User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/20/2001 Time: 09:04:45

Sample: AD26006
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/19/01 13:04 Ended: 12/19/01 13:04 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 AD26006 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-20-2001 Time: 15:18:30

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 bis(2-Ethylhexyl)phthalate at an estimated level of 0.26 ug/L.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC1901\26006A.D Vial: 4
Acq On : 20 Dec 2001 9:36 am Operator: GALOS
Sample : gc/ms unknown 10/31 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 20 10:38 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	11.87	152	1098210	20.00	ng/uL	-0.07
19) Naphthalene-d8	15.48	136	4336034	20.00	ng/uL	-0.07
31) Acenaphthene-d10	20.76	164	1886702	20.00	ng/uL	-0.07
49) Phenanthrene-d10	25.17	188	2793202	20.00	ng/uL	-0.08
59) Chrysene-d12	33.17	240	1895753	20.00	ng/uL	-0.08
68) Perylene-d12	37.26	264	957557	20.00	ng/uL	-0.08
System Monitoring Compounds						
4) 2-Fluorophenol	8.74	112	1773	0.02	ng/uL	0.00
Spiked Amount 75.000	Range 18 - 42		Recovery = 0.03%#			
5) Phenol-d5	11.17	99	608	0.01	ng/uL	0.04
Spiked Amount 75.000	Range 19 - 32		Recovery = 0.01%#			
6) 2-Chlorophenol-d4	11.39	132	1354	0.02	ng/uL	-0.03
Spiked Amount 75.000	Range 34 - 84		Recovery = 0.03%#			
7) 1,2-Dichlorobenzene-d4	12.38	152	1214	0.03	ng/uL	-0.08
Spiked Amount 50.000	Range 26 - 73		Recovery = 0.06%#			
20) Nitrobenzene-d5	13.48	82	1900	0.02	ng/uL	-0.08
Spiked Amount 50.000	Range 55 - 98		Recovery = 0.04%#			
32) 2-Fluorobiphenyl	18.76	172	3265	0.03	ng/uL	-0.08
Spiked Amount 50.000	Range 42 - 78		Recovery = 0.06%#			
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	30.04	244	2959	0.03	ng/uL	-0.08
Spiked Amount 50.000	Range 58 - 98		Recovery = 0.06%#			
Target Compounds						Qvalue
2) Pyridine	5.14	79	282	N.D.		
3) N-nitrosodimethylamine	5.49	74	977	N.D.		
8) Phenol	11.21	94	3099	N.D.		
9) bis(2-Chloroethyl)ether	11.27	93	4267	N.D.		
10) 2-Chlorophenol	11.44	128	1998	N.D.		
11) 1,3-Dichlorobenzene	11.77	146	3164	N.D.		
12) 1,4-Dichlorobenzene	11.91	146	3594	N.D.		
13) 1,2-Dichlorobenzene	12.42	146	3517	N.D.		
14) 2-Methylphenol	12.80	108	287	N.D.		
15) 2,2'-oxybis(1-Chloropropan	12.75	45	9388	N.D.		
16) 3&4-Methylphenol	13.25	108	979	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	13.14	70	2945	N.D.	
18)	Hexachloroethane	0.00	117	0	N.D.	
21)	Nitrobenzene	13.54	77	4637	N.D.	
22)	Isophorone	14.19	82	8812	N.D.	
23)	2-Nitrophenol	14.46	139	419	N.D.	
24)	2,4-Dimethylphenol	14.72	107	2341	N.D.	
25)	bis(2-Chloroethoxy)methane	14.89	93	5270	N.D.	
26)	2,4-Dichlorophenol	0.00	162	0	N.D.	
27)	1,2,4-Trichlorobenzene	15.37	180	2784	N.D.	
28)	Naphthalene	15.54	128	13137	N.D.	
29)	Hexachlorobutadiene	16.09	225	936	N.D.	
30)	4-Chloro-3-methylphenol	17.42	107	1274	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	19.02	162	5944	N.D.	
38)	Dimethylphthalate	20.13	163	7938	N.D.	
39)	2,6-Dinitrotoluene	20.33	165	364	N.D.	
40)	Acenaphthylene	20.29	152	9044	N.D.	
41)	Acenaphthene	20.85	153	6649	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	21.52	165	684	N.D.	
45)	Diethylphthalate	22.24	149	9148	N.D.	
46)	4-Chlorophenylphenylether	22.39	204	2920	N.D.	
47)	Fluorene	22.36	166	7372	N.D.	
48)	Azobenzen/ 1,2-Diphenylhyd	22.86	77	8745	N.D.	
50)	4,6-Dinitro-2-methylphenol	0.00	198	0	N.D.	
51)	N-Nitrosodiphenylamine	22.78	169	4102	N.D.	
52)	4-Bromophenyl-phenylether	23.85	248	1539	N.D.	
53)	Hexachlorobenzene	24.26	284	916	N.D.	
54)	Pentachlorophenol	0.00	266	0	N.D.	
55)	Phenanthrene	25.23	178	10532	N.D.	
56)	Anthracene	25.23	178	10532	N.D.	
57)	Di-n-butylphthalate	27.15	149	21980	N.D.	
58)	Fluoranthene	28.83	202	9284	N.D.	
60)	3,3'-Dichlorobenzidine	33.14	252	1095	N.D.	

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Response via : Initial Calibration
DataAcq Meth : NP112701

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
61)	Benzidine	0.00	184	0	N.D. d	
63)	Pyrene	29.49	202	9576	N.D.	
64)	Butylbenzylphthalate	31.64	149	7463	N.D.	
65)	Benzo(a)anthracene	33.24	228	7721	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.42	149	28165	0.26 ng/uL#	97
67)	Chrysene	33.24	228	7721	N.D.	
69)	Di-n-Octylphthalate	35.15	149	13381	N.D.	
70)	Benzo(b)fluoranthene	36.18	252	5620	N.D.	
71)	Benzo(k)fluoranthene	36.23	252	5047	N.D.	
72)	Benzo(a)pyrene	37.06	252	3973	N.D.	
73)	Indeno(1,2,3-cd)pyrene	40.98	276	280	N.D.	
74)	Dibenz(a,h)anthracene	41.01	278	2413	N.D.	
75)	Benzo(g,h,i)perylene	42.03	276	758	N.D.	

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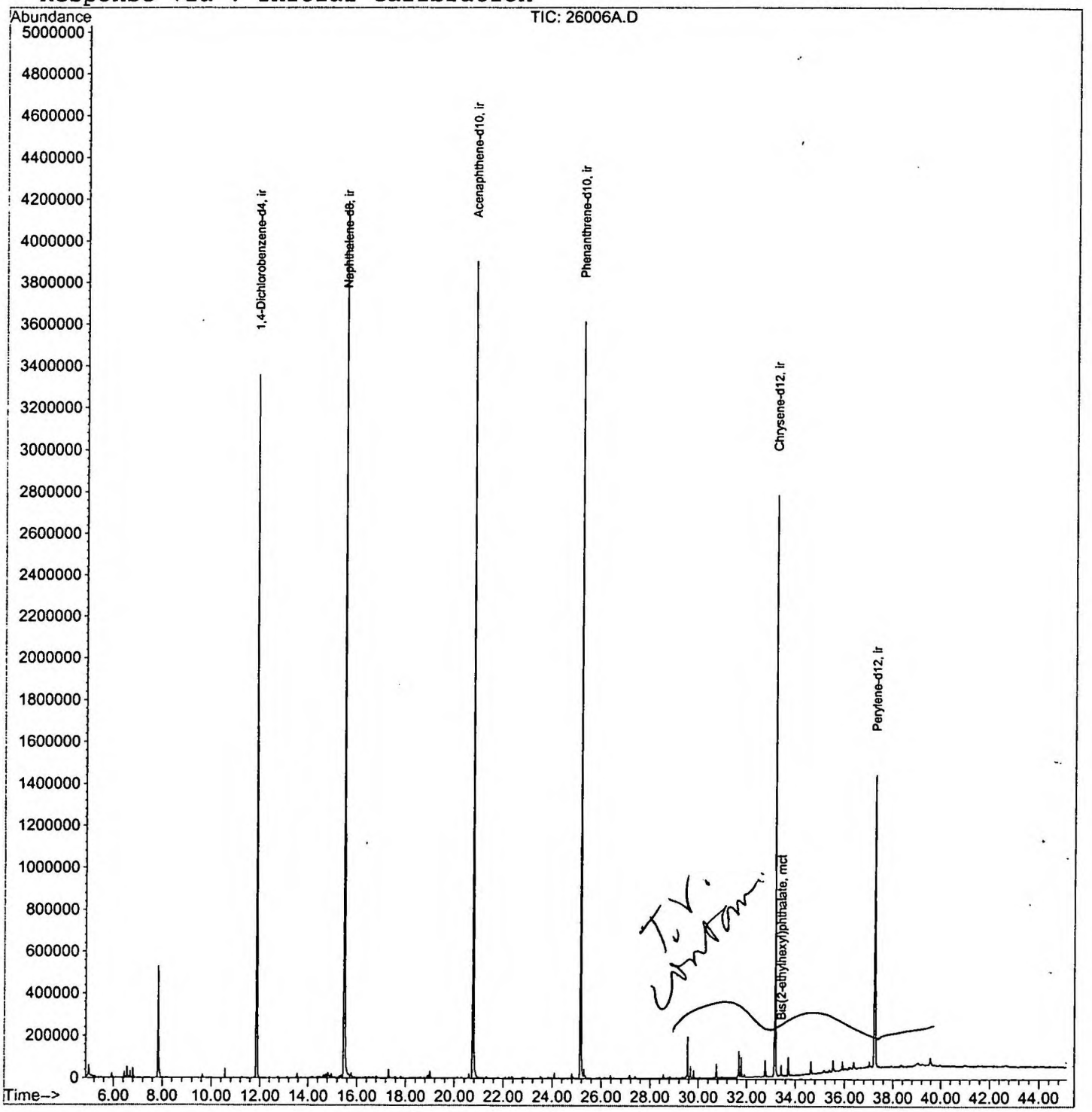
Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC1901\26006A.D Vial: 4
Acq On : 20 Dec 2001 9:36 am Operator: GALOS
Sample : gc/ms unknown 10/31 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

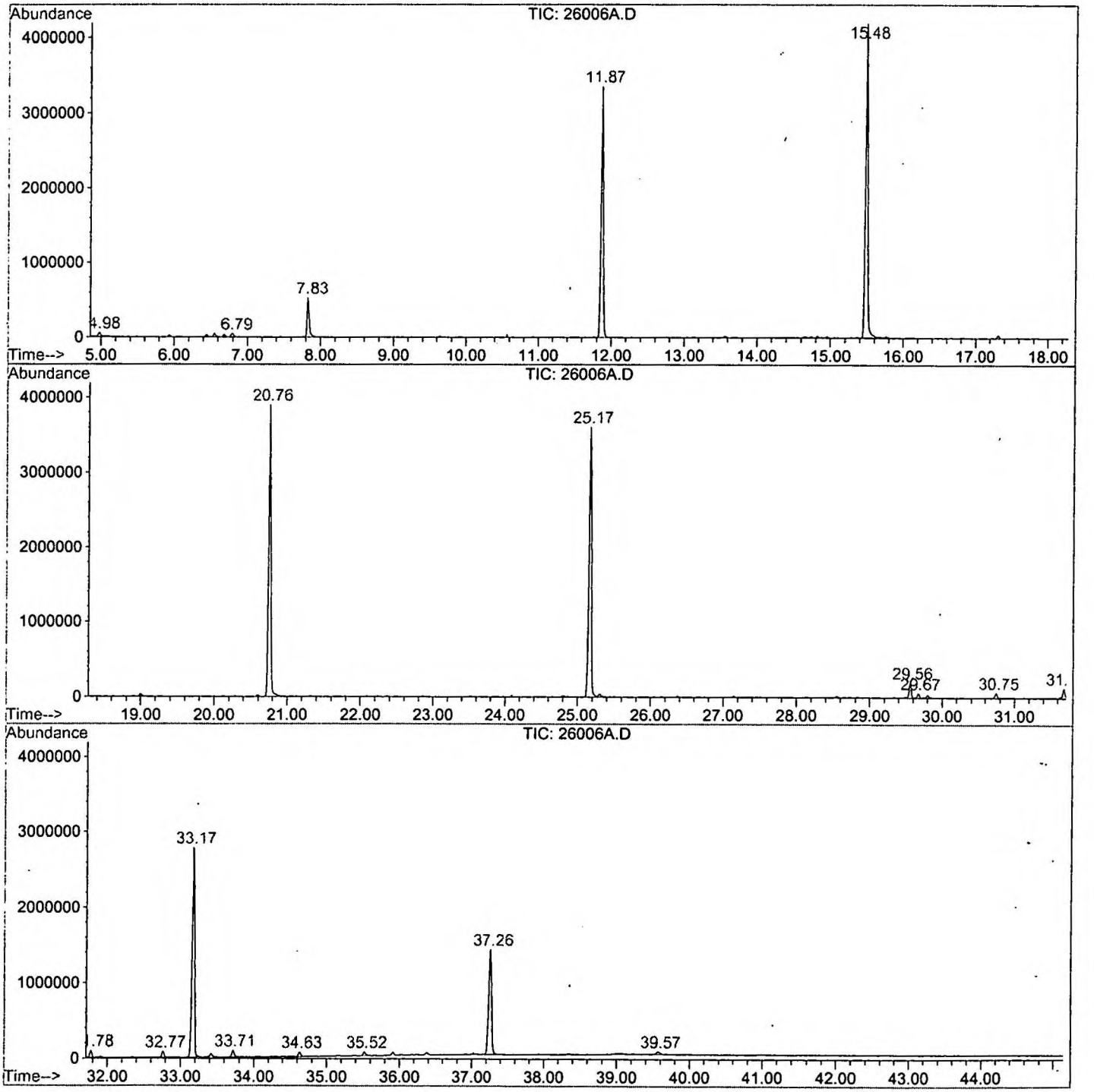
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	4.983	8	15	19	rBV	53042	113345	1.24%	0.246%
2	6.789	207	212	219	rVB2	46034	109564	1.20%	0.238%
3	7.826	321	325	341	rVB	528391	1051475	11.50%	2.283%
4	11.870	760	766	776	rBV	3358357	6833701	74.72%	14.839%
5	15.483	1154	1160	1177	rBV	4186511	9145194	100.00%	19.858%
6	20.756	1727	1735	1751	rBV	3904244	8566727	93.67%	18.602%
7	25.166	2209	2216	2227	rBV2	3610294	8075845	88.31%	17.536%
8	29.559	2691	2695	2701	rVB	198373	366067	4.00%	0.795%
9	29.669	2703	2707	2713	rVB	57083	113167	1.24%	0.246%
10	30.751	2820	2825	2829	rBV	66104	128316	1.40%	0.279%
11	31.686	2923	2927	2932	rVB	121718	225280	2.46%	0.489%
12	31.778	2932	2937	2941	rVB	93265	176982	1.94%	0.384%
13	32.768	3040	3045	3049	rBV	78946	165262	1.81%	0.359%
14	33.172	3080	3089	3095	rBV2	2776712	6468322	70.73%	14.045%
15	33.713	3143	3148	3153	rBV2	90508	201961	2.21%	0.439%
16	34.630	3242	3248	3254	rBV	64847	152842	1.67%	0.332%
17	35.519	3341	3345	3351	rBV2	51703	114969	1.26%	0.250%
18	37.262	3526	3535	3547	rVB2	1387285	3924253	42.91%	8.521%
19	39.573	3783	3787	3794	rVB8	37341	119874	1.31%	0.260%

Sum of corrected areas: 46053146

26006A.D NP112701.M Thu Dec 20 10:39:39 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC1901\26006A.D
Operator : GALOS
Acquired : 20 Dec 2001 9:36 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 10/31
Misc Info :
Vial Number: 4
Quant File :NP112701.RES (RTE Integrator)



26006A.D NP112701.M Thu Dec 20 10:39:42 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1901\26006A.D Vial: 4
Acq On : 20 Dec 2001 9:36 am Operator: GALOS
Sample : gc/ms unknown 10/31 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

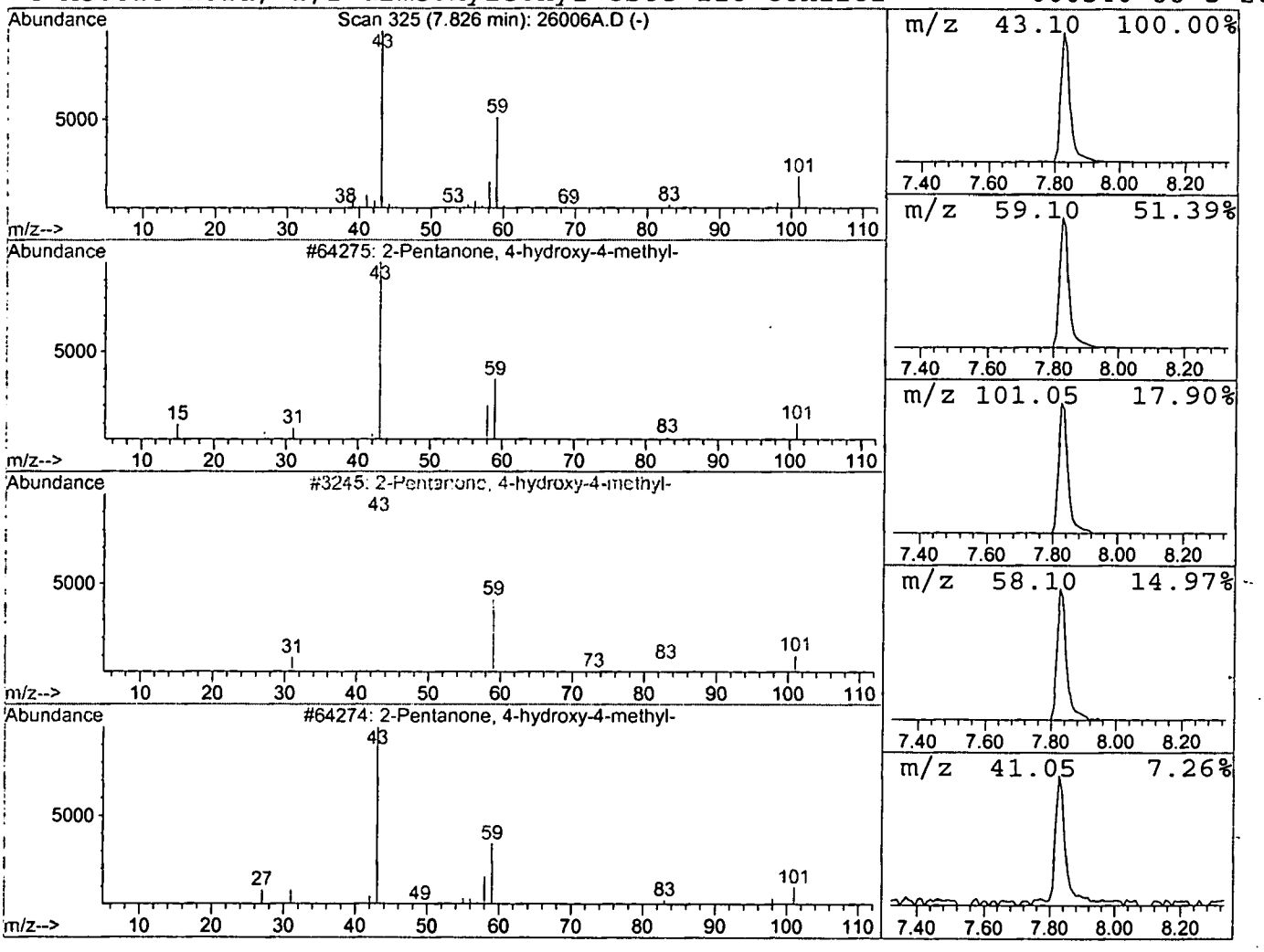
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	3.08 ng/uL	1051480	1,4-Dichlorobenzene-d4	11.87

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	36
4	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28



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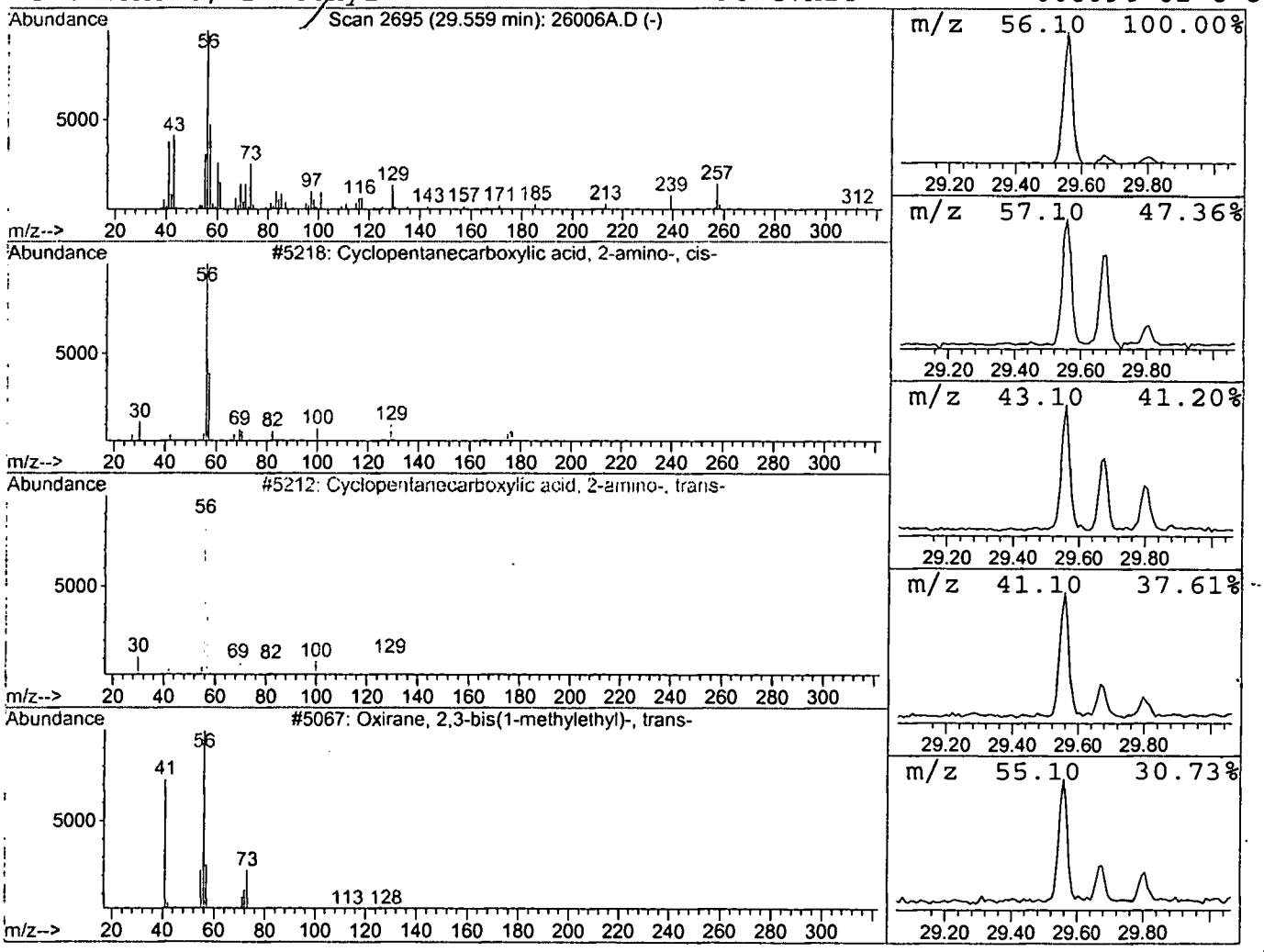
****

Peak Number 2 Cyclopentanecarboxylic acid, 2 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.56	1.13 ng/uL	366067	Chrysene-d12	33.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	037910-65-9	38
2	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	040482-05-1	38
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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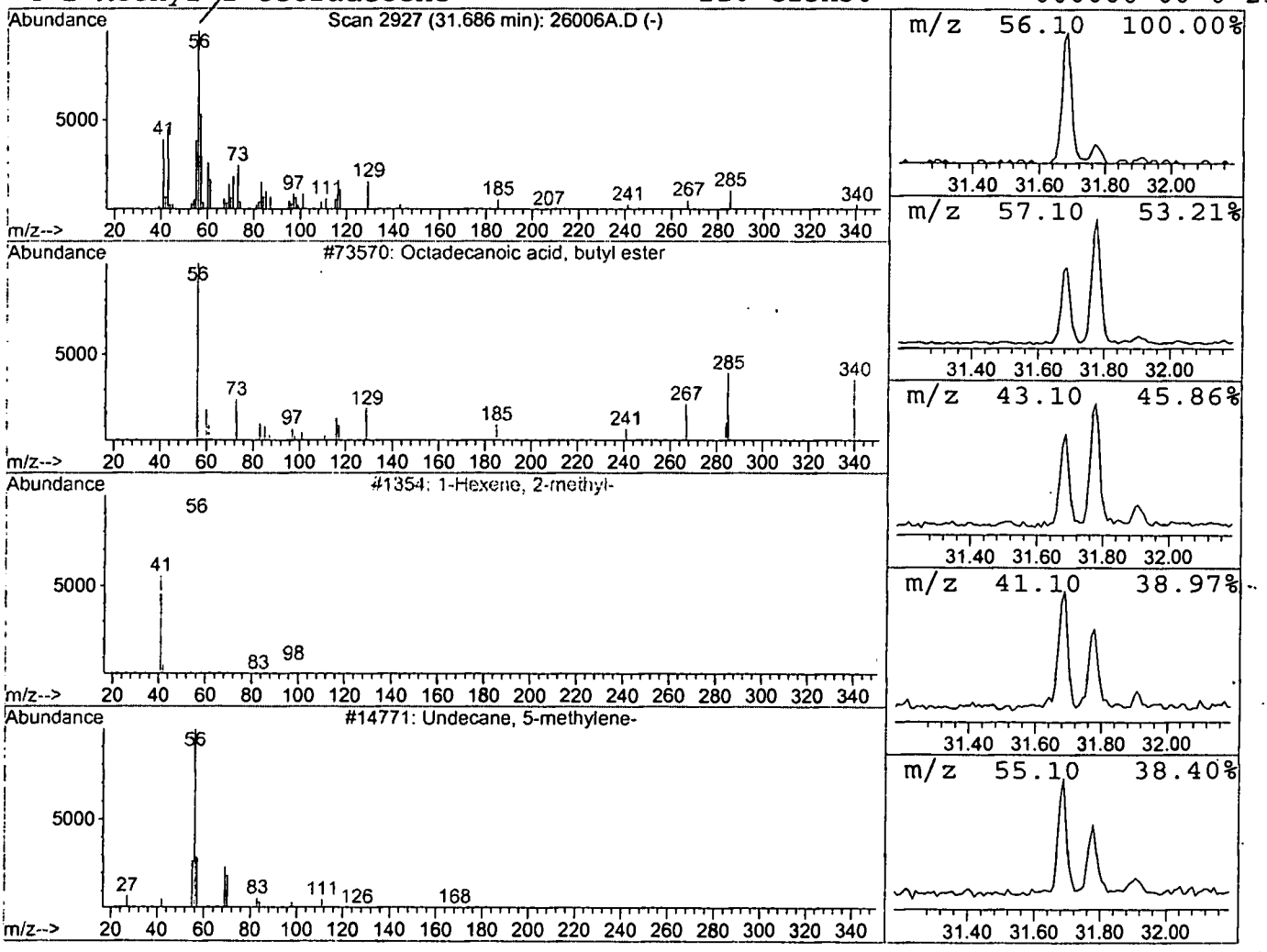
****

Peak Number 3 Octadecanoic acid, butyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.69	0.70 ng/uL	225280	Chrysene-d12	33.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	80
2	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	35
3	Undecane, 5-methylene-	168	C12H24	005698-48-6	25
4	2-Methyl-1-tetradecene	210	C15H30	000000-00-0	25



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Library : C:\DATABASE\NBS75K.L

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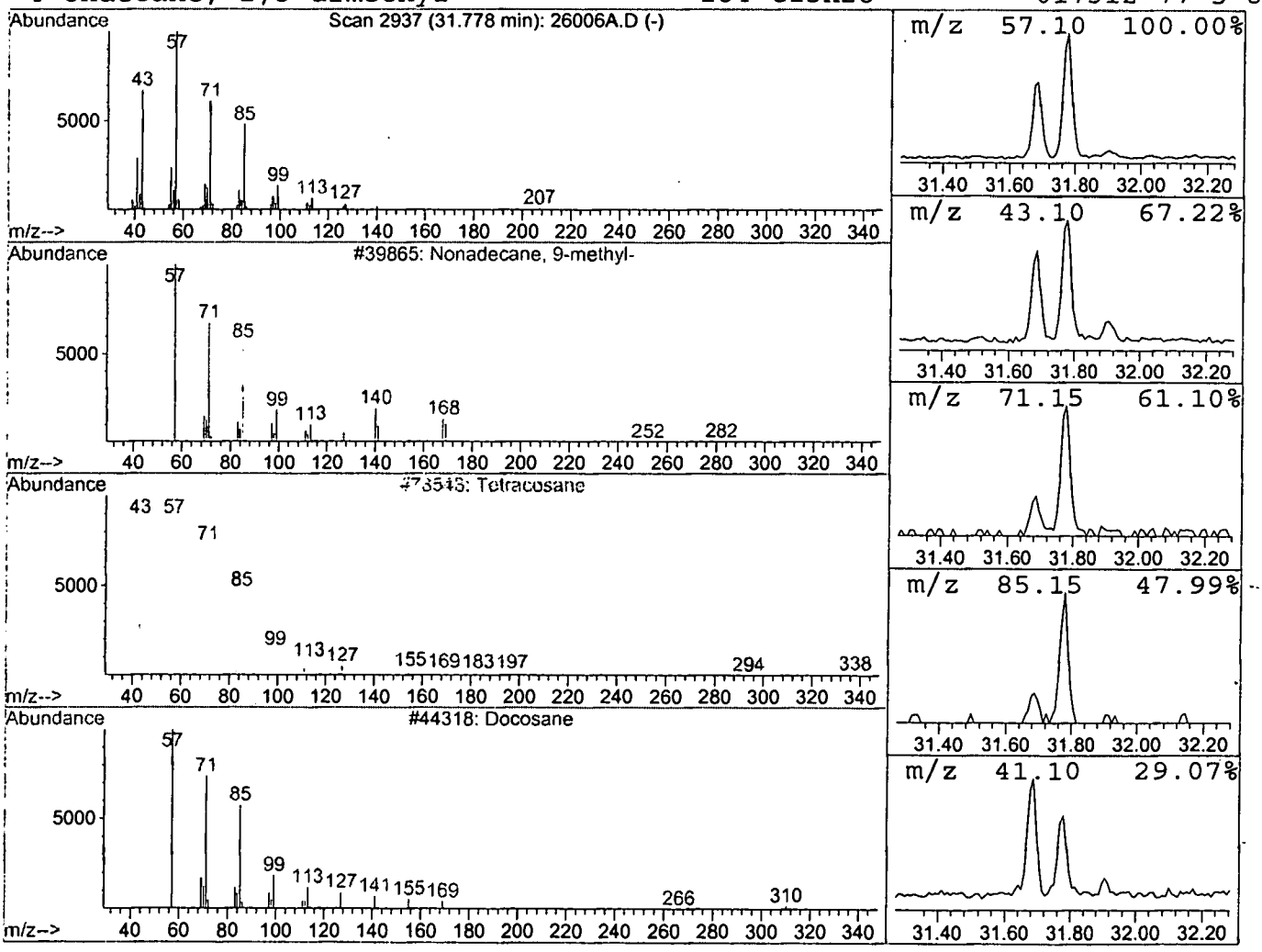
Peak Number 4 Nonadecane, 9-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.78	0.55 ng/uL	176982	Chrysene-d12	33.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	91
2	Tetracosane	338	C24H50	000646-31-1	87
3	Docosane	310	C22H46	000629-97-0	86
4	Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	86



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MS Integration Params: LSCINT.P

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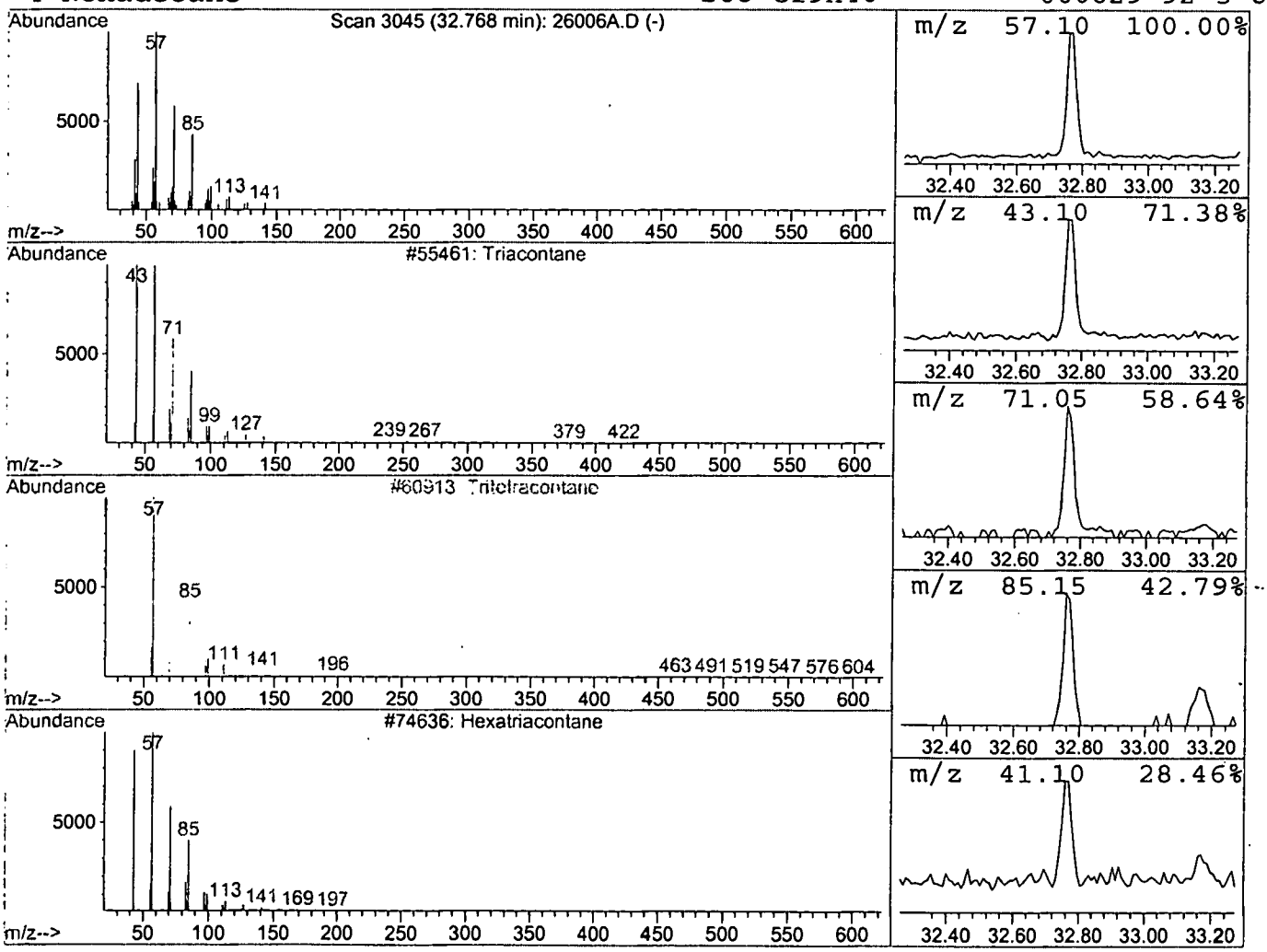
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Triacontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.77	0.51 ng/uL	165262	Chrysene-d12	33.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	91
2	Tritetracontane	605	C43H88	007098-21-7	91
3	Hexatriacontane	507	C36H74	000630-06-8	90
4	Nonadecane	268	C19H40	000629-92-5	86



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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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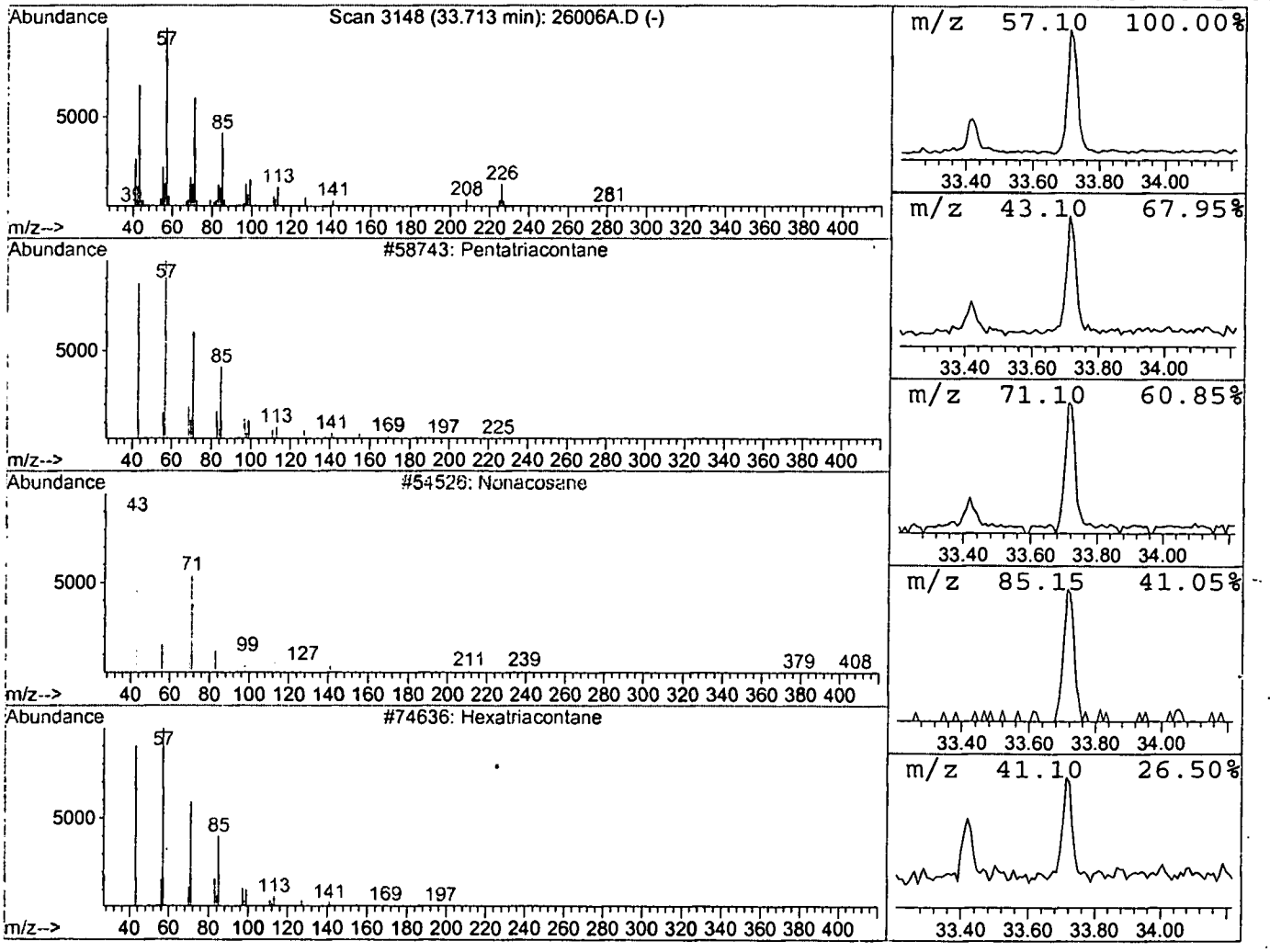
Peak Number 6 Pentatriacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.71	0.62 ng/uL	201961	Chrysene-d12	33.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentatriacontane	493	C35H72	000630-07-9	90
2	Nonacosane	408	C29H60	000630-03-5	87
3	Hexatriacontane	507	C36H74	000630-06-8	83
4	Octadecane	254	C18H38	000593-45-3	83



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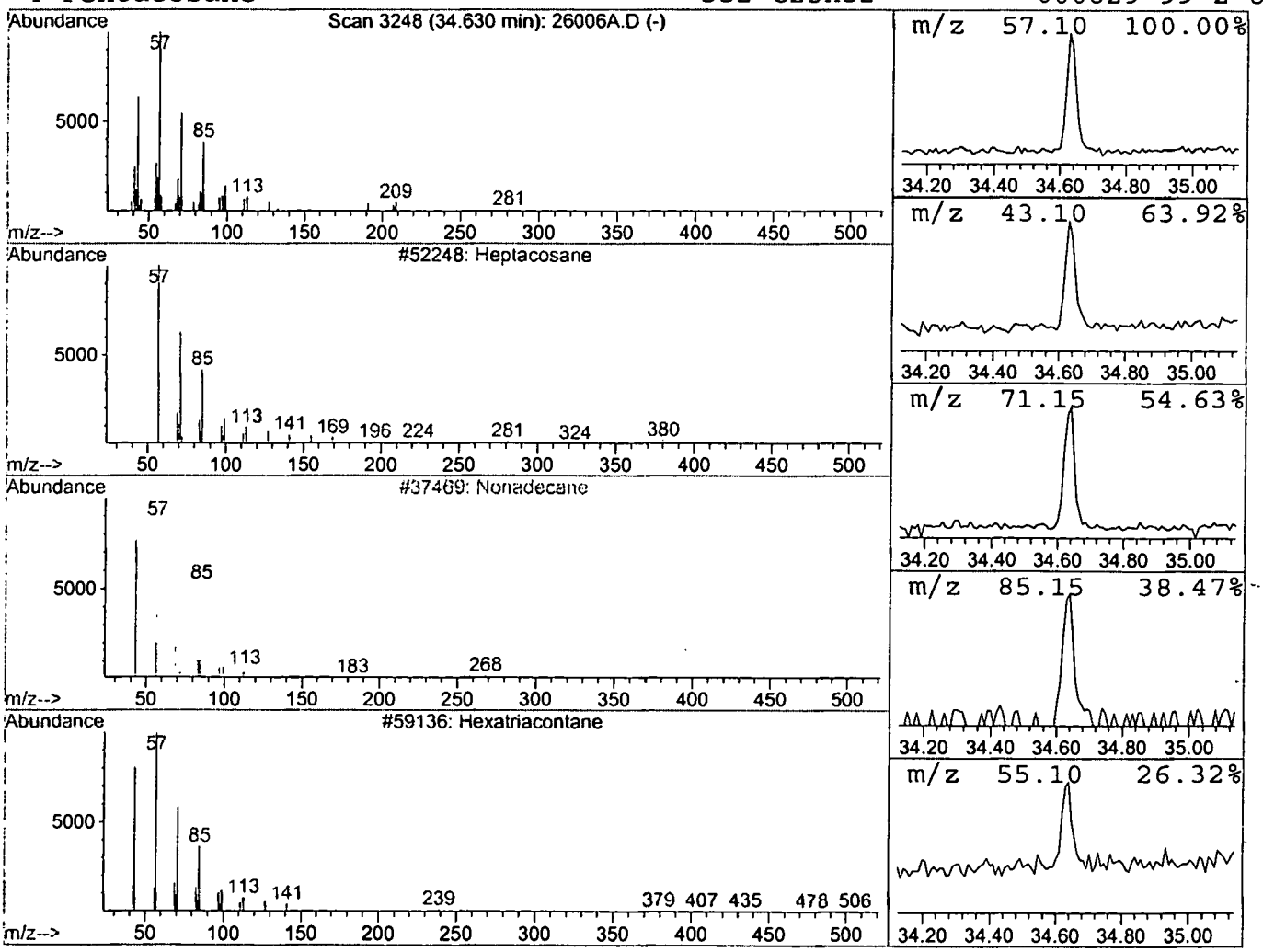
****

Peak Number 7 Heptacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.63	0.47 ng/uL	152842	Chrysene-d12	33.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	83
2	Nonadecane	268	C19H40	000629-92-5	83
3	Hexatriacontane	507	C36H74	000630-06-8	80
4	Pentacosane	352	C25H52	000629-99-2	80



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1901\26006A.D Vial: 4
Acq On : 20 Dec 2001 9:36 am Operator: GALOS
Sample : gc/ms unknown 10/31 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

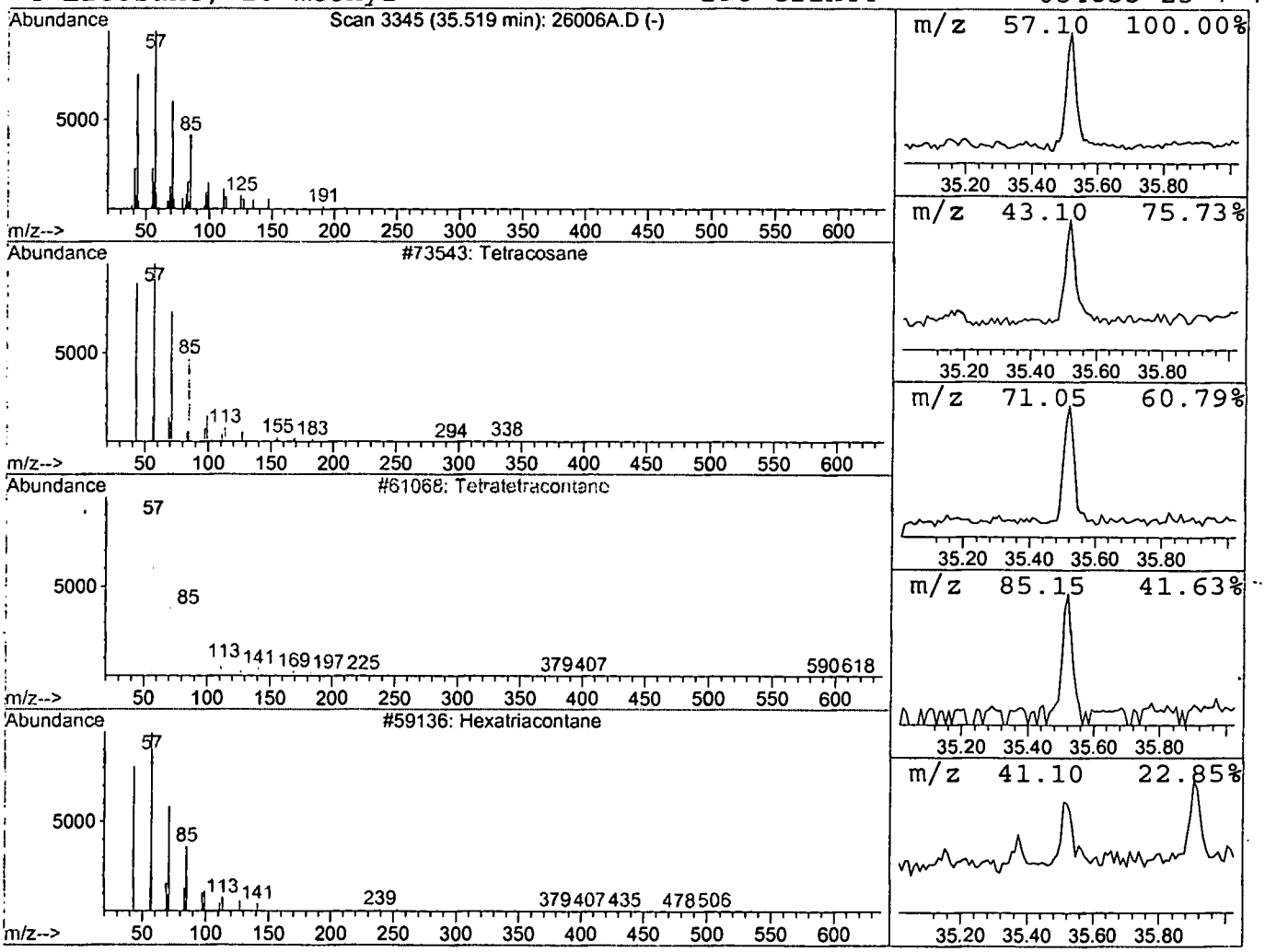
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.52	0.59 ng/uL	114969	Perylene-d12	37.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	80
2	Tetratetracontane	619	C44H90	007098-22-8	80
3	Hexatriacontane	507	C36H74	000630-06-8	80
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	72



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC1901\26006A.D Vial: 4
Acq On : 20 Dec 2001 9:36 am Operator: GALOS
Sample : gc/ms unknown 10/31 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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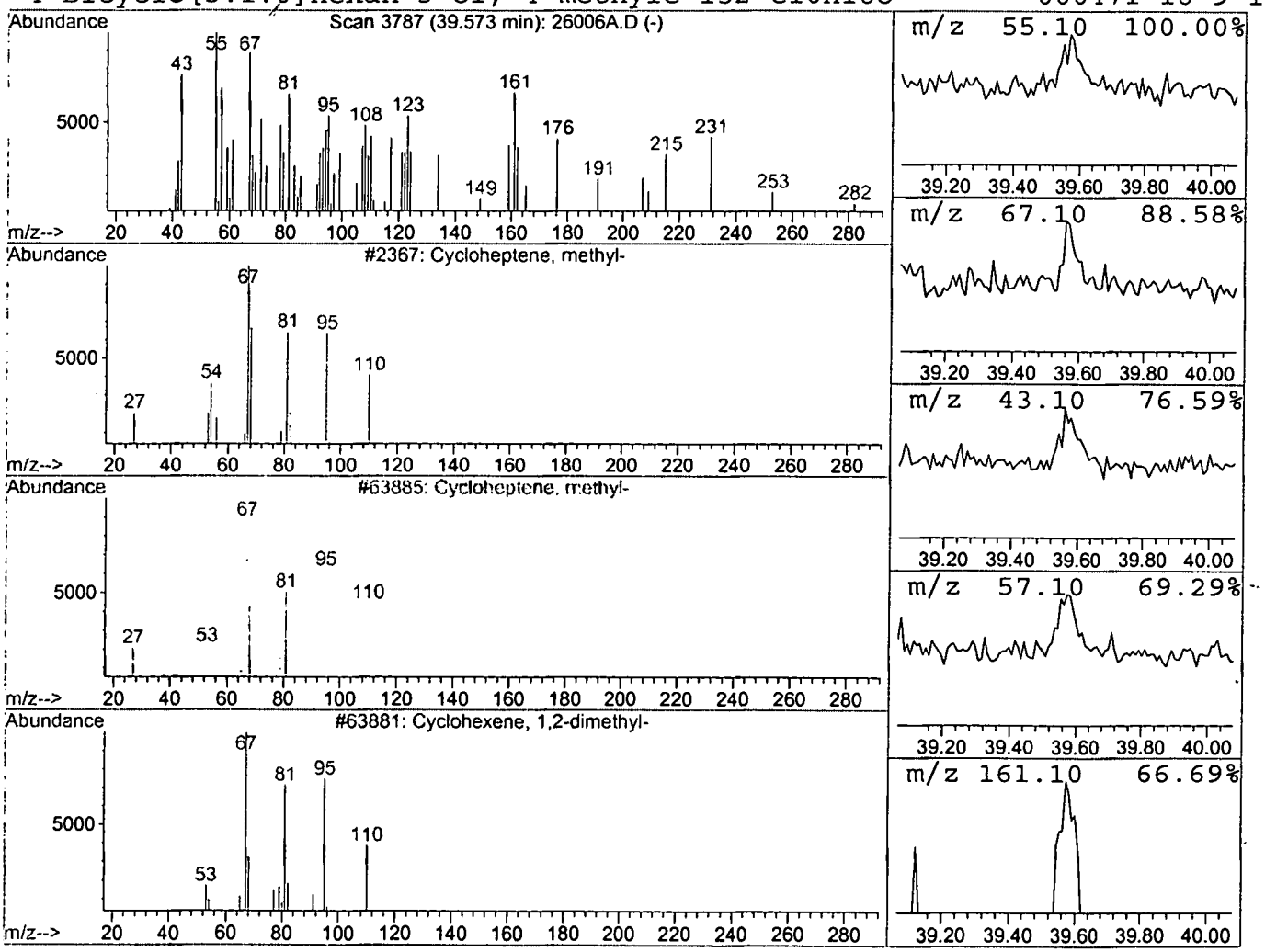
Peak Number 9 Cycloheptene, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
39.57	0.61 ng/uL	119874	Perylene-d12	37.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cycloheptene, methyl-	110	C8H14	055308-20-8	22
2	Cycloheptene, methyl-	110	C8H14	055308-20-8	22
3	Cyclohexene, 1,2-dimethyl-	110	C8H14	001674-10-8	14
4	Bicyclo[3.1.0]hexan-3-ol, 4-methyle	152	C10H16O	000471-16-9	10



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 20 Dec 2001 9:36 am
Data File: C:\HPCHEM\1\DATA\DEC1901\26006A.D
Name: gc/ms unknown 10/31
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-Pentanone, 4-hydro	7.83	3.1	ng/uL	1051480	ISTD01	11.87	6833700	20.0
Cyclopentanecarboxyl	29.56	1.1	ng/uL	366067	ISTD05	33.17	6468320	20.0
Octadecanoic acid, b	31.69	0.7	ng/uL	225280	ISTD05	33.17	6468320	20.0
Nonadecane, 9-methyl	31.78	0.5	ng/uL	176982	ISTD05	33.17	6468320	20.0
Triacontane	32.77	0.5	ng/uL	165262	ISTD05	33.17	6468320	20.0
Pentatriacontane	33.71	0.6	ng/uL	201961	ISTD05	33.17	6468320	20.0
Heptacosane	34.63	0.5	ng/uL	152842	ISTD05	33.17	6468320	20.0
Tetracosane	35.52	0.6	ng/uL	114969	ISTD06	37.26	3924250	20.0
Cycloheptene, methyl	39.57	0.6	ng/uL	119874	ISTD06	37.26	3924250	20.0

26006A.D NP112701.M Thu Dec 20 10:40:01 2001