LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Integration Parameters: LSCINT.P
Integrator: RTE
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

Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.557	69	80	88	rBV	172469	316582	2.61%	0.407%
2	3.816	115	124	130	rVV2	145475	220628	1.82%	0.283%
3	4.034	153	161	171	rVB	1200217	1653483	13.65%	2.124%
4	4.375	214	219	229	rVB	102176	155937	1.29%	0.200%
5	5.340	378	383	391	rVV	2541066	3601924	29.73%	4.627%
6	5.475	399	406	416	rBV	5953112	12116430	100.00%	15.564%
7	5.675	435	440	443	rBV	95586	139249	1.15%	0.179%
8	5.840	461	468	477	rVB	1465568	2100087	17.33%	2.698%
9	6.310	542	548	555	rBV	112141	196972	1.63%	0.253%
10	6.804	622	632	637	rBV2	130105	249348	2.06%	0.320%
11	6.869	637	643	645	rBV3	83968	147330	1.22%	0.189%
12	6.904	645	649	654	rVB2	231779	392740	3.24%	0.504%
13	6.969	654	660	661	rBV	184387	256843	2.12%	0.330%
14	6.998	661	665	670	rBV	602927	853724	7.05%	1.097%
15	7.151	685	691	696	rBV	664551	978701	8.08%	1.257%
16	7.204	696	700	705	rVB2	90480	128579	1.06%	0.165%
17	7.275	707	712	715	rBV	84510	130020	1.07%	0.167%
18	7.357	720	726	734	rVB	885227	1451014	11.98%	1.864%
19	7.439	734	740	742	rBV	473856	672459	5.55%	0.864%
20	7.463	742	744	754	rVB	324904	432855	3.57%	0.556%
21	7.816	794	804	813	rBV	741800	1356954	11.20%	1.743%
22	7.934	819	824	831	rBV2	121295	235737	1.95%	0.303%
23	8.098	844	852	860	rVB6	48773	123152	1.02%	0.158%
24	8.357	890	896	901	rVB3	129364	249714	2.06%	0.321%
25	8.451	906	912	919	rVV4	185138	420943	3.47%	0.541%
26	8.534	919	926	932	rVV	945903	1553176	12.82%	1.995%
27	8.639	936	944	949	rVV2	98297	248178	2.05%	0.319%
28	8.916	980	991	995	rBV5	129708	281862	2.33%	0.362%
29	8.981	995	1002	1007	rVV	781568	1355541	11.19%	1.741%
30	9.039	1007	1012	1022	rVB	360805	564100	4.66%	0.725%
31	9.163	1022	1033	1039	rBV7	53972	161766	1.34%	0.208%
32	9.698	1112	1124	1130	rBV	668196	1173131	9.68%	1.507%
33	9.980	1160	1172	1176	rBV2	153281	363457	3.00%	0.467%
34	10.239	1210	1216	1223	rVB	1139723	1871010	15.44%	2.403%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Integration Parameters: LSCINT.P
Integrator: RTE
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

Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Title : SVOA CALIBRATION

35	10.610	1269	1279	1284	rBV	979835	1826551	15.07%	2.346%
36	10.657	1284	1287	1295	rVB	130507	217961	1.80%	0.280%
37	10.745	1295	1302	1314	rBV	400554	827838	6.83%	1.063%
38	11.433	1412	1419	1426	rBV	127744	228534	1.89%	0.294%
39	11.927	1498	1503	1511	rVB	100828	154339	1.27%	0.198%
40	12.027	1513	1520	1526	rVV2	129392	206156	1.70%	0.265%
41	12.269	1556	1561	1569	rVB	88740	152131	1.26%	0.195%
42	12.969	1675	1680	1689	rVB2	99466	211336	1.74%	0.271%
43	13.274	1727	1732	1740	rVB2	116633	187300	1.55%	0.241%
44	13.774	1812	1817	1822	rBV	80761	135496	1.12%	0.174%
45	13.868	1827	1833	1837	rVV	1839164	2646121	21.84%	3.399%
46	13.910	1837	1840	1849	rVB	930659	1348808	11.13%	1.733%
47	14.151	1875	1881	1896	rVB	2160986	3155274	26.04%	4.053%
48	14.351	1911	1915	1923	rVB	113414	142667	1.18%	0.183%
49	14.457	1923	1933	1940	rBV2	1359467	2133774	17.61%	2.741%
50	14.686	1963	1972	1980	rBV	402050	700349	5.78%	0.900%
51	14.857	1997	2001	2008	rVB2	115552	240926	1.99%	0.309%
52	15.304	2074	2077	2082	rVB	98826	123527	1.02%	0.159%
53	15.451	2094	2102	2108	rBV	2391282	3470497	28.64%	4.458%
54	15.592	2122	2126	2131	rBV	83356	121369	1.00%	0.156%
55	15.815	2156	2164	2169	rVB	480232	694068	5.73%	0.892%
56	16.168	2219	2224	2232	rBV3	131227	314624	2.60%	0.404%
57	16.604	2295	2298	2304	rBV2	85869	129150	1.07%	0.166%
58	17.204	2395	2400	2404	rBV2	1297722	1841035	15.19%	2.365%
59	17.245	2404	2407	2411	rVV	432843	557693	4.60%	0.716%
60	17.304	2411	2417	2427	rVB2	2083305	3034496	25.04%	3.898%
61	17.980	2527	2532	2538	rBV5	131547	322831	2.66%	0.415%
62	18.103	2545	2553	2561	rBV2	890543	1688176	13.93%	2.169%
63	18.268	2577	2581	2585	rBV2	128900	198474	1.64%	0.255%
64	18.580	2630	2634	2638	rBV2	107387	174304	1.44%	0.224%
65	18.998	2702	2705	2708	rBV	130668	184653	1.52%	0.237%
66	19.262	2746	2750	2756	rVB	870445	1165968	9.62%	1.498%
67	19.321	2756	2760	2763	rBV3	130331	176302	1.46%	0.226%
68	19.380	2765	2770	2780	rBV2	705555	1050898	8.67%	1.350%
69	19.597	2802	2807	2811	rBV	2446600	3628484	29.95%	4.661%
70	20.515	2960	2963	2967	rVB	570347	589501	4.87%	0.757%
71	21.386	3106	3111	3115	rBV2	1798618	2669614	22.03%	3.429%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Integration Parameters: LSCINT.P
Integrator: RTE
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

Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Title : SVOA CALIBRATION

72	23.538	3472	3477	3482	rBV2	1834162	3223533	26.60%	4.141%
73	23.685	3497	3502	3507	rVB2	1036410	1820927	15.03%	2.339%

Sum of corrected areas: 77849311

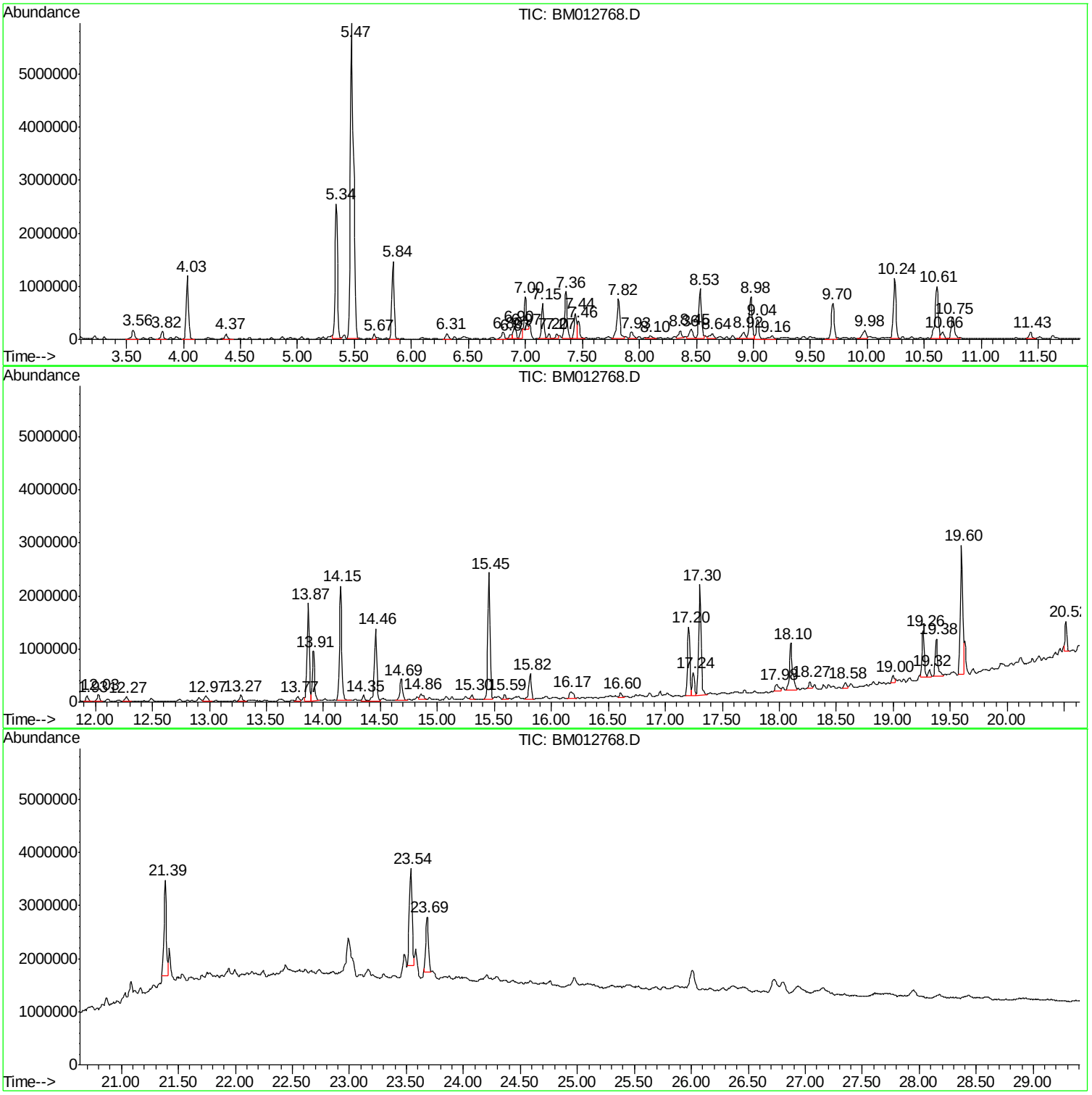
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

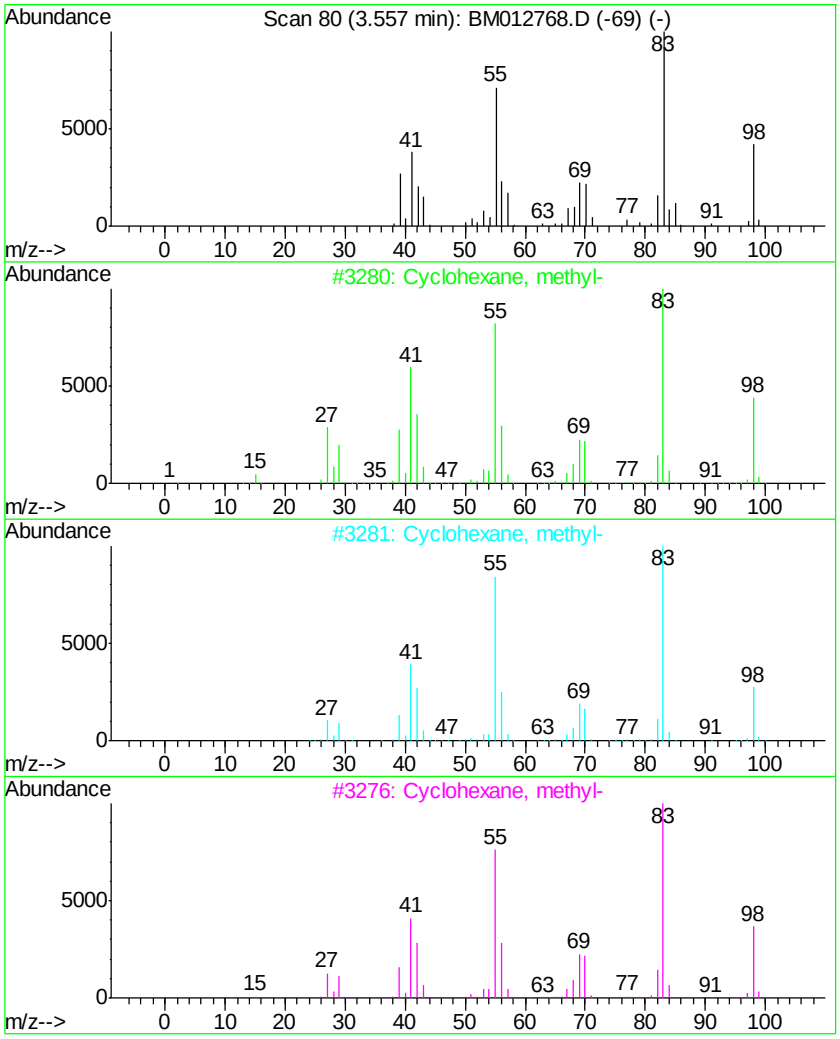
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

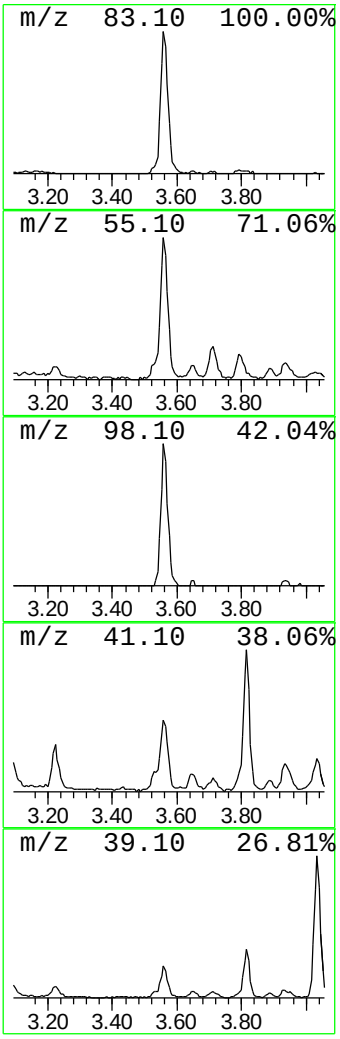
****

Peak Number 1 (DEL) Alkane: Cyclic3.56 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.56	4.67 ng/ul	316582	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, methyl-	98	C7H14	000108-87-2	94
2	Cyclohexane, methyl-	98	C7H14	000108-87-2	90
3	Cyclohexane, methyl-	98	C7H14	000108-87-2	90
4	Cyclohexane, methyl-	98	C7H14	000108-87-2	83
5	2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	53





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

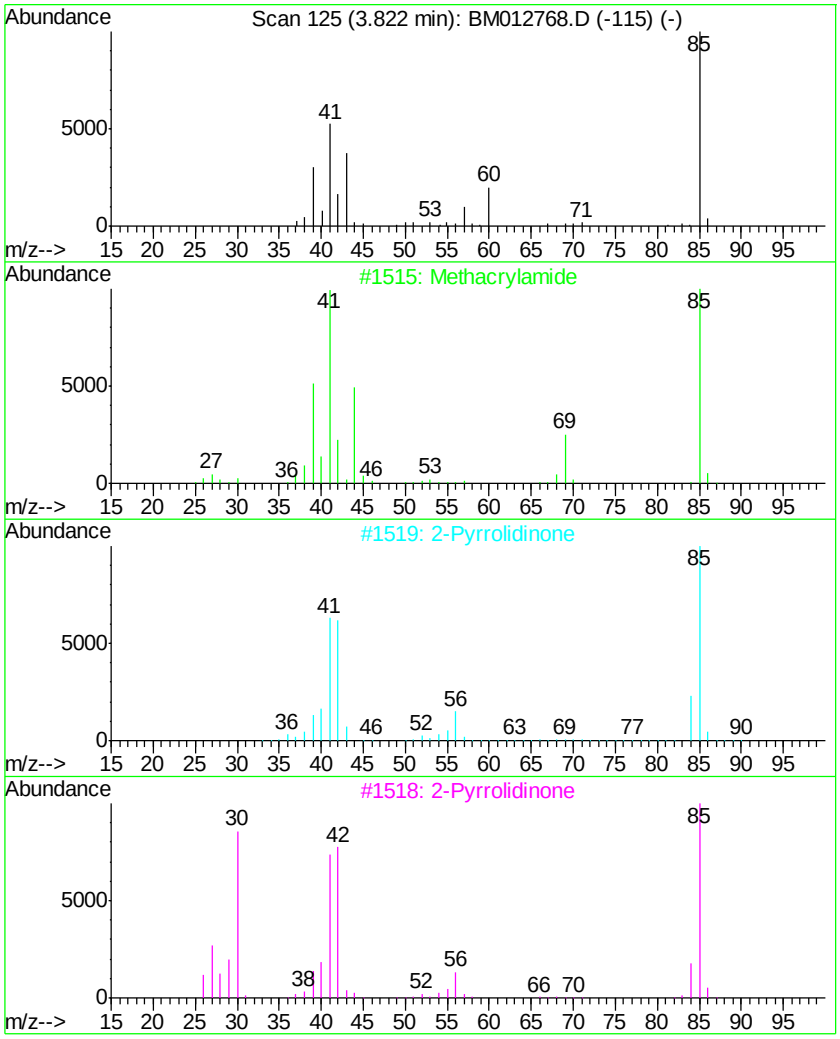
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

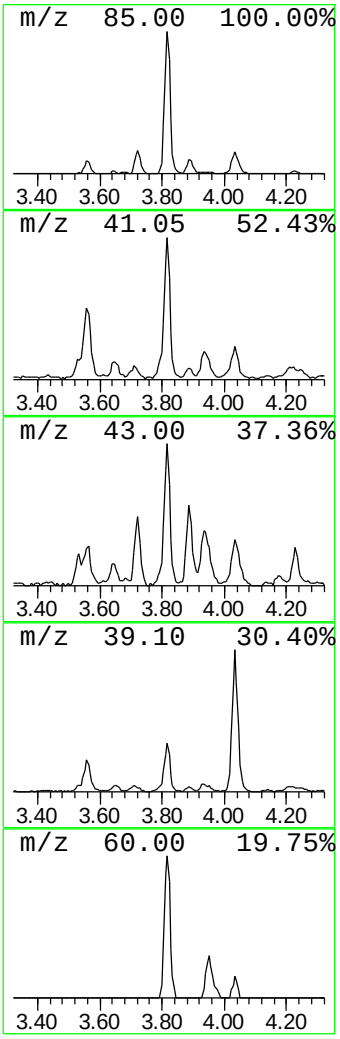
****

Peak Number 2 Methacrylamide Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.82	3.25 ng/ul	220628	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methacrylamide	85	C4H7NO	000079-39-0	50
2	2-Pyrrolidinone	85	C4H7NO	000616-45-5	42
3	2-Pyrrolidinone	85	C4H7NO	000616-45-5	40
4	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
5	Methacrylamide	85	C4H7NO	000079-39-0	9





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

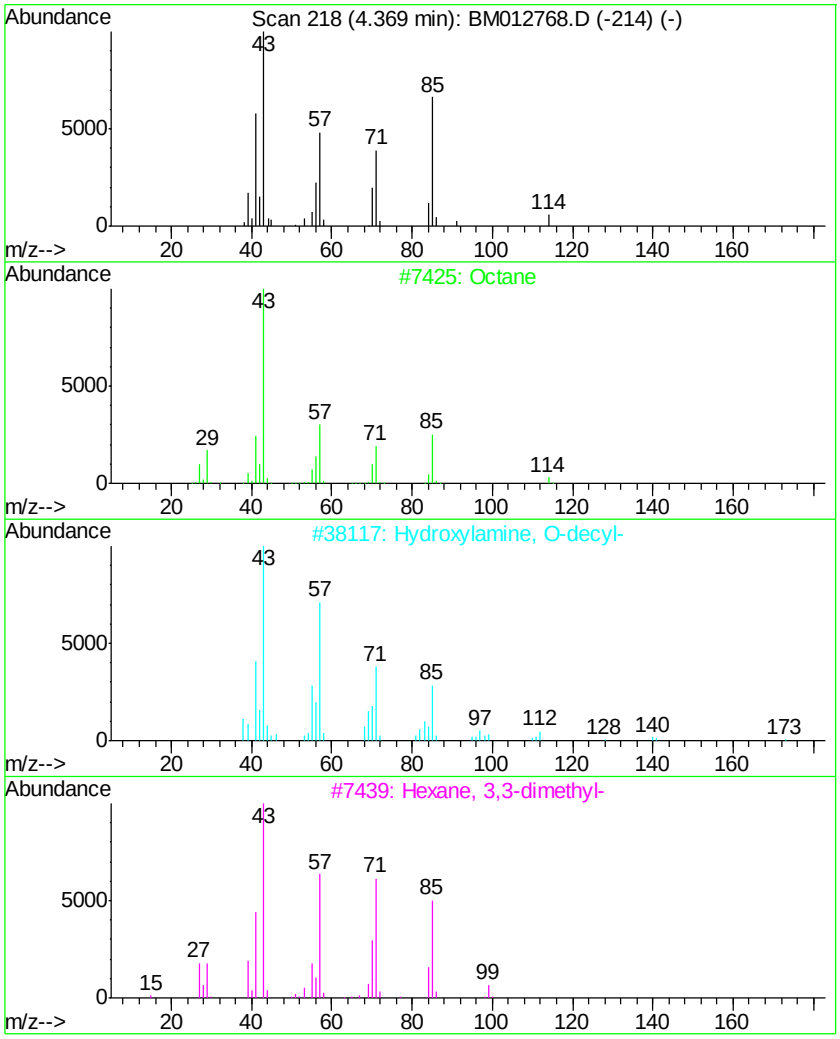
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

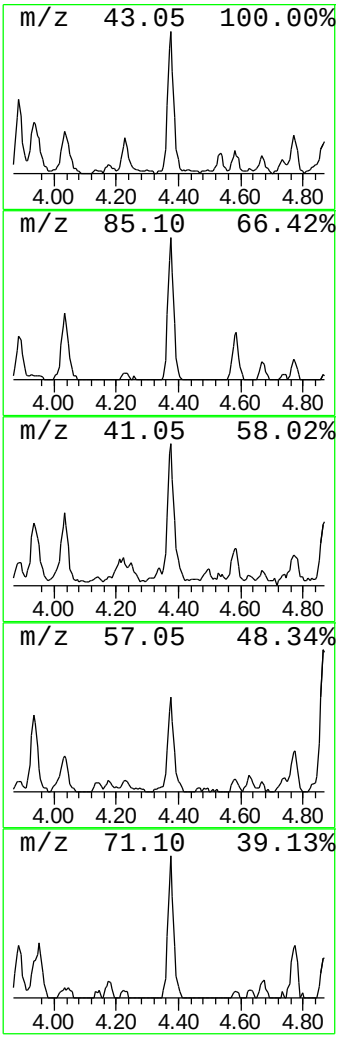
****

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	2.30 ng/ul	155937	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	83
2	Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	78
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
5	Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	59





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

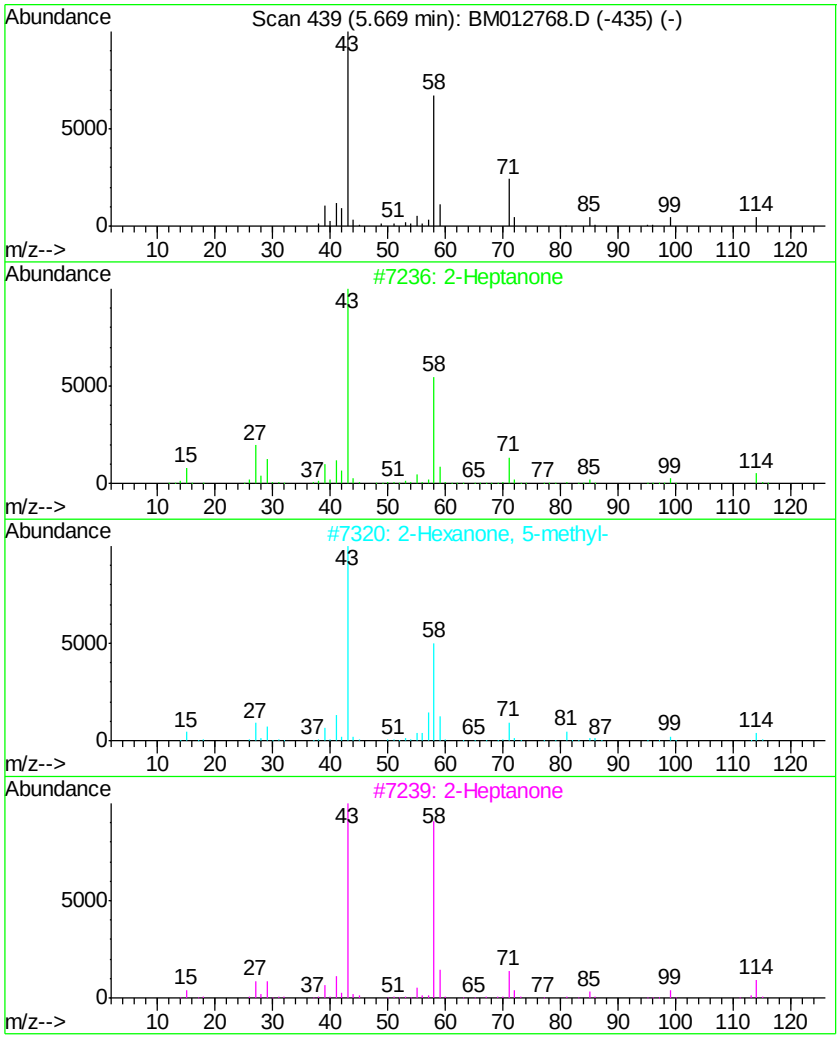
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

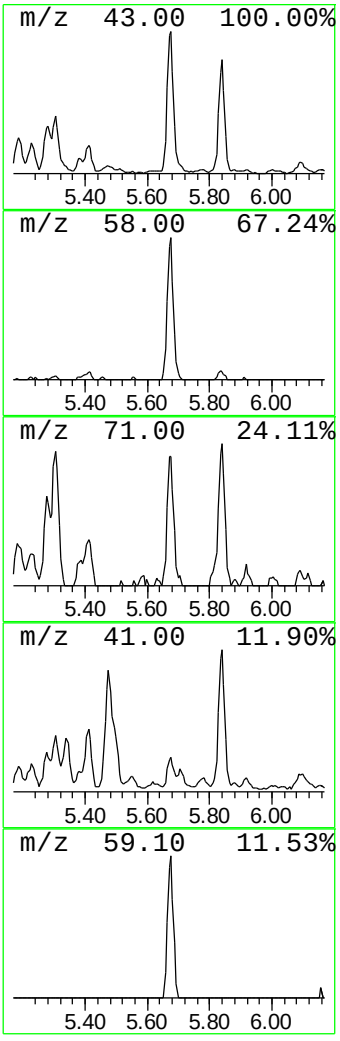
****

Peak Number 7 2-Heptanone Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.67	2.05 ng/ul	139249	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone	114	C7H14O	000110-43-0	87
2	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	80
3	2-Heptanone	114	C7H14O	000110-43-0	72
4	2-Heptanone	114	C7H14O	000110-43-0	68
5	2-Pentanone, 4,4-dimethyl-	114	C7H14O	000590-50-1	43





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

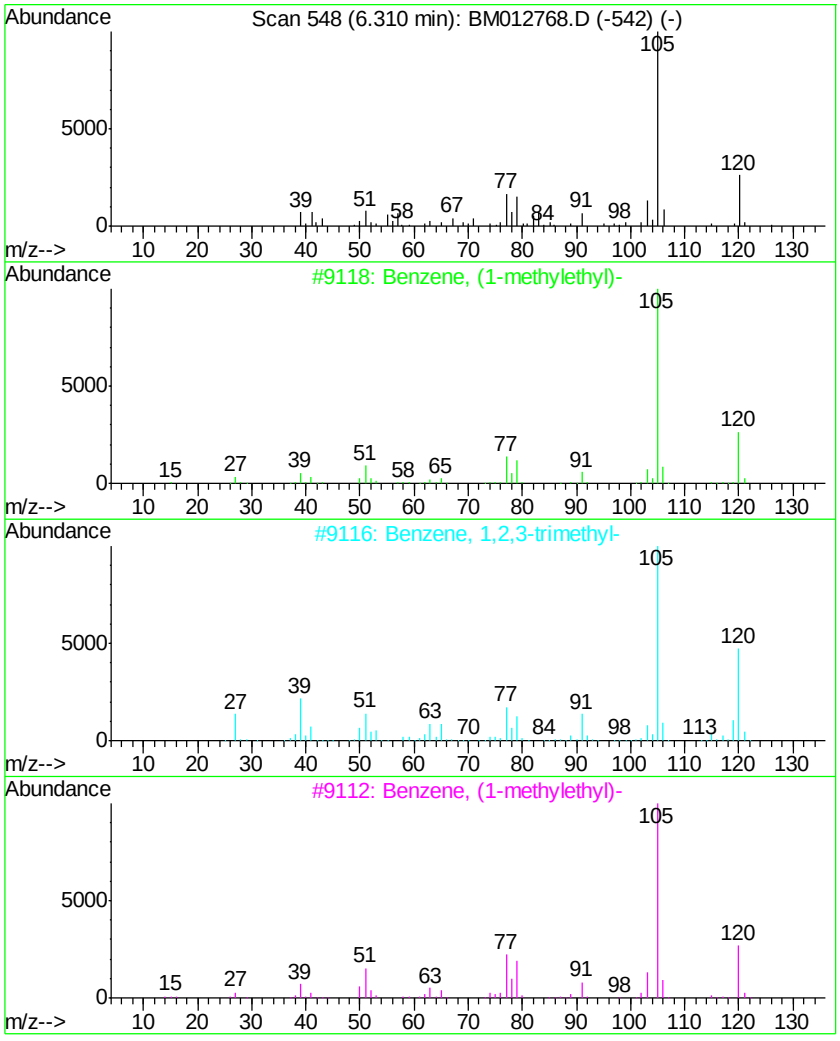
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

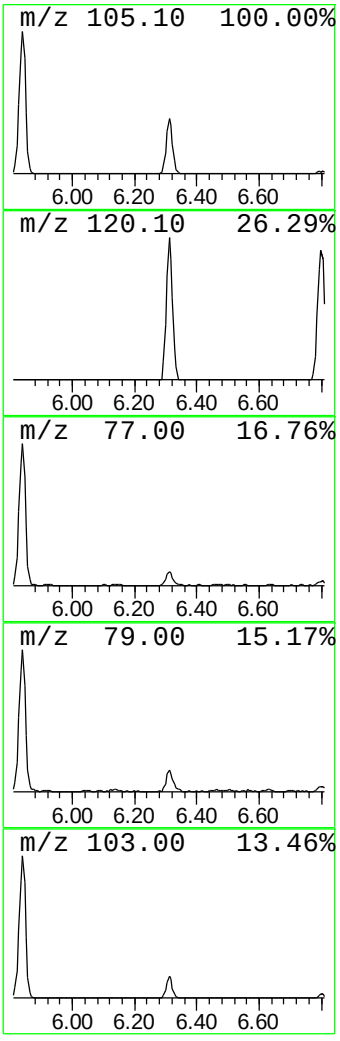
****

Peak Number 9 Benzene, (1-methylethyl)- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	2.90 ng/ul	196972	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	86





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

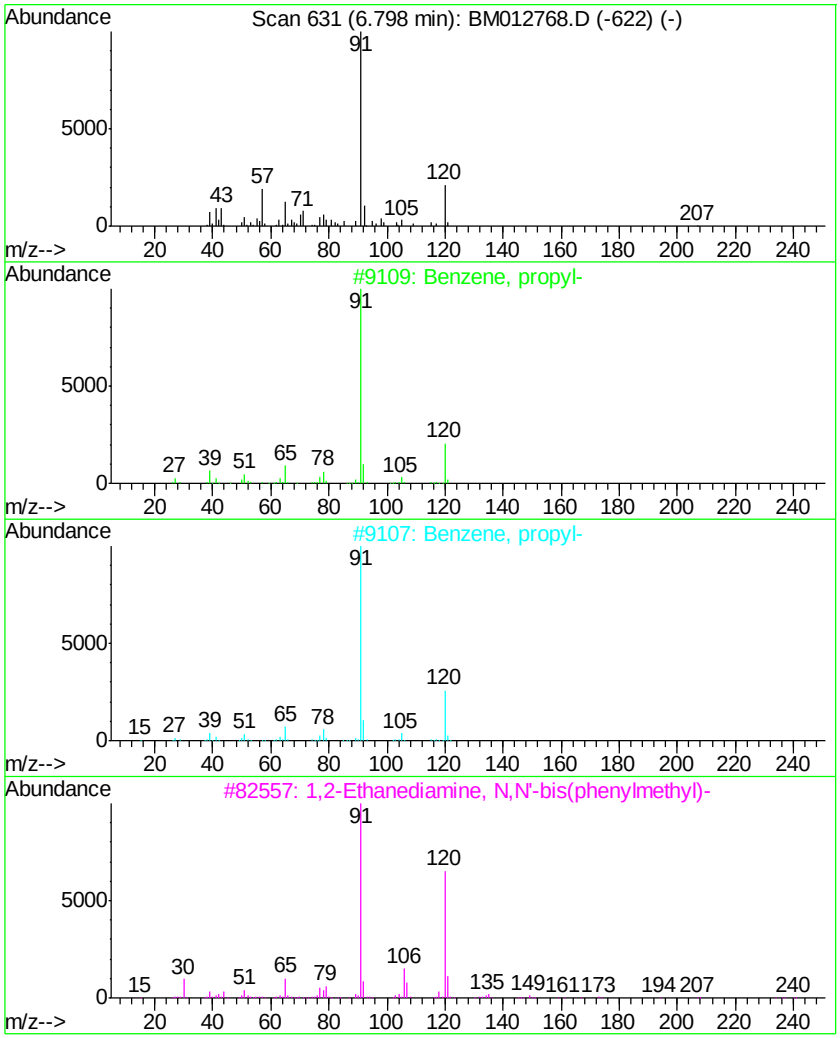
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

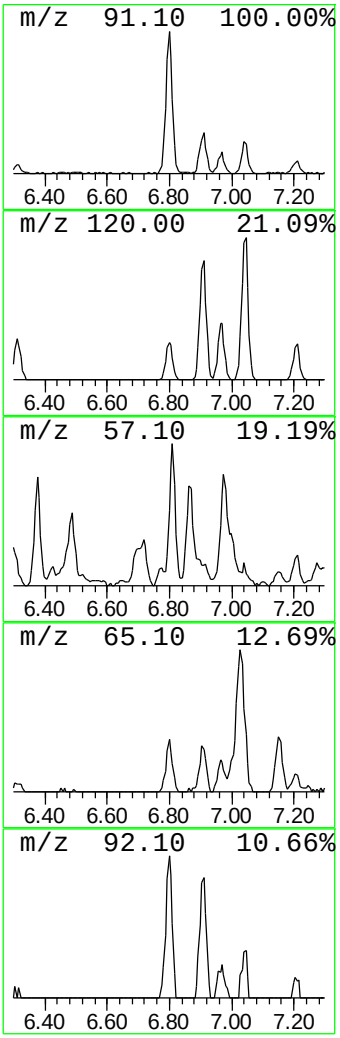
****

Peak Number 10 Benzene, propyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	3.68 ng/ul	249348	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	60
2	Benzene, propyl-	120	C9H12	000103-65-1	53
3	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	47
4	Propanenitrile, 3-[(phenylmethyl...	160	C10H12N2	000706-03-6	47
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	47





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

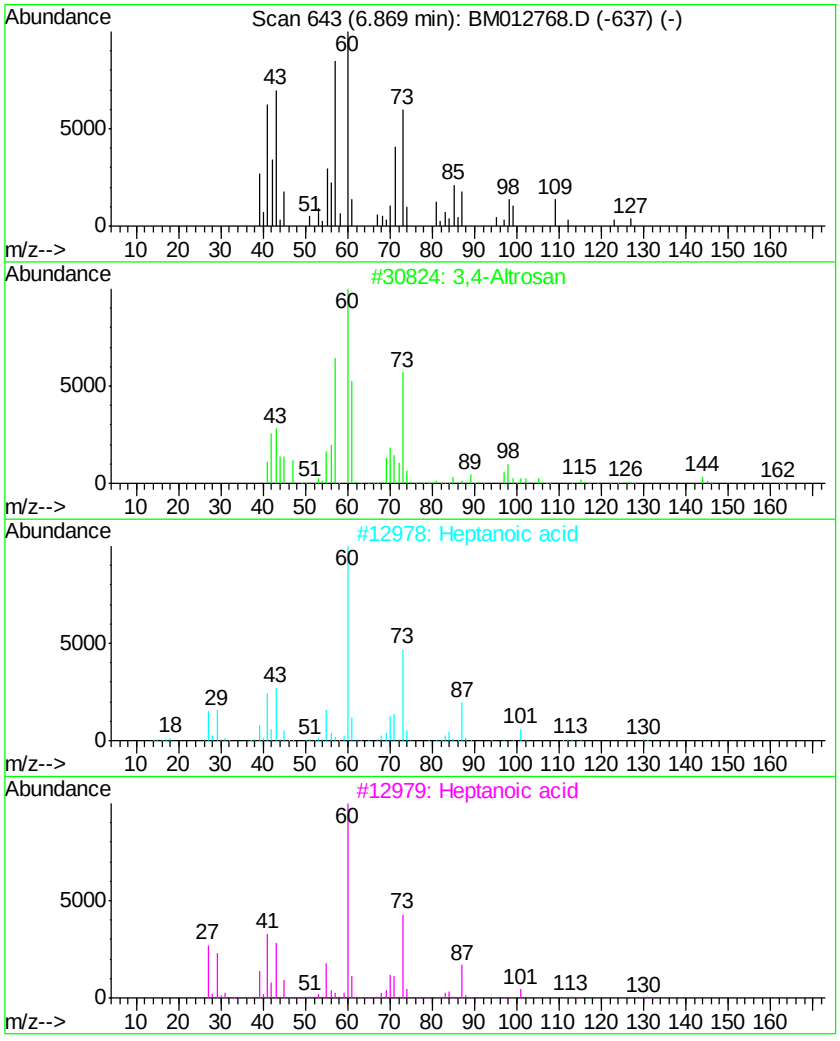
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

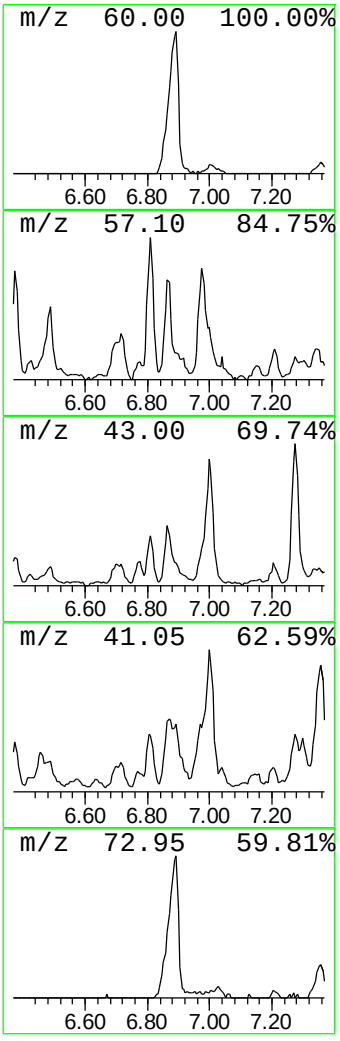
****

Peak Number 11 unknown-01 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.87	2.17 ng/ul	147330	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Altrosan	162	C6H10O5	1000129-76-4	38
2	Heptanoic acid	130	C7H14O2	000111-14-8	38
3	Heptanoic acid	130	C7H14O2	000111-14-8	37
4	Nonane, 2-methyl-	142	C10H22	000871-83-0	27
5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	27





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

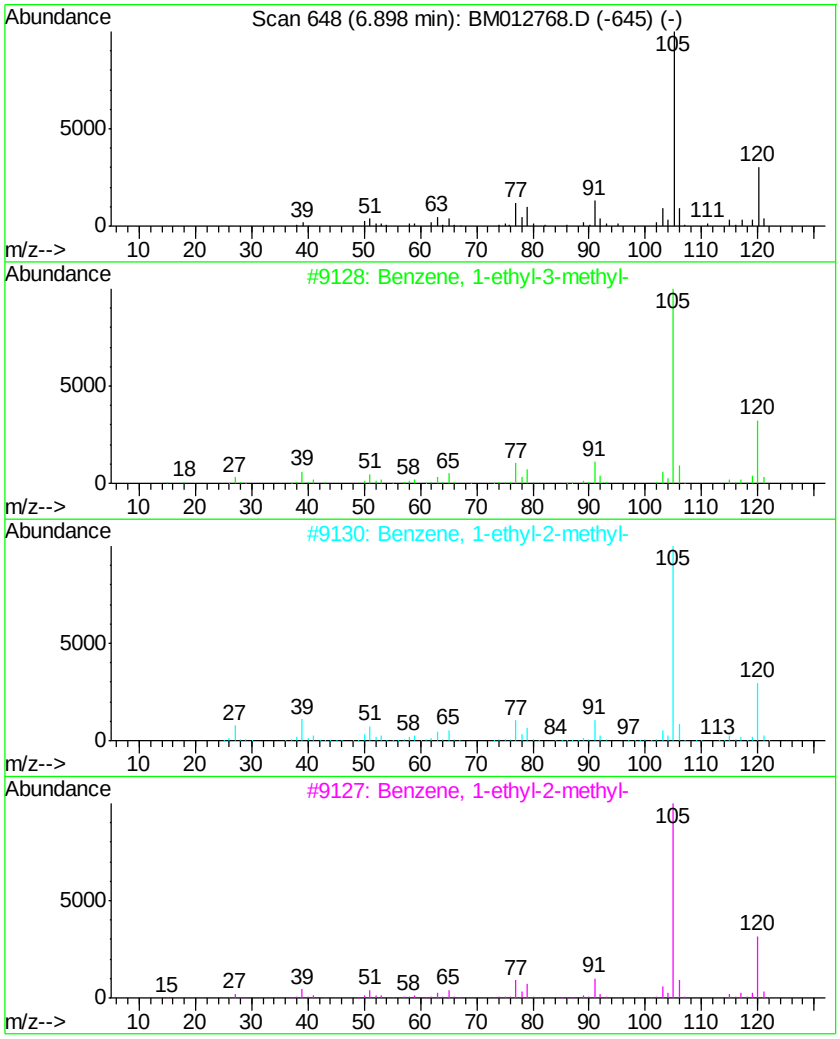
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

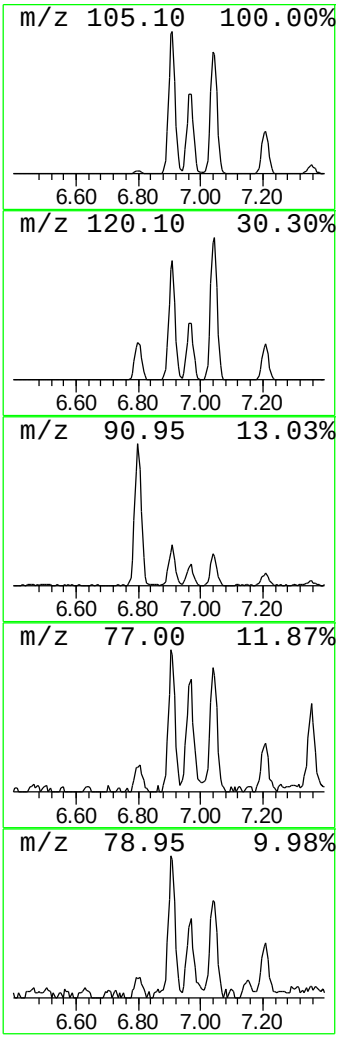
****

Peak Number 12 Benzene, 1-ethyl-3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.90	5.79 ng/ul	392740	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

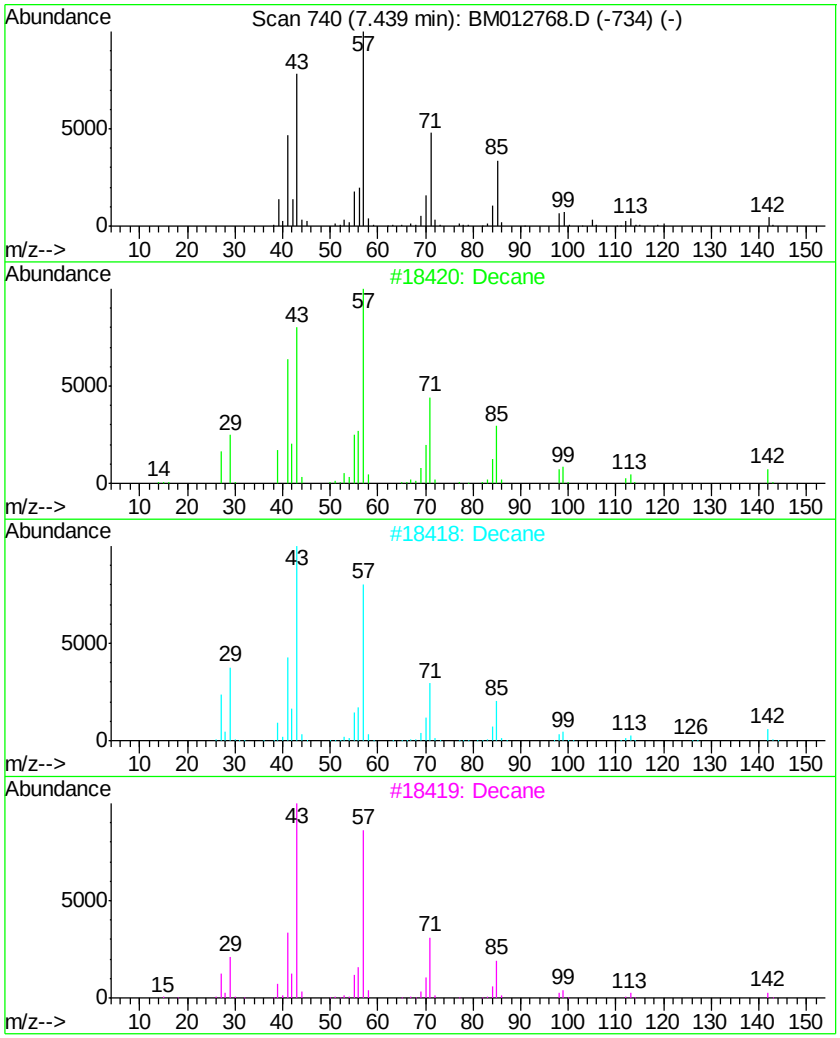
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

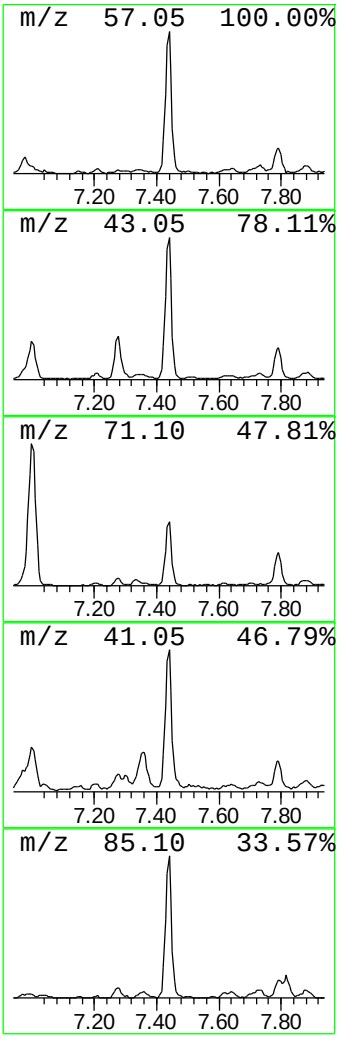
****

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.44	9.91 ng/ul	672459	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	95
2	Decane	142	C10H22	000124-18-5	95
3	Decane	142	C10H22	000124-18-5	94
4	Decane	142	C10H22	000124-18-5	87
5	Undecane	156	C11H24	001120-21-4	83





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

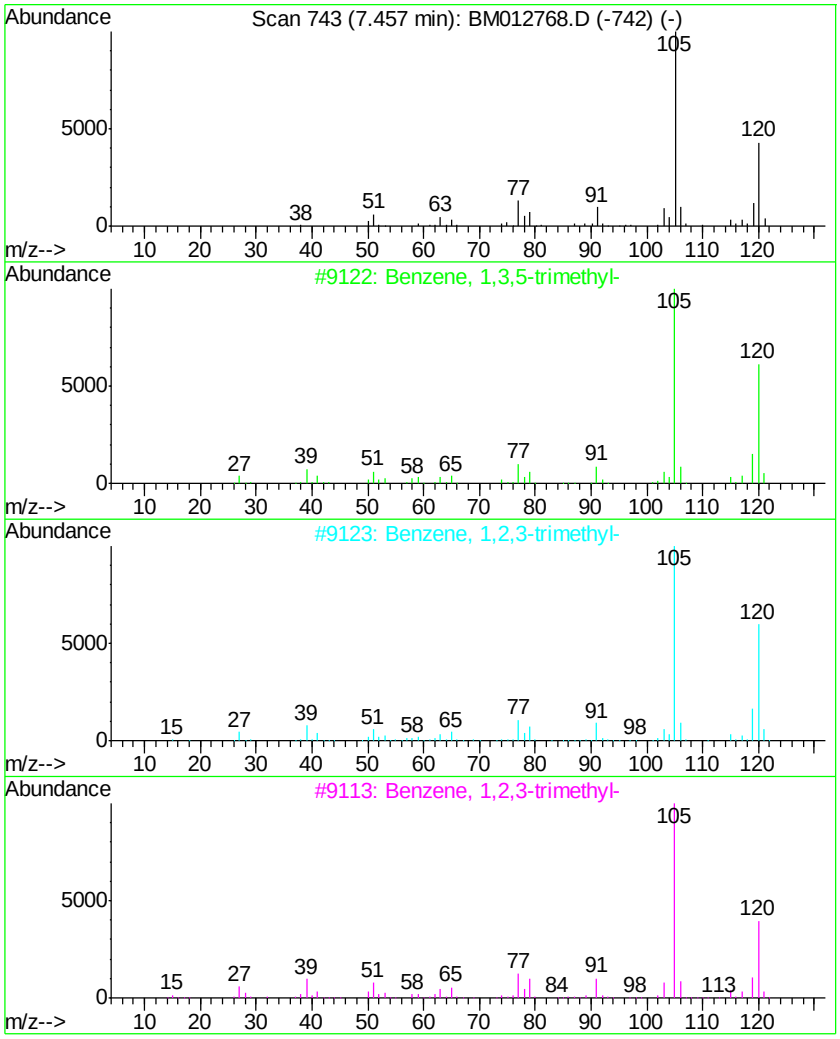
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

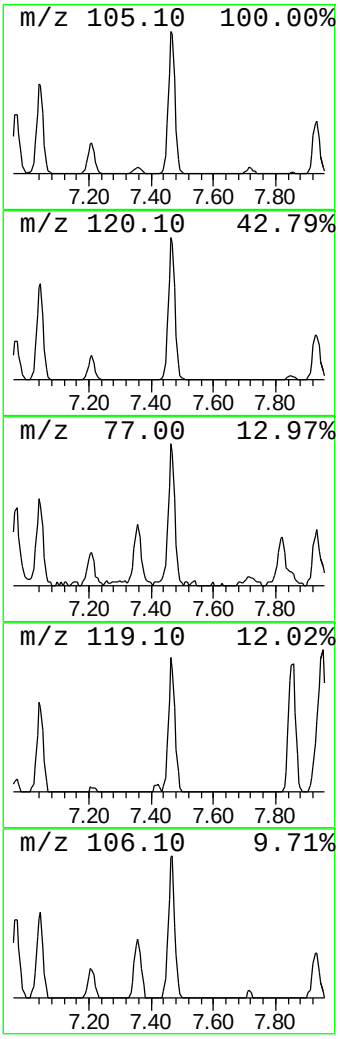
****

Peak Number 14 Benzene, 1,3,5-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.46	6.38 ng/ul	432855	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

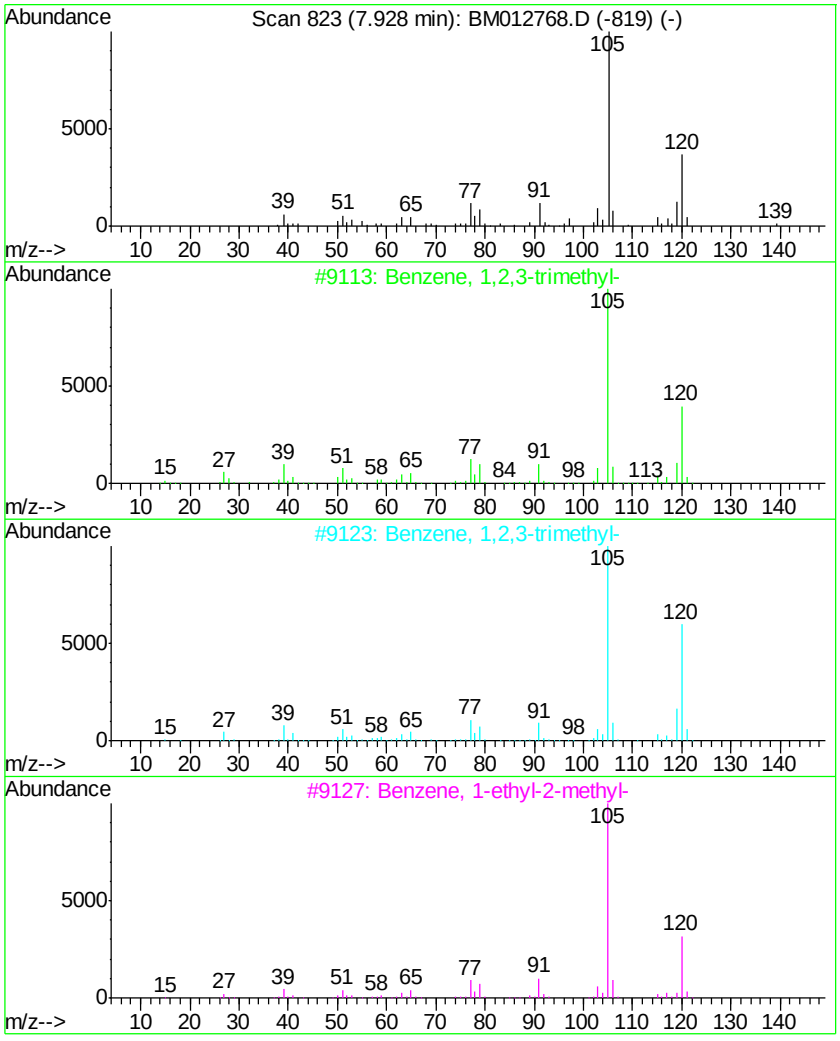
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

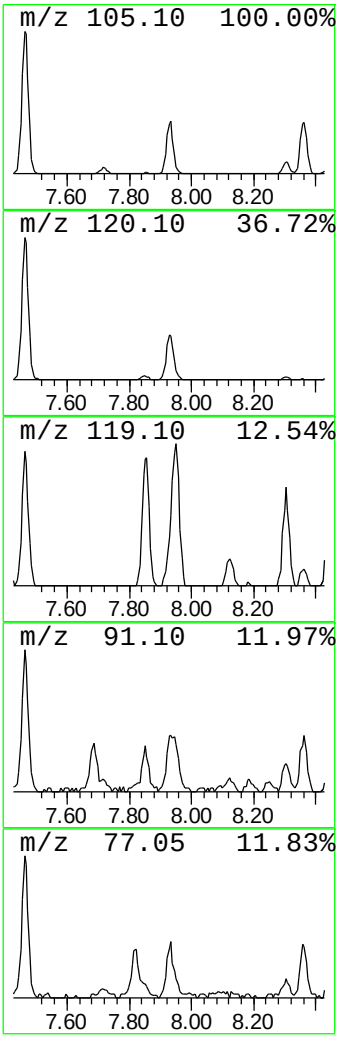
****

Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.93	3.47 ng/ul	235737	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	90





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

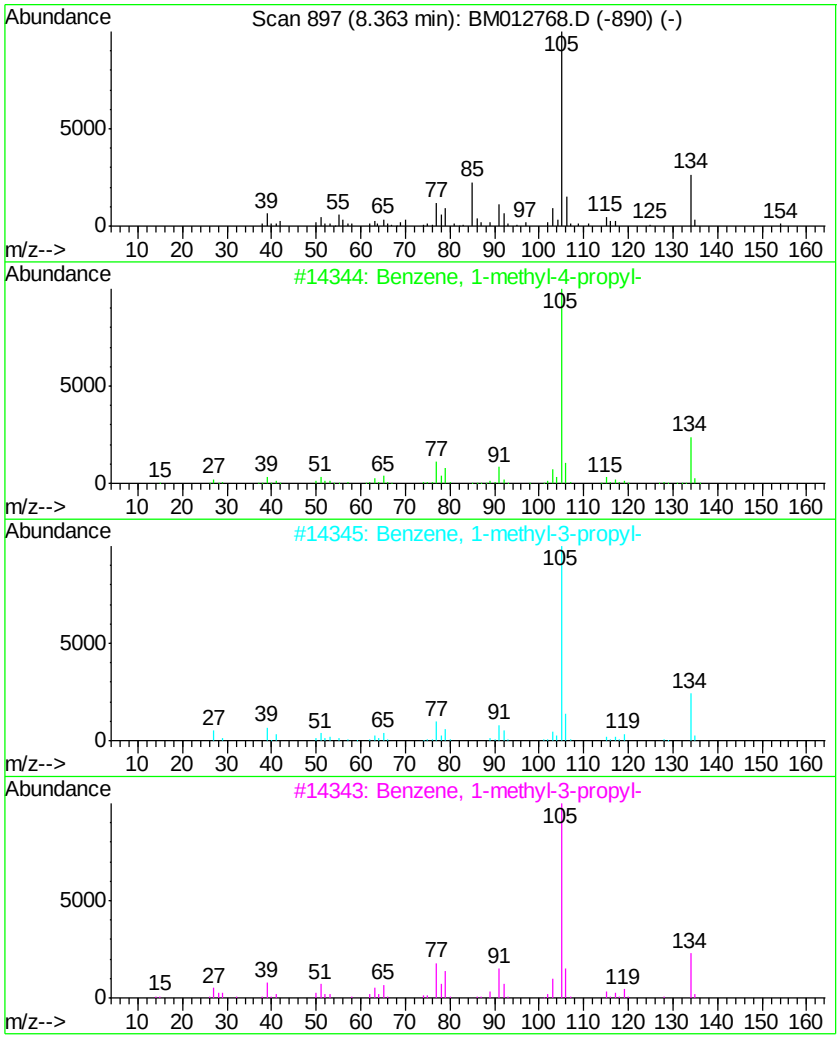
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

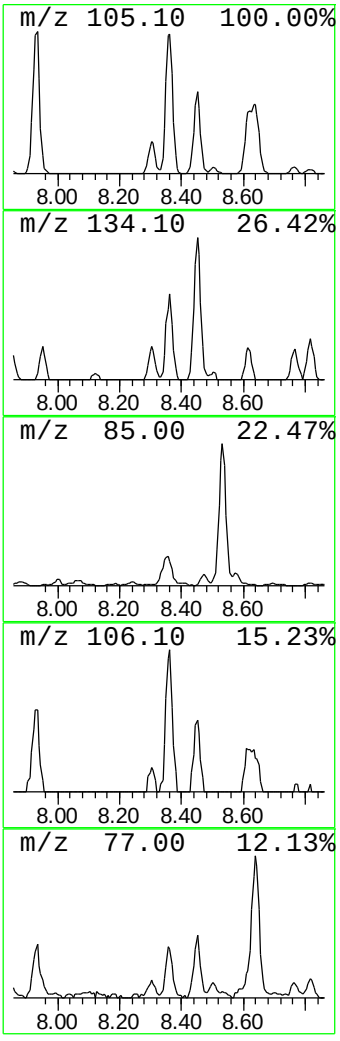
****

Peak Number 16 Benzene, 1-methyl-4-propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.36	3.68 ng/ul	249714	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	76
2	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	76
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	70
4	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	70
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	68





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

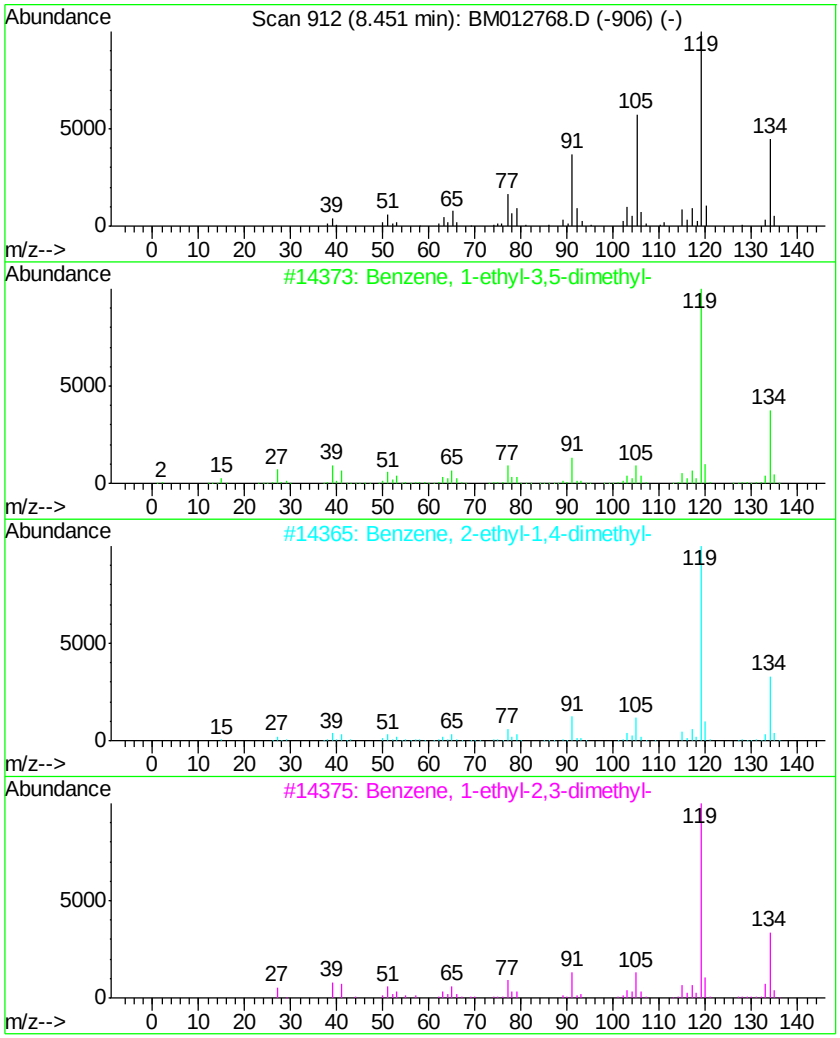
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

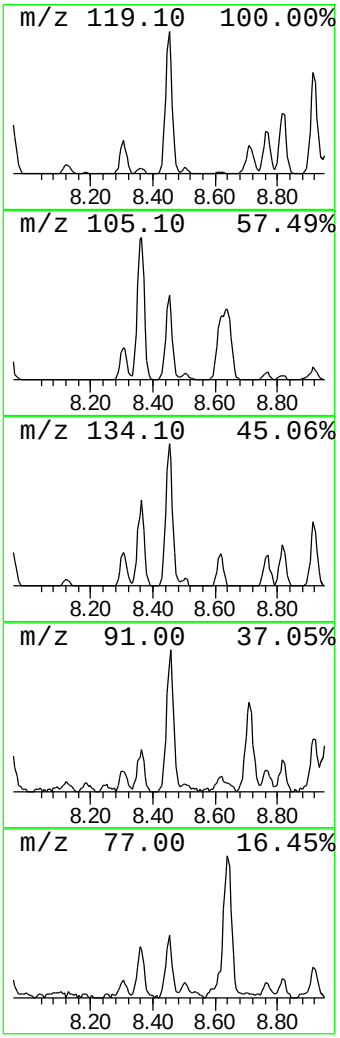
****

Peak Number 17 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.45	6.20 ng/ul	420943	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

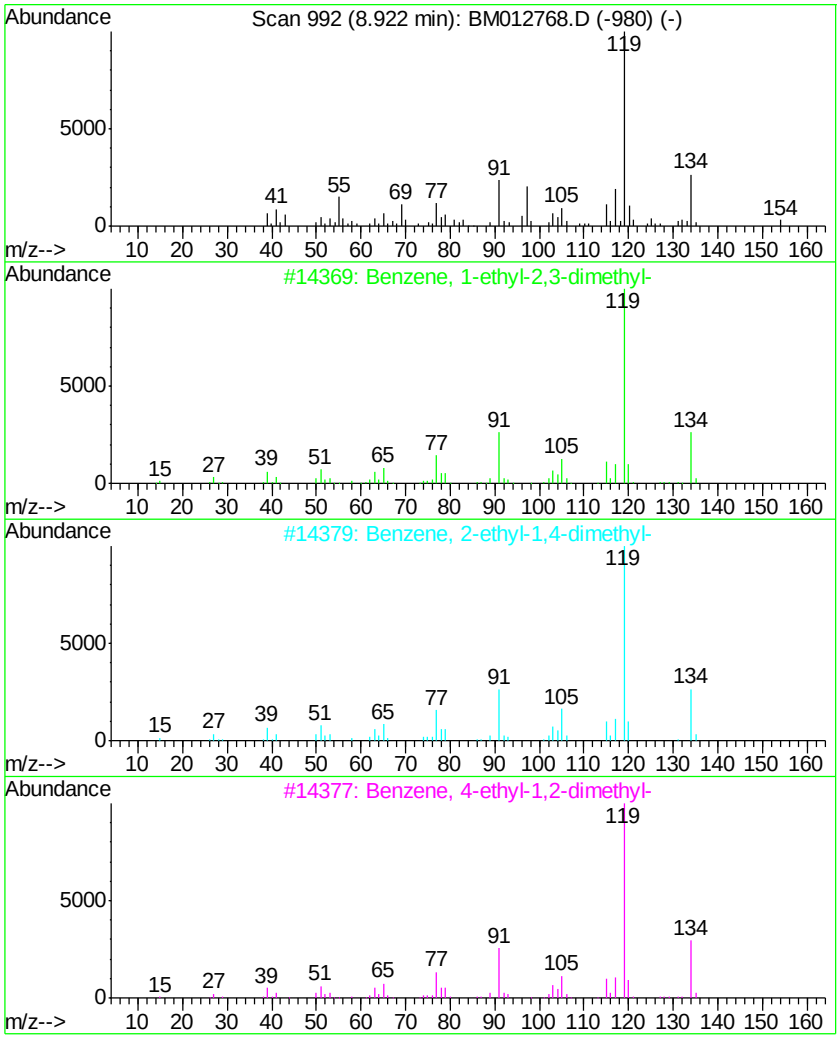
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

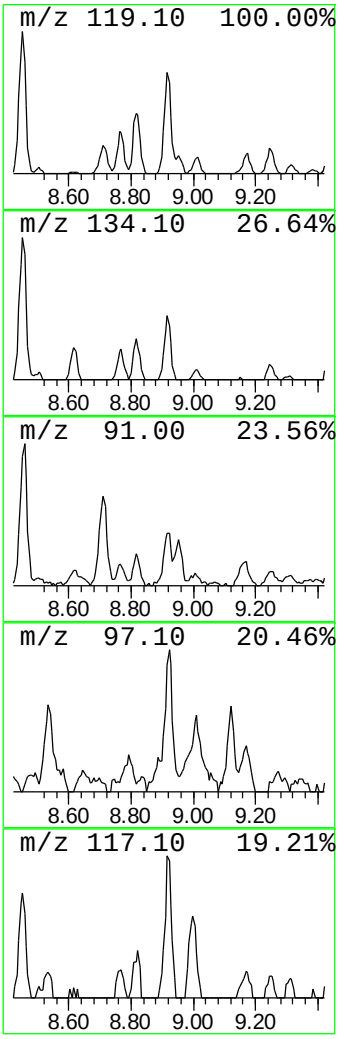
****

Peak Number 19 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.92	4.15 ng/ul	281862	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	94
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

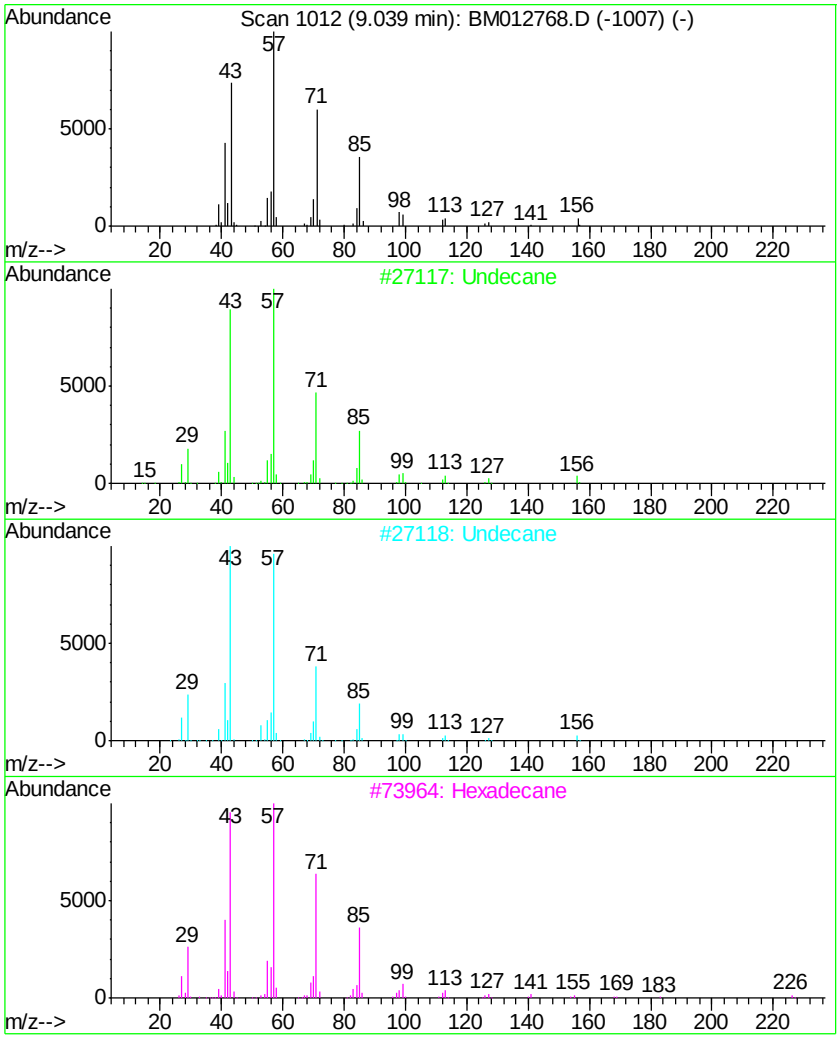
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

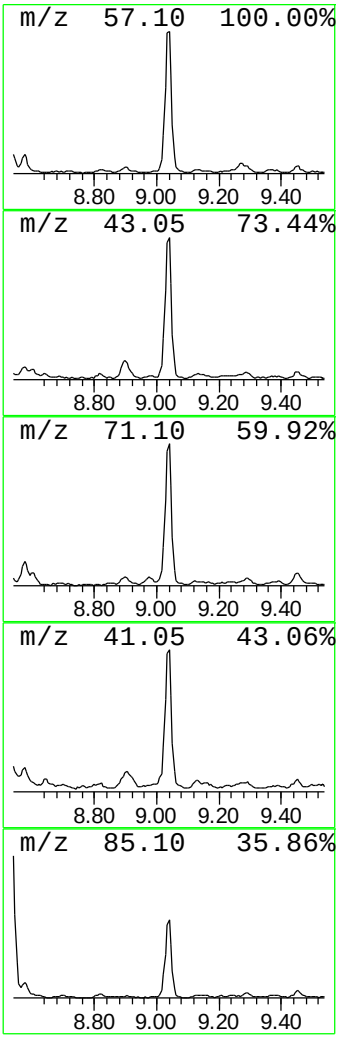
****

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.04	8.31 ng/ul	564100	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	90
2	Undecane	156	C11H24	001120-21-4	87
3	Hexadecane	226	C16H34	000544-76-3	86
4	Hexadecane	226	C16H34	000544-76-3	86
5	Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	78





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

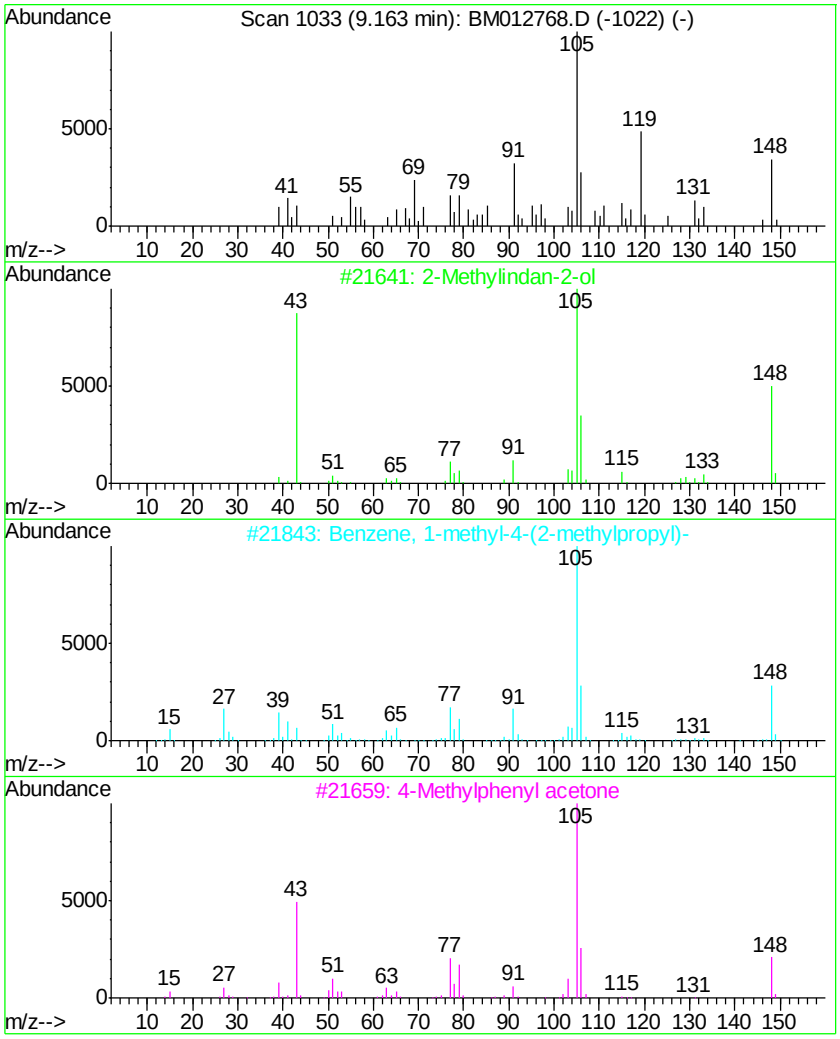
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

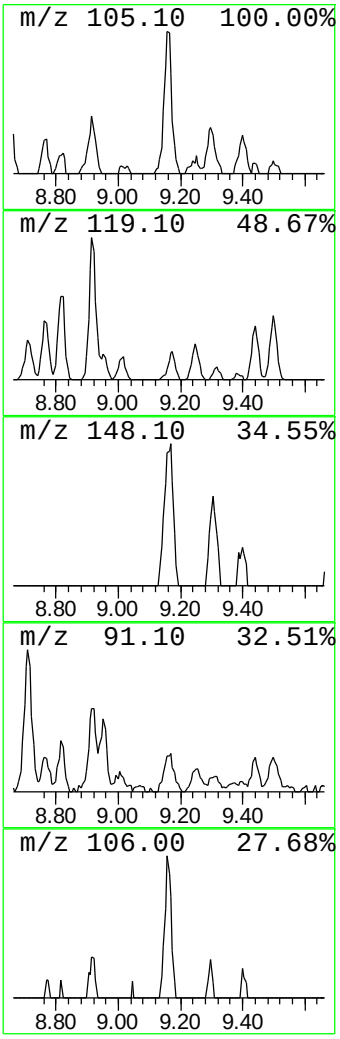
****

Peak Number 21 unknown-02 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.16	2.38 ng/ul	161766	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindan-2-ol	148	C10H12O	033223-84-6	38
2	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	38
3	4-Methylphenyl acetone	148	C10H12O	002096-86-8	30
4	Benzene, (2-methoxy-2-propenyl)-	148	C10H12O	026473-60-9	27
5	Benzene, (1,2-dimethylpropyl)-	148	C11H16	004481-30-5	22





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

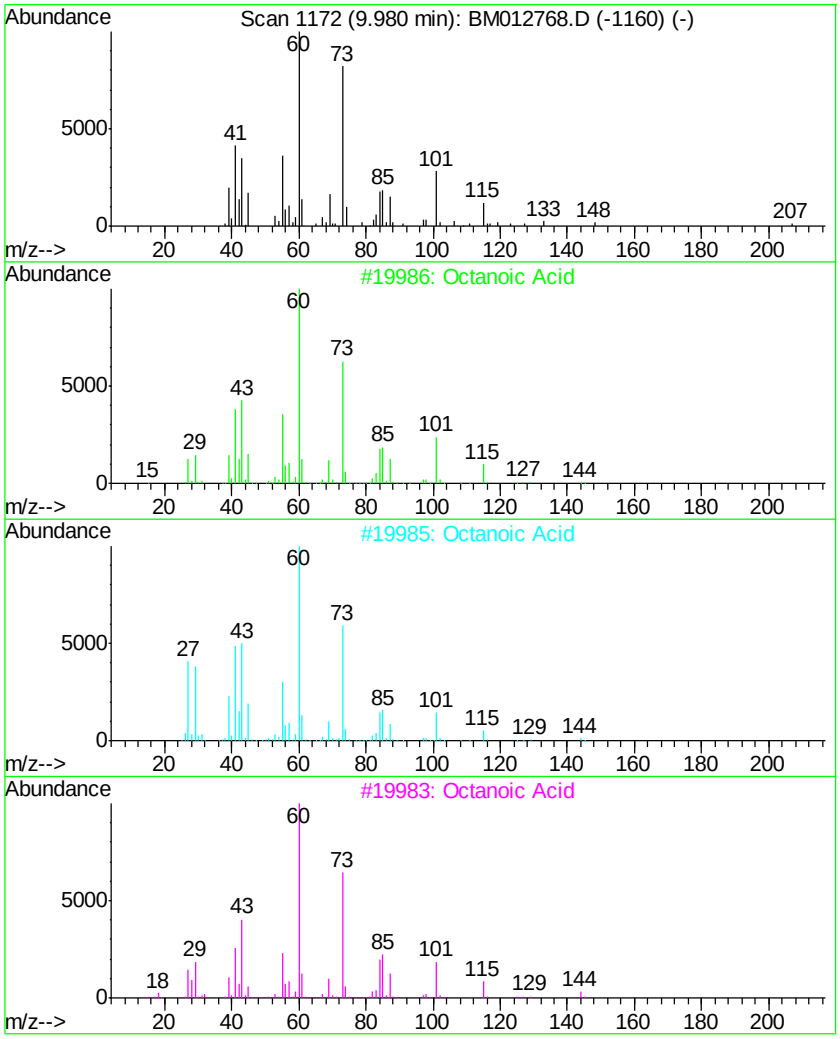
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

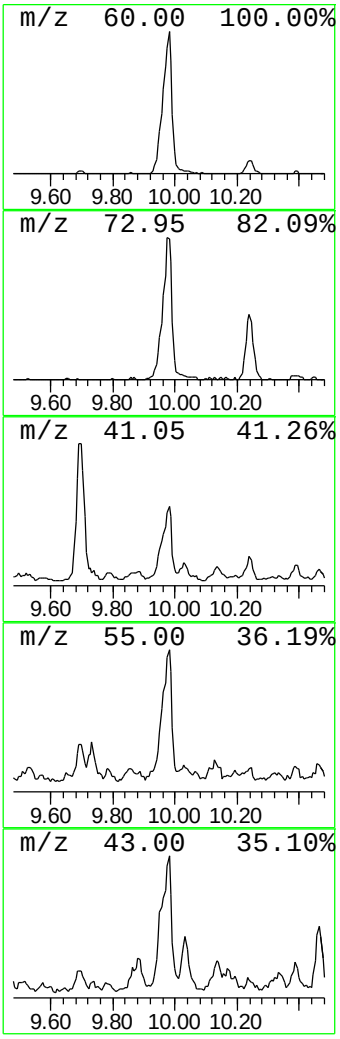
****

Peak Number 22 Octanoic Acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.98	3.98 ng/ul	363457	Naphthalene-d8	10.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic Acid	144	C8H16O2	000124-07-2	91
2	Octanoic Acid	144	C8H16O2	000124-07-2	83
3	Octanoic Acid	144	C8H16O2	000124-07-2	72
4	Hexanoic acid	116	C6H12O2	000142-62-1	53
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

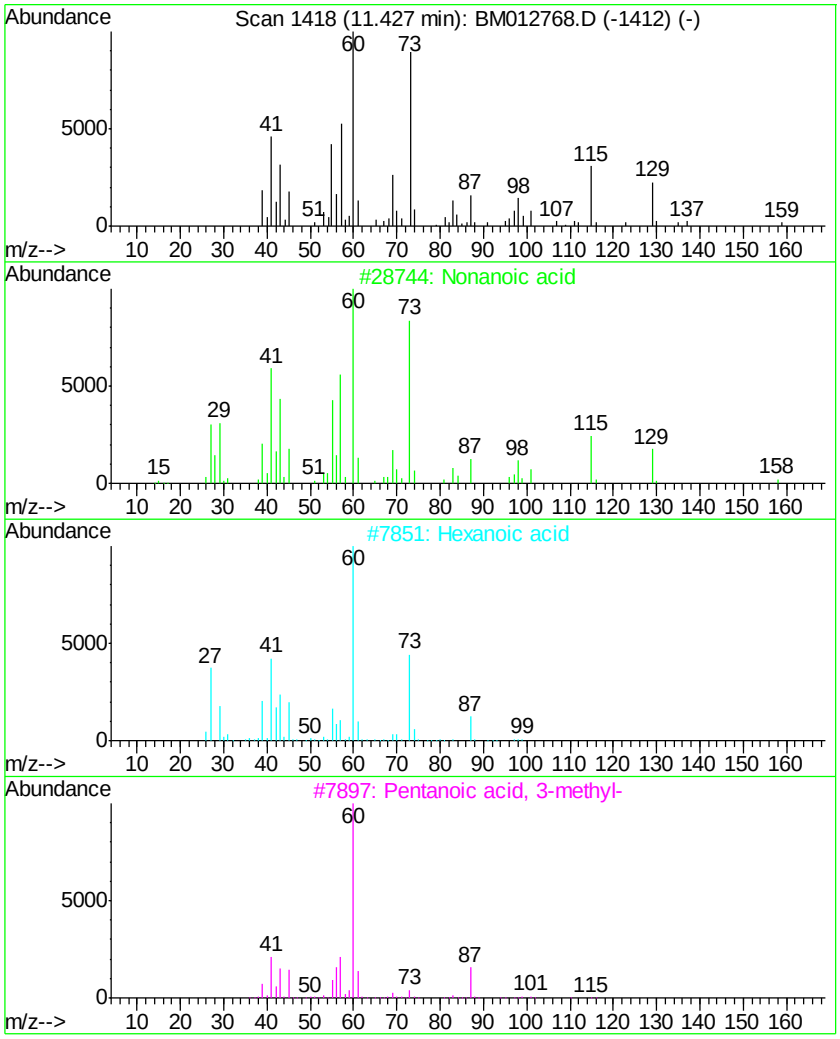
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

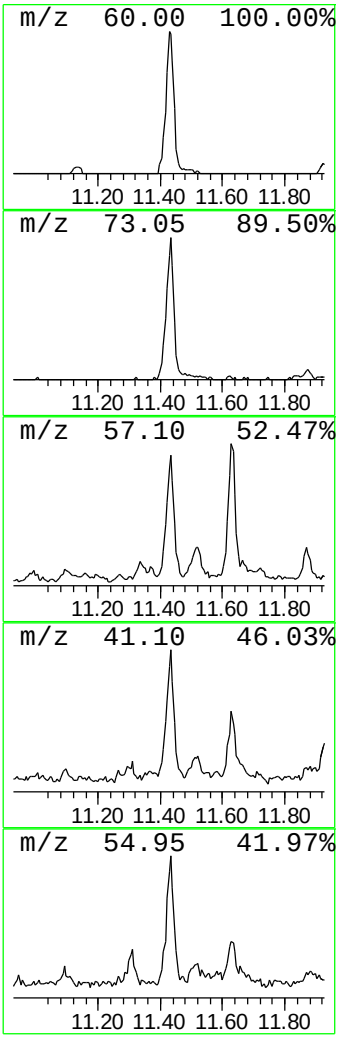
****

Peak Number 23 Nonanoic acid Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	2.50 ng/ul	228534	Naphthalene-d8	10.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	72
2	Hexanoic acid	116	C6H12O2	000142-62-1	50
3	Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	47
4	Octanoic Acid	144	C8H16O2	000124-07-2	42
5	Hexanoic acid	116	C6H12O2	000142-62-1	40





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

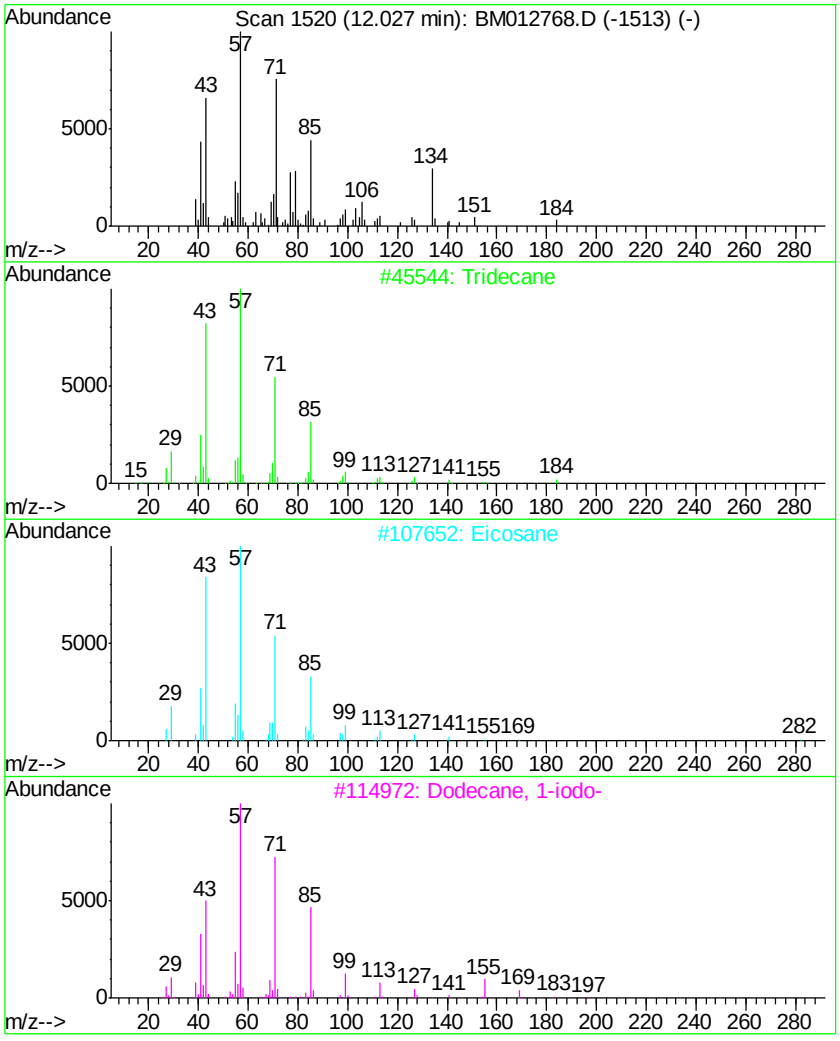
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

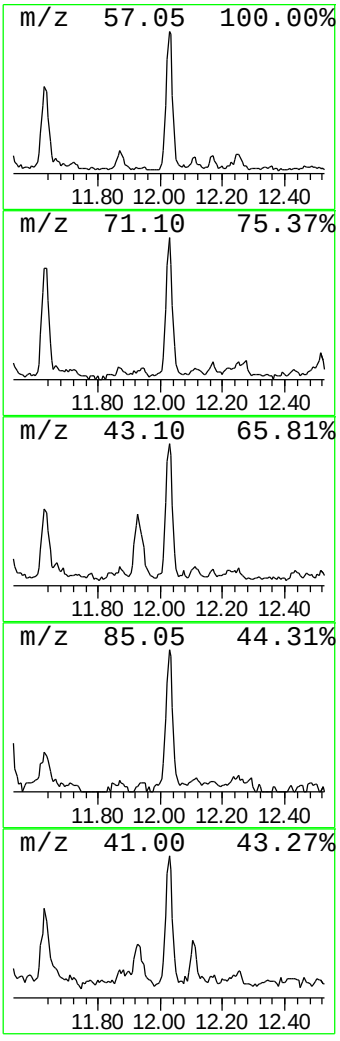
****

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	2.26 ng/ul	206156	Naphthalene-d8	10.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	60
2	Eicosane	282	C20H42	000112-95-8	53
3	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	49
4	Tetradecane	198	C14H30	000629-59-4	49
5	Dodecane	170	C12H26	000112-40-3	49





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

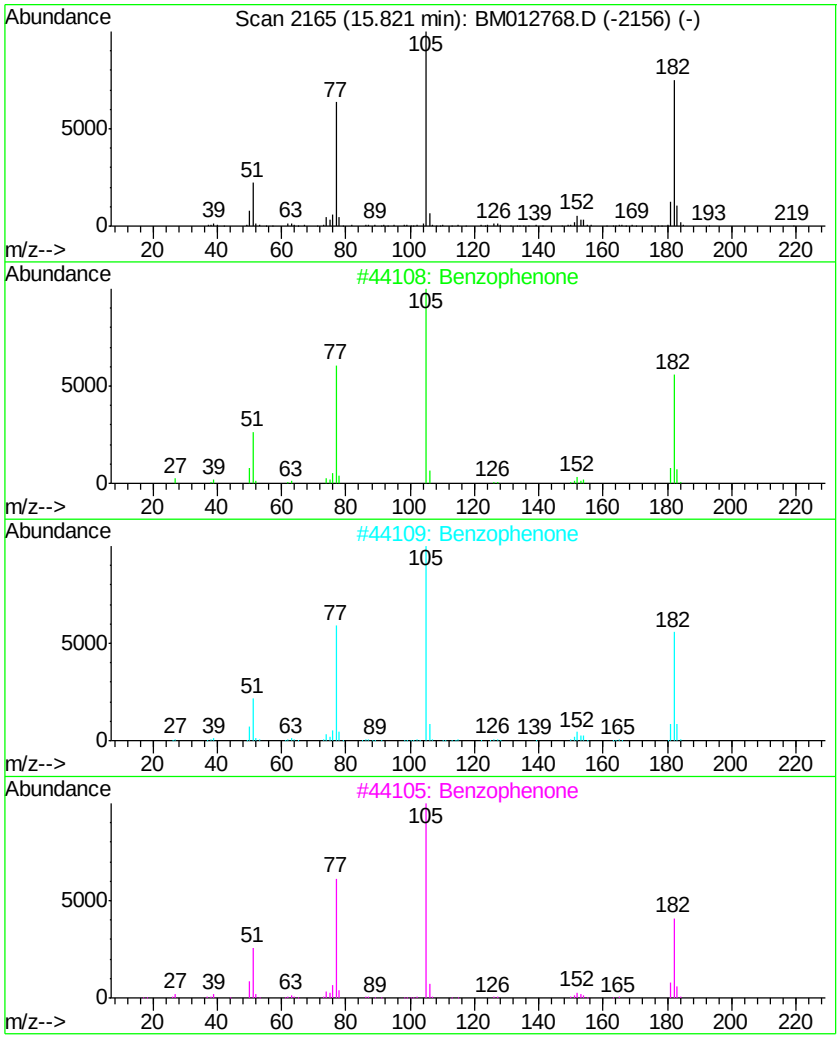
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

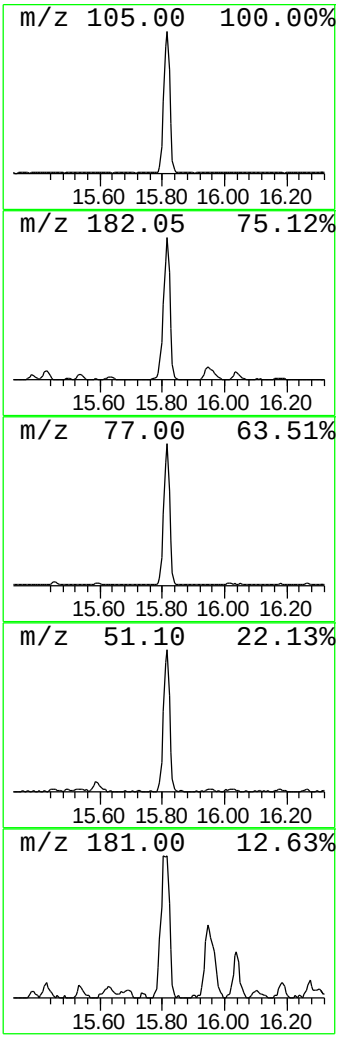
****

Peak Number 26 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	6.51 ng/ul	694068	Acenaphthene-d10	14.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	94
5	Benzophenone	182	C13H10O	000119-61-9	87





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

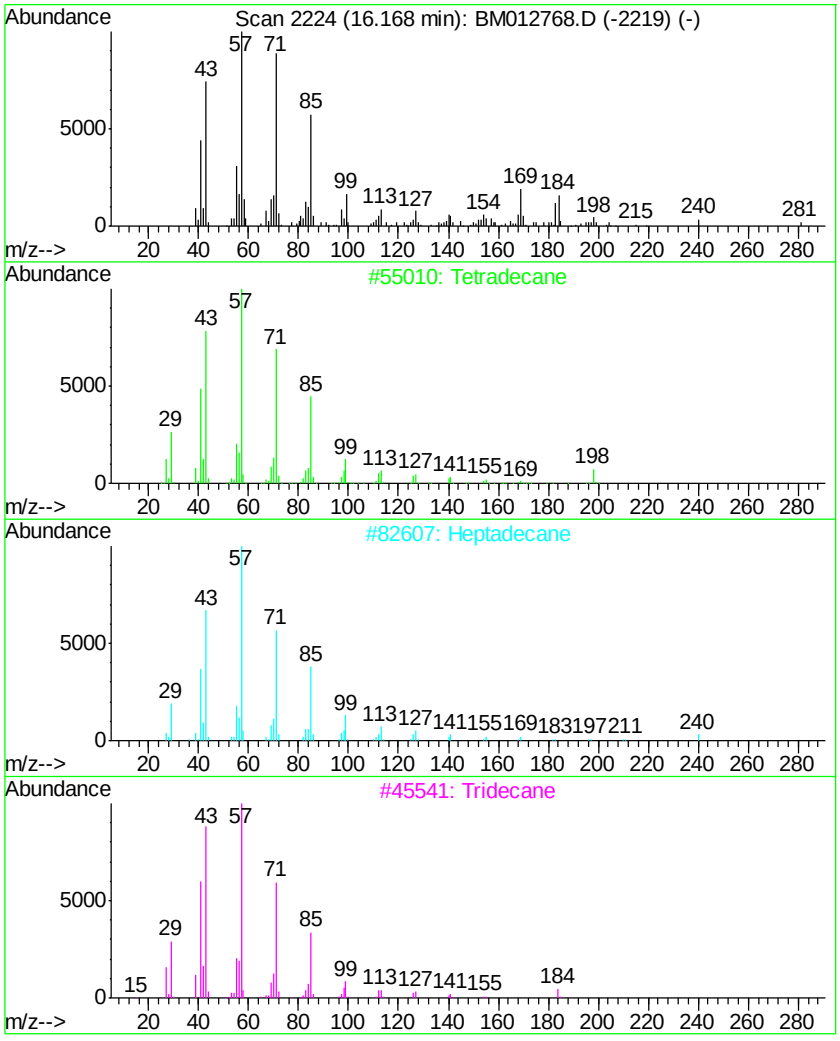
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

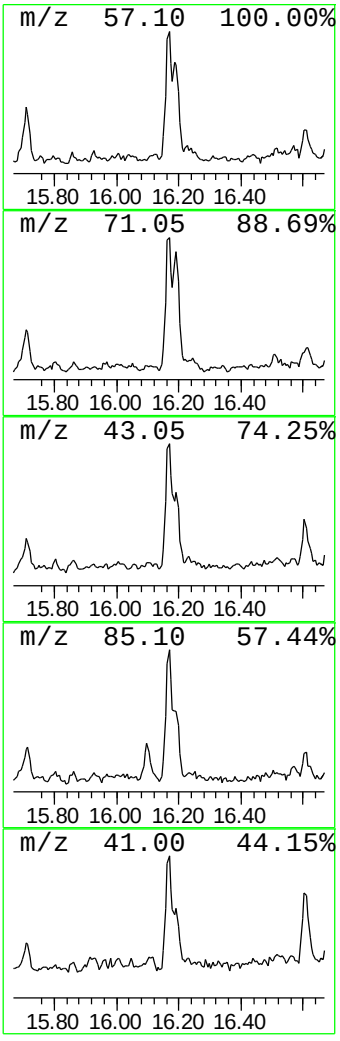
****

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.17	3.42 ng/ul	314624	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	89
2	Heptadecane	240	C17H36	000629-78-7	89
3	Tridecane	184	C13H28	000629-50-5	89
4	Heptadecane	240	C17H36	000629-78-7	87
5	Heptadecane	240	C17H36	000629-78-7	83





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

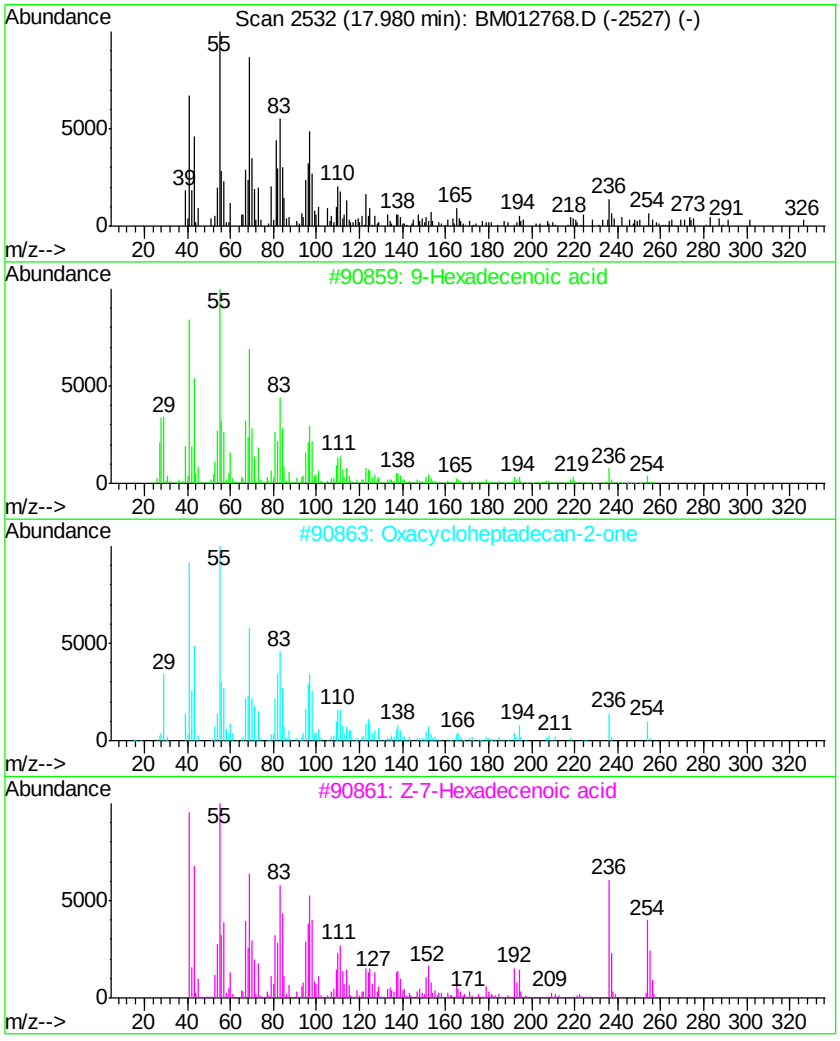
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

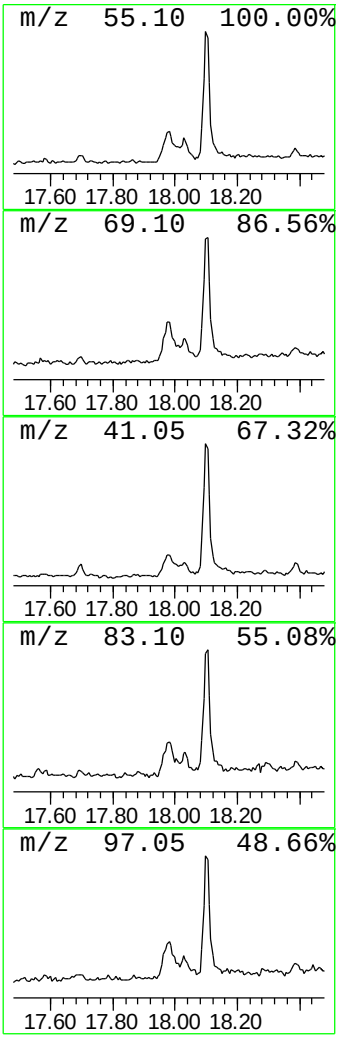
****

Peak Number 28 9-Hexadecenoic acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.98	3.51 ng/ul	322831	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Hexadecenoic acid	254	C16H30O2	002091-29-4	99
2	Oxacycloheptadecan-2-one	254	C16H30O2	000109-29-5	91
3	Z-7-Hexadecenoic acid	254	C16H30O2	1000130-90-8	90
4	1-Heptadecene	238	C17H34	006765-39-5	70
5	3-Hexadecene, (Z)-	224	C16H32	034303-81-6	64





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

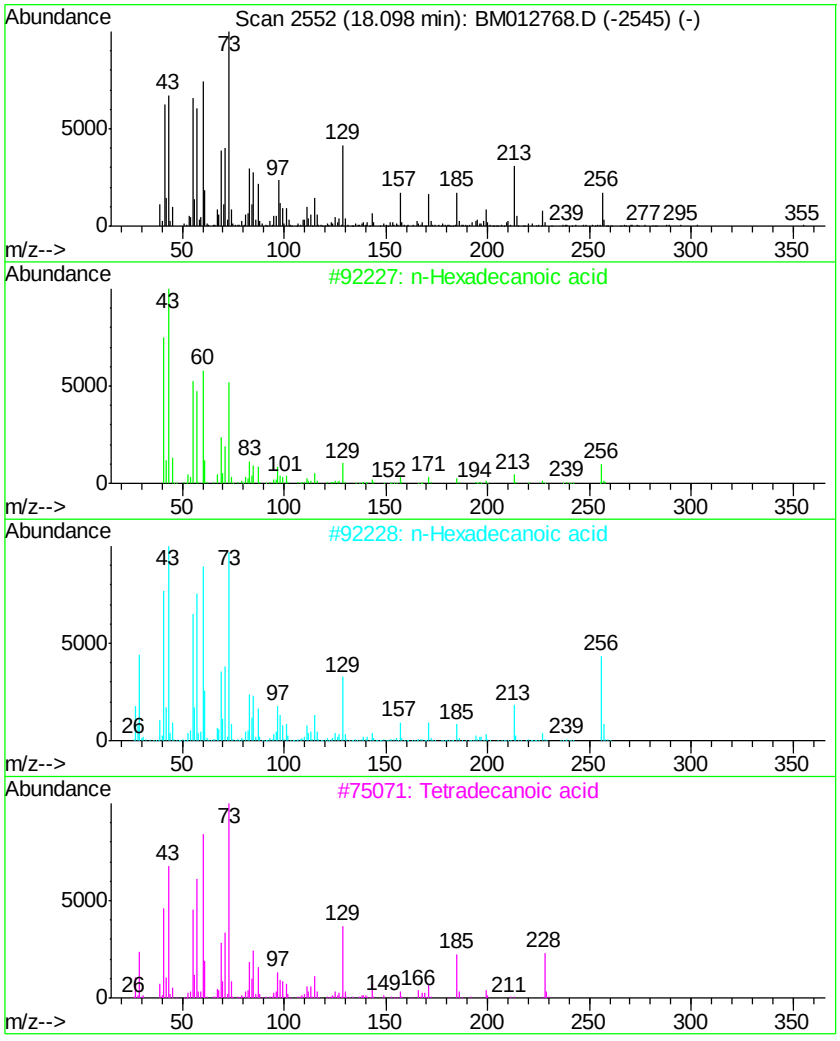
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

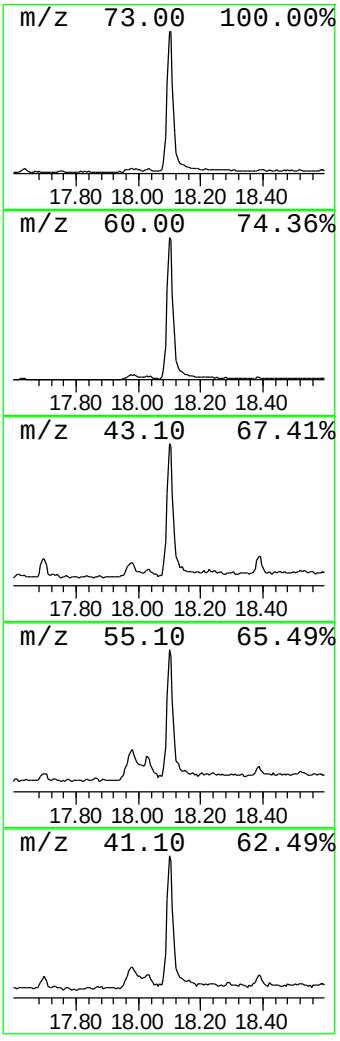
****

Peak Number 29 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.10	18.34 ng/ul	1688180	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4	Tridecanoic acid	214	C13H26O2	000638-53-9	89
5	Tridecanoic acid	214	C13H26O2	000638-53-9	86





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

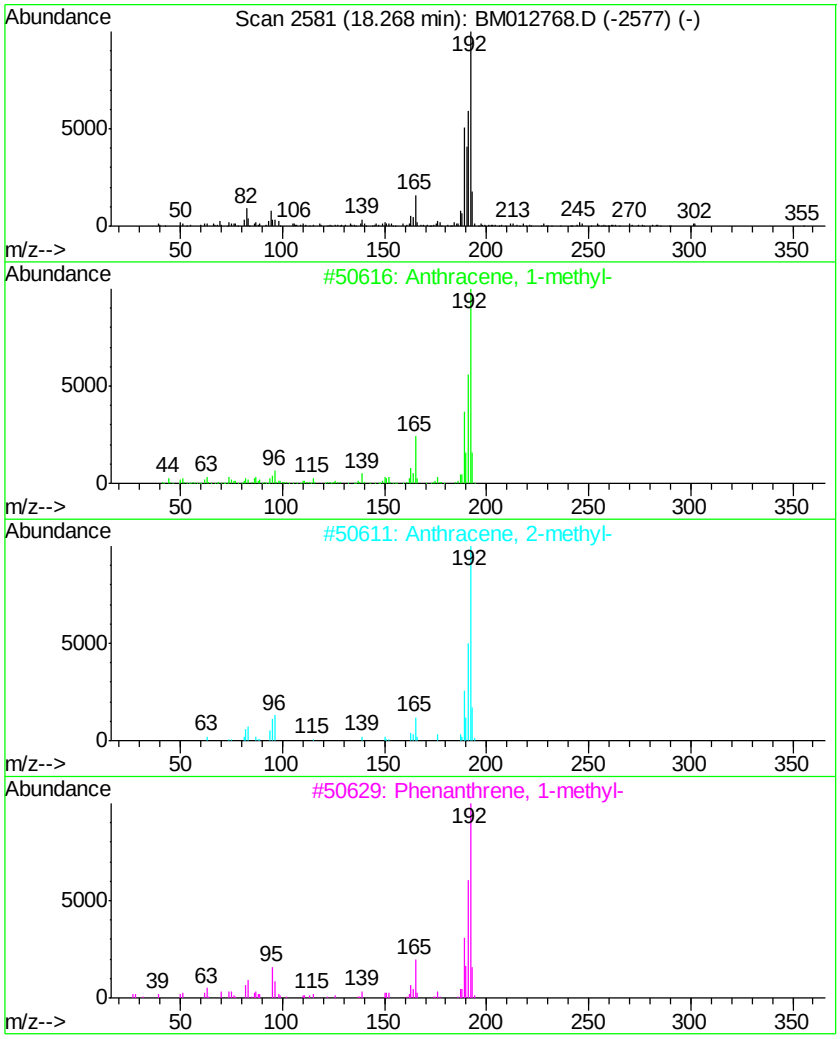
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

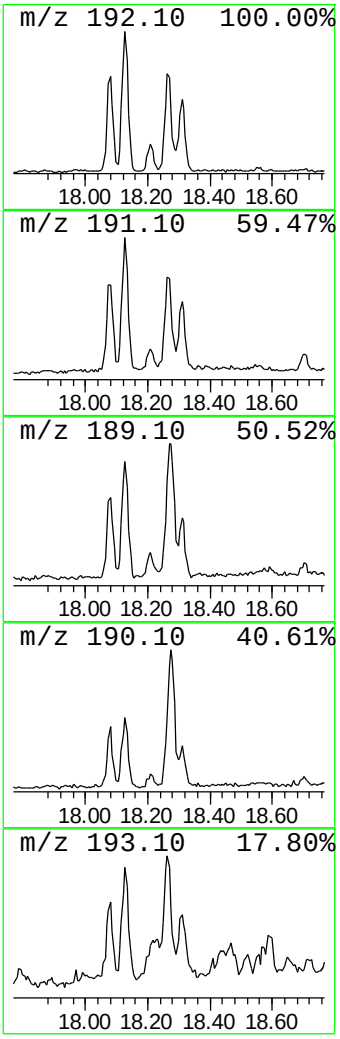
****

Peak Number 30 Anthracene, 1-methyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.27	2.16 ng/ul	198474	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	83
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	64
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	64





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

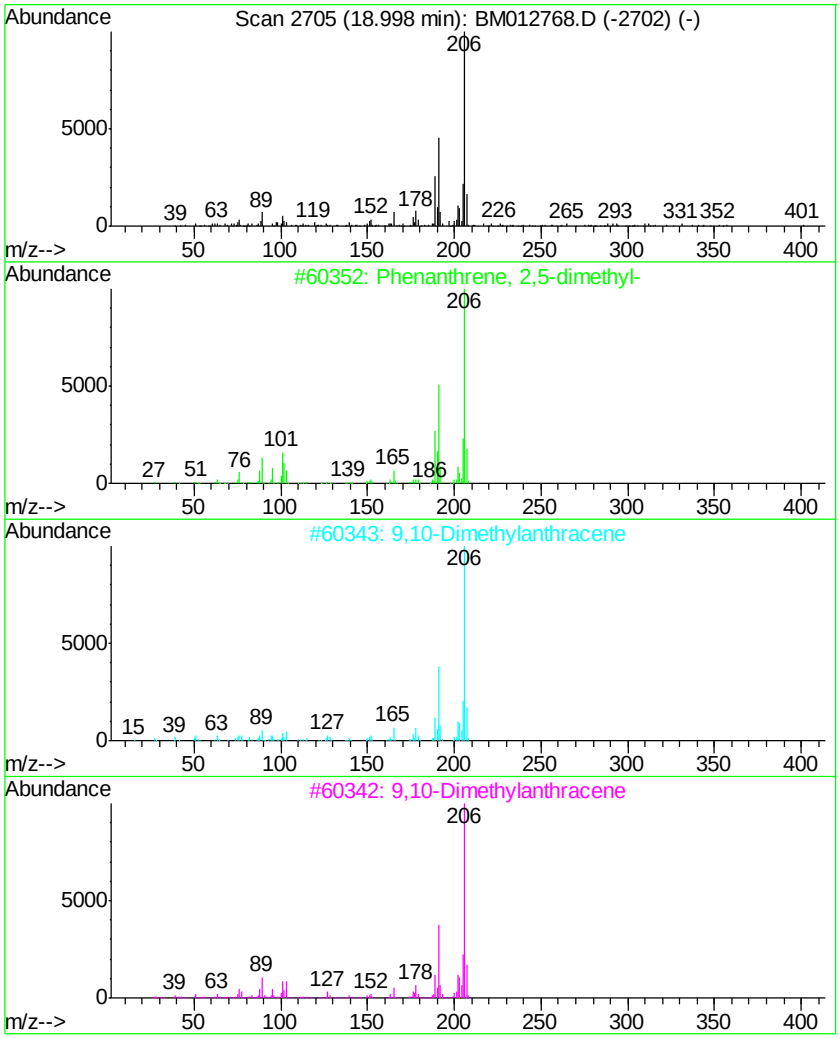
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

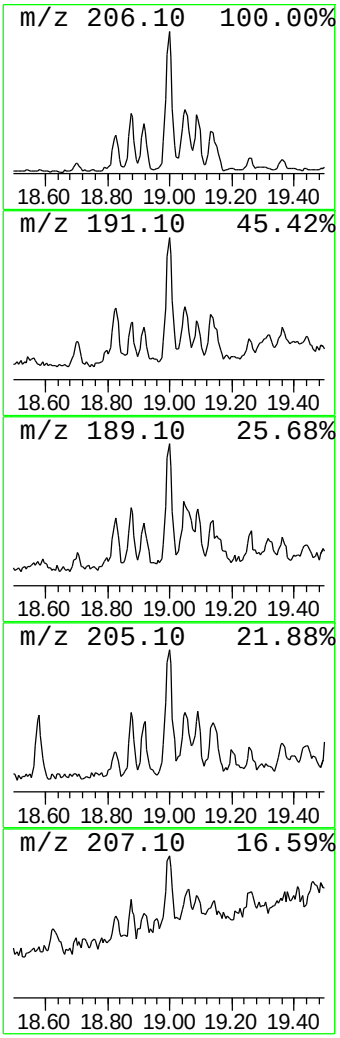
****

Peak Number 31 Phenanthrene, 2,5-dimethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.00	2.01 ng/ul	184653	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
3	9,10-Dimethylanthracene	206	C16H14	000781-43-1	90
4	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

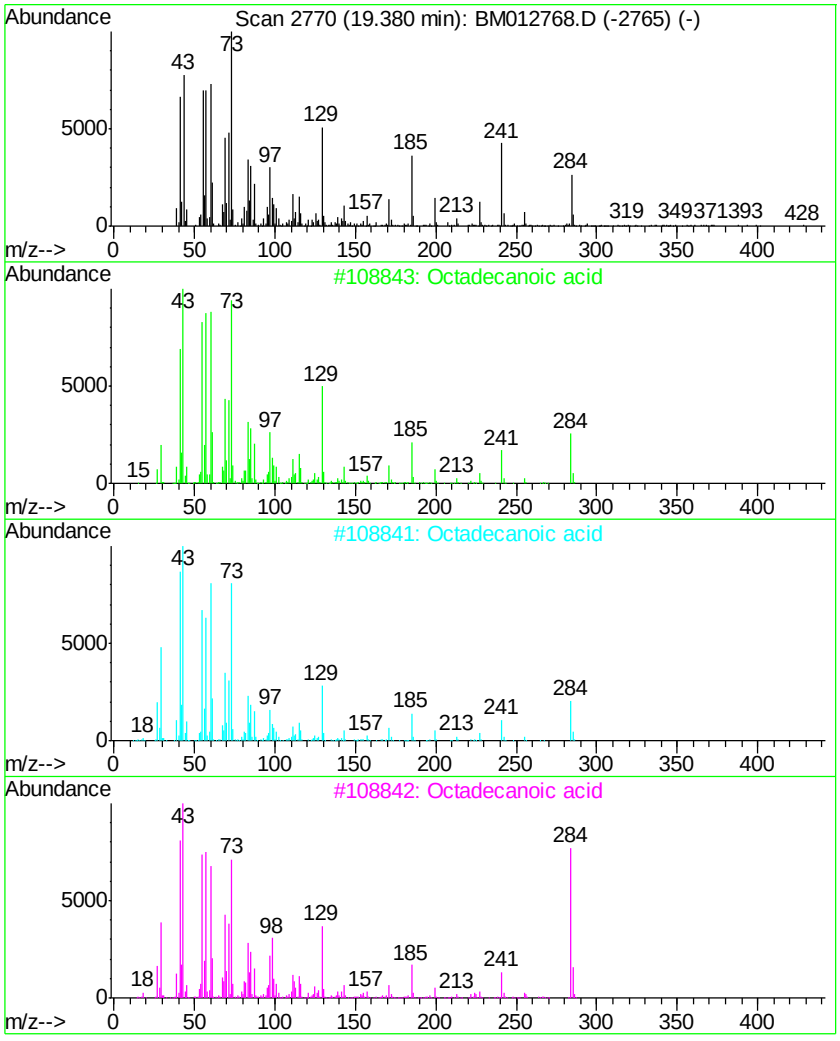
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

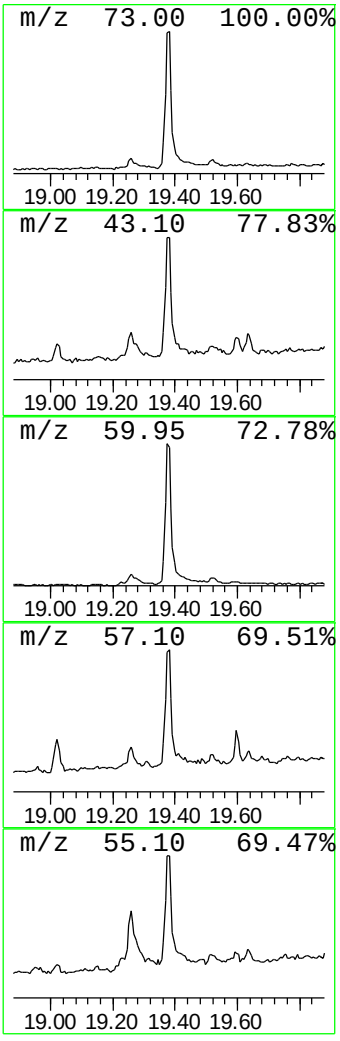
****

Peak Number 32 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.38	7.87 ng/ul	1050900	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	96
3	Octadecanoic acid	284	C18H36O2	000057-11-4	95
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
5	Octadecanoic acid	284	C18H36O2	000057-11-4	55





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

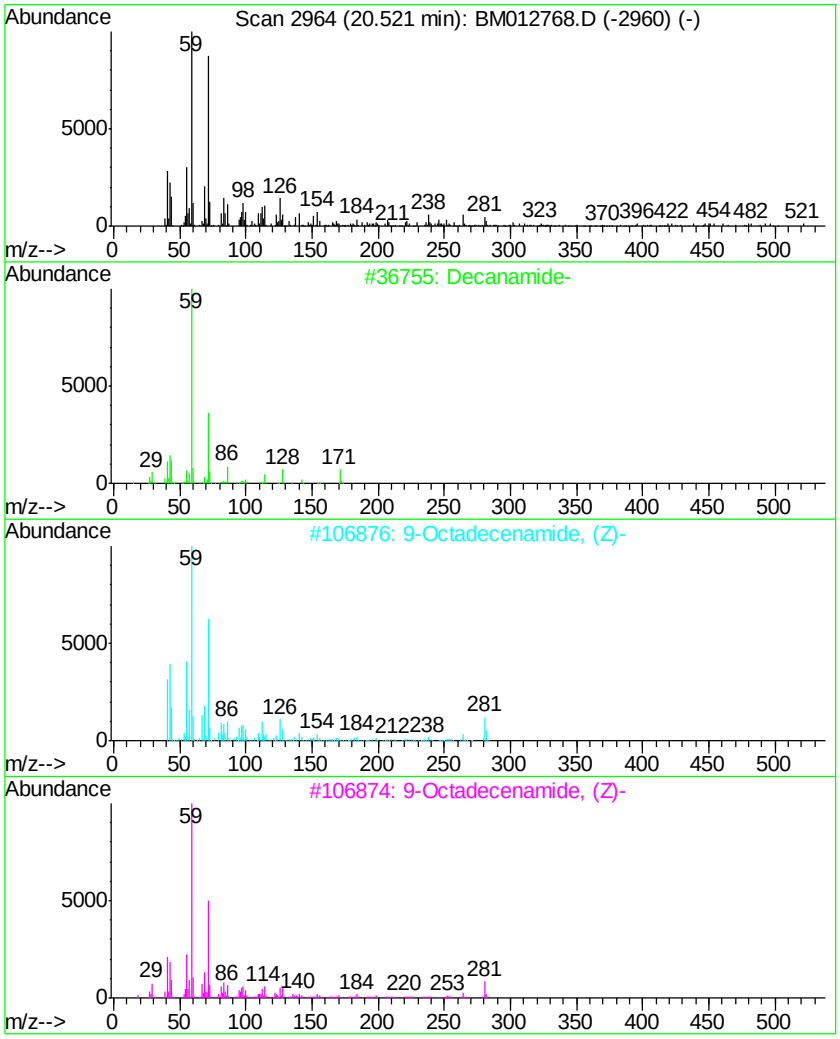
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

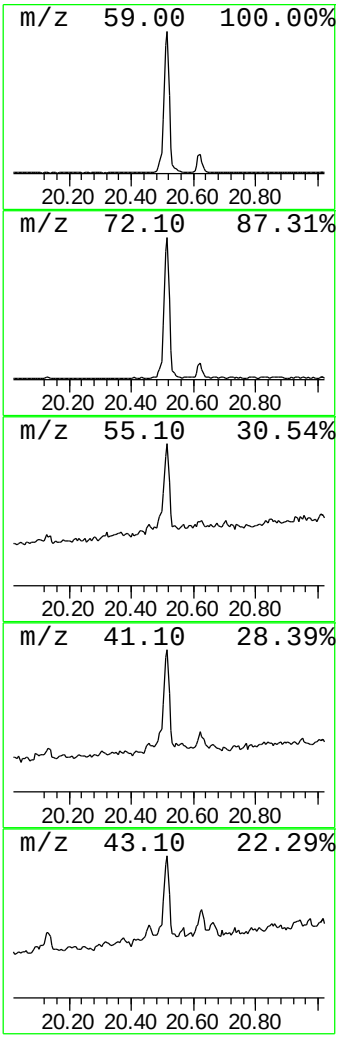
****

Peak Number 33 Decanamide- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.52	4.42 ng/ul	589501	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decanamide-	171	C10H21NO	002319-29-1	72
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	52
3	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	45
4	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	45
5	Dodecanamide	199	C12H25NO	001120-16-7	43





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110517\
Data File : BM012768.D
Acq On : 06 Nov 2017 08:45
Operator : SJ/JU
Sample : I5902-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-12(1-3)-101317

Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Cyc...	3.56	4.7	ng/ul	316582	1	7.82	1356950	20.0
Methacrylamide	3.82	3.3	ng/ul	220628	1	7.82	1356950	20.0
(DEL) Alkane: Str...	4.37	2.3	ng/ul	155937	1	7.82	1356950	20.0
2-Heptanone	5.67	2.0	ng/ul	139249	1	7.82	1356950	20.0
Benzene, (1-methy...	6.31	2.9	ng/ul	196972	1	7.82	1356950	20.0
Benzene, propyl-	6.80	3.7	ng/ul	249348	1	7.82	1356950	20.0
unknown-01	6.87	2.2	ng/ul	147330	1	7.82	1356950	20.0
Benzene, 1-ethyl-...	6.90	5.8	ng/ul	392740	1	7.82	1356950	20.0
(DEL) Alkane: Str...	7.44	9.9	ng/ul	672459	1	7.82	1356950	20.0
Benzene, 1,3,5-tr...	7.46	6.4	ng/ul	432855	1	7.82	1356950	20.0
Benzene, 1,2,3-tr...	7.93	3.5	ng/ul	235737	1	7.82	1356950	20.0
Benzene, 1-methyl...	8.36	3.7	ng/ul	249714	1	7.82	1356950	20.0
Benzene, 1-ethyl-...	8.45	6.2	ng/ul	420943	1	7.82	1356950	20.0
Benzene, 1-ethyl-...	8.92	4.2	ng/ul	281862	1	7.82	1356950	20.0
(DEL) Alkane: Str...	9.04	8.3	ng/ul	564100	1	7.82	1356950	20.0
unknown-02	9.16	2.4	ng/ul	161766	1	7.82	1356950	20.0
Octanoic Acid	9.98	4.0	ng/ul	363457	2	10.61	1826550	20.0
Nonanoic acid	11.43	2.5	ng/ul	228534	2	10.61	1826550	20.0
(DEL) Alkane: Str...	12.03	2.3	ng/ul	206156	2	10.61	1826550	20.0
Benzophenone	15.82	6.5	ng/ul	694068	3	14.46	2133770	20.0
(DEL) Alkane: Str...	16.17	3.4	ng/ul	314624	4	17.20	1841040	20.0
9-Hexadecenoic acid	17.98	3.5	ng/ul	322831	4	17.20	1841040	20.0
n-Hexadecanoic acid	18.10	18.3	ng/ul	1688180	4	17.20	1841040	20.0
Anthracene, 1-met...	18.27	2.2	ng/ul	198474	4	17.20	1841040	20.0
Phenanthrene, 2,5...	19.00	2.0	ng/ul	184653	4	17.20	1841040	20.0
Octadecanoic acid	19.38	7.9	ng/ul	1050900	5	21.39	2669610	20.0
Decanamide-	20.52	4.4	ng/ul	589501	5	21.39	2669610	20.0