LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
Data File : BM016925.D
Acq On : 01 Oct 2018 17:26
Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.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	3.193	31	35	39	rVB	1308424	1338751	21.36%	2.221%
2	7.710	795	803	811	rBV	1236132	1923269	30.68%	3.190%
3	10.492	1268	1276	1281	rBV	1591667	2698045	43.04%	4.475%
4	10.539	1281	1284	1292	rVB	256416	385734	6.15%	0.640%
5	12.157	1549	1559	1566	rBV	221529	344071	5.49%	0.571%
6	12.380	1587	1597	1603	rBV	168096	275316	4.39%	0.457%
7	13.180	1728	1733	1740	rVB	62060	88776	1.42%	0.147%
8	13.374	1760	1766	1776	rVB	48744	82895	1.32%	0.137%
9	13.510	1784	1789	1791	rBV2	72522	119702	1.91%	0.199%
10	13.668	1810	1816	1821	rBV	152359	265122	4.23%	0.440%
11	13.721	1821	1825	1833	rVB	110094	167549	2.67%	0.278%
12	13.910	1850	1857	1860	rBV	81421	124548	1.99%	0.207%
13	13.939	1860	1862	1867	rVB2	35264	42802	0.68%	0.071%
14	14.074	1877	1885	1890	rVB2	71078	116733	1.86%	0.194%
15	14.351	1923	1932	1938	rVV2	2119271	3319361	52.96%	5.506%
16	14.415	1938	1943	1951	rVV	452375	693700	11.07%	1.151%
17	14.563	1960	1968	1978	rBV	35570	81348	1.30%	0.135%
18	14.663	1978	1985	1990	rBV2	49311	85124	1.36%	0.141%
19	14.751	1994	2000	2007	rVV	460534	668875	10.67%	1.109%
20	14.827	2007	2013	2018	rVV	56916	81283	1.30%	0.135%
21	14.974	2032	2038	2042	rBV	57533	95596	1.53%	0.159%
22	15.027	2042	2047	2052	rVB	62009	78730	1.26%	0.131%
23	15.145	2059	2067	2070	rBV2	55453	107552	1.72%	0.178%
24	15.174	2070	2072	2077	rVV2	37962	54994	0.88%	0.091%
25	15.327	2096	2098	2104	rVV4	25759	44941	0.72%	0.075%
26	15.404	2104	2111	2116	rVV	490604	735837	11.74%	1.221%
27	15.527	2128	2132	2135	rVV	69390	105238	1.68%	0.175%
28	15.562	2135	2138	2143	rVV2	136704	197037	3.14%	0.327%
29	15.615	2143	2147	2150	rVV2	38711	58600	0.93%	0.097%
30	15.651	2150	2153	2157	rVV	68099	100465	1.60%	0.167%
31	15.698	2157	2161	2171	rVB	249347	360902	5.76%	0.599%
32	15.845	2179	2186	2195	rVV	245960	489920	7.82%	0.813%
33	15.933	2195	2201	2211	rVB3	58480	109077	1.74%	0.181%
34	16.080	2223	2226	2232	rVB3	28203	41872	0.67%	0.069%

LSC Area Percent Report

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Acq On : 01 Oct 2018 17:26
Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Title : SVOA CALIBRATION

35	16.215	2245	2249	2254	rVB3	22097	28580	0.46%	0.047%
36	16.409	2277	2282	2287	rVB3	107088	180359	2.88%	0.299%
37	16.456	2287	2290	2295	rVB	51470	64504	1.03%	0.107%
38	16.509	2295	2299	2301	rBV3	21238	26315	0.42%	0.044%
39	16.557	2301	2307	2312	rVB5	42198	86964	1.39%	0.144%
40	16.639	2312	2321	2324	rBV3	130033	254980	4.07%	0.423%
41	16.680	2324	2328	2331	rVB2	98261	122405	1.95%	0.203%
42	16.709	2331	2333	2338	rVB3	37137	36904	0.59%	0.061%
43	16.762	2338	2342	2348	rVB6	38632	76180	1.22%	0.126%
44	16.833	2348	2354	2356	rBV	141780	217809	3.47%	0.361%
45	16.921	2364	2369	2377	rVB	198869	321442	5.13%	0.533%
46	17.104	2393	2400	2403	rBV	2253029	3334249	53.19%	5.530%
47	17.145	2403	2407	2413	rVV	4558446	6268184	100.00%	10.397%
48	17.233	2417	2422	2427	rVV	417528	566858	9.04%	0.940%
49	17.286	2427	2431	2438	rVV2	53700	99782	1.59%	0.166%
50	17.503	2464	2468	2471	rBV5	23391	42046	0.67%	0.070%
51	17.539	2471	2474	2479	rVV	48128	84113	1.34%	0.140%
52	17.621	2484	2488	2492	rVV	88245	113752	1.81%	0.189%
53	17.680	2494	2498	2501	rVV	37797	42665	0.68%	0.071%
54	17.821	2518	2522	2528	rBV	47250	78125	1.25%	0.130%
55	17.974	2539	2548	2552	rBV	591312	867996	13.85%	1.440%
56	18.021	2552	2556	2564	rVB	829798	1119307	17.86%	1.857%
57	18.103	2564	2570	2575	rBV	83393	135880	2.17%	0.225%
58	18.168	2575	2581	2585	rVV2	566344	1046325	16.69%	1.736%
59	18.209	2585	2588	2595	rVV	369786	504827	8.05%	0.837%
60	18.298	2599	2603	2608	rBV5	24239	32387	0.52%	0.054%
61	18.480	2629	2634	2640	rBV	382030	512158	8.17%	0.850%
62	18.603	2651	2655	2658	rBV	57534	75565	1.21%	0.125%
63	18.727	2672	2676	2680	rVV	124824	177002	2.82%	0.294%
64	18.780	2680	2685	2688	rVV	123812	176086	2.81%	0.292%
65	18.815	2688	2691	2695	rVB	94812	115735	1.85%	0.192%
66	18.898	2700	2705	2710	rVV	226162	359909	5.74%	0.597%
67	18.956	2710	2715	2718	rVV3	135980	277066	4.42%	0.460%
68	18.992	2718	2721	2725	rVV	124442	156932	2.50%	0.260%
69	19.039	2725	2729	2738	rVV3	104050	225417	3.60%	0.374%
70	19.162	2744	2750	2757	rBV	4289099	5564257	88.77%	9.229%
71	19.227	2757	2761	2764	rVV	82319	114814	1.83%	0.190%

LSC Area Percent Report

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Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Title : SVOA CALIBRATION

72	19.262	2764	2767	2773	rVB5	89502	132208	2.11%	0.219%
73	19.356	2778	2783	2787	rBV3	26270	40713	0.65%	0.068%
74	19.403	2787	2791	2795	rBV	88061	129624	2.07%	0.215%
75	19.445	2795	2798	2801	rVB	26628	29312	0.47%	0.049%
76	19.497	2801	2807	2808	rBV	728186	992271	15.83%	1.646%
77	19.527	2808	2812	2820	rVB	2090047	2912324	46.46%	4.831%
78	19.597	2820	2824	2831	rVB2	259230	396769	6.33%	0.658%
79	19.703	2838	2842	2850	rVB	261424	365391	5.83%	0.606%
80	19.762	2850	2852	2858	rVB6	36626	46132	0.74%	0.077%
81	19.856	2862	2868	2877	rBV2	250184	561841	8.96%	0.932%
82	19.939	2878	2882	2885	rBV3	39166	64362	1.03%	0.107%
83	19.986	2885	2890	2892	rBV2	197643	319739	5.10%	0.530%
84	20.021	2892	2896	2901	rVB	404818	607765	9.70%	1.008%
85	20.121	2909	2913	2917	rBV	126012	167268	2.67%	0.277%
86	20.174	2917	2922	2927	rBV2	245057	490993	7.83%	0.814%
87	20.221	2927	2930	2933	rVB2	105634	122628	1.96%	0.203%
88	20.262	2933	2937	2942	rVB2	105953	158198	2.52%	0.262%
89	20.321	2943	2947	2950	rBV	88197	110387	1.76%	0.183%
90	20.368	2950	2955	2959	rBV3	135393	228505	3.65%	0.379%
91	20.586	2987	2992	2994	rBV4	51967	104986	1.67%	0.174%
92	20.939	3048	3052	3055	rBV	192649	242714	3.87%	0.403%
93	20.974	3055	3058	3062	rVV	332527	435284	6.94%	0.722%
94	21.009	3062	3064	3069	rVB	142895	170219	2.72%	0.282%
95	21.286	3104	3111	3115	rBV2	2769846	5324976	84.95%	8.832%
96	21.327	3115	3118	3126	rVB	1582428	2031574	32.41%	3.370%
97	21.839	3202	3205	3209	rVB	199913	257844	4.11%	0.428%
98	21.886	3211	3213	3218	rVB	195739	259259	4.14%	0.430%
99	22.856	3373	3378	3383	rBV	652600	1373715	21.92%	2.279%
100	23.521	3485	3491	3499	rBV	1718250	3219967	51.37%	5.341%

Sum of corrected areas: 60288612

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Instrument :
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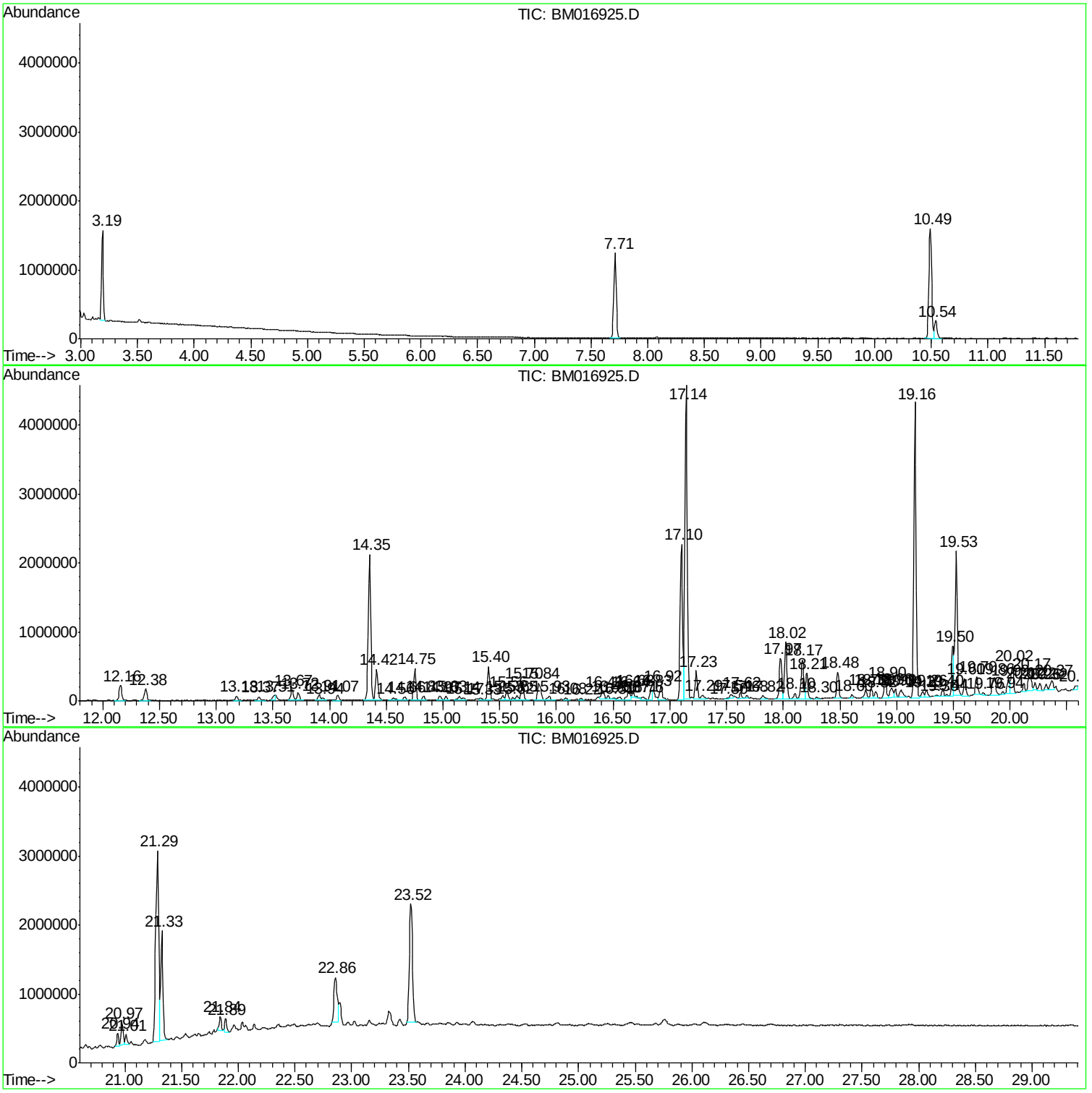
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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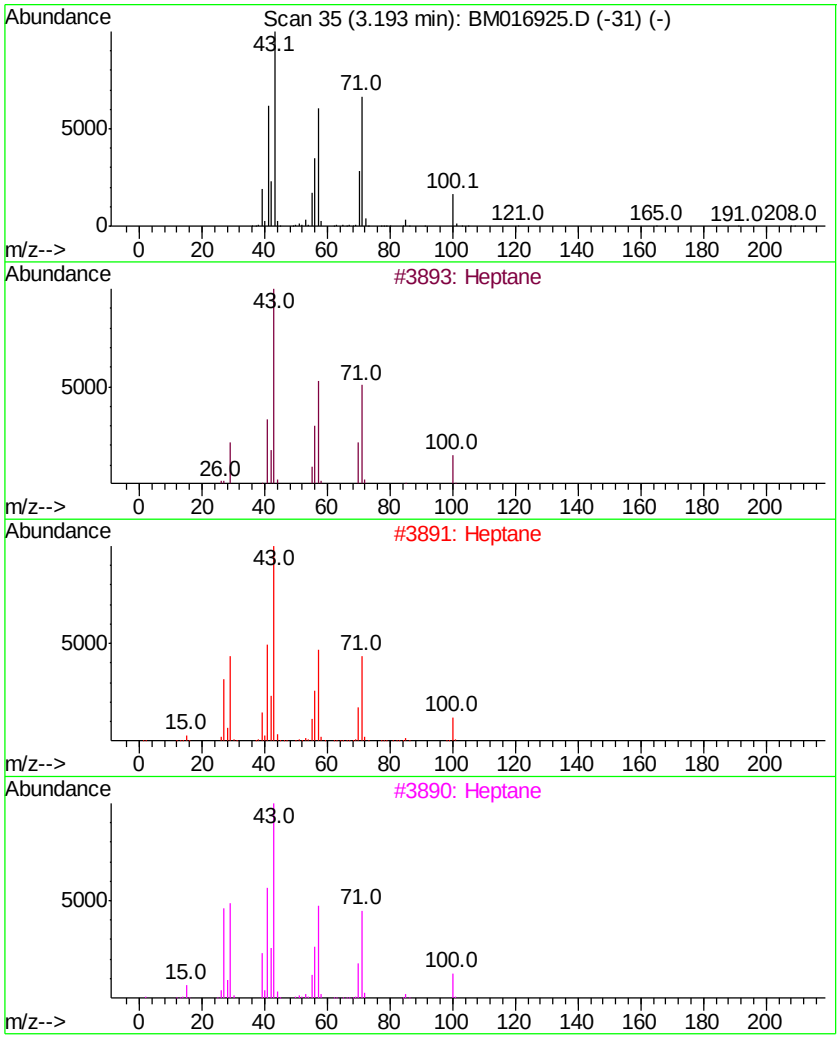
Quant Title : SVOA CALIBRATION

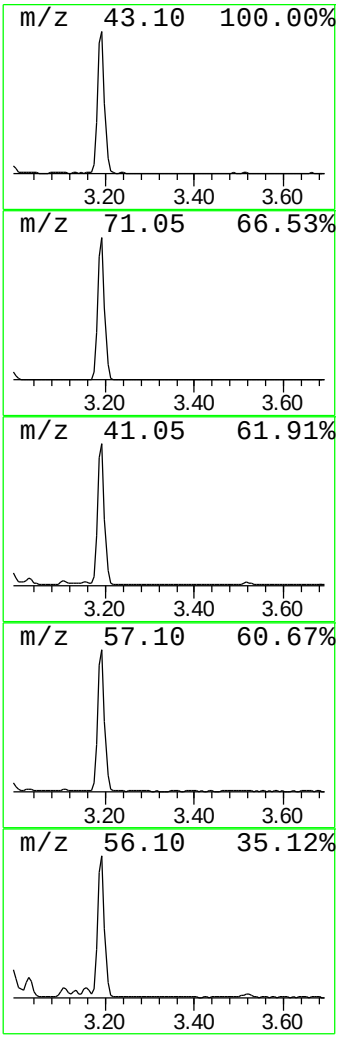
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.19	13.92 ng/ul	1338750	1,4-Dichlorobenzene-d4	7.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	91
2	Heptane	100	C7H16	000142-82-5	91
3	Heptane	100	C7H16	000142-82-5	91
4	Heptane	100	C7H16	000142-82-5	87
5	Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	43





Library Search Compound Report

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Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

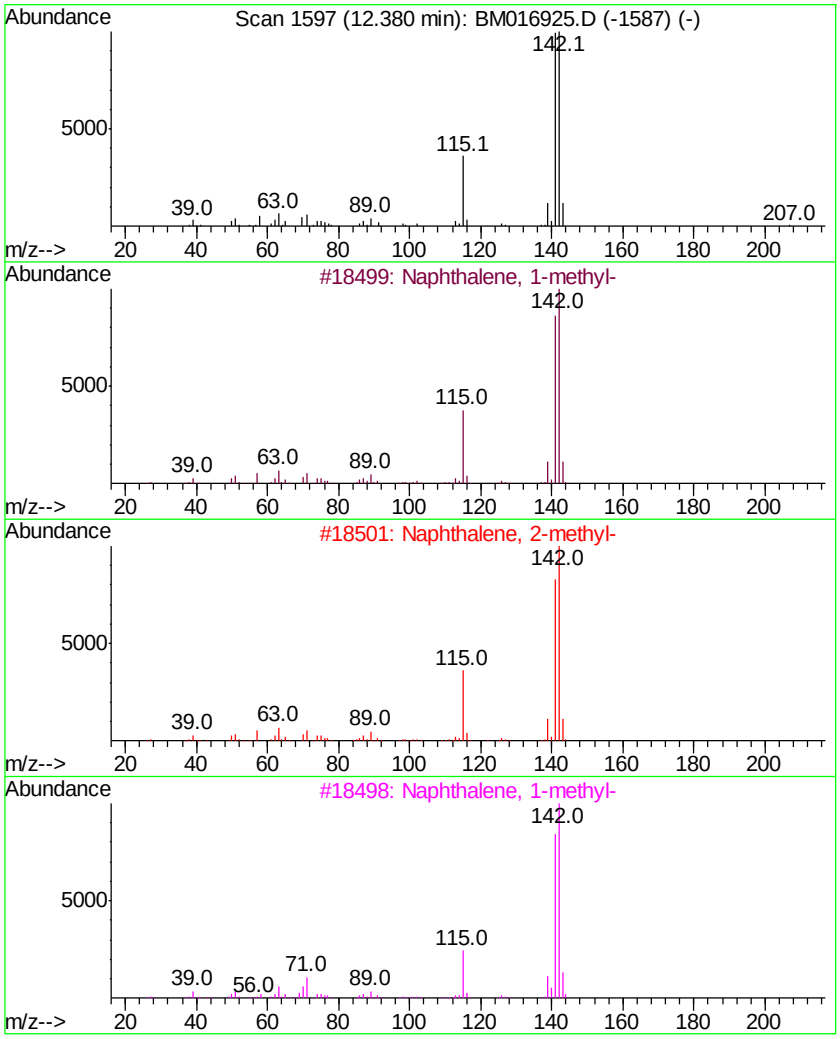
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

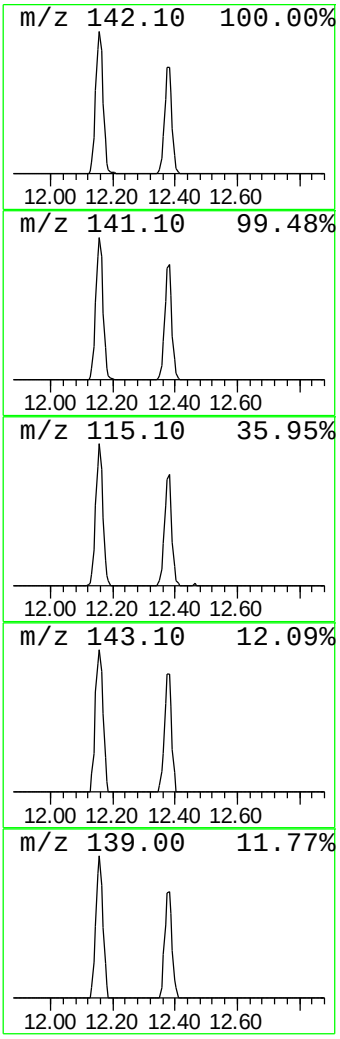
****

Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.38	2.04 ng/ul	275316	Naphthalene-d8	10.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91





Library Search Compound Report

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Quant Title : SVOA CALIBRATION

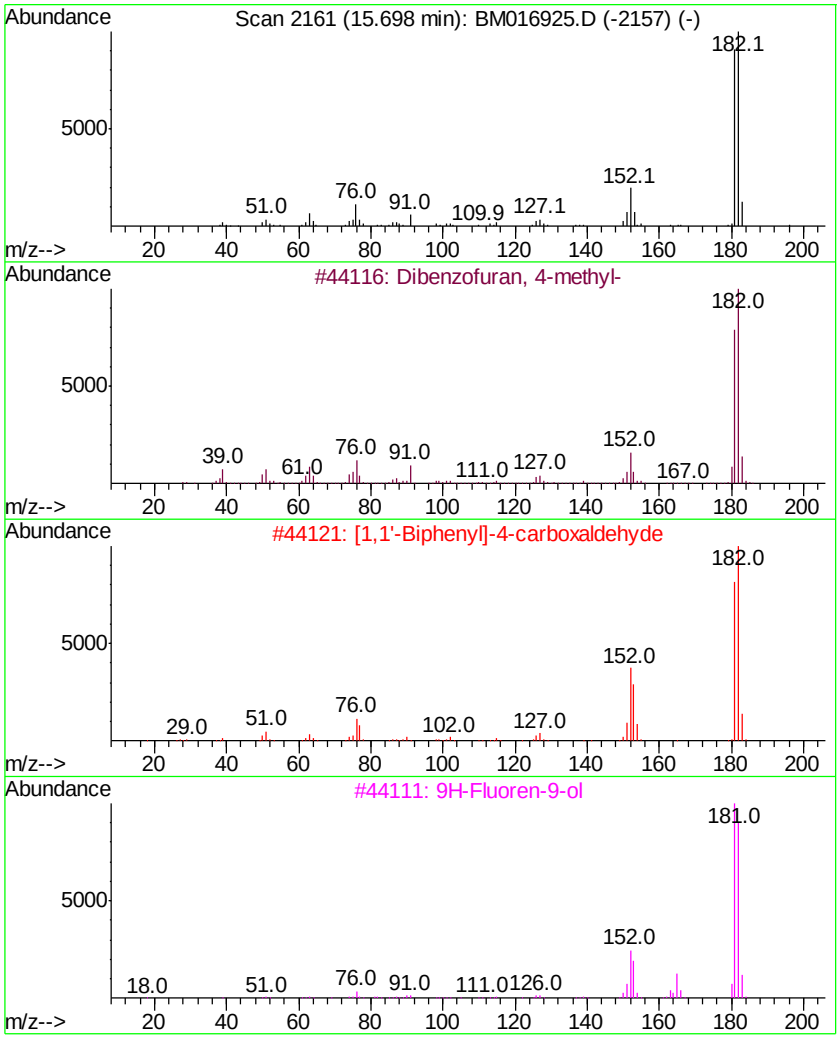
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

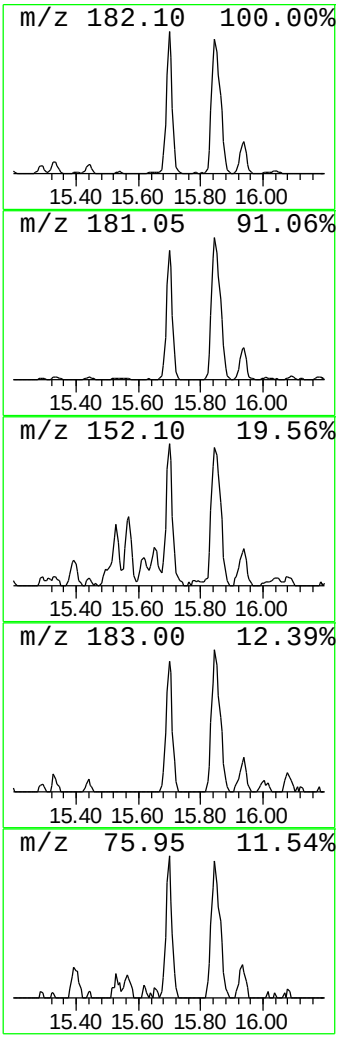
****

Peak Number 3 Dibenzofuran, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	2.17 ng/ul	360902	Acenaphthene-d10	14.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	90
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78
4	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	72
5	3-(2-Naphthyl)acrylaldehyde	182	C13H10O	020884-05-3	72





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

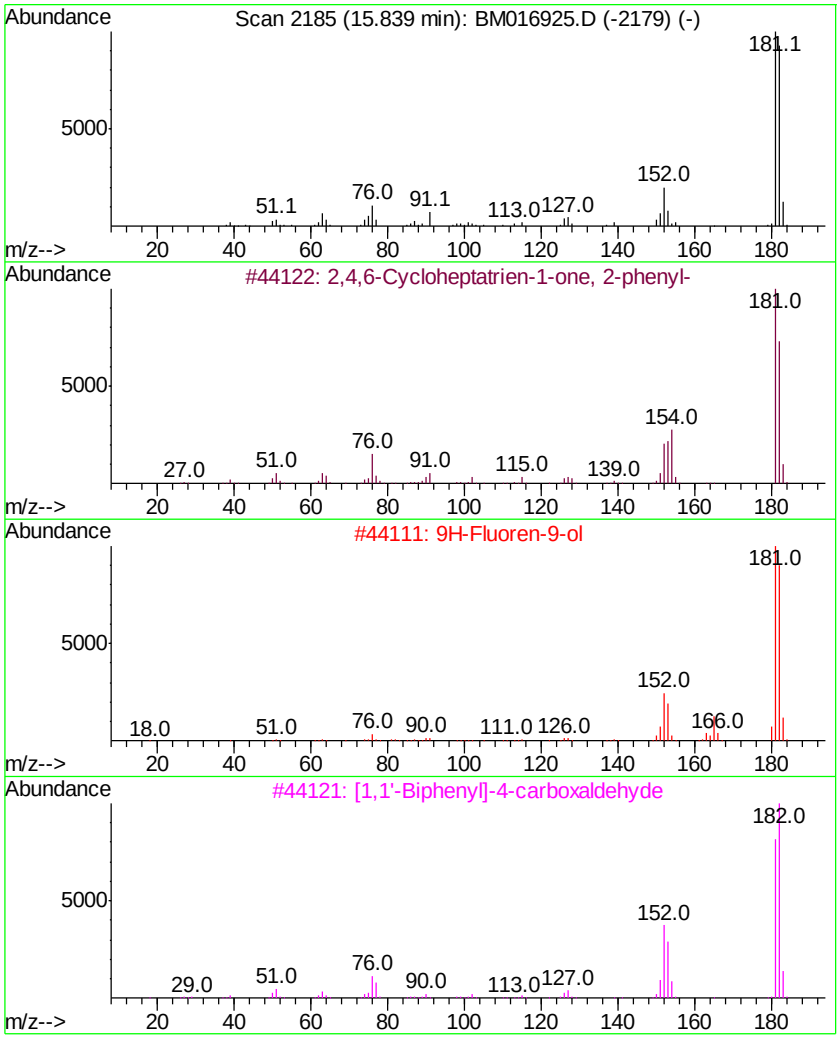
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

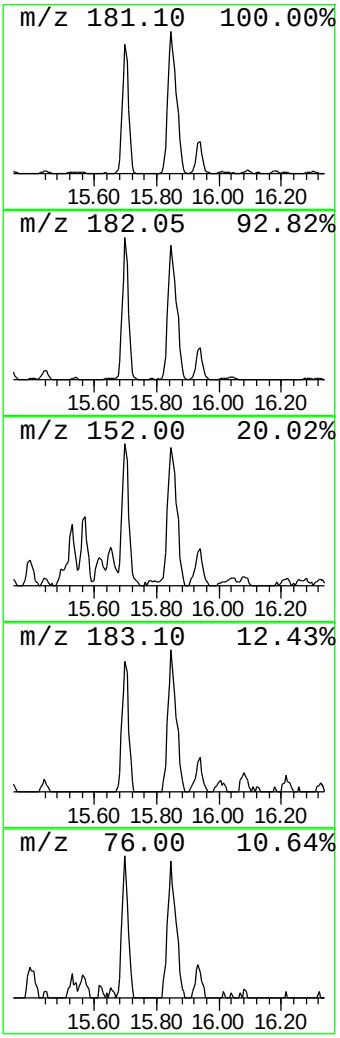
****

Peak Number 4 2,4,6-Cycloheptatrien-1-one... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.84	2.94 ng/ul	489920	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	87
2	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
3	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	83
4	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	81
5	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78





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ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

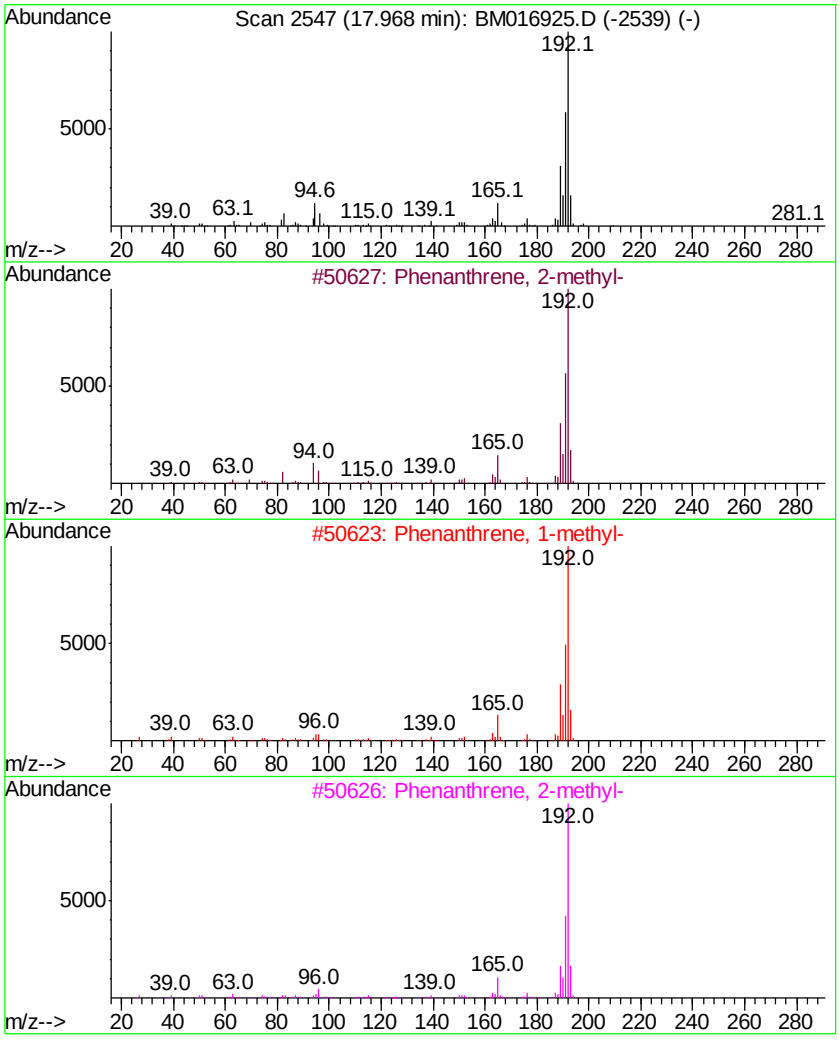
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

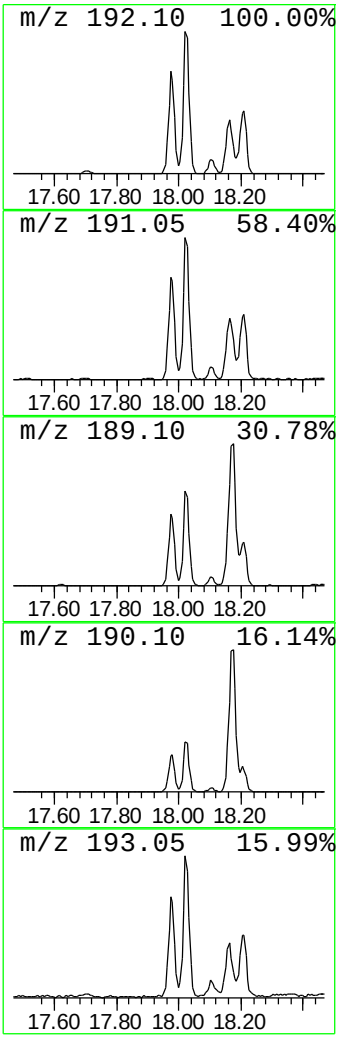
****

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	5.21 ng/ul	867996	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95





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Acq On : 01 Oct 2018 17:26
Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

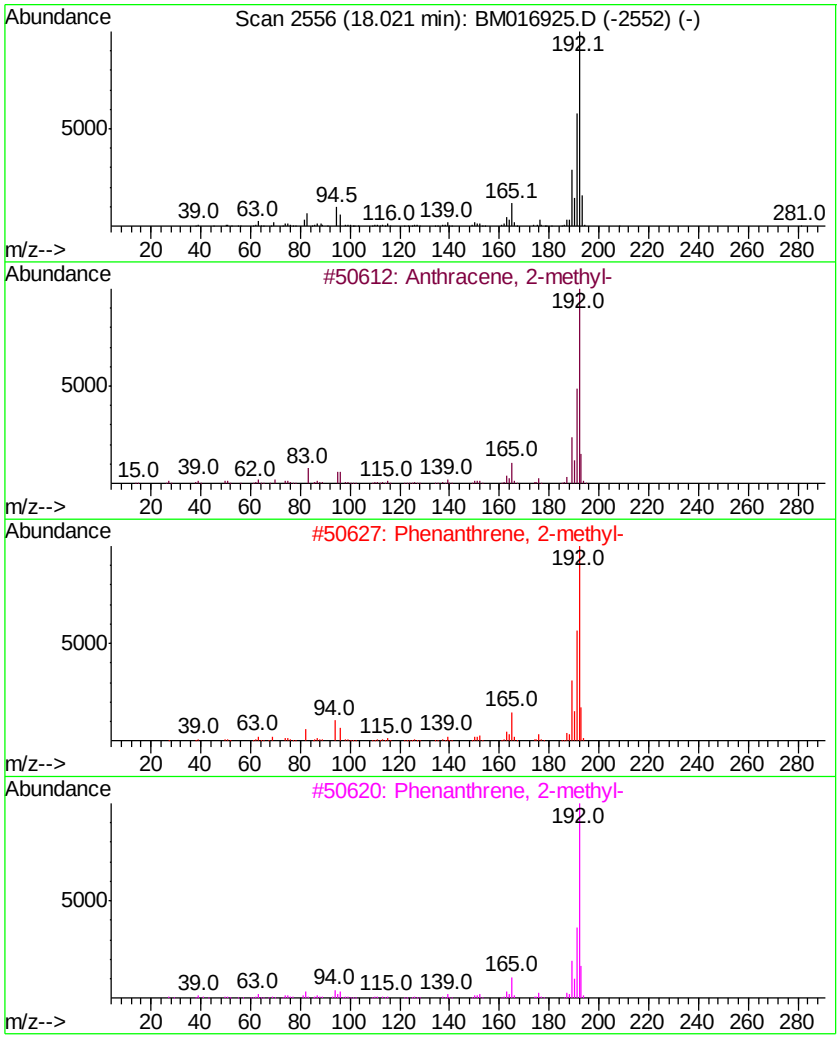
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

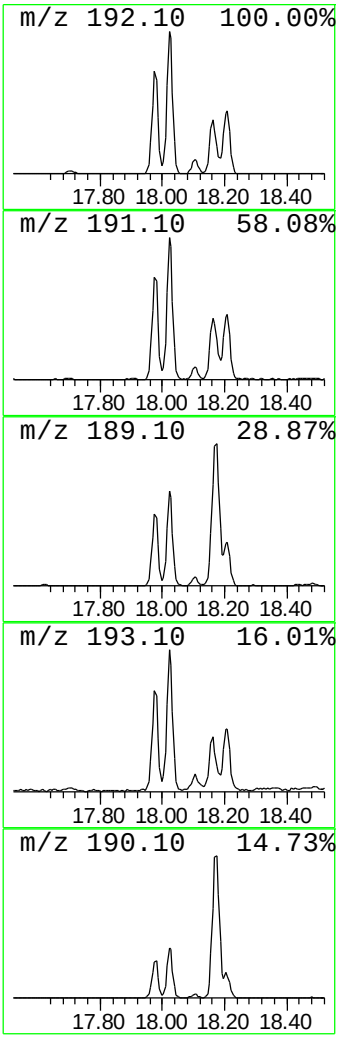
****

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.02	6.71 ng/ul	1119310	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94





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Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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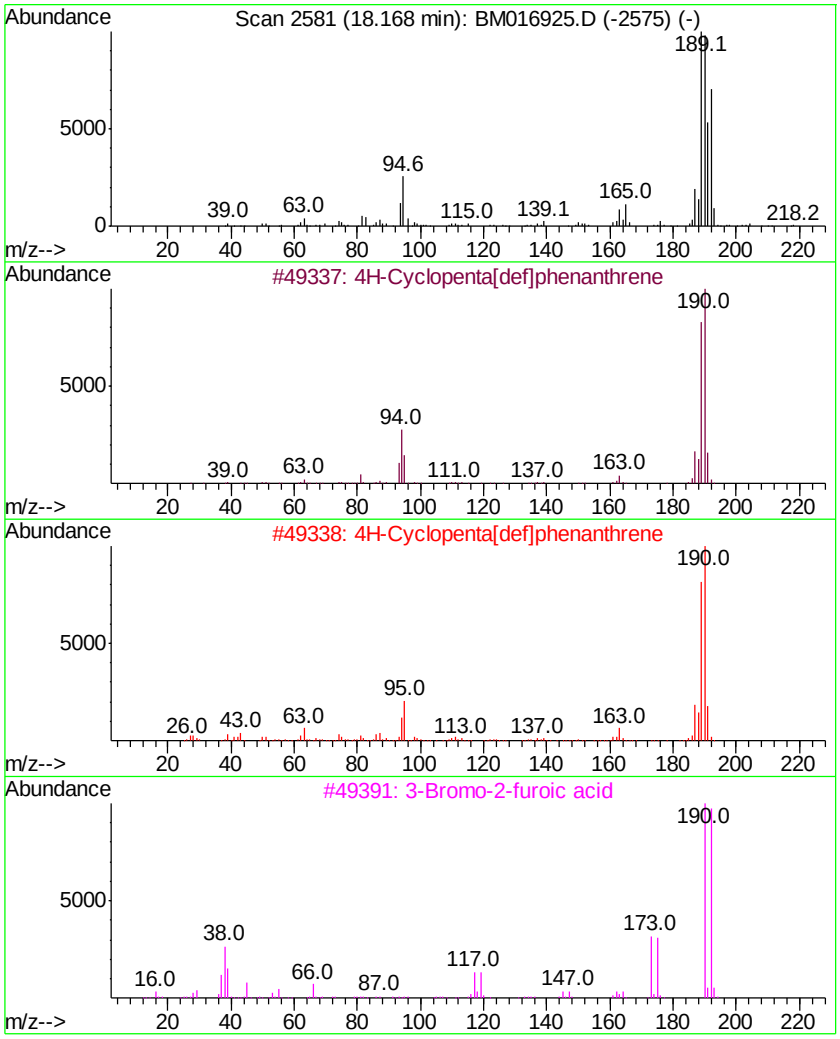
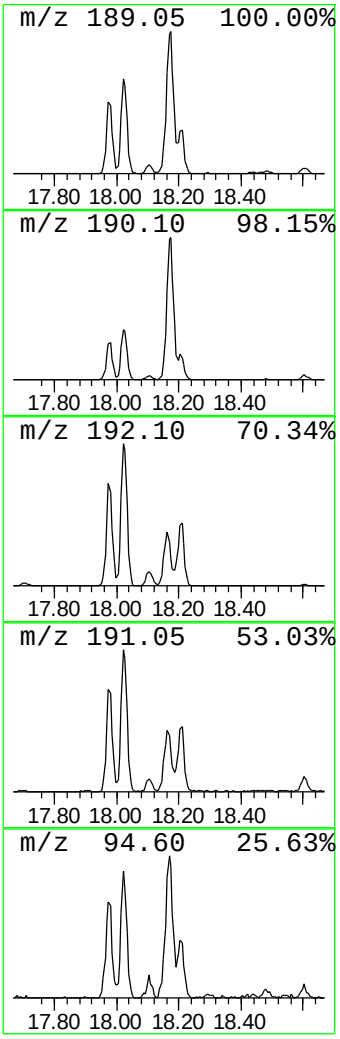
Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	6.28 ng/ul	1046330	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	92
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	41
3	3-Bromo-2-furoic acid	190	C5H3BrO3	014903-90-3	27
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	18
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	15





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Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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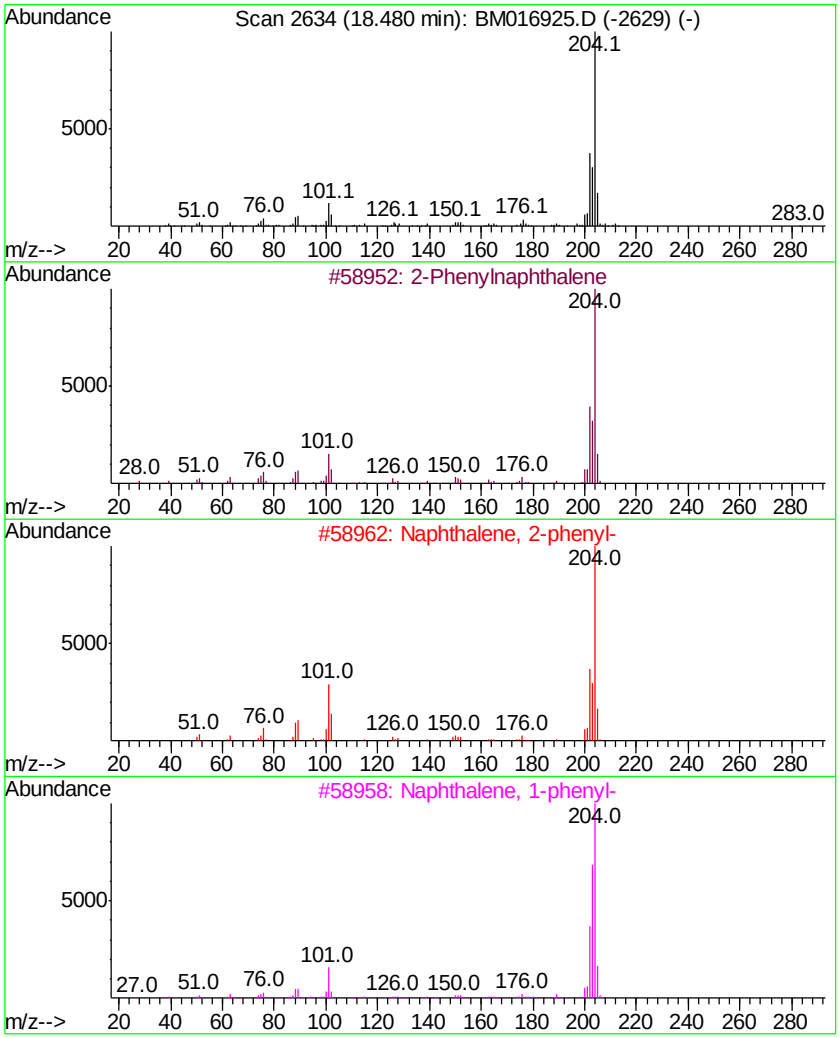
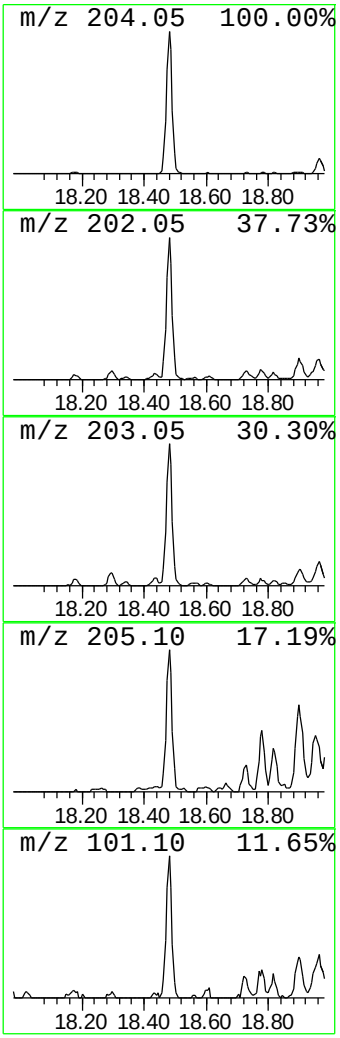
Peak Number 9 2-Phenylnaphthalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	3.07 ng/ul	512158	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenylnaphthalene	204	C16H12	035465-71-5	96
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	94
3	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	92
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	90
5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
Data File : BM016925.D
Acq On : 01 Oct 2018 17:26
Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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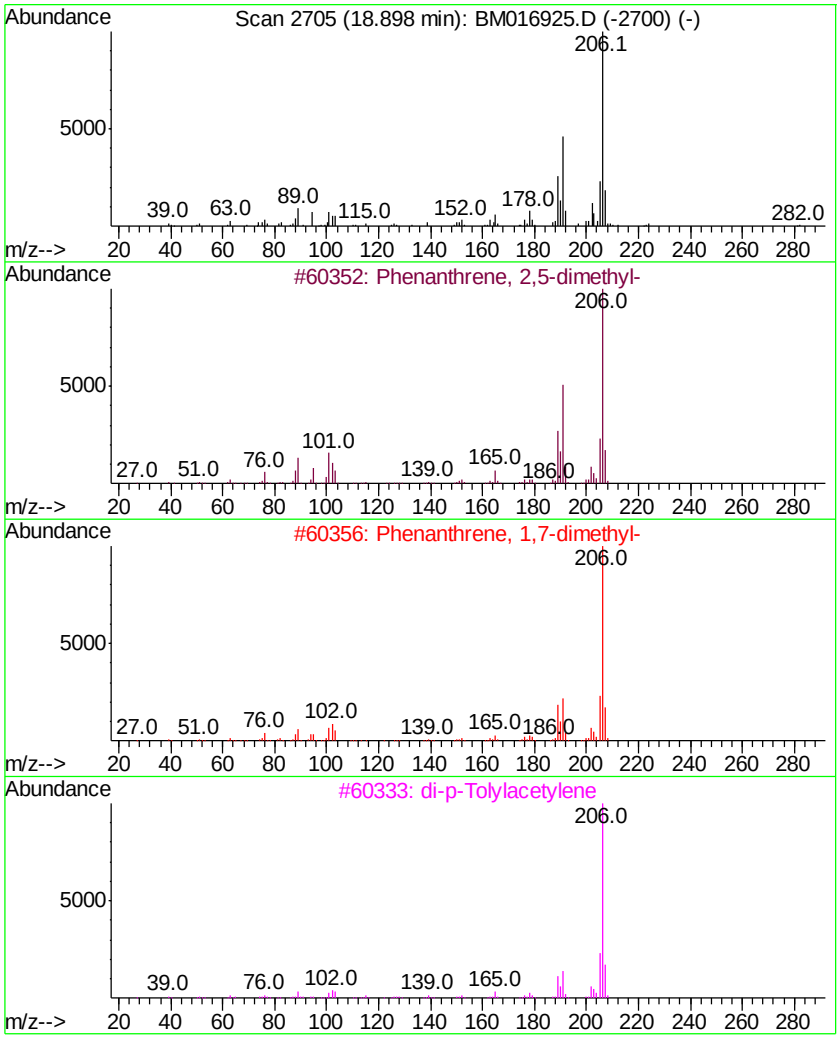
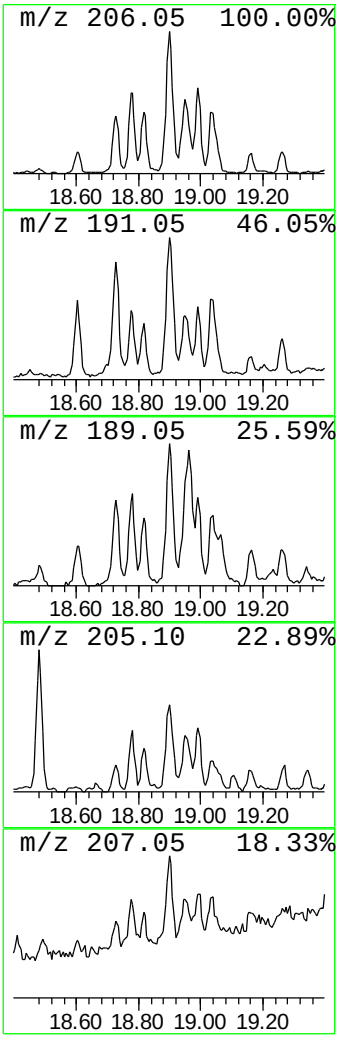
Peak Number 10 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.90	2.16 ng/ul	359909	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
4	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
5	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91





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Data File : BM016925.D
Acq On : 01 Oct 2018 17:26
Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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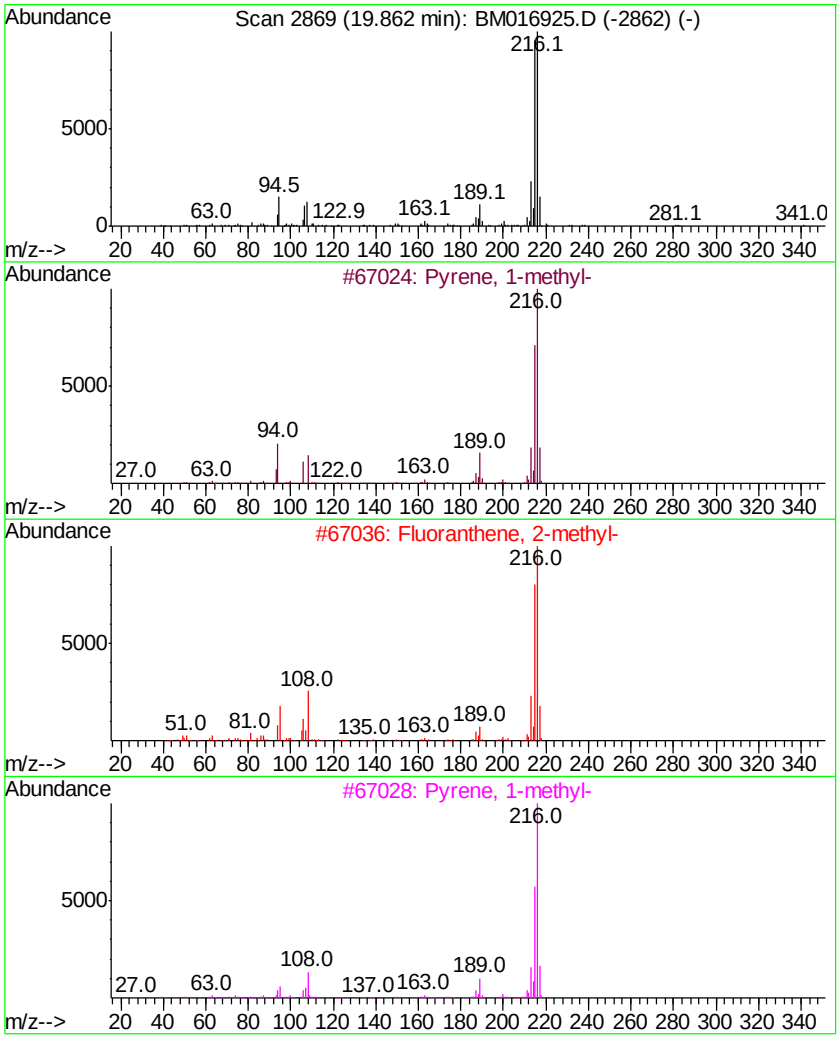
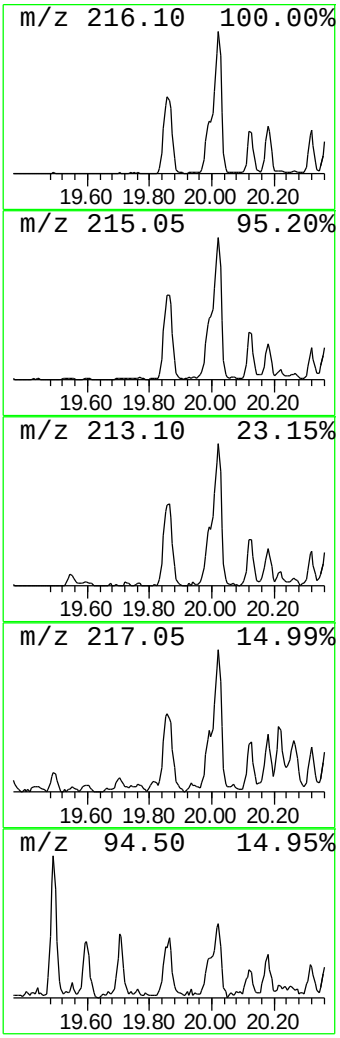
Peak Number 11 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.86	2.11 ng/ul	561841	Chrysene-d12	21.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	92
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	92
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	92





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Data File : BM016925.D
Acq On : 01 Oct 2018 17:26
Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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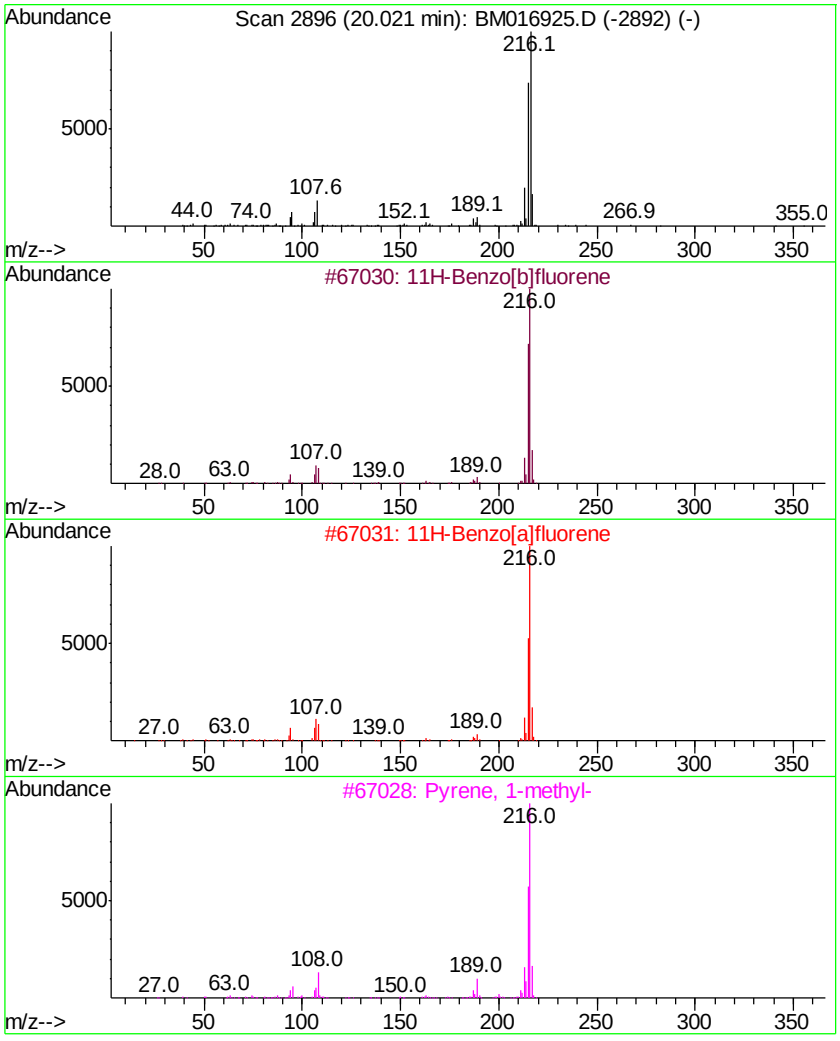
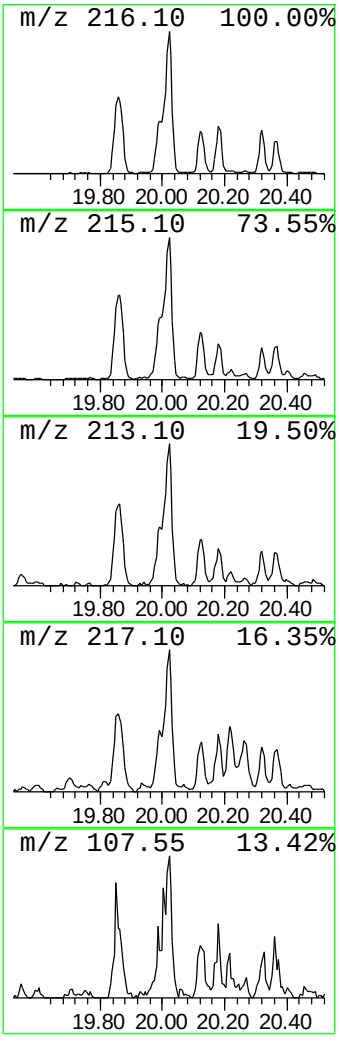
Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.02	2.28 ng/ul	607765	Chrysene-d12	21.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	97
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	96
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100118\
Data File : BM016925.D
Acq On : 01 Oct 2018 17:26
Operator : MJ/SJ
Sample : J5122-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BC0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.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: Str...	3.19	13.9	ng/ul	1338750	1	7.71	1923270	20.0
Naphthalene, 1-me...	12.38	2.0	ng/ul	275316	2	10.49	2698050	20.0
Dibenzofuran, 4-m...	15.70	2.2	ng/ul	360902	3	14.35	3319360	20.0
2,4,6-Cycloheptat...	15.84	2.9	ng/ul	489920	4	17.10	3334250	20.0
Phenanthrene, 2-m...	17.97	5.2	ng/ul	867996	4	17.10	3334250	20.0
Anthracene, 2-met...	18.02	6.7	ng/ul	1119310	4	17.10	3334250	20.0
4H-Cyclopenta[def...	18.17	6.3	ng/ul	1046330	4	17.10	3334250	20.0
2-Phenylnaphthalene	18.48	3.1	ng/ul	512158	4	17.10	3334250	20.0
Phenanthrene, 2,5...	18.90	2.2	ng/ul	359909	4	17.10	3334250	20.0
Pyrene, 1-methyl-	19.86	2.1	ng/ul	561841	5	21.29	5324980	20.0
11H-Benzo[b]fluorene	20.02	2.3	ng/ul	607765	5	21.29	5324980	20.0