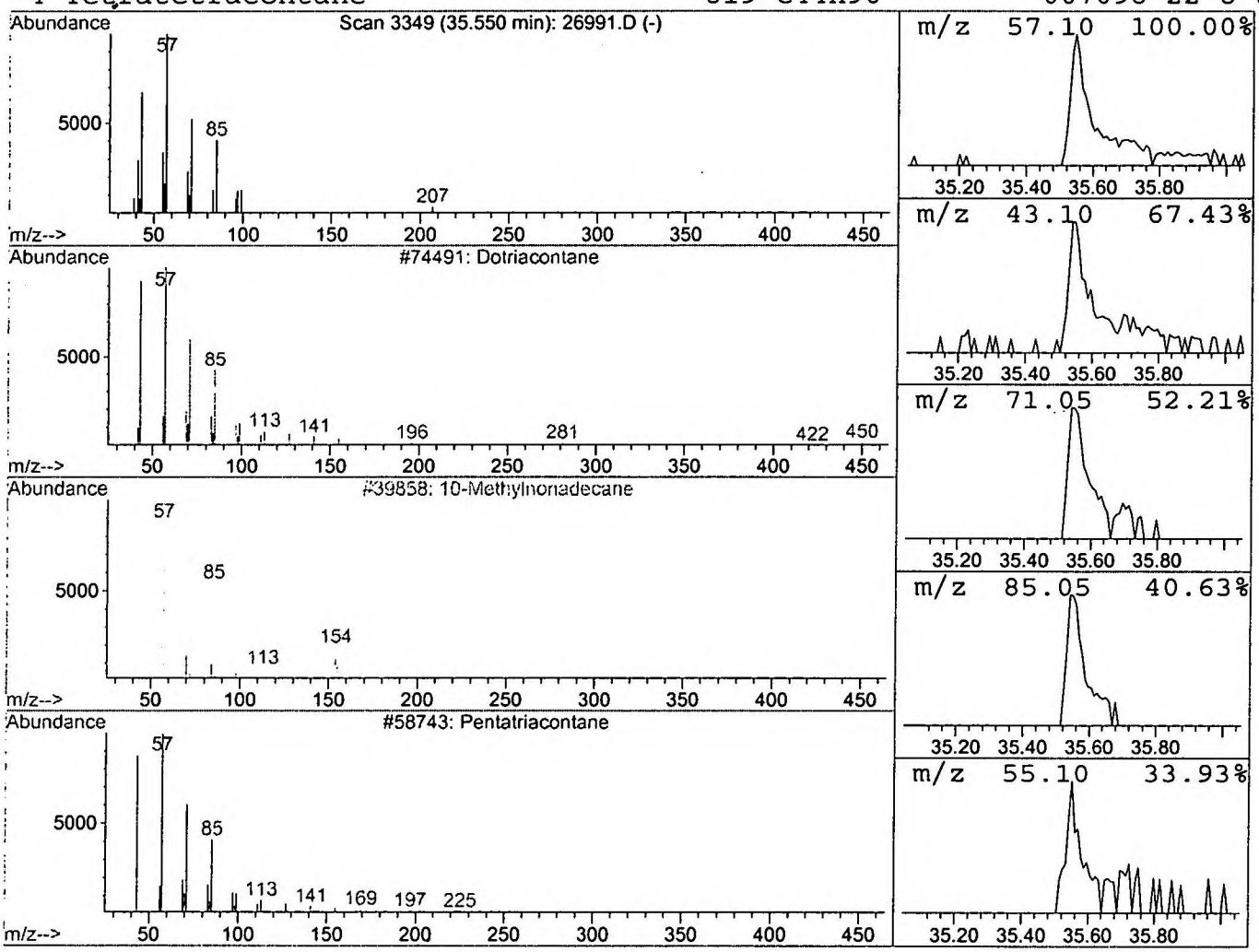
Peak Number 10 Dotriacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.55	0.47 ng/uL	114035	Perylene-d12	37.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dotriacontane	451	C32H66	000544-85-4	83
2	10-Methylnonadecane	282	C20H42	000000-00-0	83
3	Pentatriacontane	493	C35H72	000630-07-9	83
4	Tetratetracontane	619	C44H90	007098-22-8	83



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 7 Dec 2001 12:43 am
Data File: C:\HPCHEM\1\DATA\DEC0601\26991.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-	5.03	0.5	ng/uL	168172	ISTD01	11.89	6497670	20.0
2-Pentanone, 4-hydro	7.85	4.0	ng/uL	1287410	ISTD01	11.89	6497670	20.0
Cyclopentanecarboxyl	29.59	0.9	ng/uL	262827	ISTD05	33.18	5774340	20.0
Eicosane	30.77	0.4	ng/uL	120174	ISTD05	33.18	5774340	20.0
1-Hexene, 5-methyl-	31.72	0.6	ng/uL	170499	ISTD05	33.18	5774340	20.0
Heptadecane	31.81	1.0	ng/uL	295236	ISTD05	33.18	5774340	20.0
Tetracosane	32.80	0.8	ng/uL	241724	ISTD05	33.18	5774340	20.0
Hexadecane	33.75	0.9	ng/uL	252486	ISTD05	33.18	5774340	20.0
10-Methylnonadecane	34.66	0.5	ng/uL	154643	ISTD05	33.18	5774340	20.0
Dotriacontane	35.55	0.5	ng/uL	114035	ISTD06	37.27	4832750	20.0

26991.D NP112701.M Tue Dec 11 14:54:40 2001