LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

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-BM100317.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	4.034	156	161	169	rVB	130240	183336	1.94%	0.204%
2	4.969	314	320	329	rVB2	55340	108529	1.15%	0.121%
3	5.363	383	387	395	rVB	76194	113758	1.20%	0.127%
4	5.499	404	410	424	rVB	206715	451086	4.77%	0.502%
5	5.851	464	470	479	rVB3	182388	442246	4.68%	0.492%
6	6.846	634	639	645	rBV3	68409	110239	1.17%	0.123%
7	6.951	652	657	662	rBV2	76475	117180	1.24%	0.130%
8	7.028	662	670	677	rBV	815160	1480464	15.66%	1.648%
9	7.198	691	699	705	rBV	643538	1054055	11.15%	1.173%
10	7.398	727	733	741	rBV	929470	1498850	15.85%	1.669%
11	7.475	741	746	749	rVV	174577	263424	2.79%	0.293%
12	7.510	749	752	759	rVB	179125	270231	2.86%	0.301%
13	7.869	802	813	821	rBV	1141988	1894680	20.04%	2.109%
14	8.110	851	854	866	rVB4	51119	120983	1.28%	0.135%
15	8.410	899	905	915	rVB3	59201	142221	1.50%	0.158%
16	8.504	915	921	926	rBV2	73658	134328	1.42%	0.150%
17	8.569	926	932	946	rVV	951279	1656575	17.52%	1.844%
18	8.969	989	1000	1004	rBV3	51830	104073	1.10%	0.116%
19	9.034	1004	1011	1016	rVV	820622	1376810	14.56%	1.533%
20	9.081	1016	1019	1025	rVB	143918	211571	2.24%	0.236%
21	9.757	1120	1134	1145	rBV	702980	1294415	13.69%	1.441%
22	10.292	1218	1225	1234	rBV	1004353	1795111	18.99%	1.998%
23	10.375	1234	1239	1248	rVB3	54191	110610	1.17%	0.123%
24	10.634	1276	1283	1284	rBV3	66309	101922	1.08%	0.113%
25	10.675	1284	1290	1296	rVV	1486149	2584308	27.33%	2.877%
26	10.722	1296	1298	1306	rVB	99163	141660	1.50%	0.158%
27	10.816	1306	1314	1330	rBV	300102	673042	7.12%	0.749%
28	11.486	1414	1428	1440	rBV	1048864	1894856	20.04%	2.109%
29	12.081	1524	1529	1539	rVB	64812	104199	1.10%	0.116%
30	13.322	1734	1740	1749	rBV2	71329	140832	1.49%	0.157%
31	13.916	1832	1841	1845	rVV	1760631	2527273	26.73%	2.813%
32	13.963	1845	1849	1856	rVB	794060	1179661	12.48%	1.313%
33	14.033	1856	1861	1864	rBV3	70620	94814	1.00%	0.106%
34	14.204	1884	1890	1895	rBV	1924345	3027252	32.02%	3.370%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
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Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

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-BM100317.M
Title : SVOA CALIBRATION

35	14.392	1917	1922	1931	rVB	118505	174455	1.85%	0.194%
36	14.516	1932	1943	1950	rVV3	2088641	4001536	42.32%	4.455%
37	14.574	1950	1953	1962	rVB4	64424	122004	1.29%	0.136%
38	14.710	1970	1976	1987	rBV	289368	629809	6.66%	0.701%
39	14.910	2005	2010	2019	rVB2	98064	186184	1.97%	0.207%
40	15.345	2079	2084	2090	rVB	129967	173809	1.84%	0.193%
41	15.504	2104	2111	2117	rBV	2334738	3407565	36.04%	3.793%
42	15.627	2127	2132	2139	rBV	496419	732328	7.75%	0.815%
43	16.204	2226	2230	2234	rBV	175314	281577	2.98%	0.313%
44	16.333	2247	2252	2255	rBV	78676	125742	1.33%	0.140%
45	16.386	2258	2261	2267	rVB2	63462	113790	1.20%	0.127%
46	16.645	2301	2305	2311	rVB4	92037	172005	1.82%	0.191%
47	16.715	2313	2317	2321	rVV2	78977	102870	1.09%	0.115%
48	16.992	2359	2364	2370	rBV4	111796	184075	1.95%	0.205%
49	17.133	2385	2388	2396	rBV9	40896	98164	1.04%	0.109%
50	17.257	2403	2409	2413	rBV	2098723	2976222	31.48%	3.313%
51	17.298	2413	2416	2420	rVB	913254	1116636	11.81%	1.243%
52	17.357	2420	2426	2430	rBV	2067201	3025942	32.00%	3.369%
53	18.133	2553	2558	2562	rBV	1491706	2006933	21.23%	2.234%
54	18.174	2562	2565	2570	rVB	242790	371157	3.93%	0.413%
55	18.292	2581	2585	2594	rBV3	578830	965229	10.21%	1.075%
56	18.398	2599	2603	2605	rVV	482823	675920	7.15%	0.752%
57	18.445	2605	2611	2616	rVB	872700	1739112	18.39%	1.936%
58	18.627	2639	2642	2646	rBV	128434	147766	1.56%	0.164%
59	18.674	2648	2650	2654	rVB3	119363	141589	1.50%	0.158%
60	18.980	2698	2702	2707	rVB	379077	544888	5.76%	0.607%
61	19.045	2709	2713	2719	rVV2	171759	325778	3.45%	0.363%
62	19.115	2721	2725	2732	rVB2	537650	869602	9.20%	0.968%
63	19.309	2752	2758	2763	rVB	1874403	2577219	27.26%	2.869%
64	19.409	2771	2775	2787	rBV	1618190	2305980	24.39%	2.567%
65	19.645	2810	2815	2818	rBV	2760140	4171400	44.12%	4.644%
66	19.674	2818	2820	2827	rVB	1660168	1911751	20.22%	2.128%
67	20.556	2966	2970	2976	rVB	7879280	9454710	100.00%	10.525%
68	21.121	3063	3066	3069	rVB2	613885	679253	7.18%	0.756%
69	21.427	3112	3118	3122	rBV2	2819102	4774091	50.49%	5.315%
70	21.462	3122	3124	3130	rVB	939564	1129162	11.94%	1.257%
71	23.062	3391	3396	3414	rVB2	1510218	4095554	43.32%	4.559%

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Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

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 >
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Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Title : SVOA CALIBRATION

72	23.609	3484	3489	3494	rBV2	1746475	3171517	33.54%	3.531%
73	23.756	3509	3514	3520	rBV	1569705	2741819	29.00%	3.052%

Sum of corrected areas: 89829737

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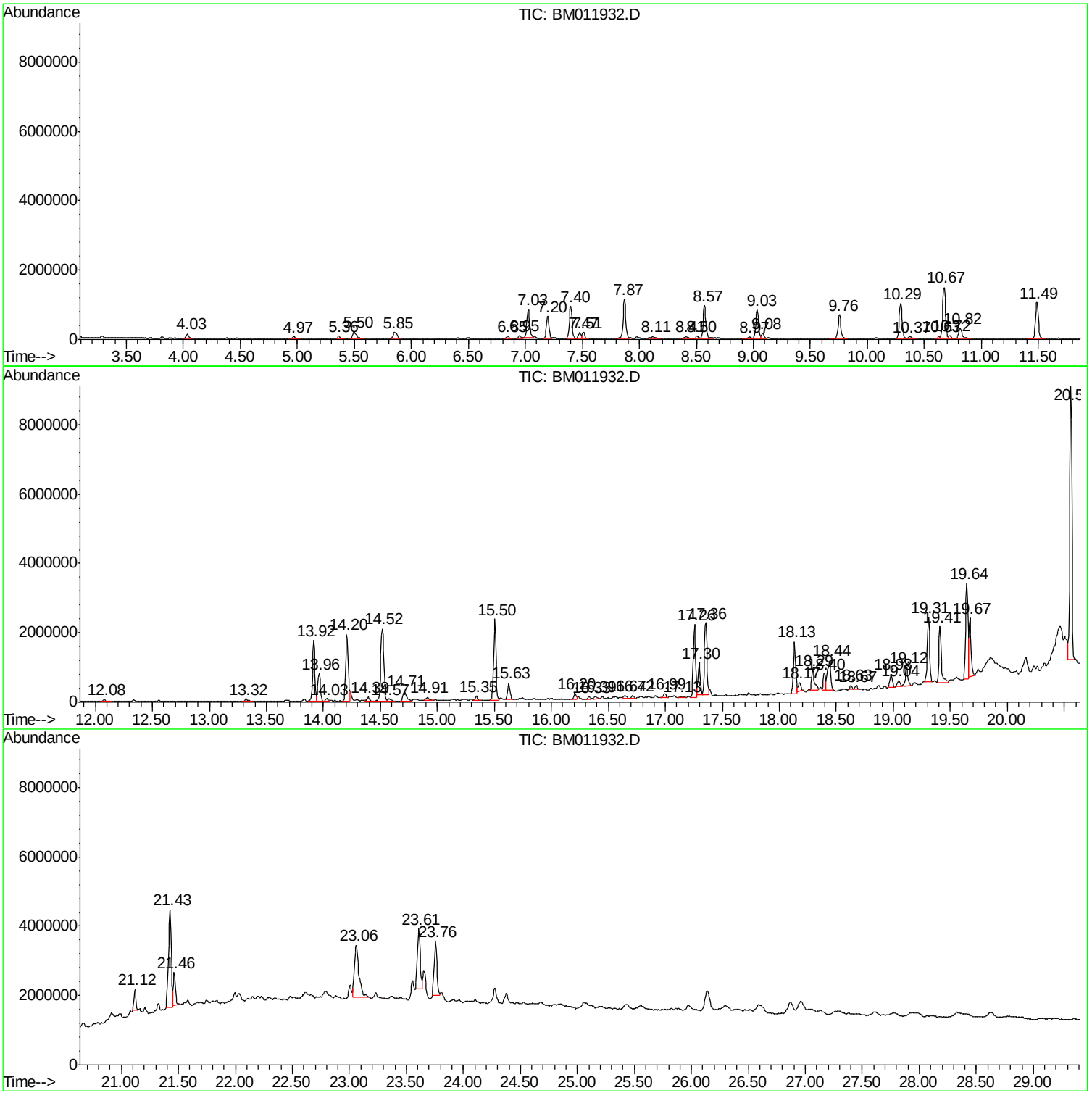
Instrument :
BNA_M
ClientSampleId :
B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
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Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

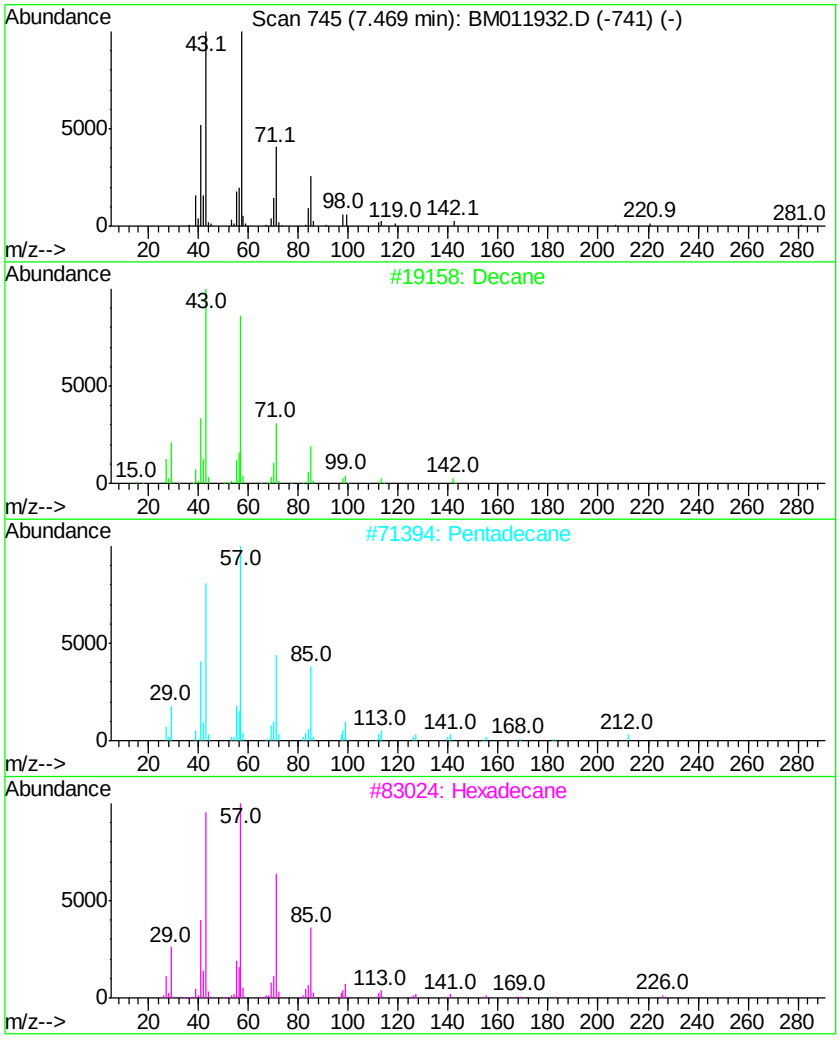
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

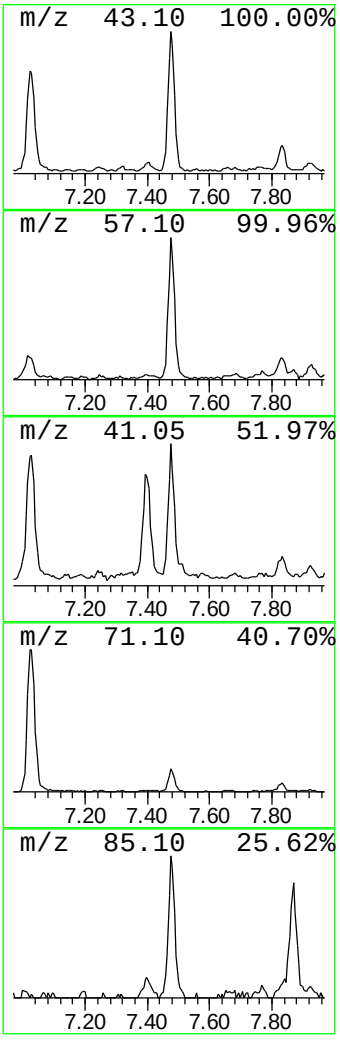
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	2.78 ng/ul	263424	1,4-Dichlorobenzene-d4	7.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	93
2	Pentadecane	212	C15H32	000629-62-9	83
3	Hexadecane	226	C16H34	000544-76-3	83
4	Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	80
5	Tridecane	184	C13H28	000629-50-5	78





Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

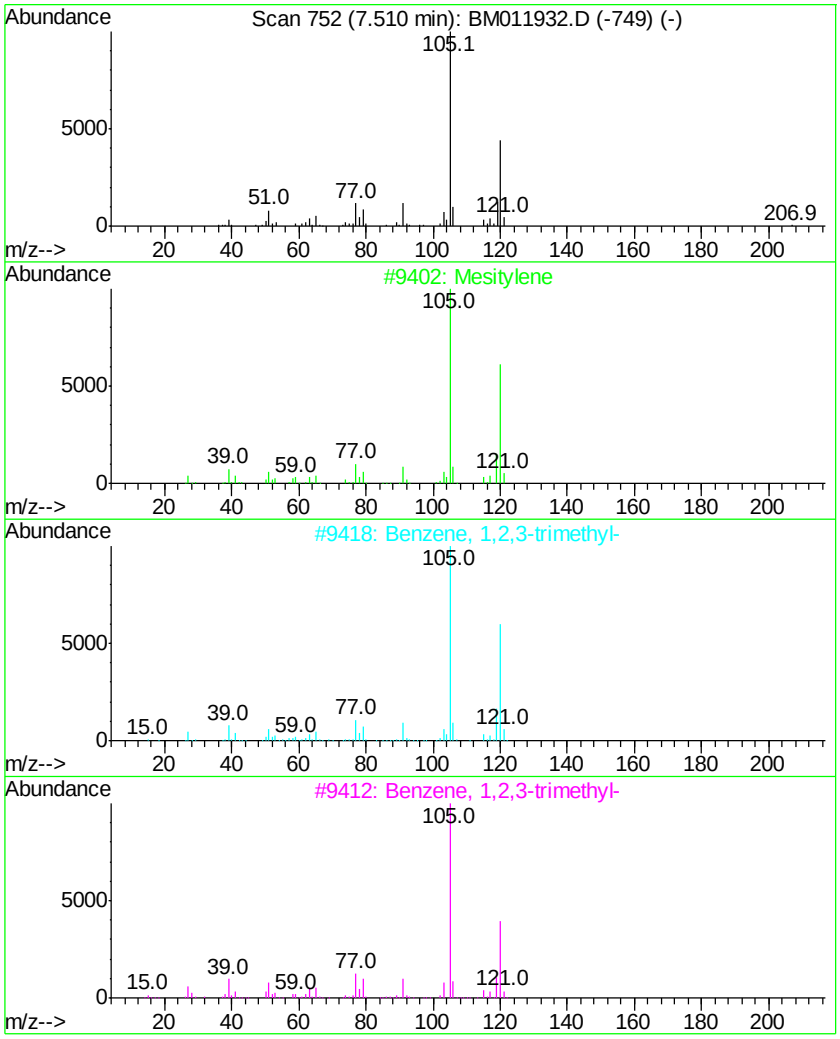
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

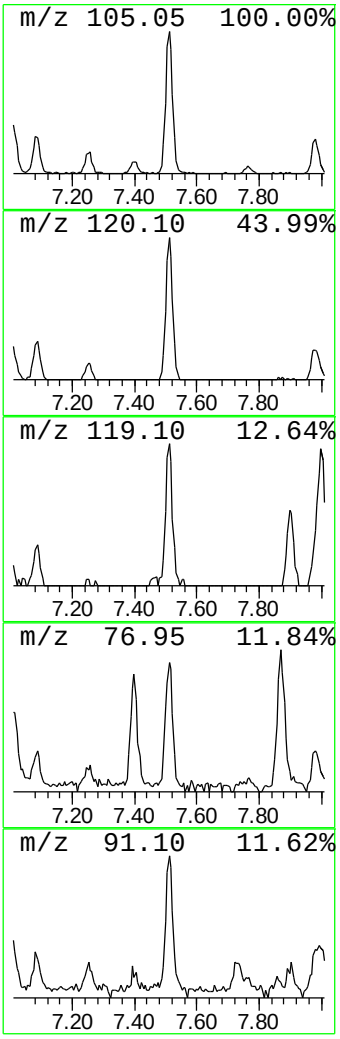
****

Peak Number 4 Mesitylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.51	2.85 ng/ul	270231	1,4-Dichlorobenzene-d4	7.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

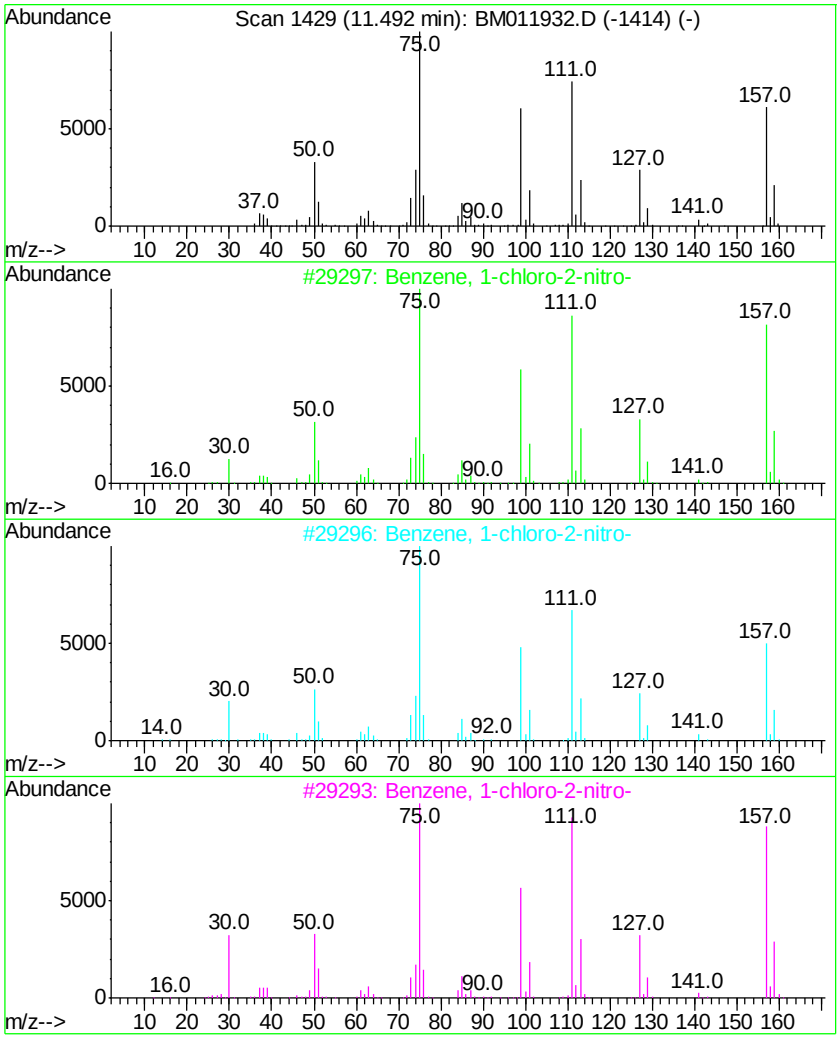
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

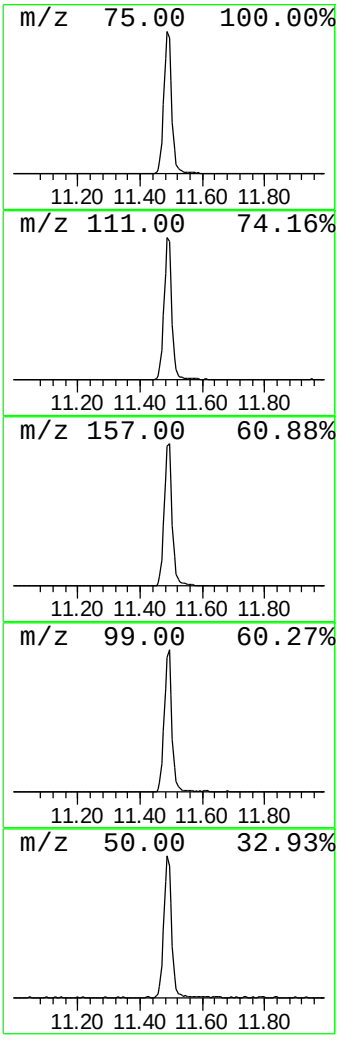
****

Peak Number 5 Benzene, 1-chloro-2-nitro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	14.66 ng/ul	1894860	Naphthalene-d8	10.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96
5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95





Library Search Compound Report

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Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

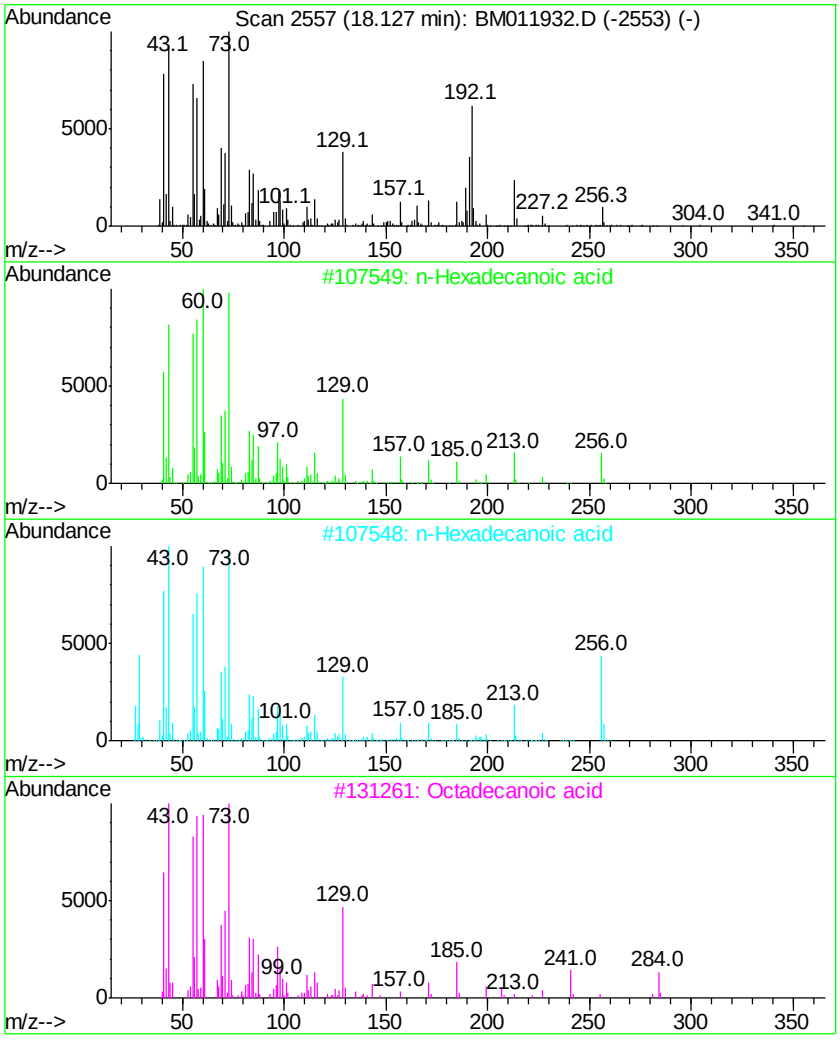
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

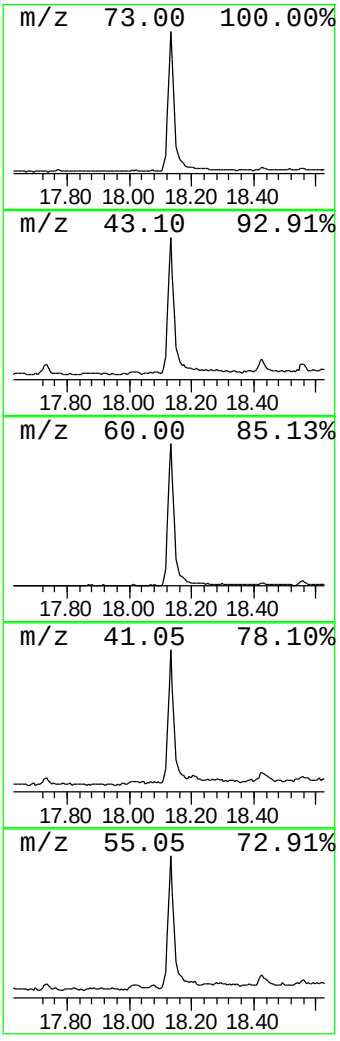
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	13.49 ng/ul	2006930	Phenanthrene-d10	17.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	Octadecanoic acid	284	C18H36O2	000057-11-4	92
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5	Tridecanoic acid	214	C13H26O2	000638-53-9	86





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
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Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

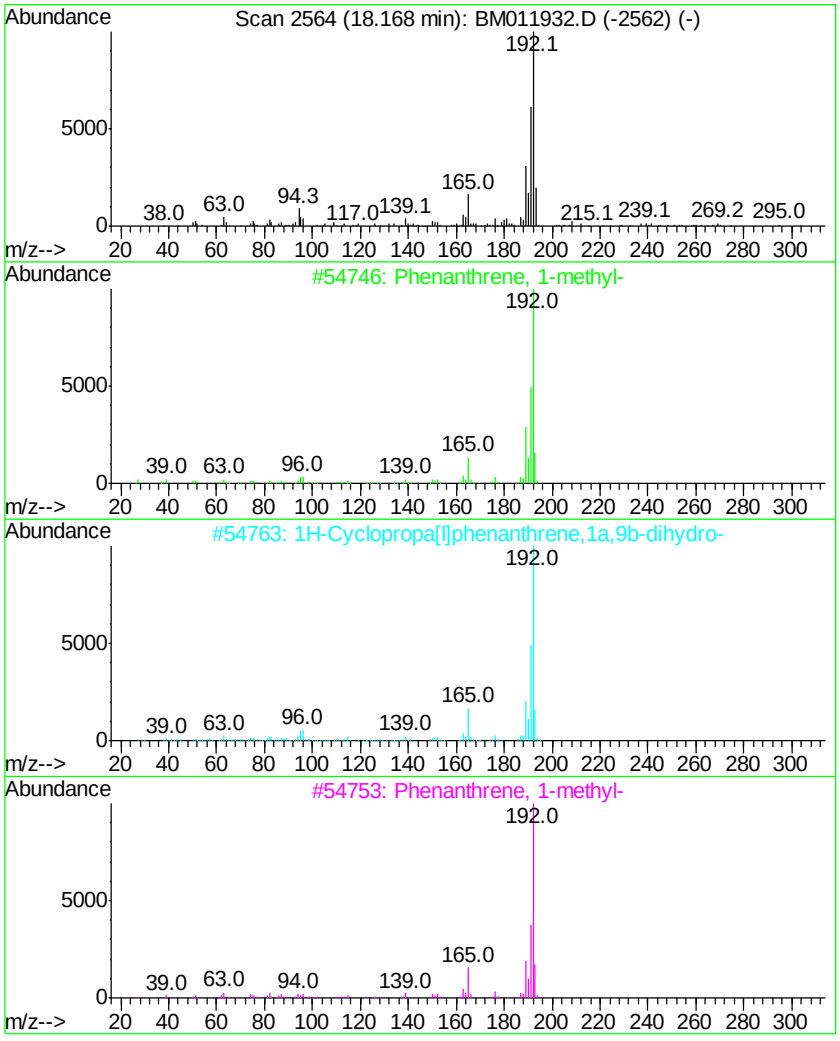
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

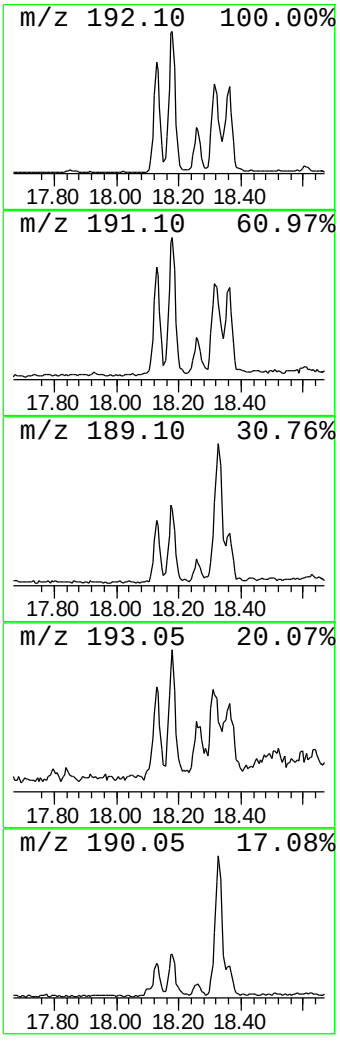
****

Peak Number 7 Phenanthrene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	2.49 ng/ul	371157	Phenanthrene-d10	17.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94





Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

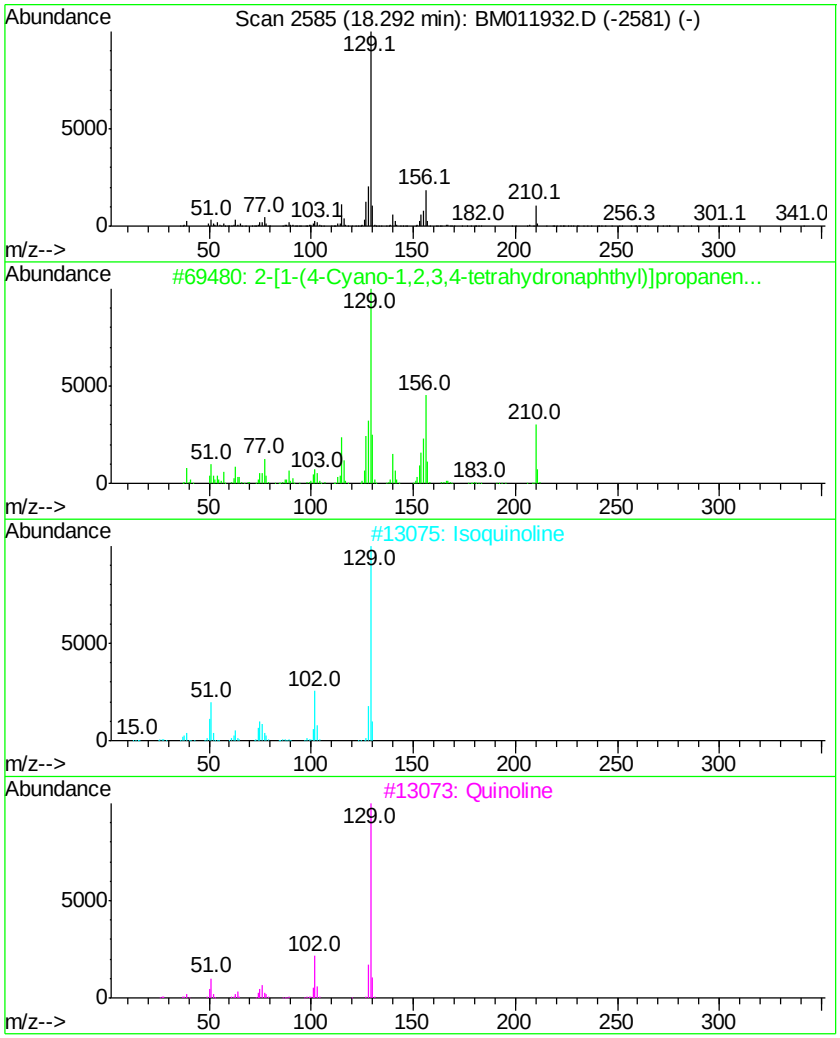
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

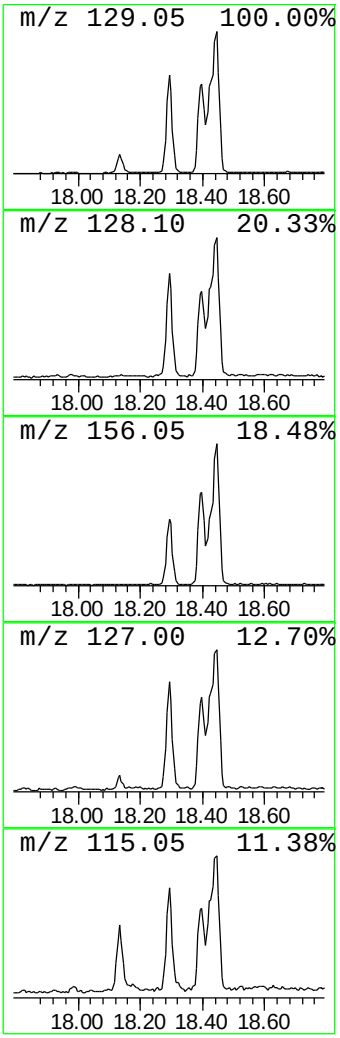
****

Peak Number 8 2-[1-(4-Cyano-1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.29	6.49 ng/ul	965229	Phenanthrene-d10	17.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-39-3	70
2	Isoquinoline	129	C9H7N	000119-65-3	42
3	Quinoline	129	C9H7N	000091-22-5	40
4	Isoquinoline	129	C9H7N	000119-65-3	40
5	3-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-40-6	38





Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

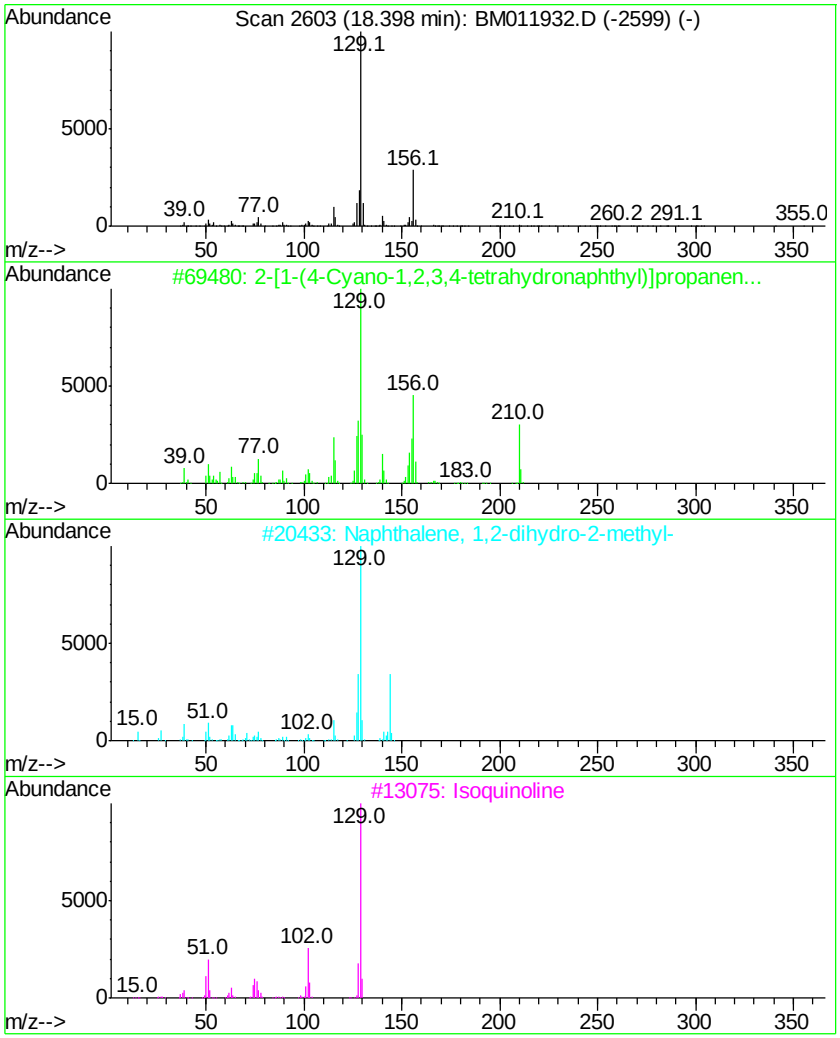
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

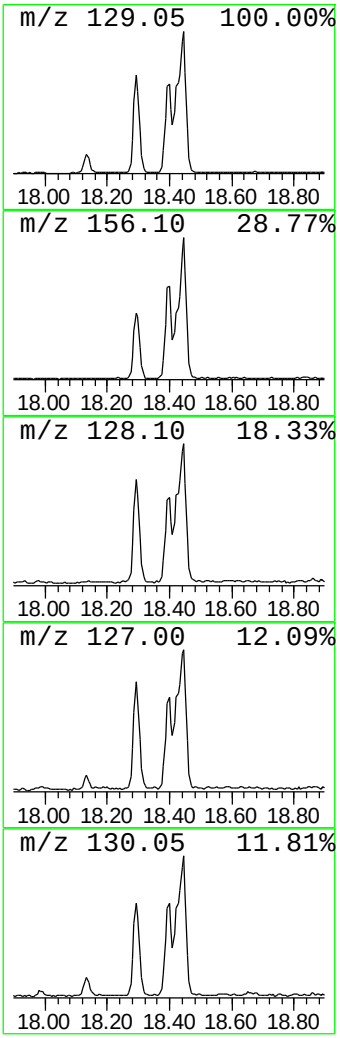
****

Peak Number 9 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.40	4.54 ng/ul	675920	Phenanthrene-d10	17.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-39-3	43
2	Naphthalene, 1,2-dihydro-2-methyl-	144	C11H12	021564-79-4	40
3	Isoquinoline	129	C9H7N	000119-65-3	40
4	Isoquinoline	129	C9H7N	000119-65-3	38
5	3-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-40-6	38





Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

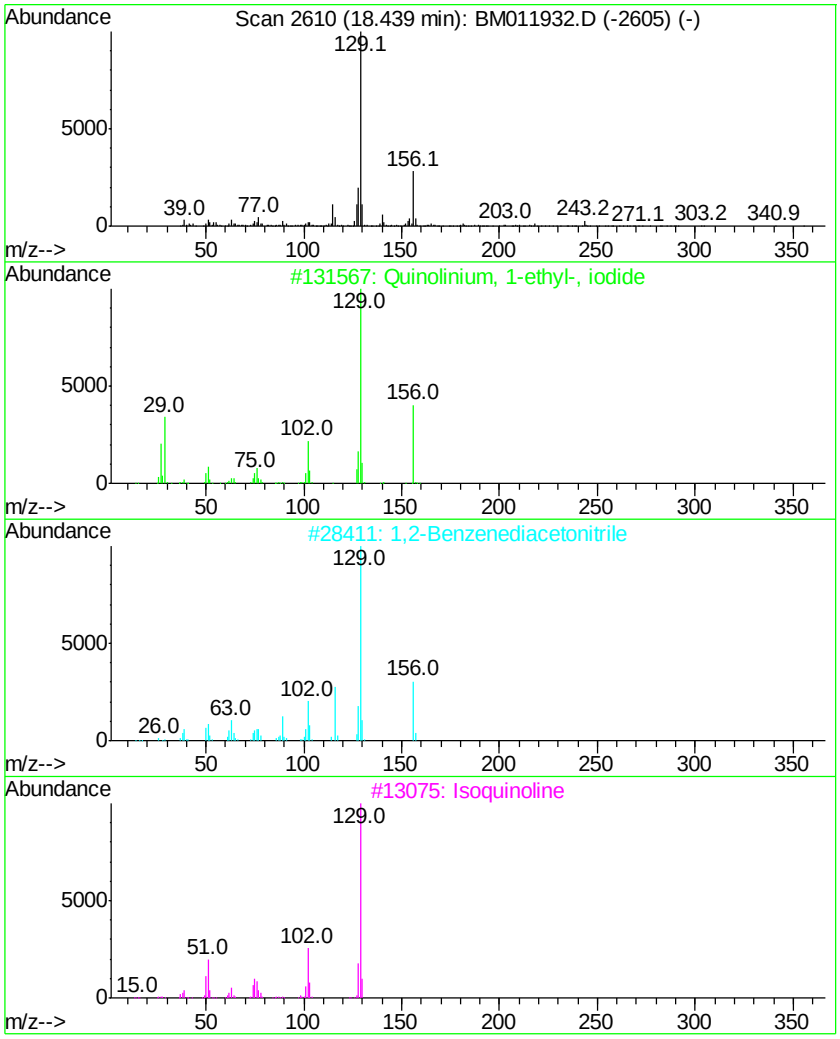
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

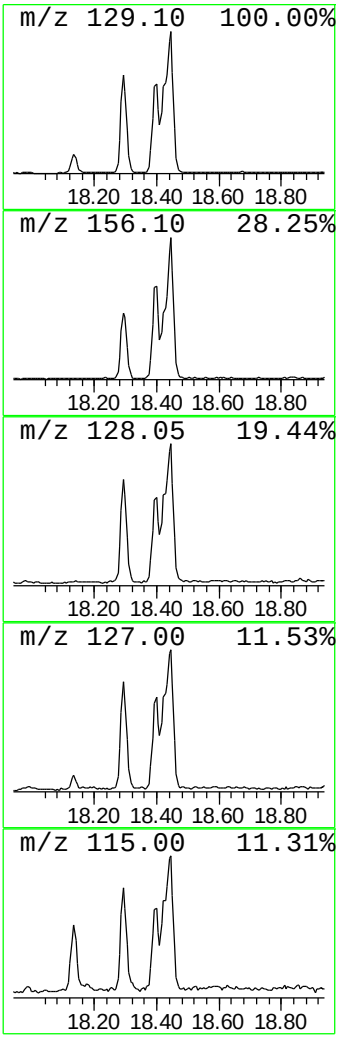
****

Peak Number 10 Quinolinium, 1-ethyl-, iodide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.44	11.69 ng/ul	1739110	Phenanthrene-d10	17.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinolinium, 1-ethyl-, iodide	285	C11H12IN	000634-35-5	64
2	1,2-Benzenediacetonitrile	156	C10H8N2	000613-73-0	59
3	Isoquinoline	129	C9H7N	000119-65-3	52
4	2-Propenenitrile, 3-phenyl-, (E)-	129	C9H7N	001885-38-7	50
5	Quinolinium, 1-ethyl-, iodide	285	C11H12IN	000634-35-5	50





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Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

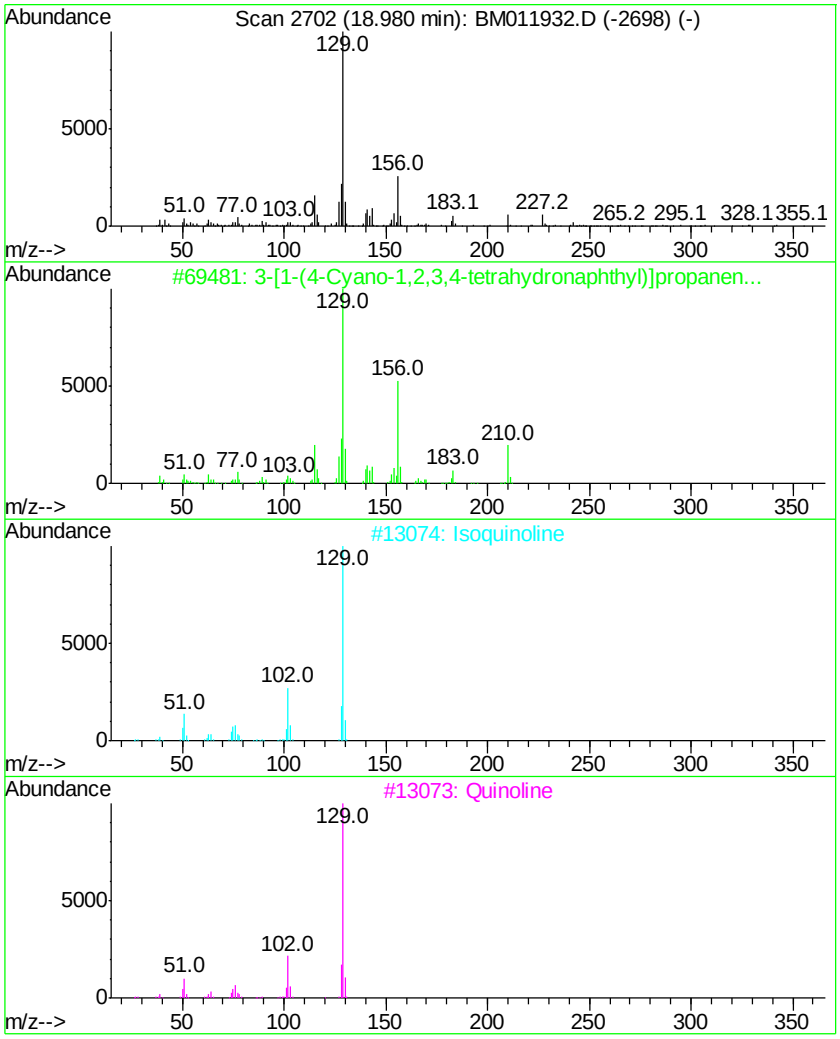
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

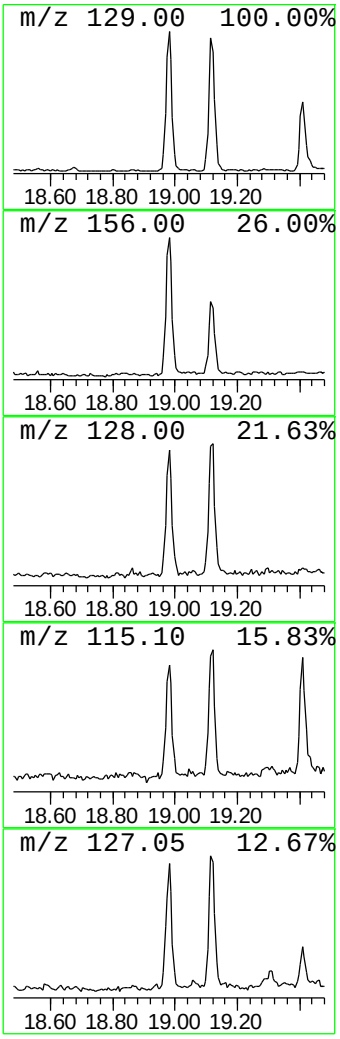
****

Peak Number 11 3-[1-(4-Cyano-1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.98	3.66 ng/ul	544888	Phenanthrene-d10	17.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-40-6	97
2	Isoquinoline	129	C9H7N	000119-65-3	52
3	Quinoline	129	C9H7N	000091-22-5	52
4	Quinoline	129	C9H7N	000091-22-5	50
5	2-Propenenitrile, 3-phenyl-, (E)-	129	C9H7N	001885-38-7	50





Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
Operator : SJ/JU
Sample : I5449-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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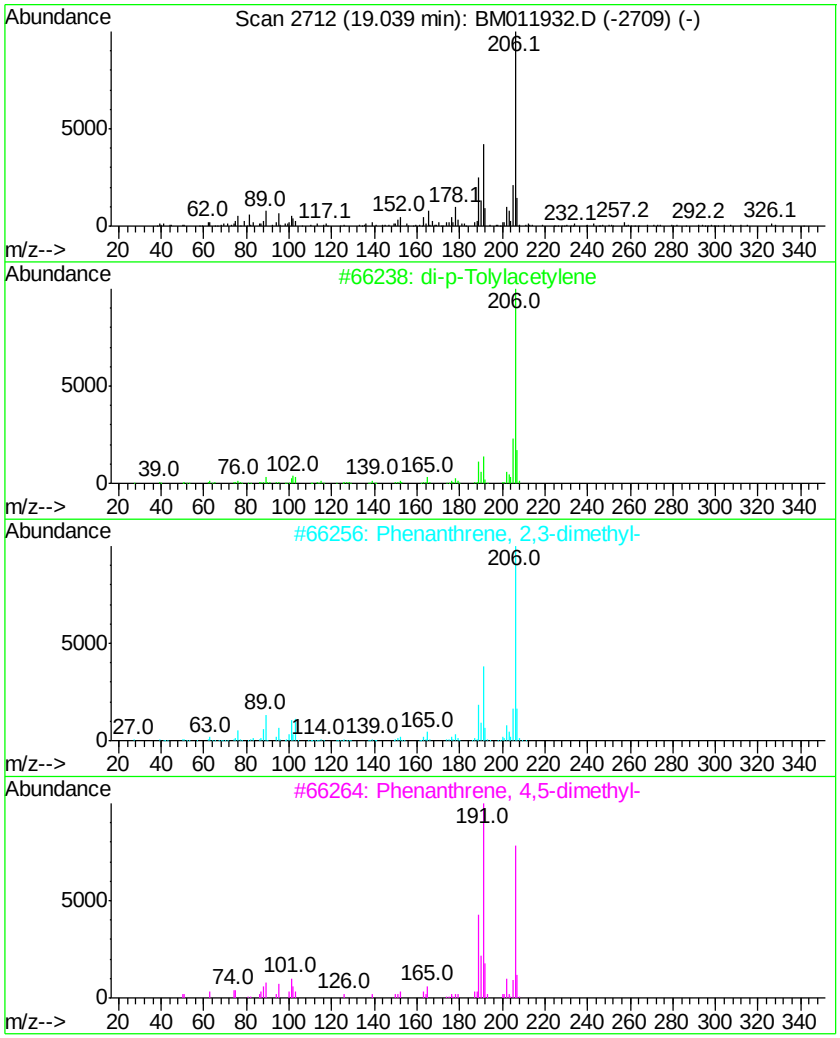
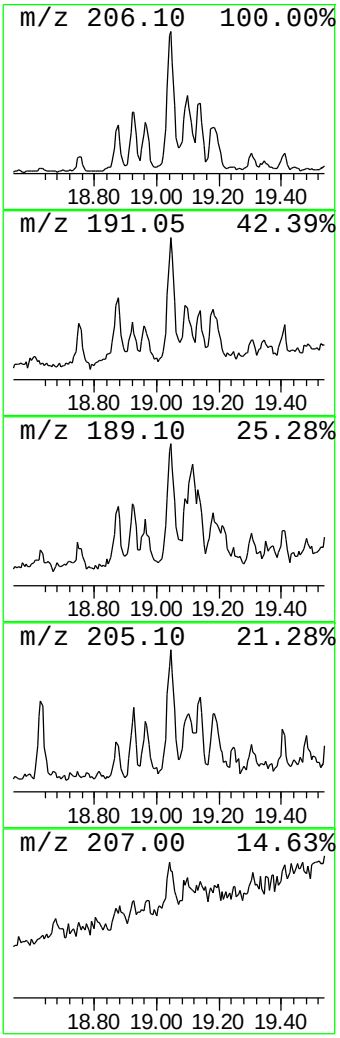
Peak Number 12 di-p-Tolylacetylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.04	2.19 ng/ul	325778	Phenanthrene-d10	17.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
2	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	81
3	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	74
4	1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	74
5	9,10-Dimethylanthracene	206	C16H14	000781-43-1	72





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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
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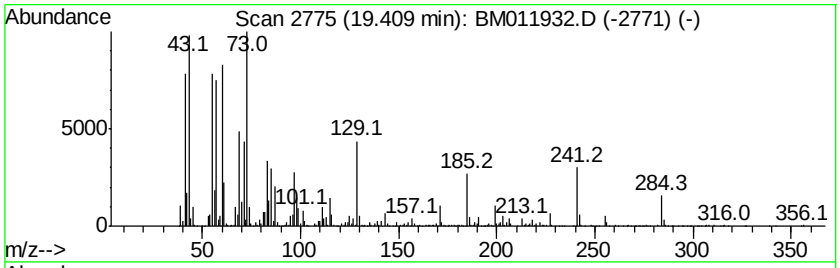
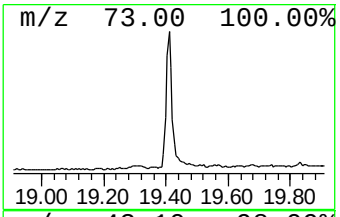
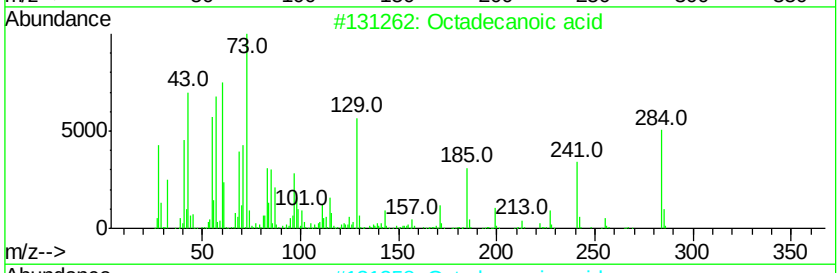
Peak Number 14 Octadecanoic acid Concentration Rank 4

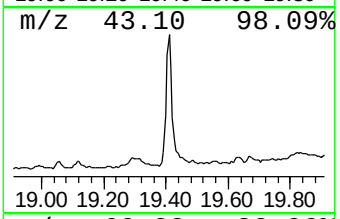
R.T.	EstConc	Area	Relative to ISTD	R.T.
19.41	9.66 ng/ul	2305980	Chrysene-d12	21.43

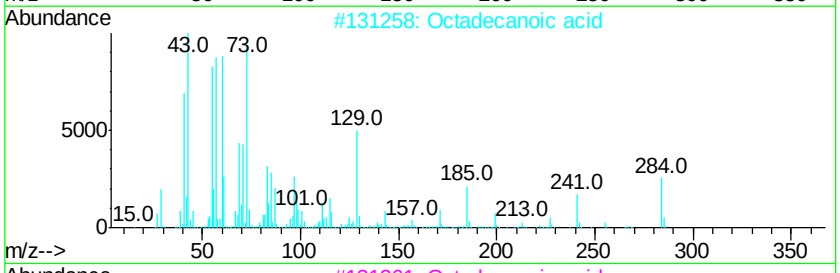
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	98
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	93

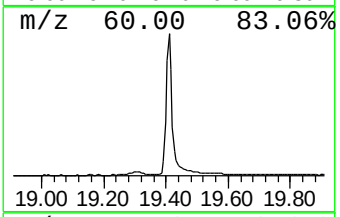


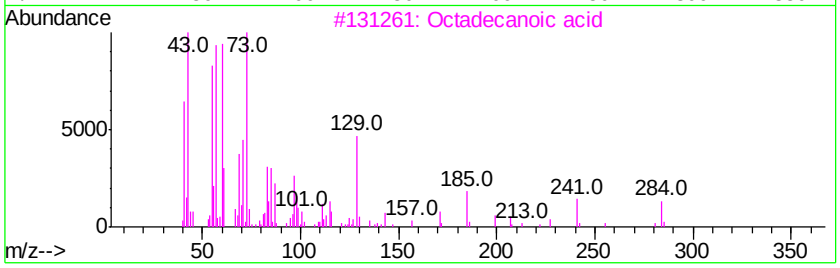


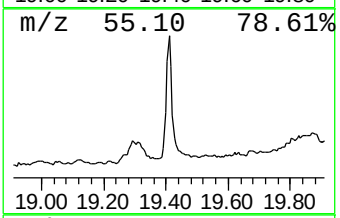


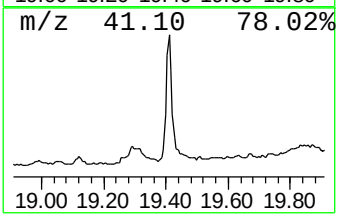












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Quant Title : SVOA CALIBRATION

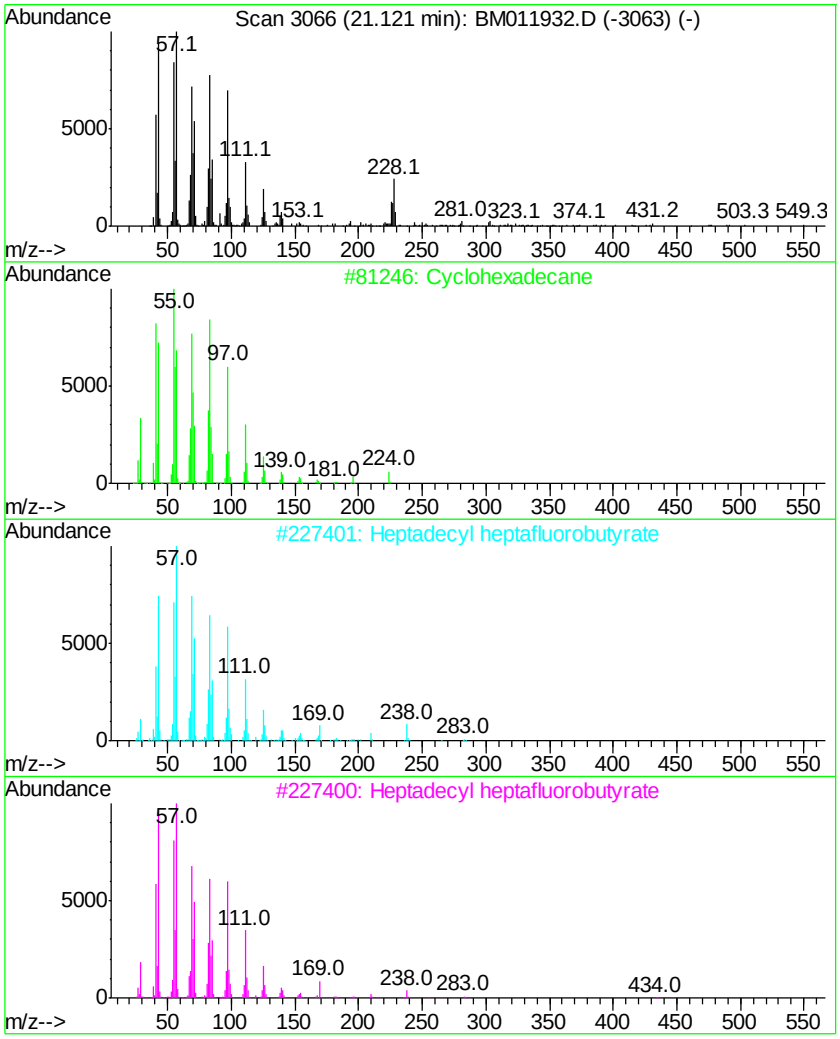
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

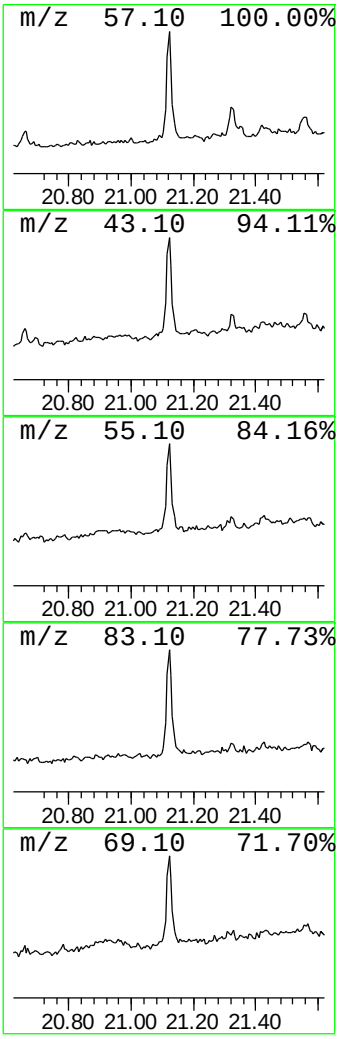
****

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.12	2.85 ng/ul	679253	Chrysene-d12	21.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	94
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
4	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM100517\
Data File : BM011932.D
Acq On : 05 Oct 2017 19:32
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Misc :
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ClientSampleId : B-60(5-7)-092617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Str...	7.47	2.8	ng/ul	263424	1	7.87	1894680	20.0
Mesitylene	7.51	2.9	ng/ul	270231	1	7.87	1894680	20.0
Benzene, 1-chloro...	11.49	14.7	ng/ul	1894860	2	10.67	2584310	20.0
n-Hexadecanoic acid	18.13	13.5	ng/ul	2006930	4	17.26	2976220	20.0
Phenanthrene, 1-m...	18.17	2.5	ng/ul	371157	4	17.26	2976220	20.0
2-[1-(4-Cyano-1,2...	18.29	6.5	ng/ul	965229	4	17.26	2976220	20.0
unknown-01	18.40	4.5	ng/ul	675920	4	17.26	2976220	20.0
Quinolinium, 1-et...	18.44	11.7	ng/ul	1739110	4	17.26	2976220	20.0
3-[1-(4-Cyano-1,2...	18.98	3.7	ng/ul	544888	4	17.26	2976220	20.0
di-p-Tolylacetylene	19.04	2.2	ng/ul	325778	4	17.26	2976220	20.0
Octadecanoic acid	19.41	9.7	ng/ul	2305980	5	21.43	4774090	20.0
(DEL) Alkane: Str...	21.12	2.9	ng/ul	679253	5	21.43	4774090	20.0