

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
 Data File : BM030551.D
 Acq On : 23 Jun 2021 13:49
 Operator : CG/JU
 Sample : M2662-11DL 10X
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGDF5DL

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\SFAM-EPA-BM062221.M
 Title : SVOA CALIBRATION

Signal : TIC: BM030551.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.016	153	158	165	rVB	25467	39316	0.82%	0.079%
2	5.457	399	403	412	rBV2	19637	39812	0.83%	0.080%
3	6.975	651	661	673	rVB2	70832	141677	2.97%	0.286%
4	7.140	684	689	699	rBV	53866	88483	1.86%	0.179%
5	7.340	717	723	732	rVB	69680	116964	2.45%	0.236%
6	7.804	795	802	811	rBV	906108	1427413	29.93%	2.884%
7	8.510	916	922	930	rBV	91264	155899	3.27%	0.315%
8	8.963	992	999	1006	rVV	63630	113463	2.38%	0.229%
9	9.681	1115	1121	1128	rBV	50214	89062	1.87%	0.180%
10	10.222	1206	1213	1227	rBV	74036	151601	3.18%	0.306%
11	10.592	1269	1276	1283	rBV	1240499	2093936	43.91%	4.231%
12	10.639	1283	1284	1291	rVB	34230	43442	0.91%	0.088%
13	10.739	1295	1301	1313	rBV2	63049	140785	2.95%	0.284%
14	13.269	1724	1731	1739	rBV3	19852	34981	0.73%	0.071%
15	13.598	1782	1787	1797	rBV3	21723	59036	1.24%	0.119%
16	13.757	1809	1814	1819	rBV	41982	67115	1.41%	0.136%
17	13.845	1825	1829	1835	rVV	171366	262781	5.51%	0.531%
18	13.992	1846	1854	1857	rBV2	29133	43912	0.92%	0.089%
19	14.127	1871	1877	1881	rBV	196173	332204	6.97%	0.671%
20	14.433	1920	1929	1936	rBV	1879764	2832833	59.40%	5.725%
21	14.498	1936	1940	1948	rVB2	93232	159122	3.34%	0.322%
22	14.651	1959	1966	1975	rBV4	33544	85466	1.79%	0.173%
23	14.833	1990	1997	2005	rVV	154787	231539	4.86%	0.468%
24	14.910	2005	2010	2015	rVV	47144	65047	1.36%	0.131%
25	15.051	2026	2034	2039	rBV2	48812	98059	2.06%	0.198%
26	15.104	2039	2043	2050	rVB	47826	66834	1.40%	0.135%
27	15.227	2055	2064	2066	rBV2	37415	76379	1.60%	0.154%
28	15.298	2072	2076	2082	rVB4	20273	38245	0.80%	0.077%
29	15.427	2090	2098	2102	rVV	279609	431766	9.05%	0.873%
30	15.480	2102	2107	2116	rVV2	258156	422787	8.87%	0.854%
31	15.604	2125	2128	2131	rVV5	22487	35790	0.75%	0.072%
32	15.645	2131	2135	2139	rVV4	22077	36880	0.77%	0.075%
33	15.692	2139	2143	2147	rVV2	30534	49369	1.04%	0.100%
34	15.774	2152	2157	2164	rVV	166965	256256	5.37%	0.518%
35	15.845	2164	2169	2173	rVB4	27503	35327	0.74%	0.071%
36	15.921	2173	2182	2191	rVB2	189661	426701	8.95%	0.862%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

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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\SFAM-EPA-BM062221.M
 Title : SVOA CALIBRATION

37	16.010	2191	2197	2205	rVB2	39384	70538	1.48%	0.143%
38	16.163	2216	2223	2229	rVB3	42077	83234	1.75%	0.168%
39	16.480	2269	2277	2282	rVV2	126381	268829	5.64%	0.543%
40	16.533	2282	2286	2291	rVV	50828	85589	1.79%	0.173%
41	16.633	2297	2303	2308	rVV2	61122	110543	2.32%	0.223%
42	16.710	2308	2316	2320	rVV2	114168	208178	4.37%	0.421%
43	16.751	2320	2323	2326	rVV	101071	145904	3.06%	0.295%
44	16.780	2326	2328	2333	rVV	49270	75007	1.57%	0.152%
45	16.839	2333	2338	2344	rVV3	54567	112547	2.36%	0.227%
46	16.904	2344	2349	2356	rVV	109768	234087	4.91%	0.473%
47	16.962	2356	2359	2360	rVV2	31577	39789	0.83%	0.080%
48	16.992	2360	2364	2378	rVB	163608	293172	6.15%	0.592%
49	17.174	2388	2395	2398	rBV	2029681	2935476	61.56%	5.932%
50	17.215	2398	2402	2407	rVV	2650882	3686438	77.30%	7.450%
51	17.274	2407	2412	2413	rVV	272821	368854	7.73%	0.745%
52	17.304	2413	2417	2423	rVV	1145178	1672722	35.08%	3.380%
53	17.362	2423	2427	2433	rVV3	62583	127518	2.67%	0.258%
54	17.445	2438	2441	2445	rVV2	34370	53896	1.13%	0.109%
55	17.615	2466	2470	2476	rVV2	45302	100089	2.10%	0.202%
56	17.692	2480	2483	2490	rVV	68981	98652	2.07%	0.199%
57	17.757	2490	2494	2503	rVB6	33087	76315	1.60%	0.154%
58	17.892	2513	2517	2526	rVB2	34649	69055	1.45%	0.140%
59	18.045	2533	2543	2547	rBV	366933	563692	11.82%	1.139%
60	18.092	2547	2551	2558	rVB	469196	690546	14.48%	1.395%
61	18.174	2559	2565	2571	rVV	292864	427171	8.96%	0.863%
62	18.245	2571	2577	2581	rVV2	616676	1106446	23.20%	2.236%
63	18.280	2581	2583	2591	rVV	230544	269153	5.64%	0.544%
64	18.357	2593	2596	2600	rVV3	26513	35571	0.75%	0.072%
65	18.545	2624	2628	2633	rVV	273922	357652	7.50%	0.723%
66	18.792	2666	2670	2674	rBV2	58359	79732	1.67%	0.161%
67	18.845	2675	2679	2682	rBV	72569	90096	1.89%	0.182%
68	18.880	2682	2685	2689	rVB	69219	90087	1.89%	0.182%
69	18.962	2693	2699	2703	rBV2	168482	313561	6.58%	0.634%
70	19.027	2706	2710	2713	rBV2	81644	111951	2.35%	0.226%
71	19.098	2719	2722	2726	rBV3	71609	109347	2.29%	0.221%
72	19.227	2738	2744	2750	rBV	3741282	4767854	99.98%	9.635%
73	19.292	2751	2755	2758	rVV	48022	63749	1.34%	0.129%
74	19.321	2758	2760	2765	rVB2	57348	71394	1.50%	0.144%
75	19.374	2765	2769	2774	rBV	273733	380807	7.99%	0.770%
76	19.468	2780	2785	2788	rBV3	51171	97775	2.05%	0.198%

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 Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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 Title : SVOA CALIBRATION

77	19.556	2795	2800	2802	rBV3	726914	1011150	21.20%	2.043%
78	19.586	2802	2805	2814	rVV	1249117	1930445	40.48%	3.901%
79	19.662	2814	2818	2824	rVB2	151716	249401	5.23%	0.504%
80	19.768	2832	2836	2843	rVB	204684	293736	6.16%	0.594%
81	19.827	2843	2846	2853	rVB3	57346	93389	1.96%	0.189%
82	19.909	2854	2860	2870	rBV3	182107	475882	9.98%	0.962%
83	20.045	2878	2883	2885	rBV3	141381	231971	4.86%	0.469%
84	20.080	2885	2889	2895	rVB	545577	781536	16.39%	1.579%
85	20.180	2902	2906	2910	rBV	535183	668316	14.01%	1.351%
86	20.239	2910	2916	2921	rVB2	199591	375821	7.88%	0.759%
87	20.321	2926	2930	2936	rVV5	77538	161690	3.39%	0.327%
88	20.380	2937	2940	2943	rVV	66665	84370	1.77%	0.170%
89	20.427	2944	2948	2952	rVB3	105773	150226	3.15%	0.304%
90	20.550	2966	2969	2974	rVB	190900	214802	4.50%	0.434%
91	20.786	3006	3009	3014	rBV	75198	111297	2.33%	0.225%
92	20.992	3041	3044	3047	rBV	104232	128590	2.70%	0.260%
93	21.027	3047	3050	3054	rBV	197331	239296	5.02%	0.484%
94	21.339	3097	3103	3107	rBV2	2507133	4768709	100.00%	9.637%
95	21.380	3107	3110	3119	rVB	1099736	1424903	29.88%	2.879%
96	21.497	3127	3130	3135	rBV3	80707	149903	3.14%	0.303%
97	21.545	3136	3138	3142	rVB	86974	110340	2.31%	0.223%
98	21.892	3195	3197	3201	rVB	186067	223589	4.69%	0.452%
99	22.933	3367	3374	3378	rBV	674259	1615735	33.88%	3.265%
100	23.609	3483	3489	3498	rBV2	1407715	2657238	55.72%	5.370%

Sum of corrected areas: 49485633

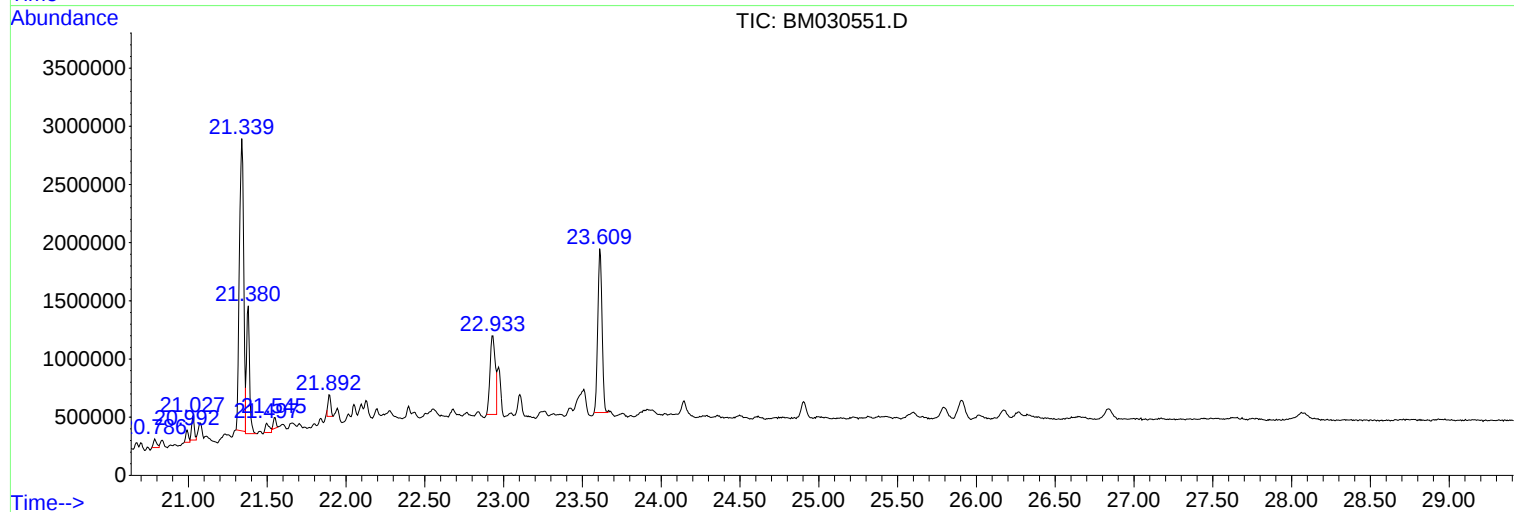
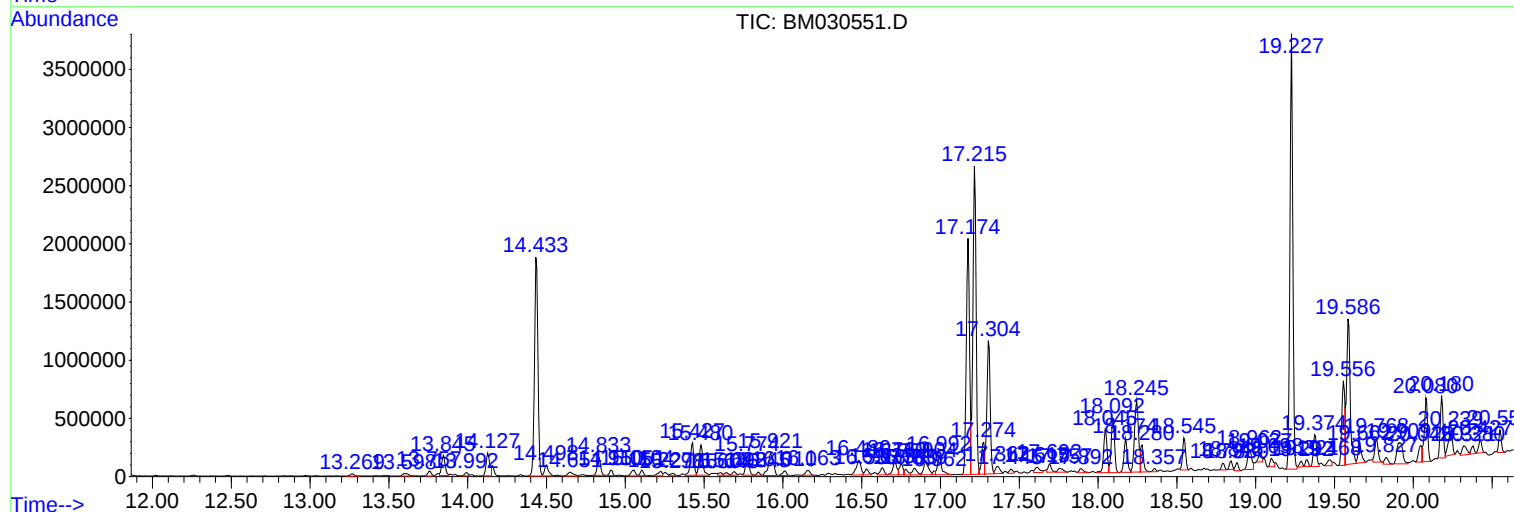
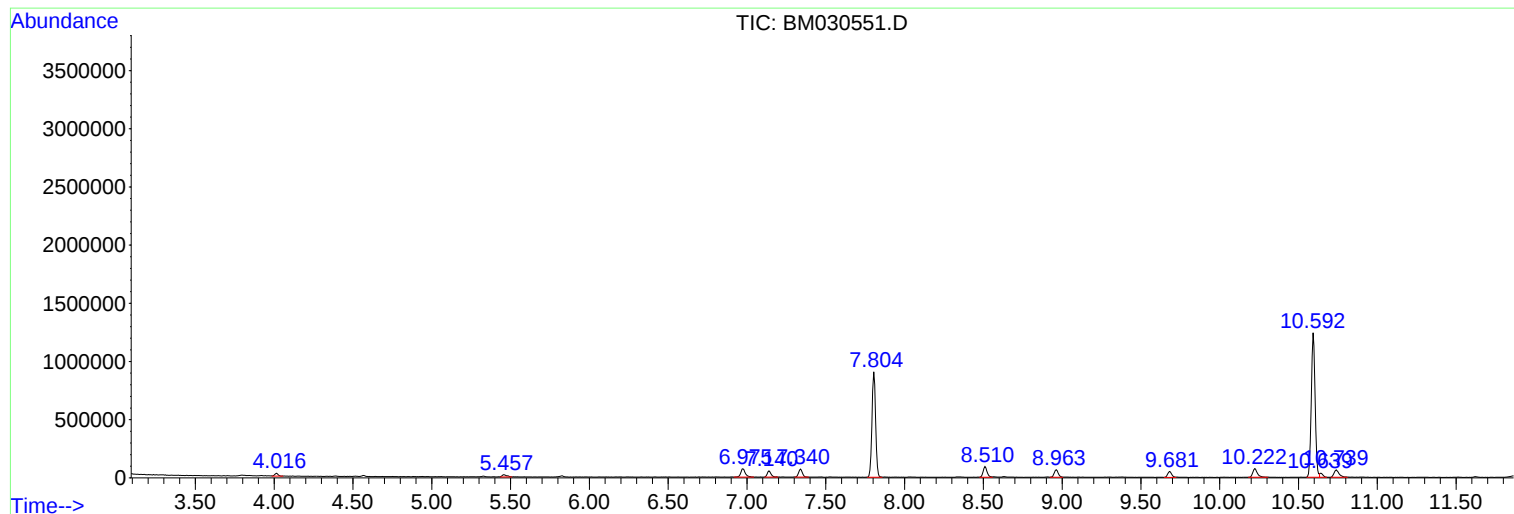
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L

TIC Integration Parameters: LSCINT.P



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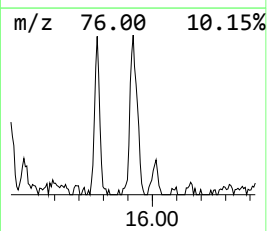
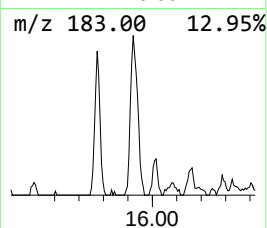
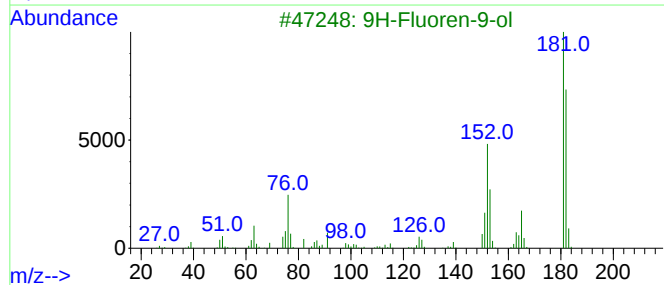
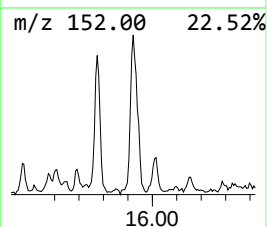
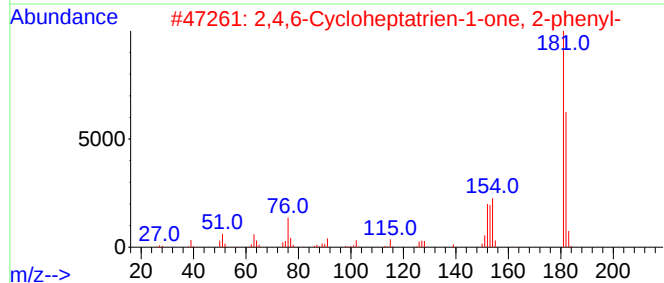
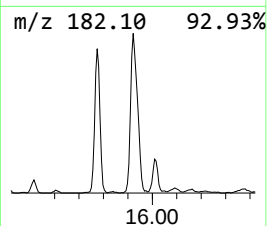
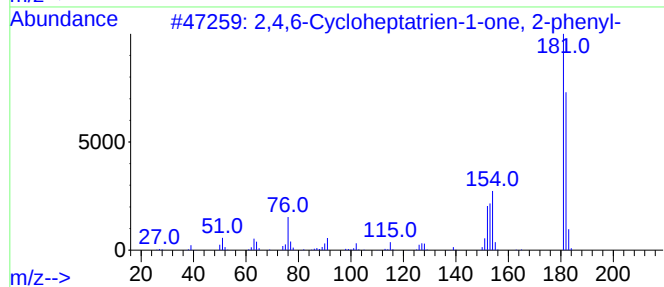
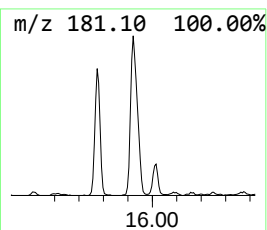
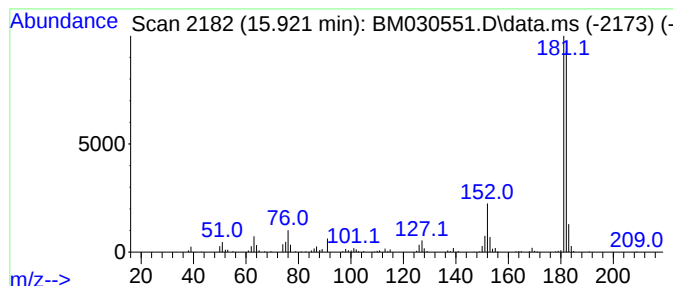
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.920	2.91 ng/ul	426701	Phenanthrene-d10	17.174

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			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	87
3			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
4			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
5			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	83



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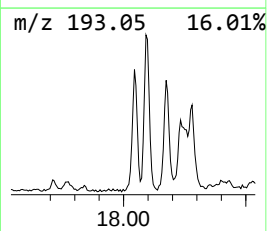
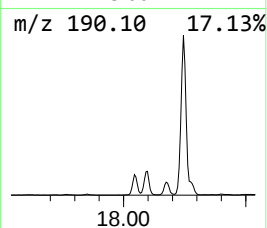
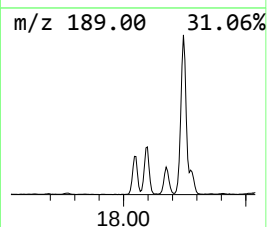
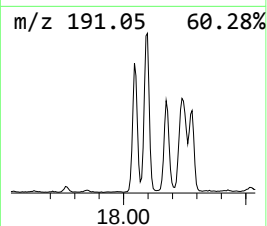
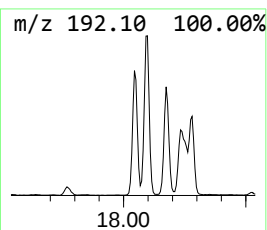
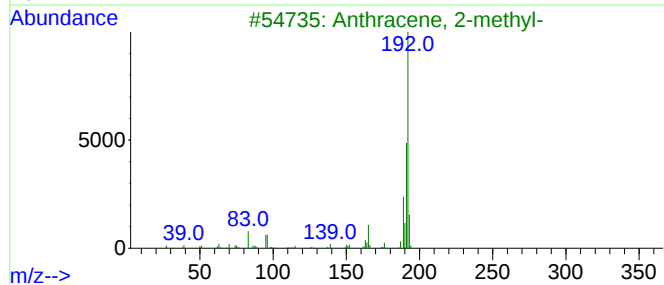
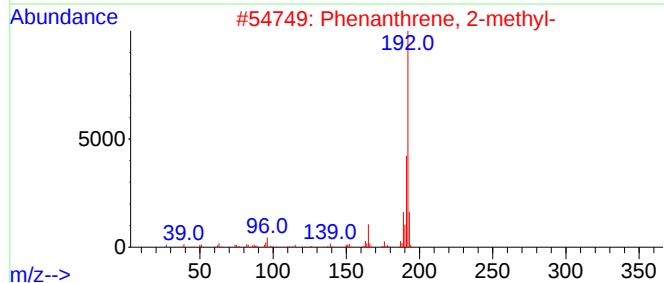
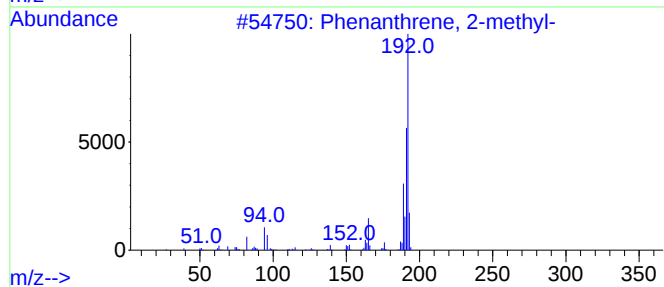
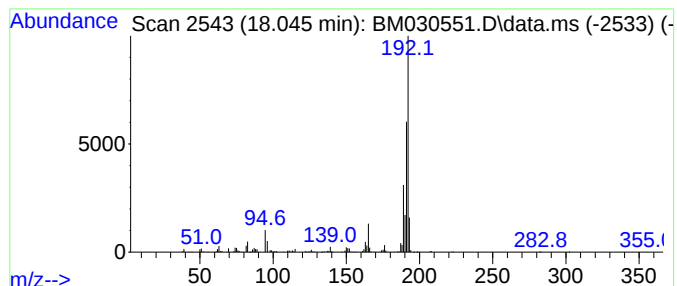
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Anthracene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.040	3.84 ng/ul	563692	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



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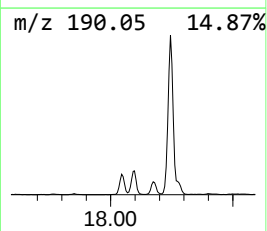
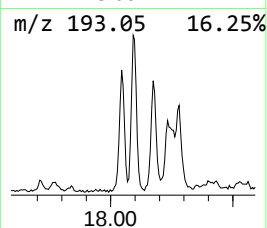
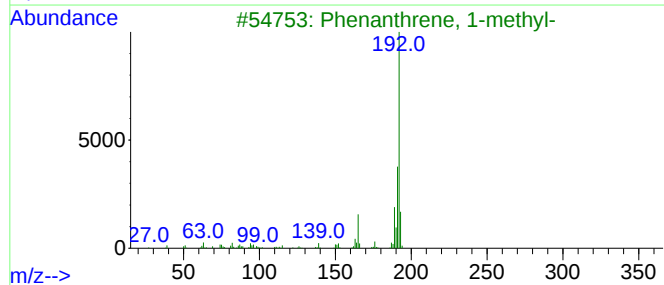
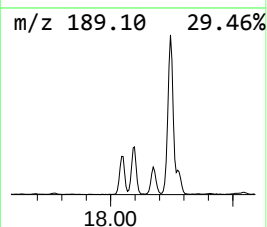
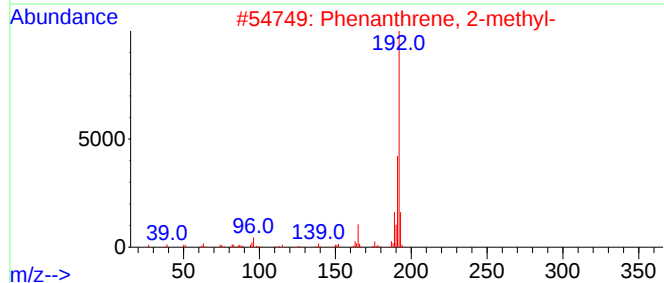
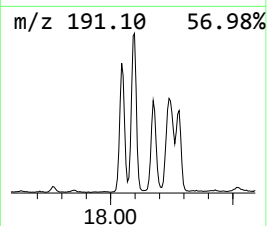
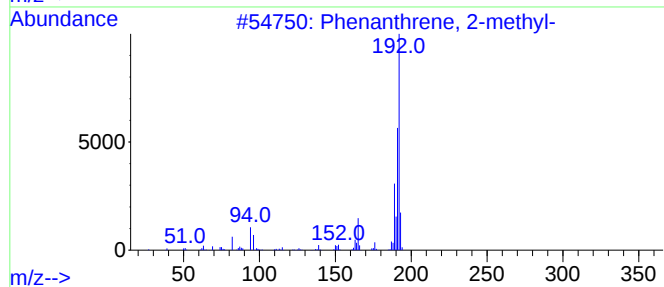
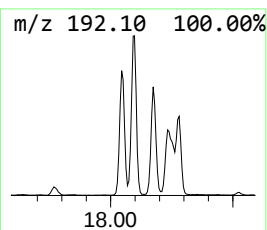
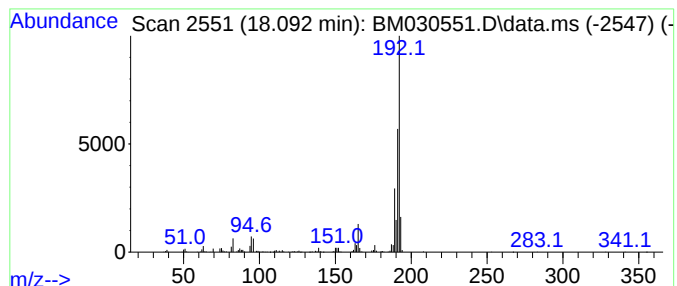
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.090	4.70 ng/ul	690546	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
Data File : BM030551.D
Acq On : 23 Jun 2021 13:49
Operator : CG/JU
Sample : M2662-11DL 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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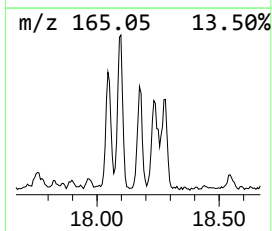
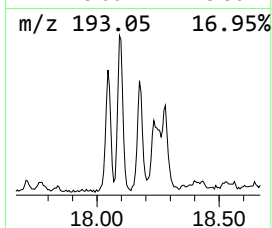
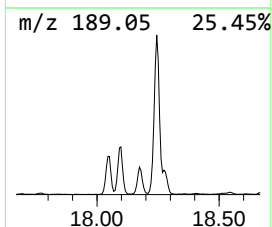
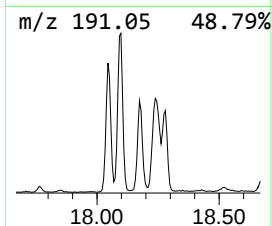
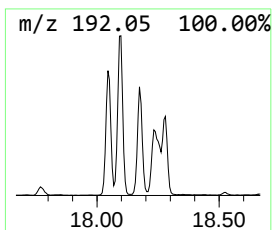
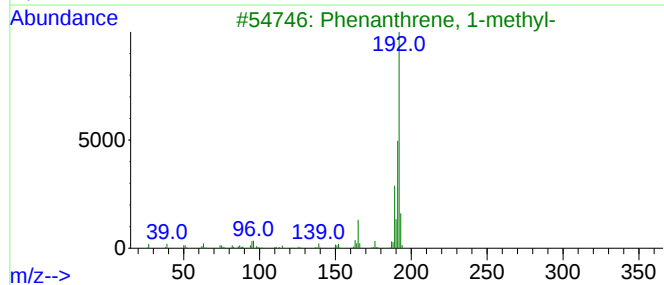
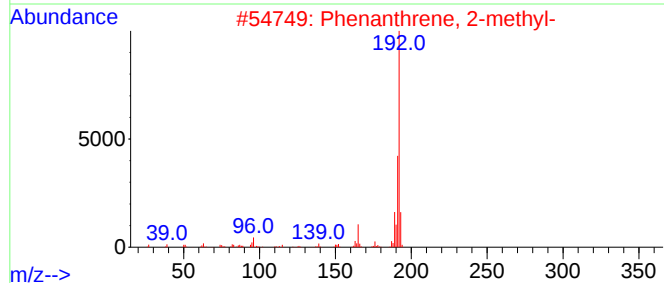
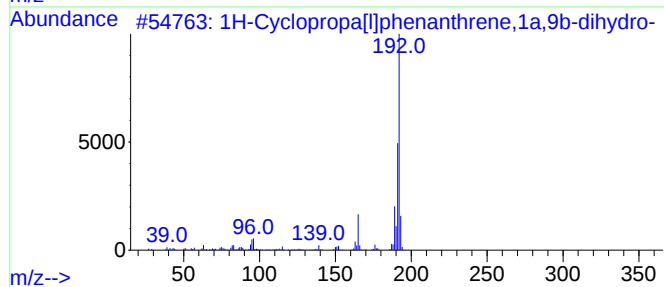
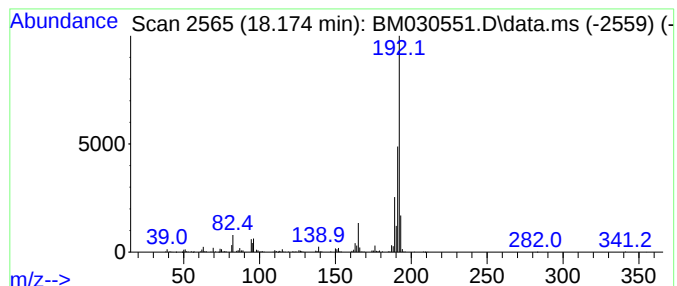
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1H-Cyclopropa[1]phenanthren... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.170	2.91 ng/ul	427171	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	98
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
Data File : BM030551.D
Acq On : 23 Jun 2021 13:49
Operator : CG/JU
Sample : M2662-11DL 10X
Misc :
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ClientSampleId :
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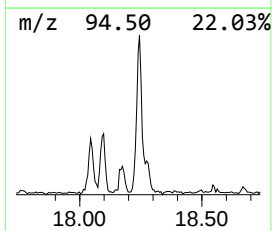
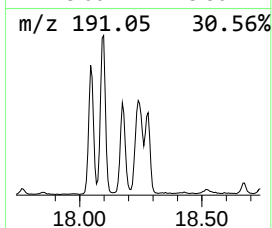
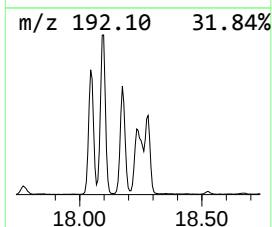
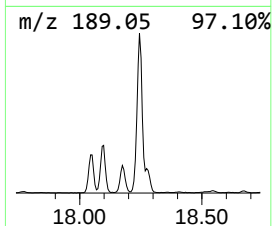
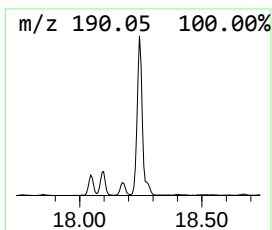
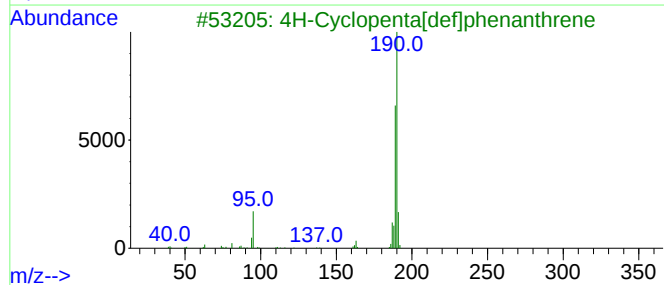
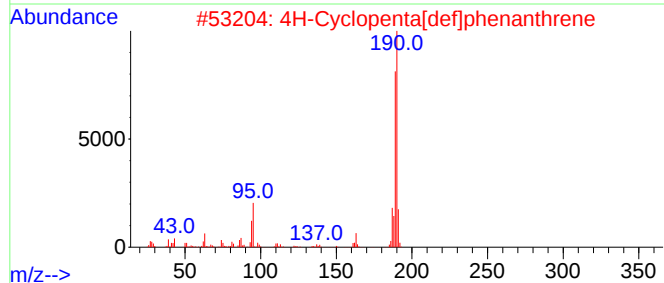
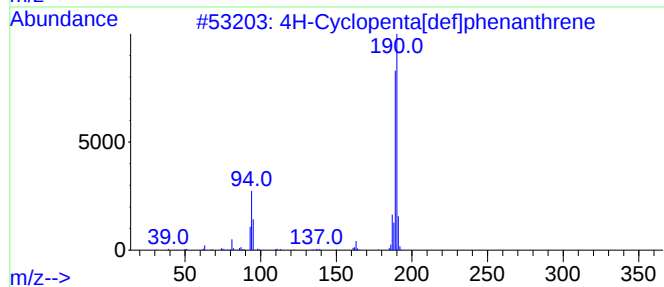
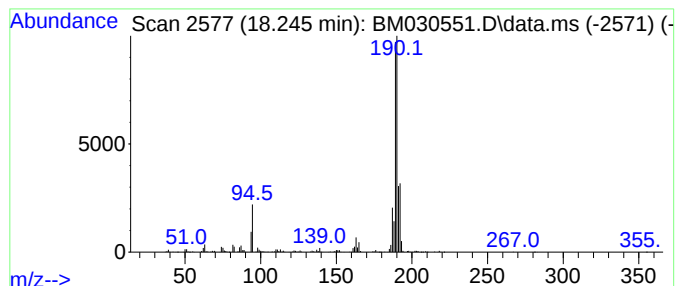
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.240	7.54 ng/ul	1106450	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
5			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
Data File : BM030551.D
Acq On : 23 Jun 2021 13:49
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Misc :
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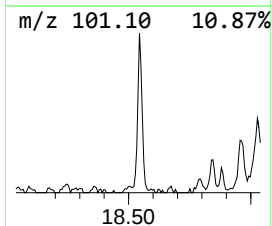
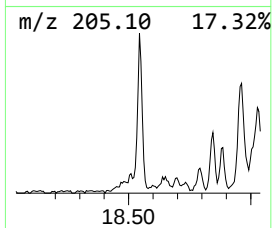
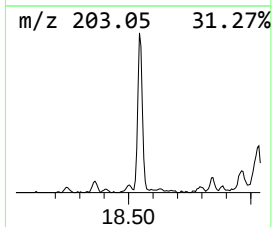
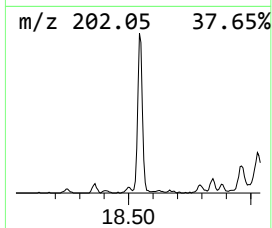
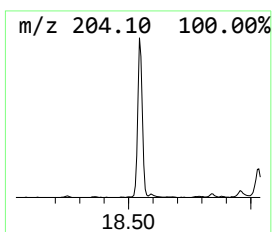
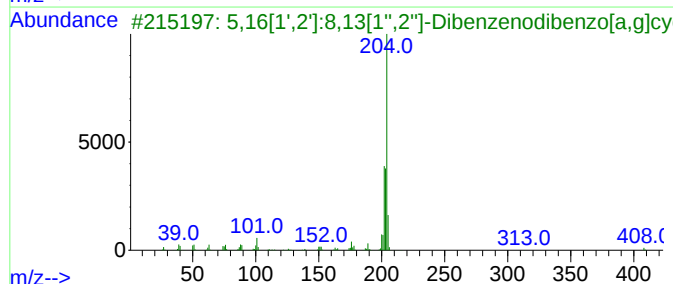
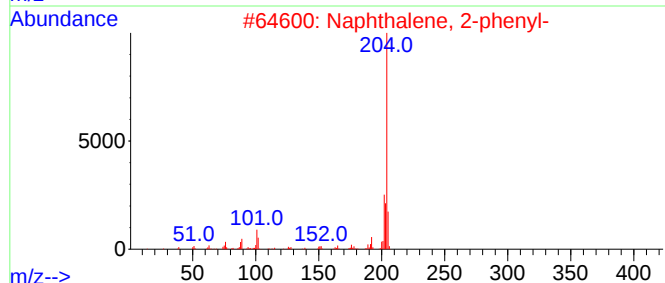
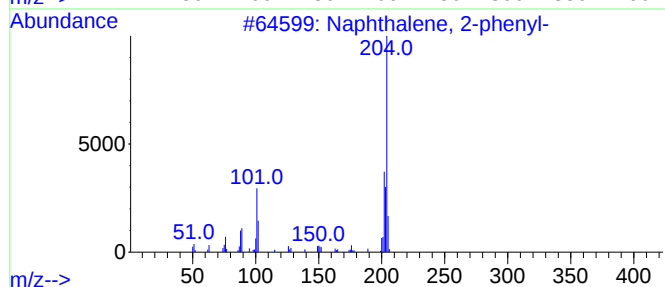
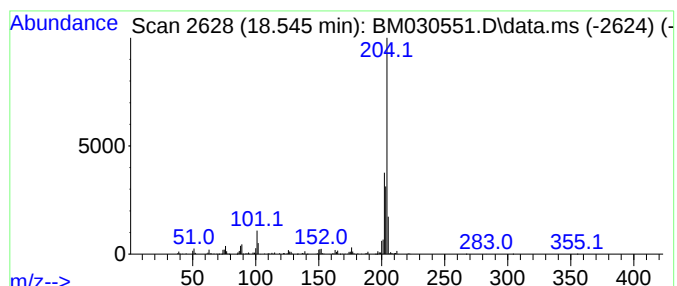
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 2-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.540	2.44 ng/ul	357652	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	90
2			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	87
3			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	74
5			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
Data File : BM030551.D
Acq On : 23 Jun 2021 13:49
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Sample : M2662-11DL 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

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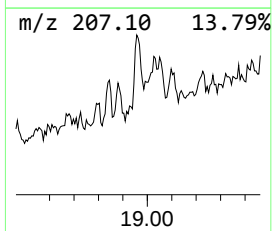
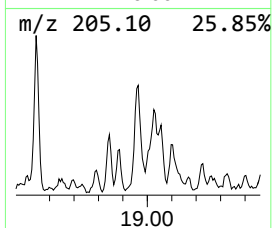
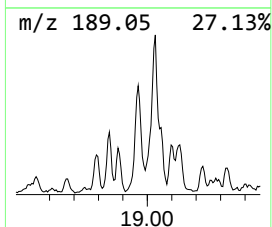
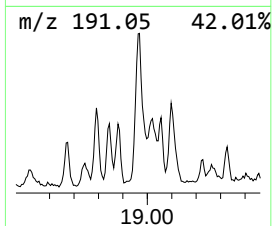
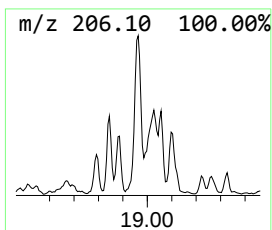
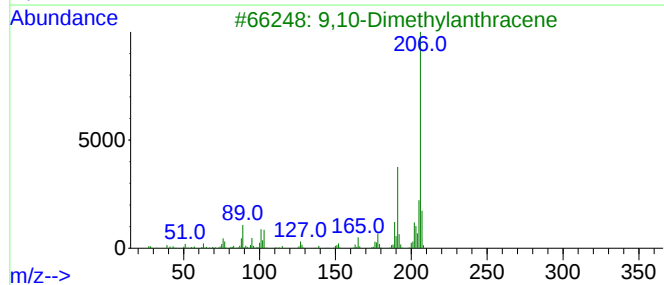
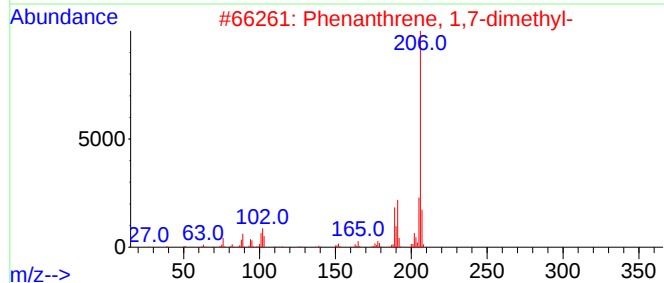
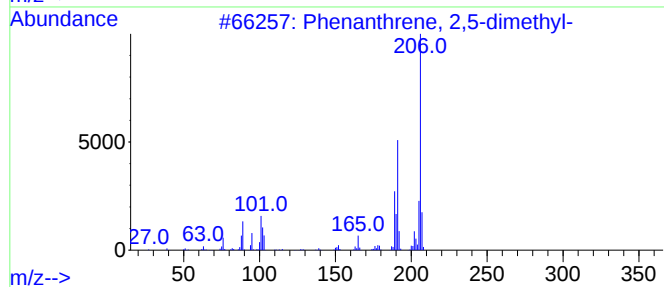
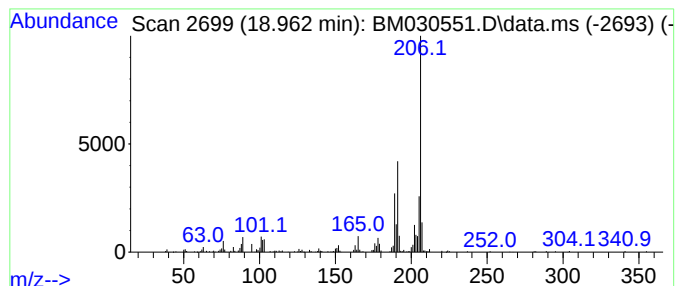
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.960	2.14 ng/ul	313561	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
3			9,10-Dimethylantracene	206	C16H14	000781-43-1	91
4			9,10-Dimethylantracene	206	C16H14	000781-43-1	90
5			di-p-Tolylacetylene	206	C16H14	002789-88-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
Data File : BM030551.D
Acq On : 23 Jun 2021 13:49
Operator : CG/JU
Sample : M2662-11DL 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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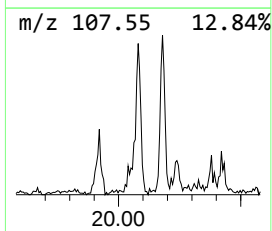
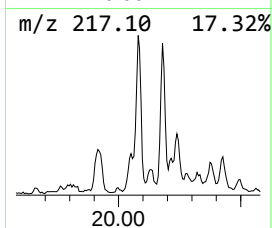
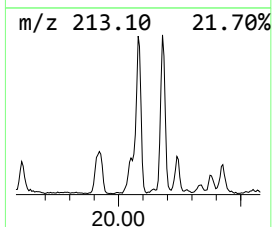
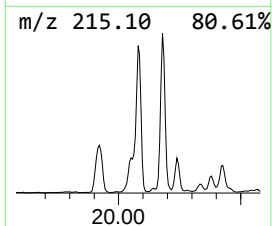
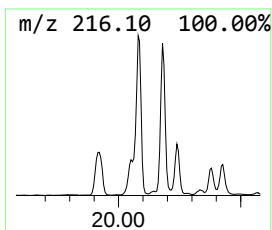
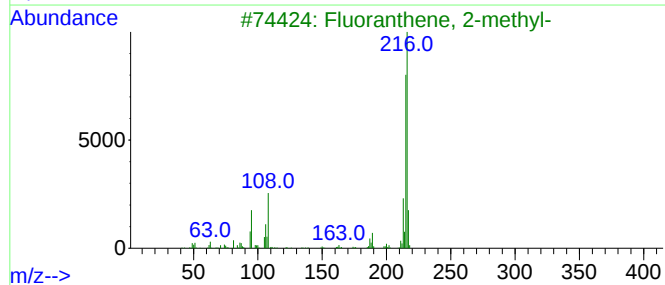
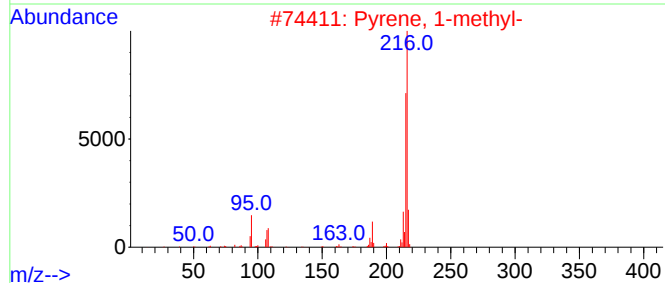
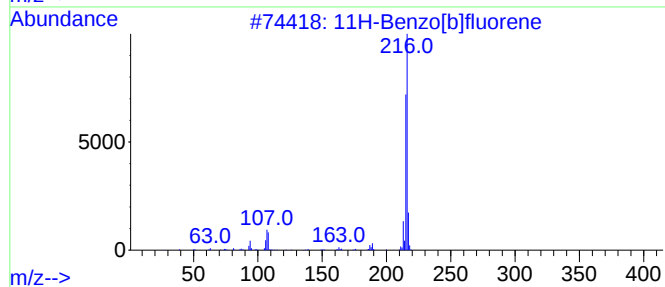
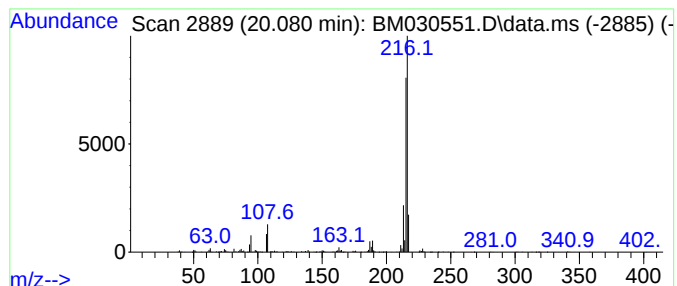
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.080	3.28 ng/ul	781536	Chrysene-d12	21.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
Data File : BM030551.D
Acq On : 23 Jun 2021 13:49
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Misc :
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Instrument :
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ClientSampleId :
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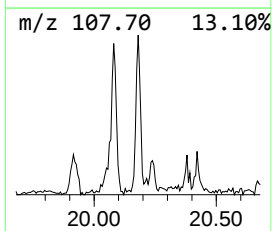
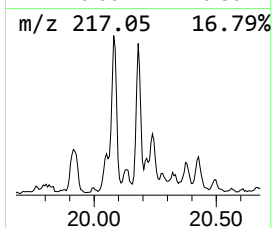
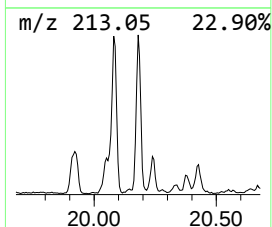
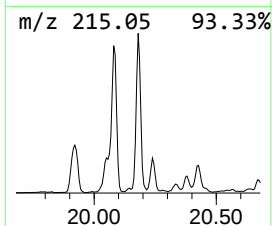
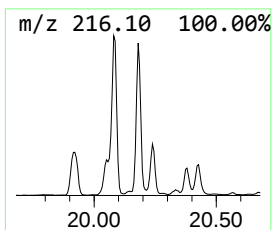
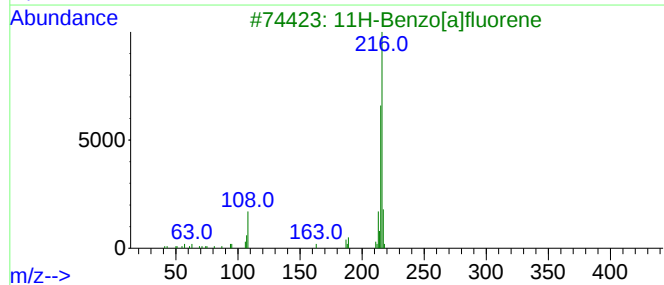
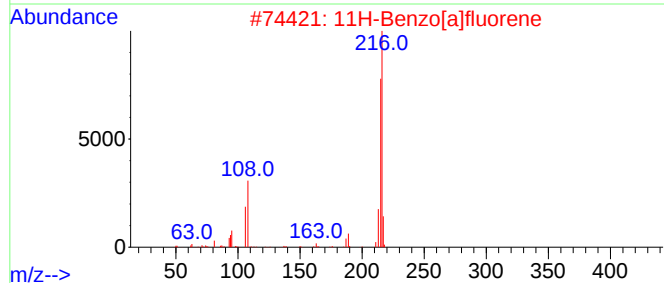
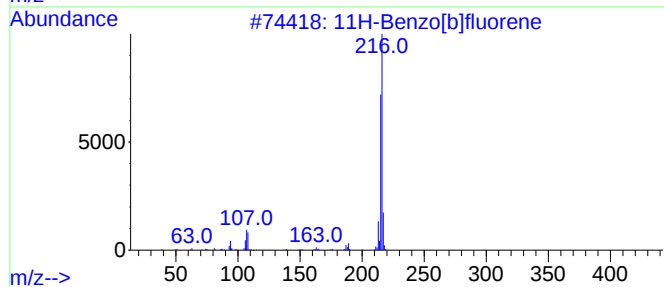
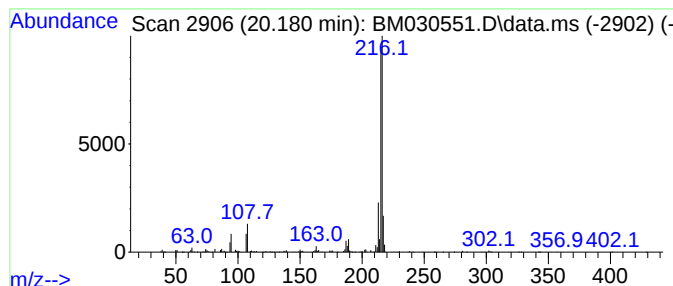
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 11H-Benzo[a]fluorene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.180	2.80 ng/ul	668316	Chrysene-d12	21.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	92
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062221\
Data File : BM030551.D
Acq On : 23 Jun 2021 13:49
Operator : CG/JU
Sample : M2662-11DL 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Anthracene, 2-m...	18.040	3.8	ng/ul	563692	4	17.174	2935480	20.0
Phenanthrene, 2...	18.090	4.7	ng/ul	690546	4	17.174	2935480	20.0
1H-Cyclopropa[l...	18.170	2.9	ng/ul	427171	4	17.174	2935480	20.0
4H-Cyclopenta[d...	18.240	7.5	ng/ul	1106450	4	17.174	2935480	20.0
Naphthalene, 2-...	18.540	2.4	ng/ul	357652	4	17.174	2935480	20.0
Phenanthrene, 2...	18.960	2.1	ng/ul	313561	4	17.174	2935480	20.0
11H-Benzo[b]flu...	20.080	3.3	ng/ul	781536	5	21.345	4768710	20.0
11H-Benzo[a]flu...	20.180	2.8	ng/ul	668316	5	21.345	4768710	20.0