

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
 Data File : BM034989.D
 Acq On : 11 May 2022 17:31
 Operator : CG/JU
 Sample : N2603-19
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4A0

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-BM051022.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034989.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.781	114	118	132	rBV	232826	372337	3.74%	0.284%
2	3.887	132	136	143	rVV	77073	105350	1.06%	0.080%
3	7.081	672	679	691	rBV	905673	1438540	14.44%	1.096%
4	7.251	701	708	717	rBV	704912	1091025	10.95%	0.831%
5	7.451	735	742	755	rVV	956446	1466611	14.72%	1.117%
6	7.922	814	822	831	rBV	776057	1233911	12.39%	0.940%
7	8.622	934	941	950	rBV	991571	1629757	16.36%	1.241%
8	9.075	1005	1018	1028	rBV	896622	1471697	14.77%	1.121%
9	9.798	1132	1141	1150	rBV	707314	1194964	11.99%	0.910%
10	10.340	1225	1233	1245	rBV	1082299	1746842	17.53%	1.331%
11	10.504	1256	1261	1269	rBV	69006	113869	1.14%	0.087%
12	10.722	1290	1298	1303	rBV	1042129	1782793	17.90%	1.358%
13	10.769	1303	1306	1313	rVV	213146	348432	3.50%	0.265%
14	10.851	1313	1320	1339	rVV	712154	1279812	12.85%	0.975%
15	12.381	1574	1580	1587	rVB	104428	160512	1.61%	0.122%
16	12.598	1612	1617	1623	rVB	75145	113405	1.14%	0.086%
17	13.392	1740	1752	1760	rBV2	85525	143871	1.44%	0.110%
18	13.963	1839	1849	1855	rBV	1844097	2524876	25.34%	1.923%
19	14.245	1887	1897	1910	rVV	2019820	3043829	30.55%	2.319%
20	14.551	1940	1949	1955	rBV	1436842	2115770	21.24%	1.612%
21	14.616	1955	1960	1967	rVB	525530	753511	7.56%	0.574%
22	14.739	1975	1981	1997	rBV	569137	943790	9.47%	0.719%
23	14.951	2011	2017	2024	rVB	324058	518743	5.21%	0.395%
24	15.410	2092	2095	2100	rVB	97724	121603	1.22%	0.093%
25	15.545	2110	2118	2123	rBV	2544314	3704997	37.19%	2.822%
26	15.598	2123	2127	2132	rVV	593662	800879	8.04%	0.610%
27	15.657	2132	2137	2142	rVV	648935	868344	8.72%	0.661%
28	15.763	2151	2155	2161	rBV2	62262	117165	1.18%	0.089%
29	15.892	2173	2177	2186	rVB	119791	208009	2.09%	0.158%
30	16.033	2197	2201	2209	rVV	130707	256375	2.57%	0.195%
31	16.245	2231	2237	2241	rBV	881975	1276519	12.81%	0.972%
32	16.298	2244	2246	2256	rVB2	94489	145425	1.46%	0.111%
33	16.604	2286	2298	2303	rBV3	87290	263764	2.65%	0.201%
34	16.945	2351	2356	2364	rVB	141412	226312	2.27%	0.172%
35	17.063	2372	2376	2380	rVV2	186208	295323	2.96%	0.225%
36	17.110	2380	2384	2395	rVB	289399	484191	4.86%	0.369%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
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37	17.292	2409	2415	2418	rBV	1440476	2043039	20.51%	1.556%
38	17.339	2418	2423	2427	rVV	5415253	7652663	76.82%	5.829%
39	17.392	2427	2432	2435	rVV	2638478	3759416	37.74%	2.864%
40	17.427	2435	2438	2444	rVB	1440289	1912816	19.20%	1.457%

41	17.486	2444	2448	2453	rVB2	85425	137133	1.38%	0.104%
42	17.692	2475	2483	2488	rBV	523187	847416	8.51%	0.645%
43	17.810	2498	2503	2508	rBV2	167896	255820	2.57%	0.195%
44	18.168	2560	2564	2566	rVV	376965	528487	5.30%	0.403%
45	18.198	2566	2569	2570	rVV	444440	561657	5.64%	0.428%

46	18.215	2570	2572	2577	rVV	597578	752876	7.56%	0.573%
47	18.268	2577	2581	2583	rVV	106512	158955	1.60%	0.121%
48	18.298	2583	2586	2591	rVV	199113	276615	2.78%	0.211%
49	18.368	2591	2598	2601	rVV3	778521	1308934	13.14%	0.997%
50	18.398	2601	2603	2609	rVB	246035	298713	3.00%	0.228%

51	18.498	2617	2620	2626	rVB2	89474	150920	1.51%	0.115%
52	18.668	2645	2649	2652	rVV	351187	450880	4.53%	0.343%
53	18.704	2652	2655	2661	rVB	240767	310288	3.11%	0.236%
54	18.962	2696	2699	2703	rVV	79076	108070	1.08%	0.082%
55	19.086	2715	2720	2724	rBV2	140664	217675	2.18%	0.166%

56	19.127	2724	2727	2729	rBV	93352	112280	1.13%	0.086%
57	19.221	2739	2743	2751	rBV2	128693	261870	2.63%	0.199%
58	19.351	2757	2765	2772	rVB	7574690	9962404	100.00%	7.589%
59	19.468	2782	2785	2793	rVB2	216909	302612	3.04%	0.231%
60	19.592	2802	2806	2810	rVB	143187	183806	1.84%	0.140%

61	19.680	2815	2821	2824	rBV2	3929955	6239571	62.63%	4.753%
62	19.709	2824	2826	2834	rVV	5733833	6927084	69.53%	5.276%
63	19.780	2834	2838	2843	rVB	185627	266260	2.67%	0.203%
64	19.886	2852	2856	2862	rBV	321727	442495	4.44%	0.337%
65	19.945	2863	2866	2872	rVB	200975	283360	2.84%	0.216%

66	20.033	2875	2881	2887	rBV3	254590	686719	6.89%	0.523%
67	20.086	2887	2890	2894	rVB	118987	144339	1.45%	0.110%
68	20.168	2900	2904	2906	rBV3	164427	245410	2.46%	0.187%
69	20.204	2906	2910	2914	rVV	828289	1044092	10.48%	0.795%
70	20.304	2923	2927	2931	rBV	703062	856881	8.60%	0.653%

71	20.362	2931	2937	2941	rVV2	347501	585176	5.87%	0.446%
72	20.404	2941	2944	2947	rVB	138633	166343	1.67%	0.127%
73	20.439	2948	2950	2956	rVB2	83308	117229	1.18%	0.089%
74	20.504	2957	2961	2964	rBV	173683	226304	2.27%	0.172%
75	20.545	2965	2968	2972	rVV2	187697	258229	2.59%	0.197%

76	20.609	2976	2979	2983	rVB2	231706	242899	2.44%	0.185%
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77	20.909	3027	3030	3032	rBV2	86518	113668	1.14%	0.087%
78	20.951	3033	3037	3043	rVB	255810	415262	4.17%	0.316%
79	21.115	3060	3065	3069	rBV2	419295	620045	6.22%	0.472%
80	21.156	3069	3072	3075	rVV	454535	545211	5.47%	0.415%
81	21.192	3075	3078	3083	rVV2	877378	1334084	13.39%	1.016%
82	21.245	3083	3087	3095	rVB2	358095	569780	5.72%	0.434%
83	21.456	3115	3123	3129	rBV2	4318659	7899269	79.29%	6.017%
84	21.509	3129	3132	3141	rVB	2816473	3640804	36.55%	2.773%
85	21.633	3148	3153	3158	rBV2	622902	903362	9.07%	0.688%
86	21.739	3168	3171	3175	rVB2	246945	326414	3.28%	0.249%
87	21.792	3176	3180	3186	rVV2	443317	664881	6.67%	0.506%
88	22.045	3220	3223	3227	rVB	318249	430802	4.32%	0.328%
89	22.098	3229	3232	3238	rVB	305981	441714	4.43%	0.336%
90	22.256	3255	3259	3261	rBV2	219024	301353	3.02%	0.230%
91	22.292	3262	3265	3270	rVB4	319814	451809	4.54%	0.344%
92	23.139	3402	3409	3414	rBV	2449501	6066118	60.89%	4.621%
93	23.315	3435	3439	3450	rVB	434636	782036	7.85%	0.596%
94	23.650	3490	3496	3500	rBV	1260634	2444001	24.53%	1.862%
95	23.703	3500	3505	3509	rVV2	1937936	3872697	38.87%	2.950%
96	23.756	3509	3514	3522	rVB	2219475	4346862	43.63%	3.311%
97	23.856	3525	3531	3536	rVV2	1016293	2174510	21.83%	1.656%
98	23.909	3536	3540	3550	rVB	601308	1254651	12.59%	0.956%
99	26.327	3942	3951	3963	rVB2	1176241	3695046	37.09%	2.815%
100	27.085	4072	4080	4100	rVB	650216	2337544	23.46%	1.781%

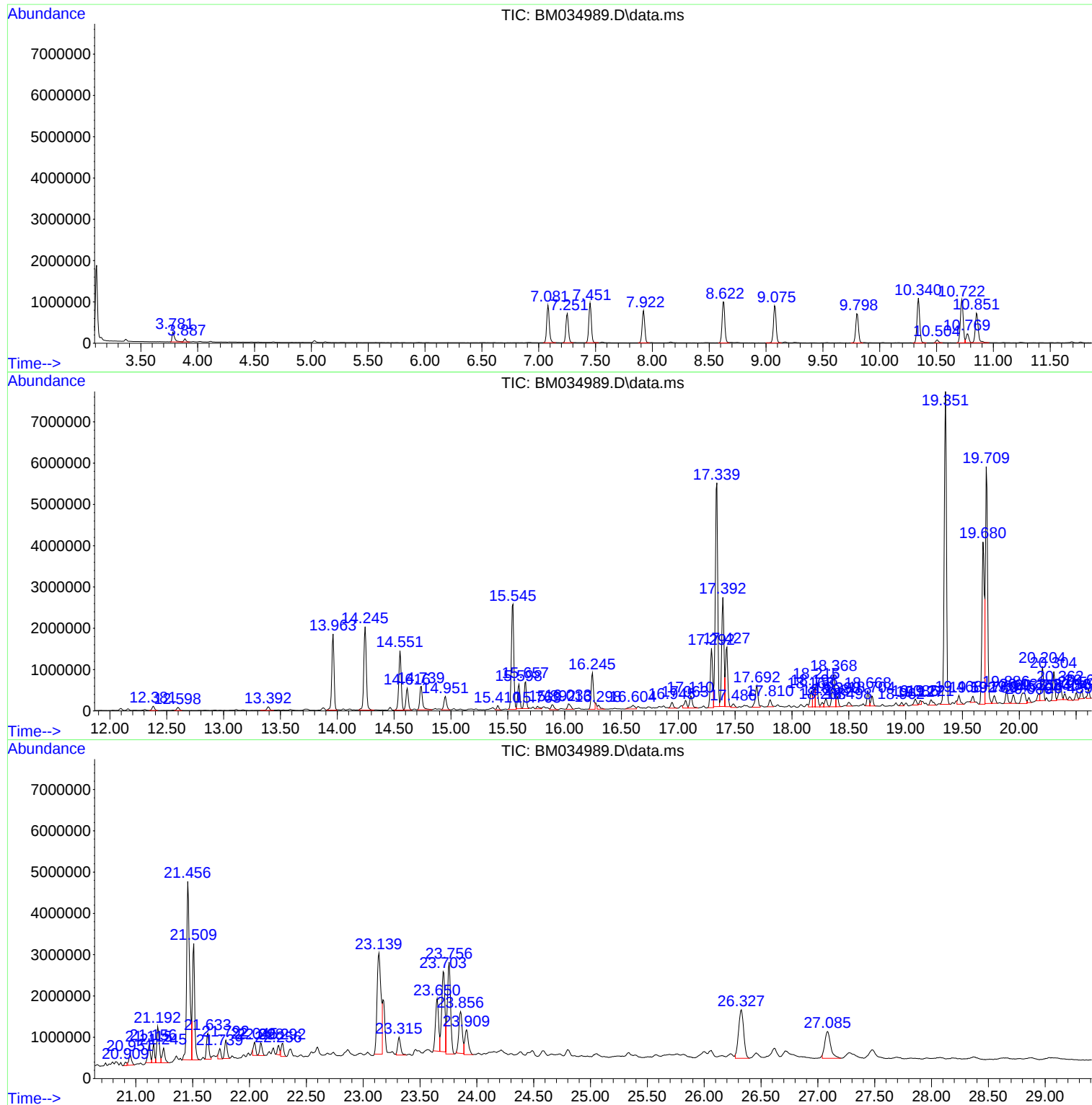
Sum of corrected areas: 131282812

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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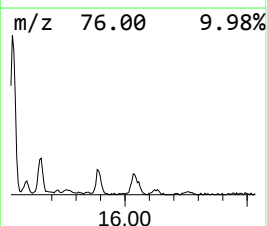
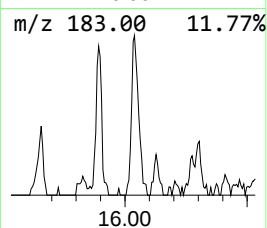
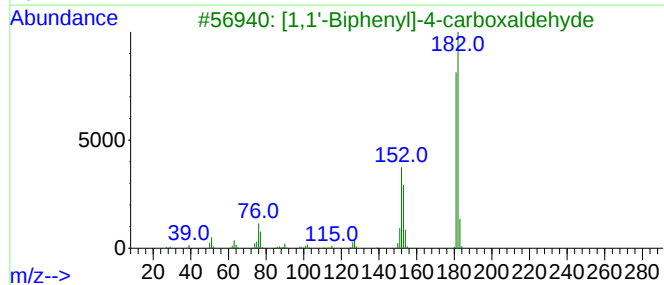
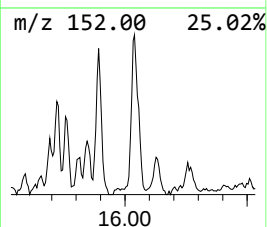
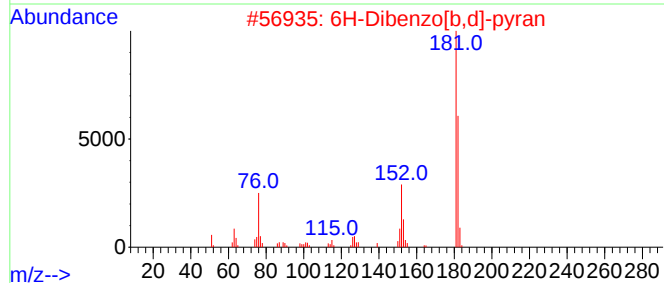
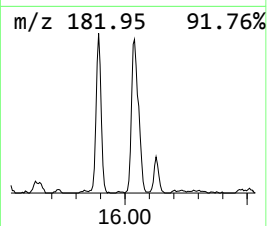
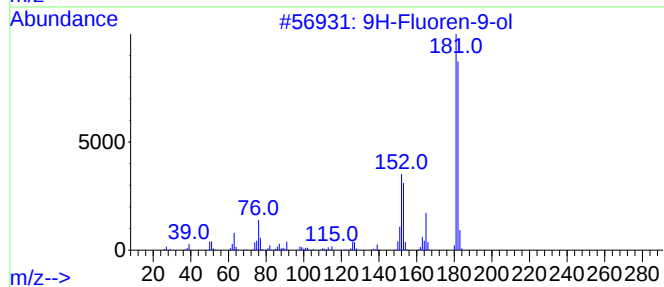
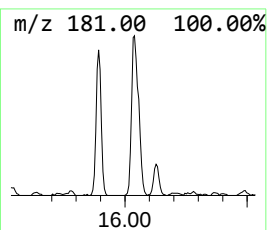
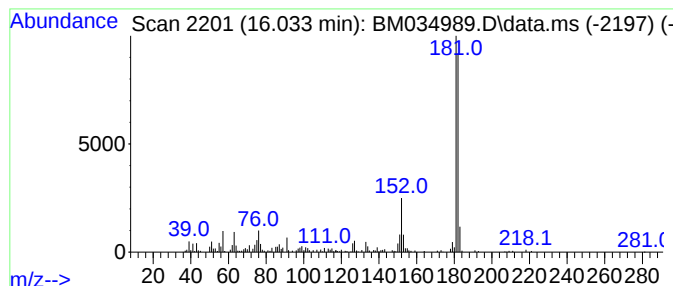
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 9H-Fluoren-9-ol Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.033	2.51 ng/ul	256375	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	93
2			6H-Dibenzo[b,d]-pyran	182	C13H10O	000229-95-8	90
3			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
4			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
5			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86



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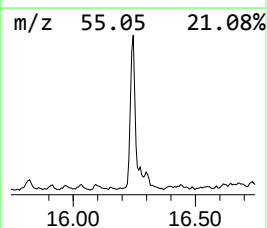
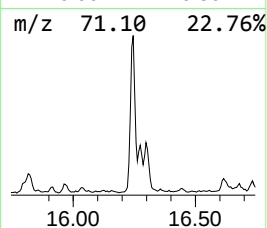
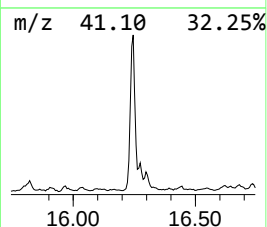
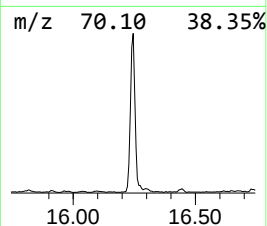
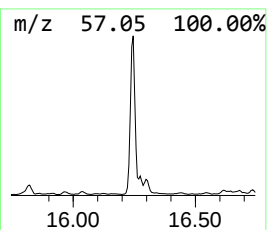
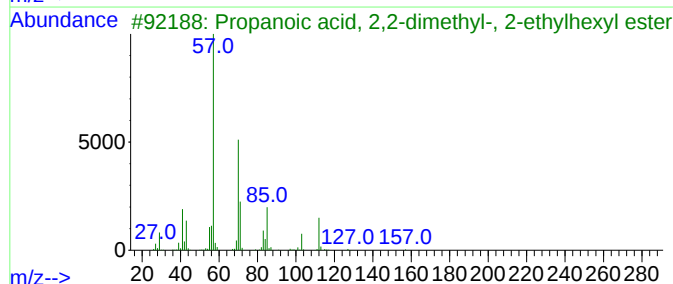
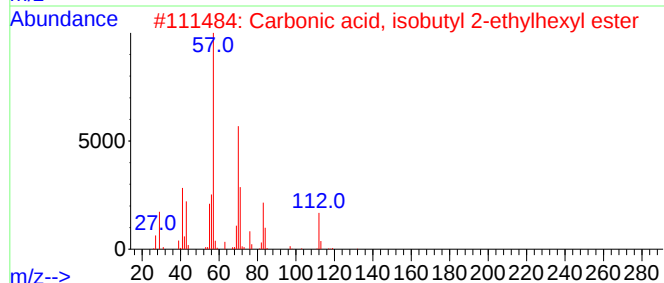
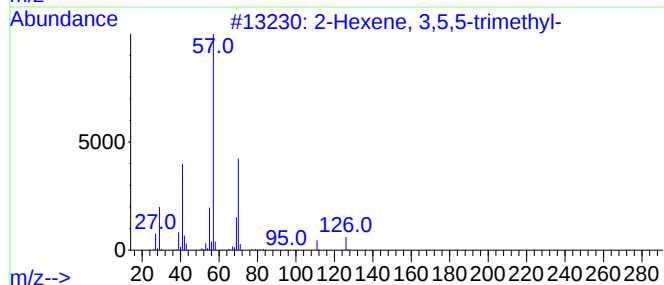
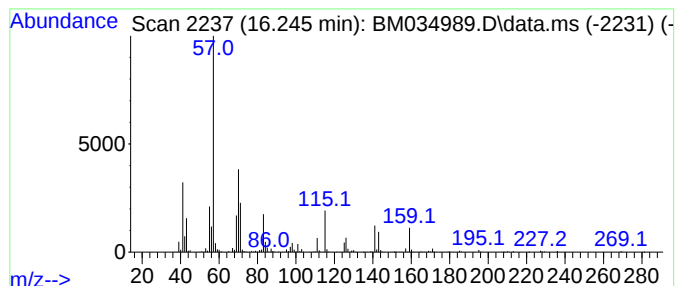
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.245	12.50 ng/ul	1276520	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexene, 3,5,5-trimethyl-		126	C9H18		026456-76-8	43
2	Carbonic acid, isobutyl 2-ethylh...		230	C13H26O3		1000383-13-3	37
3	Propanoic acid, 2,2-dimethyl-, 2...		214	C13H26O2		016387-18-1	37
4	2-Propyl-1-pentanol		130	C8H18O		058175-57-8	35
5	3,5,5-Trimethylhexyl propionate		200	C12H24O2		1000429-26-0	35



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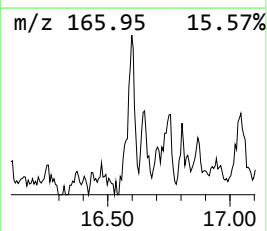
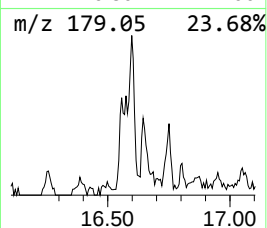
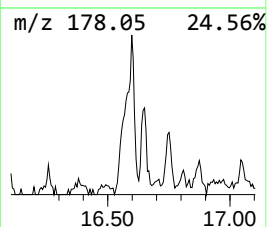
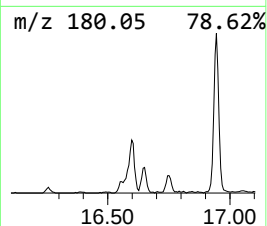
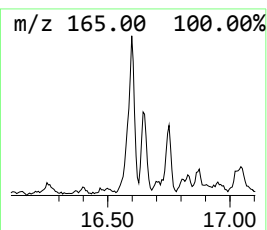
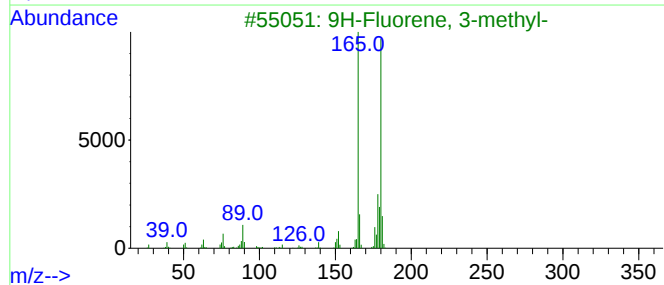
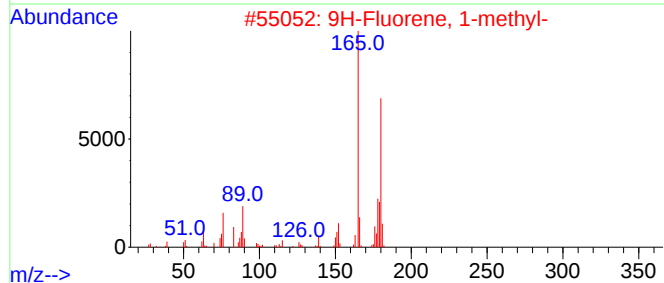
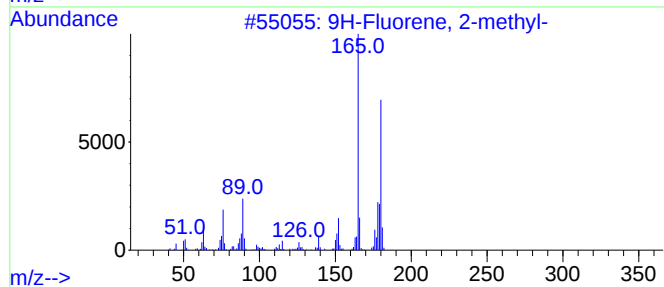
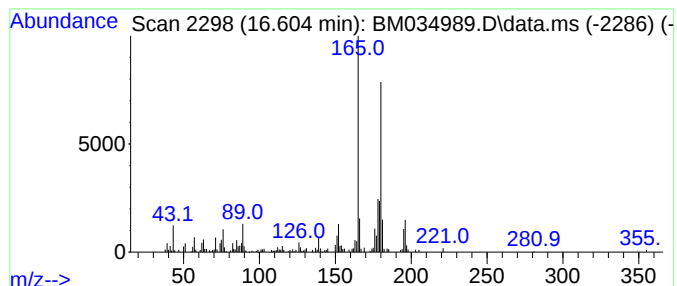
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 9H-Fluorene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.604	2.58 ng/ul	263764	Phenanthrene-d10	17.292

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	96
2		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	95
3		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	94
4		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	90
5		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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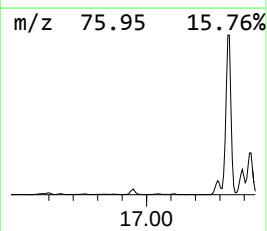
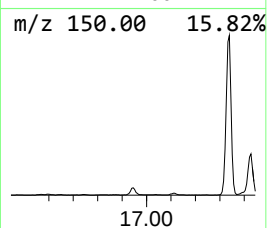
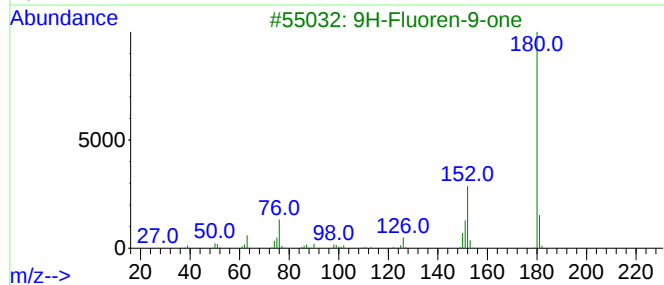
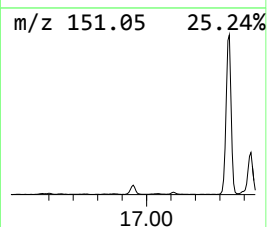
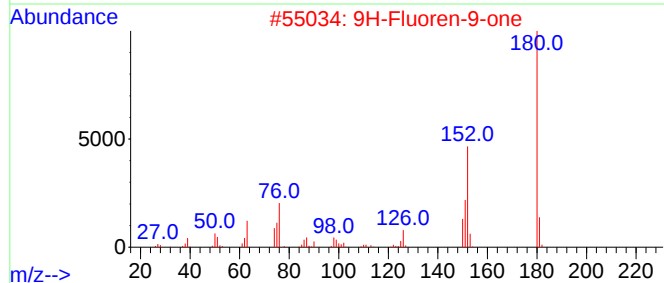
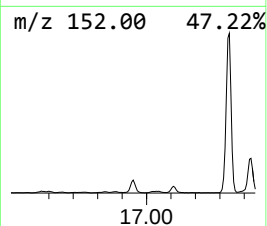
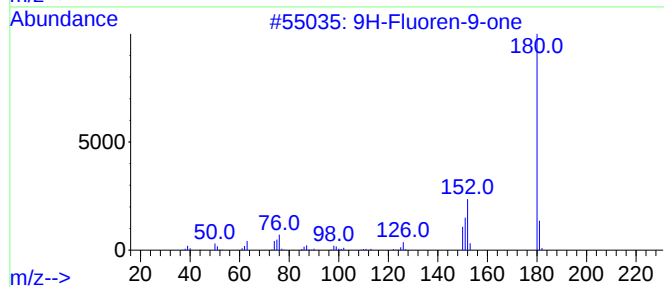
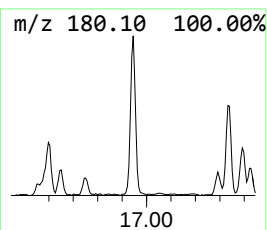
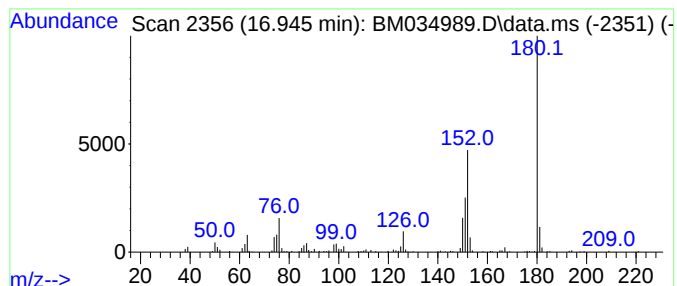
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9H-Fluoren-9-one Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.945	2.22 ng/ul	226312	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one			180	C13H8O	000486-25-9	97
2	9H-Fluoren-9-one			180	C13H8O	000486-25-9	94
3	9H-Fluoren-9-one			180	C13H8O	000486-25-9	94
4	9H-Fluoren-9-one			180	C13H8O	000486-25-9	94
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
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ClientSampleId :
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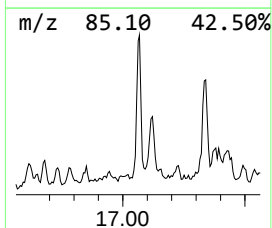
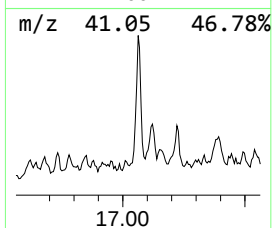
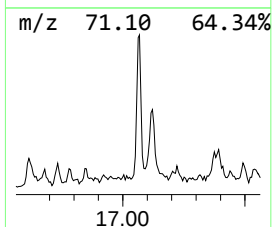
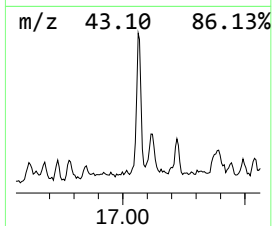
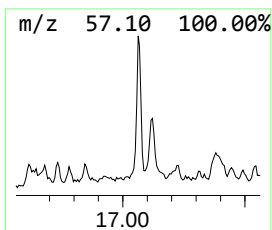
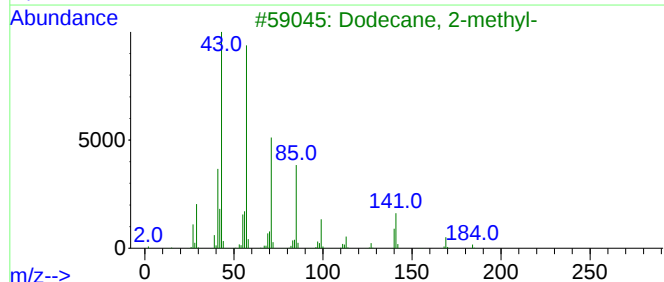
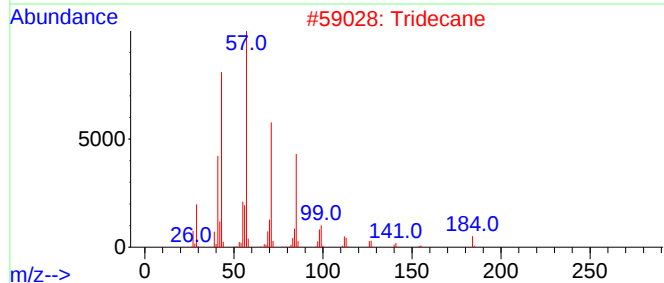
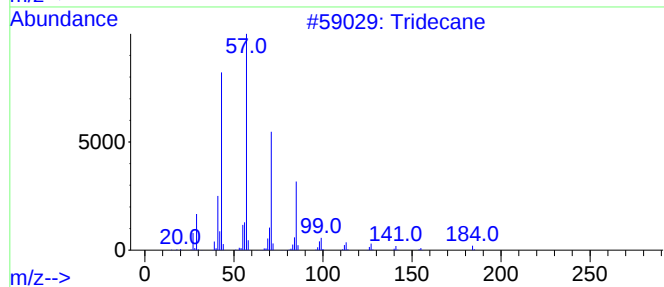
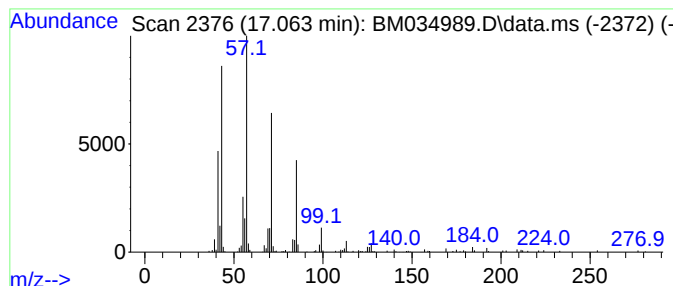
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.063	2.89 ng/ul	295323	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Tridecane	184	C13H28	000629-50-5	90
3			Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
4			Heptacosane	380	C27H56	000593-49-7	90
5			Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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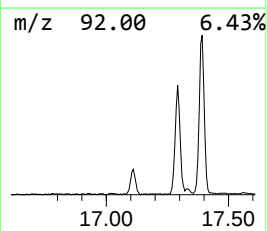
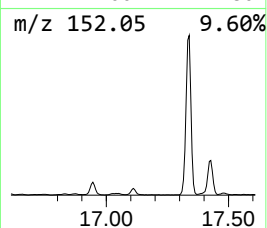
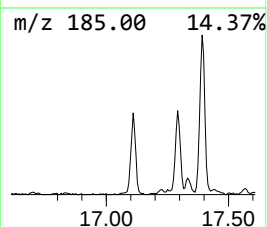
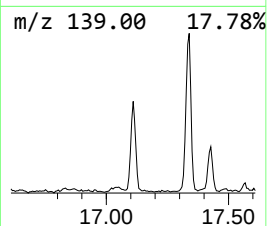
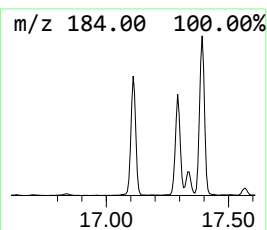
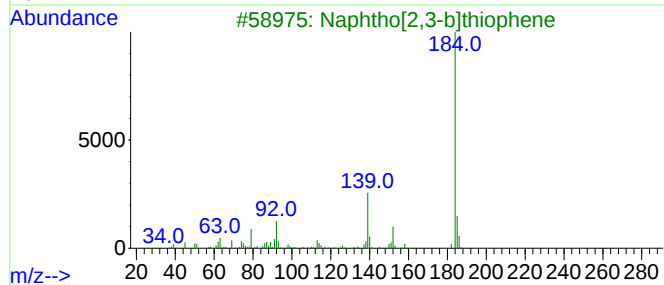
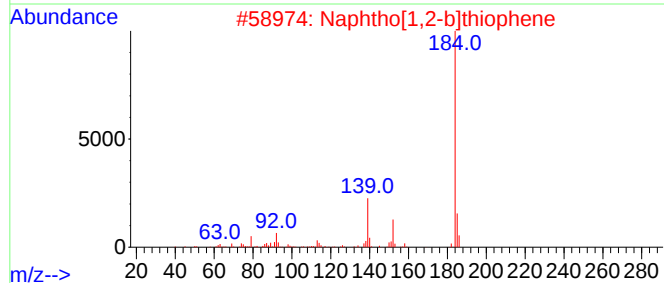
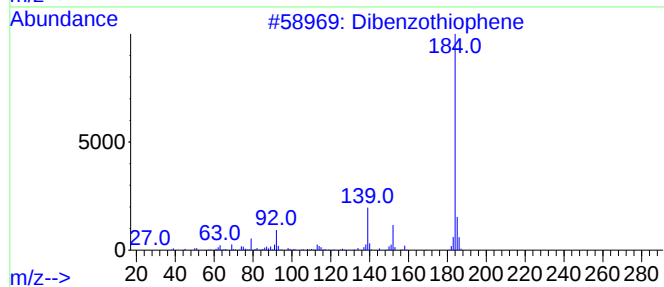
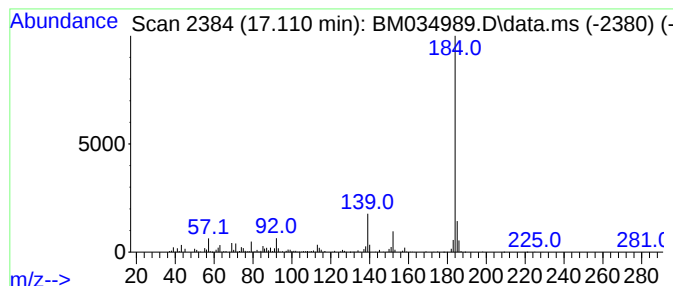
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Dibenzothiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.110	4.74 ng/ul	484191	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	97
3			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	97
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	97
5			Dibenzothiophene	184	C12H8S	000132-65-0	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
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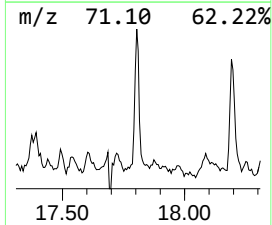
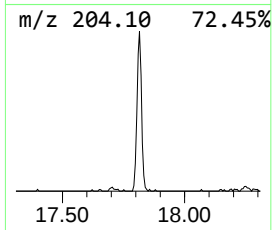
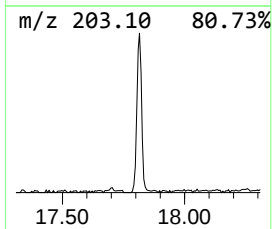
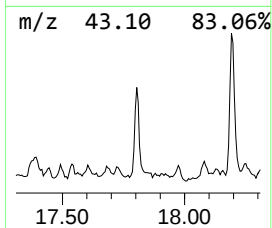
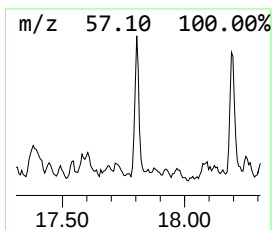
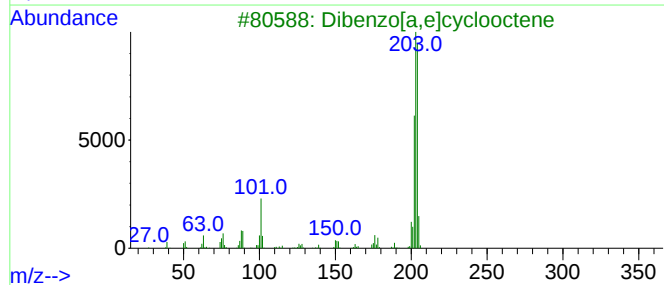
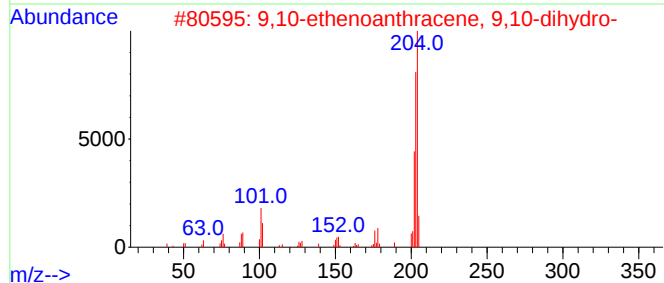
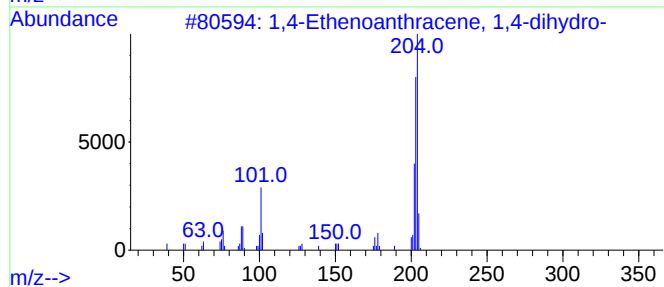
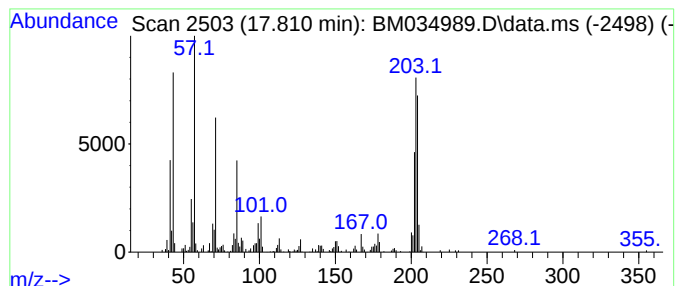
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1,4-Ethenoanthracene, 1,4-d... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.810	2.50 ng/ul	255820	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Ethenoanthracene, 1,4-dihydro-	204	C16H12	027765-96-4	64
2			9,10-ethenoanthracene, 9,10-dihy...	204	C16H12	1000400-47-8	64
3			Dibenzo[a,e]cyclooctene	204	C16H12	000262-89-5	62
4			Pyrene, 4,5-dihydro-	204	C16H12	006628-98-4	62
5			Nonadecane	268	C19H40	000629-92-5	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
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Acq On : 11 May 2022 17:31
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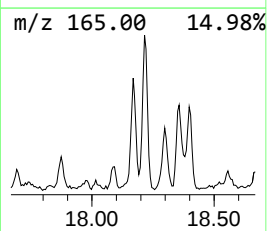
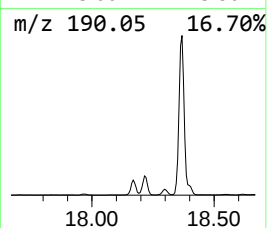
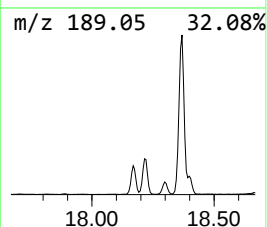
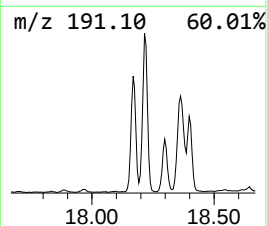
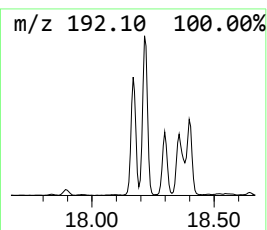
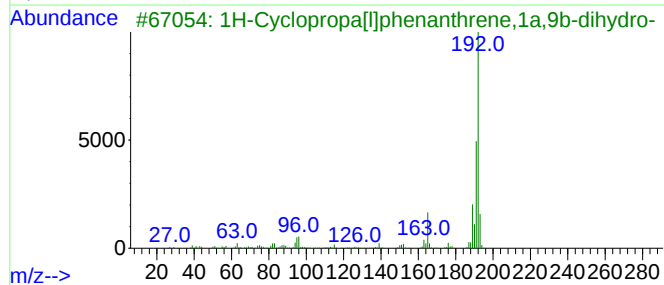
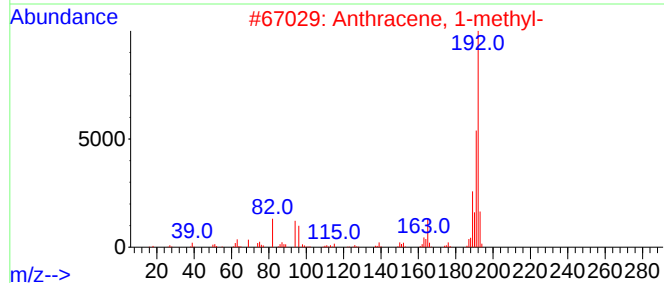
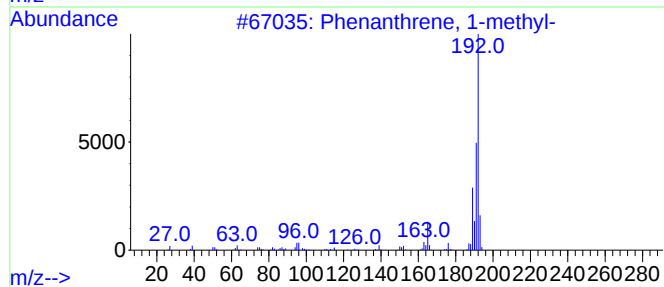
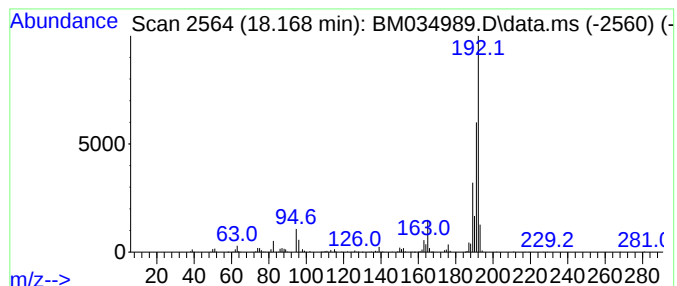
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.168	5.17 ng/ul	528487	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	91
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
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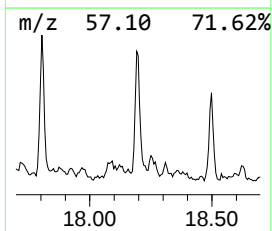
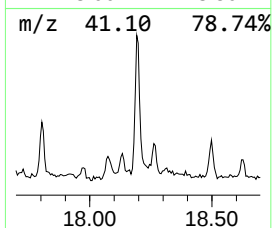
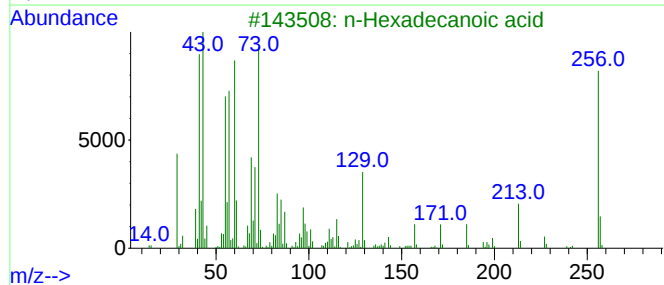
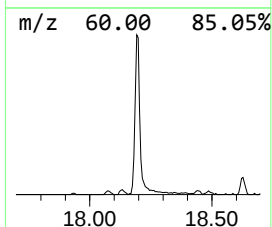
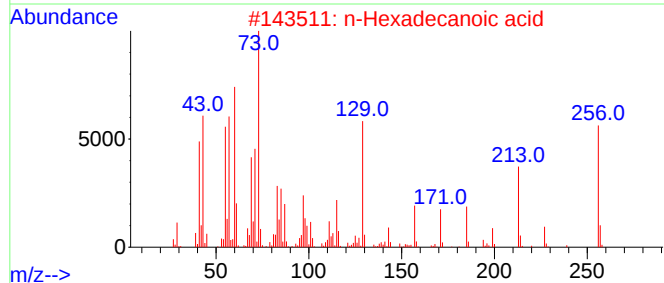
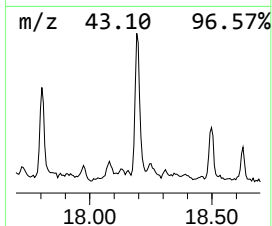
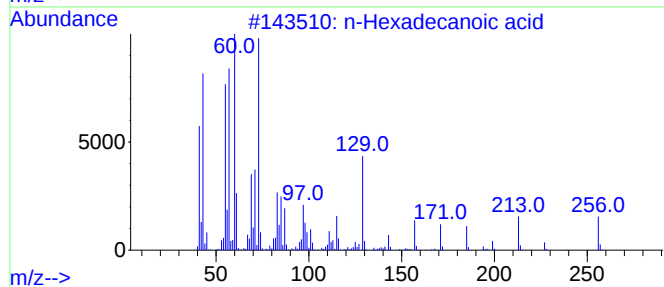
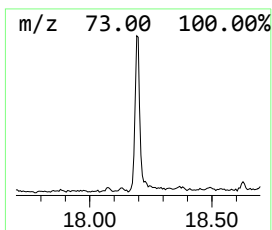
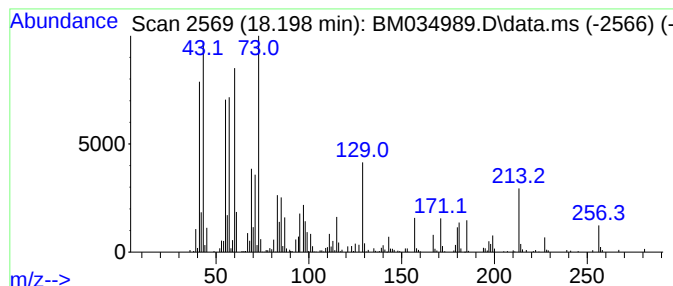
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.198	5.50 ng/ul	561657	Phenanthrene-d10	17.292

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tridecanoic acid	214	C13H26O2	000638-53-9	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
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ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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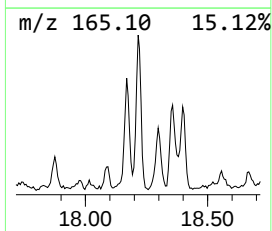
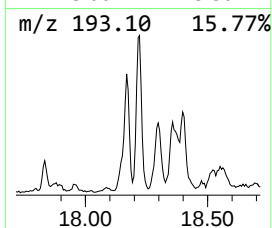
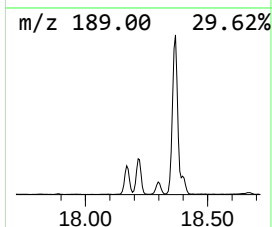
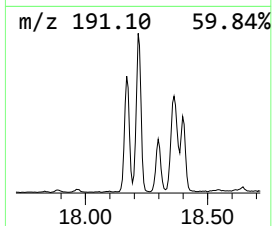
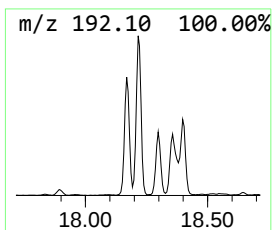
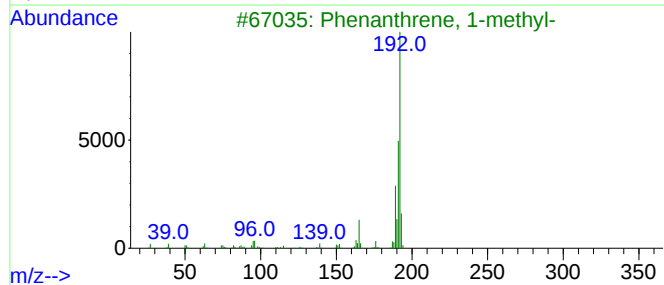
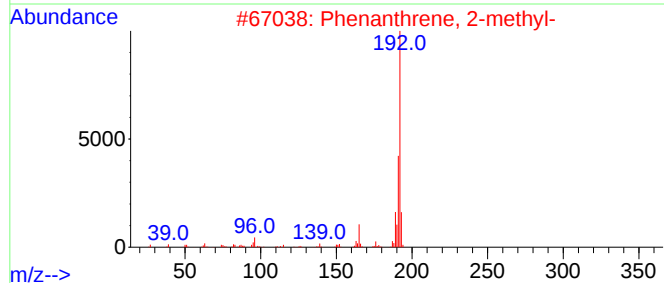
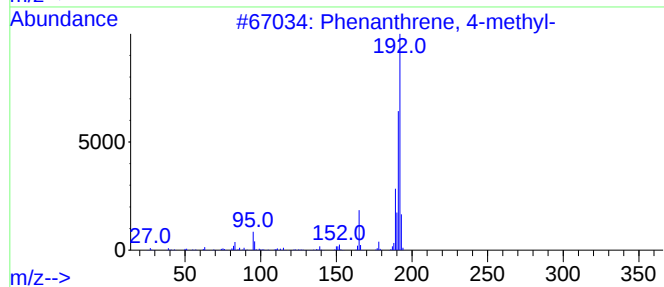
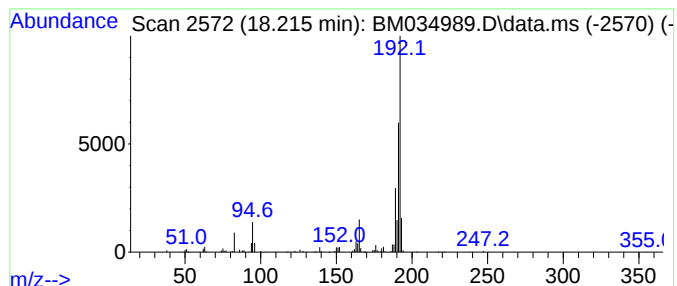
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Phenanthrene, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.215	7.37 ng/ul	752876	Phenanthrene-d10	17.292

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
Misc :
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Instrument :
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ClientSampleId :
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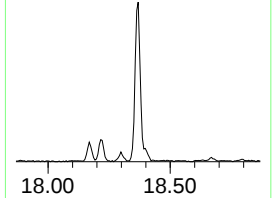
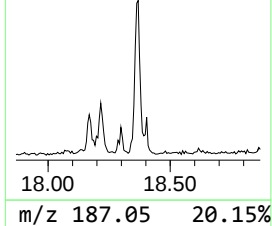
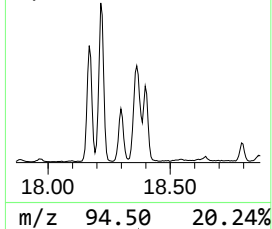
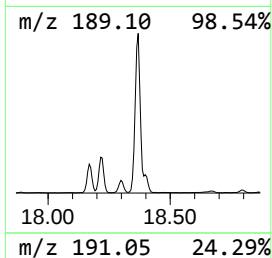
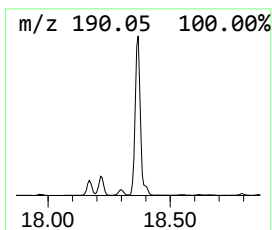
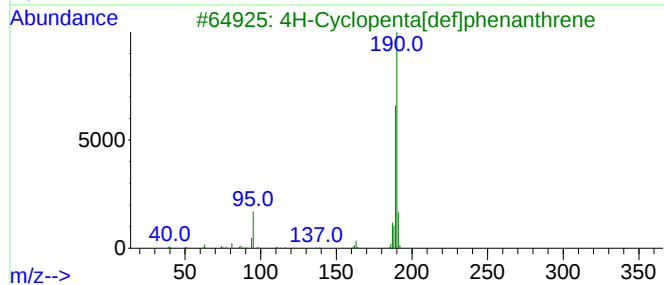
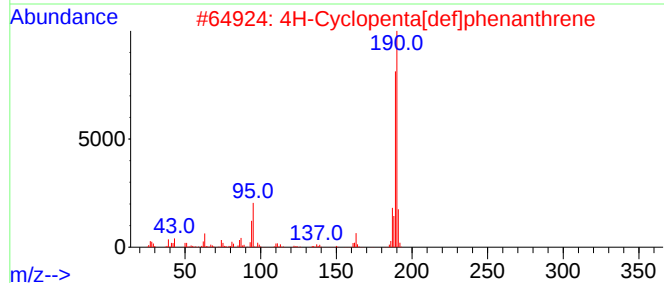
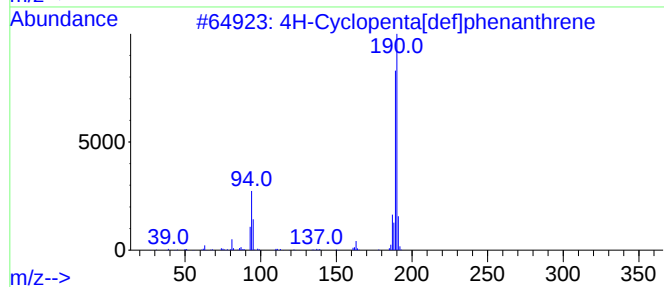
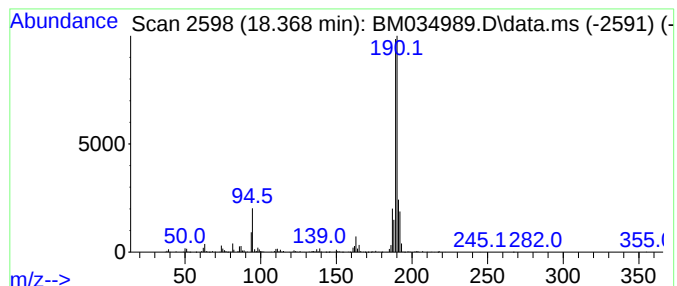
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.368	12.81 ng/ul	1308930	Phenanthrene-d10	17.292

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	90
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	80
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
5			Methyl diselenide	190	C2H6Se2	007101-31-7	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
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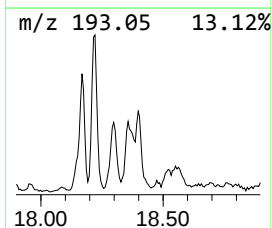
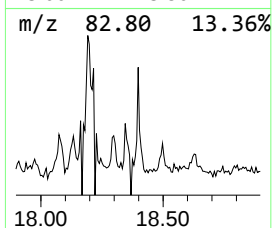
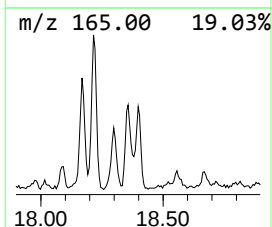
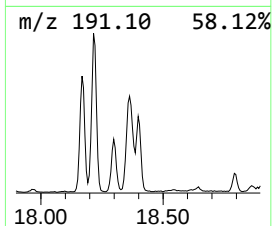
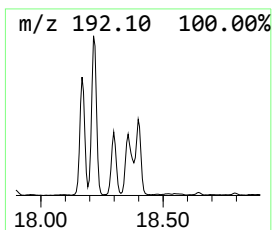
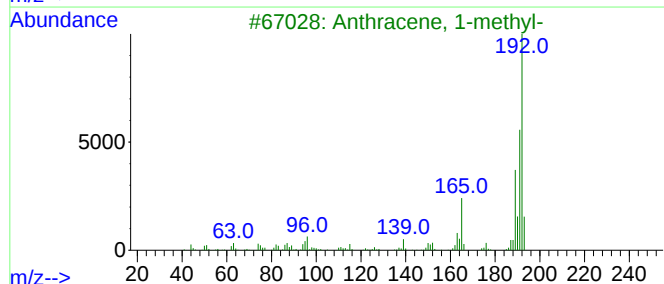
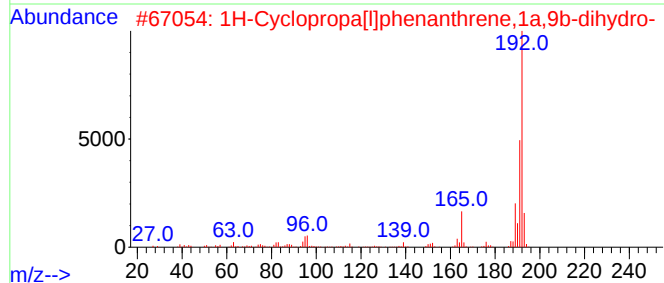
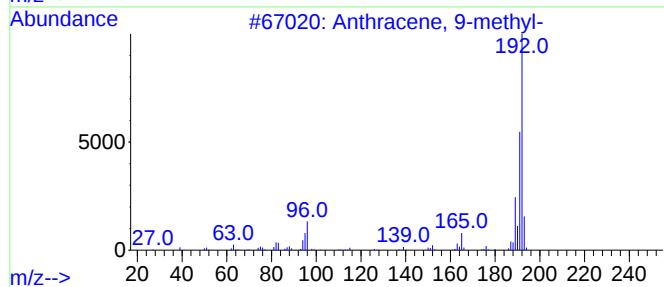
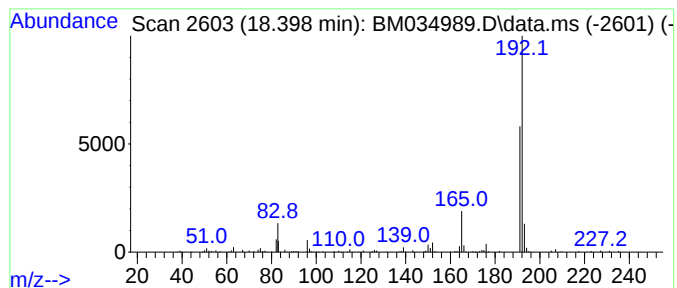
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Anthracene, 9-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	2.92 ng/ul	298713	Phenanthrene-d10	17.292

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
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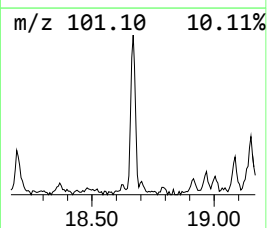
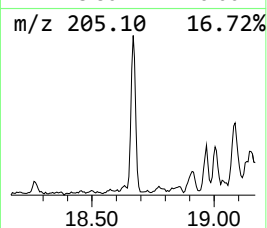
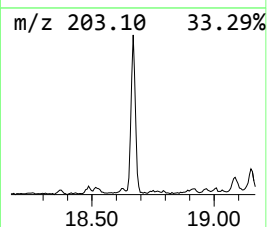
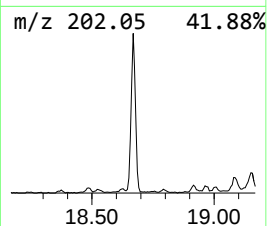
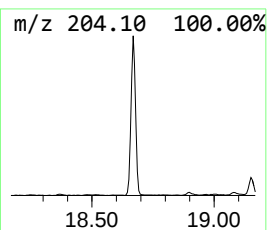
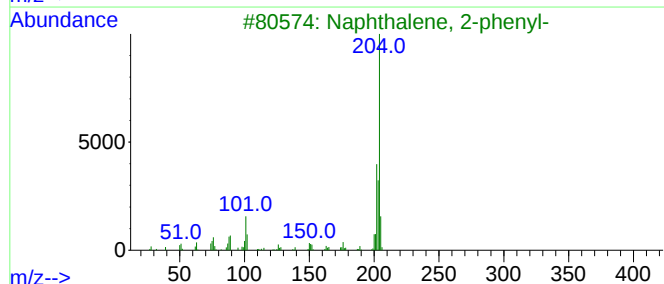
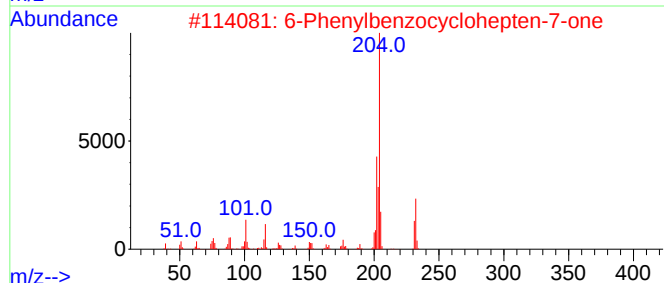
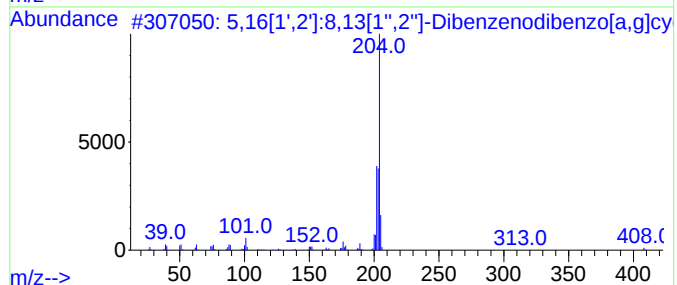
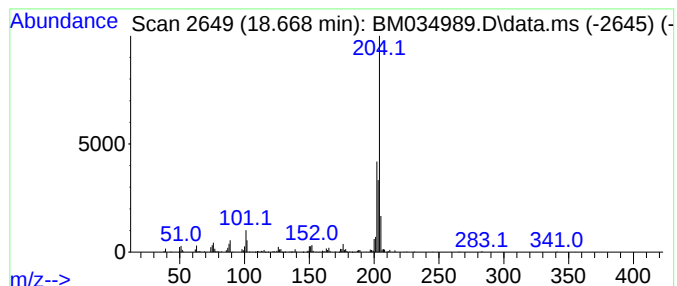
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.668	4.41 ng/ul	450880	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91		
2	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	90		
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86		
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83		
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
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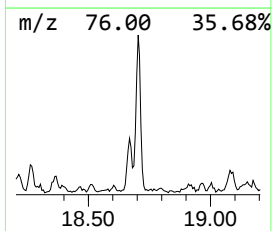
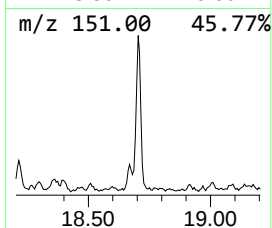
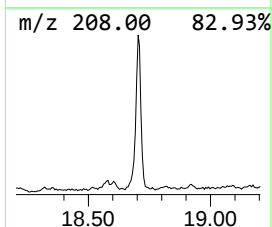
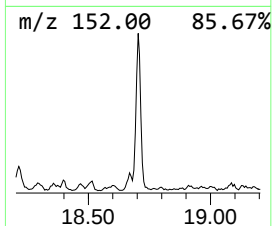
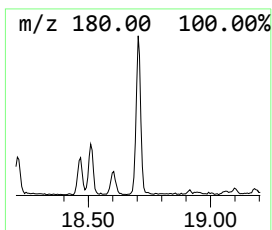
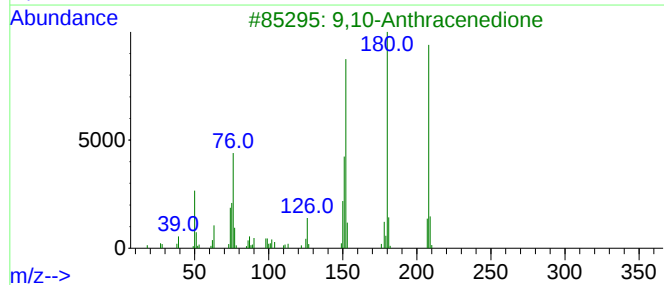
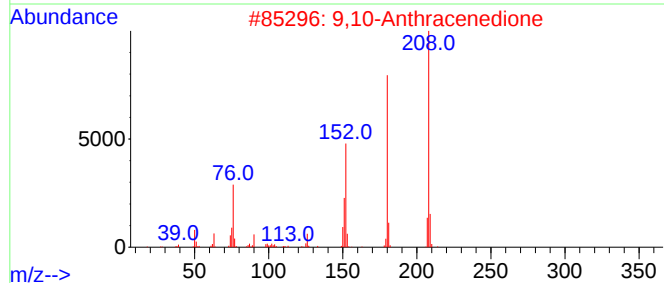
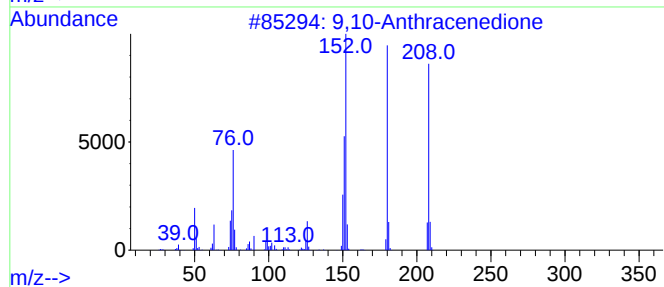
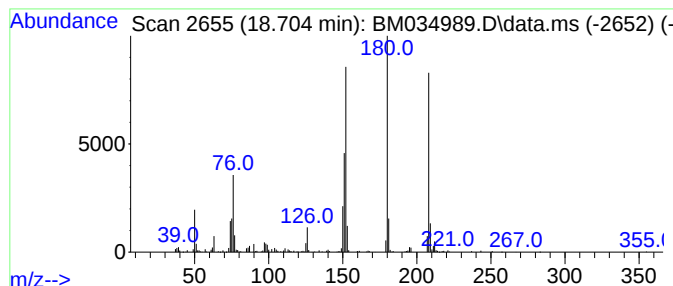
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.704	3.04 ng/ul	310288	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
5			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
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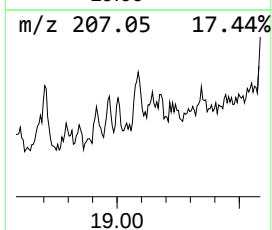
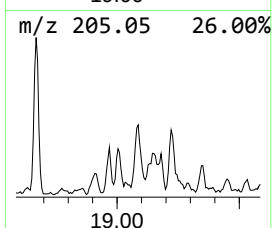
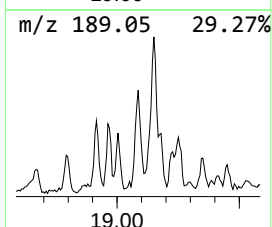
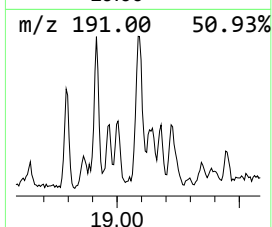
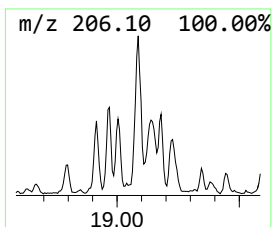
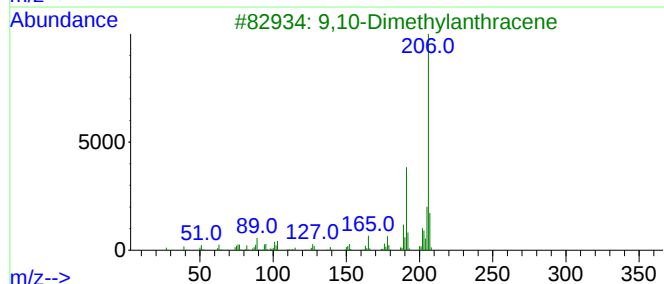
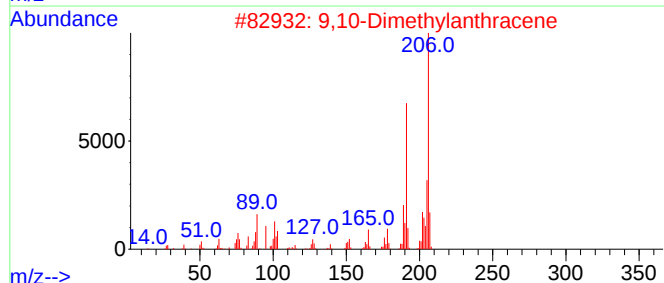
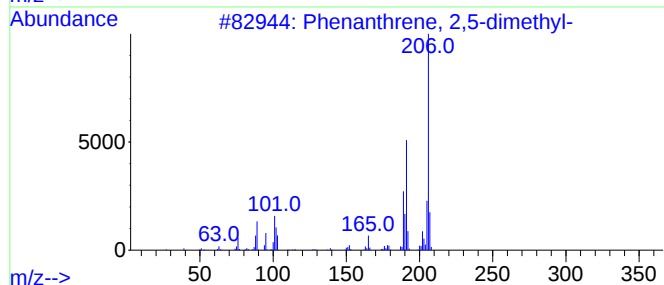
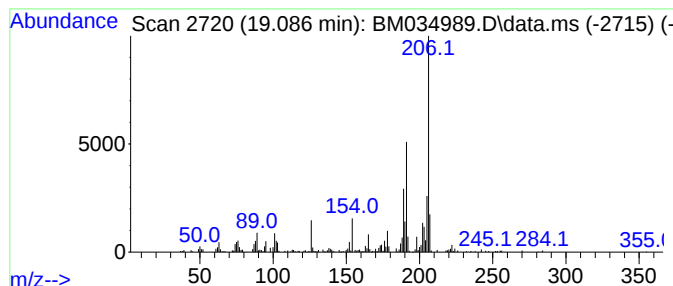
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Phenanthrene, 2,5-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.086	2.13 ng/ul	217675	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	90
3			9,10-Dimethylantracene	206	C16H14	000781-43-1	89
4			9,10-Dimethylantracene	206	C16H14	000781-43-1	89
5			di-p-Tolylacetylene	206	C16H14	002789-88-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
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Acq On : 11 May 2022 17:31
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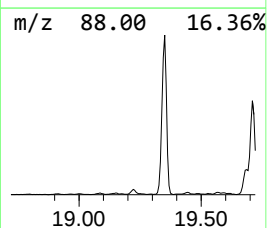
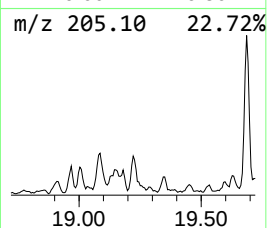
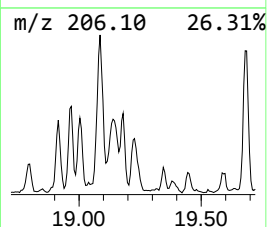
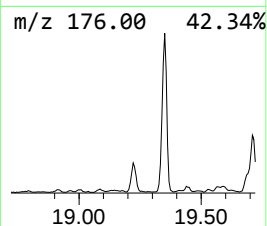
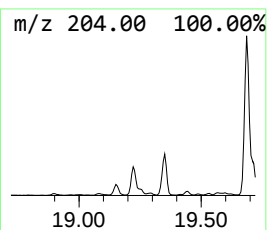
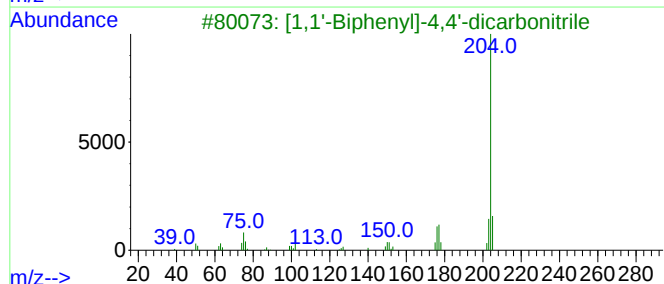
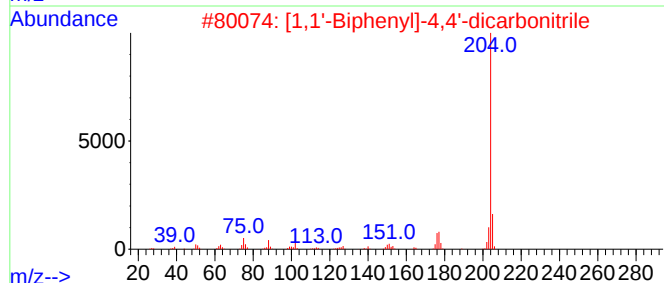
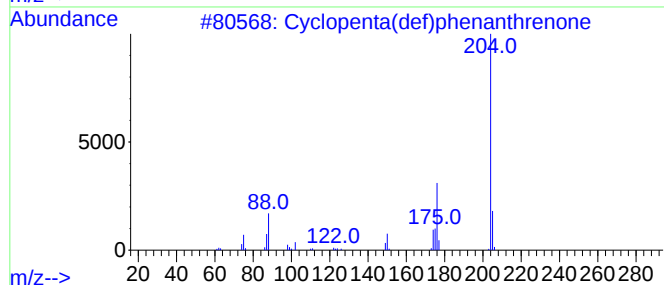
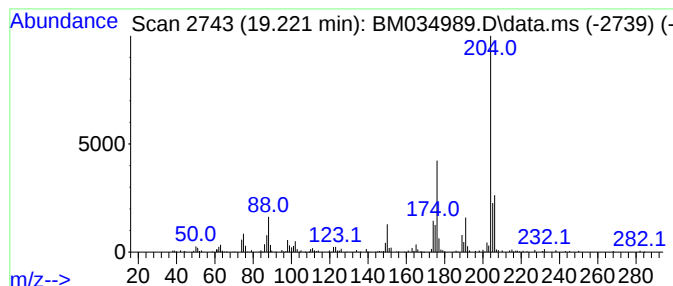
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Cyclopenta(def)phenanthrenone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.221	2.56 ng/ul	261870	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	94
2			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	50
3			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	49
4			[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	46
5			L-Rhamnose, (R,R,S,S)-, 4TMS der...	452	C18H44O5Si4	108392-01-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
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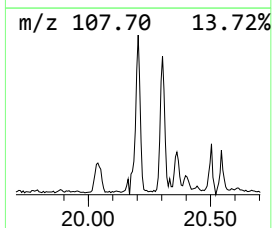
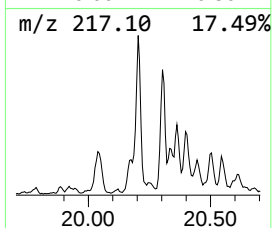
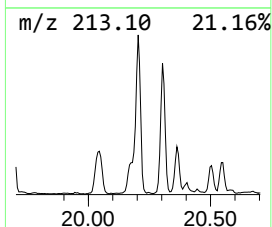
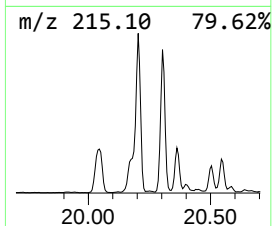
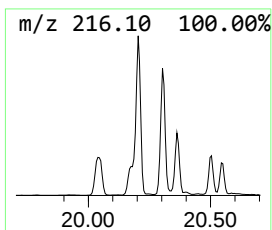
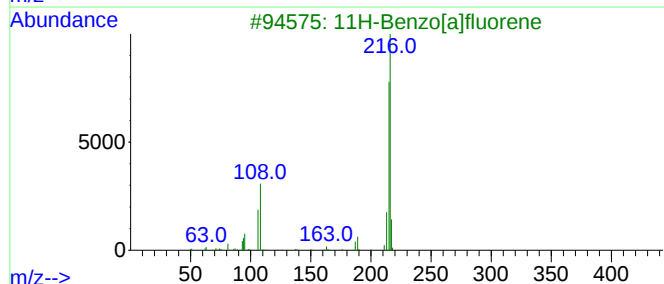
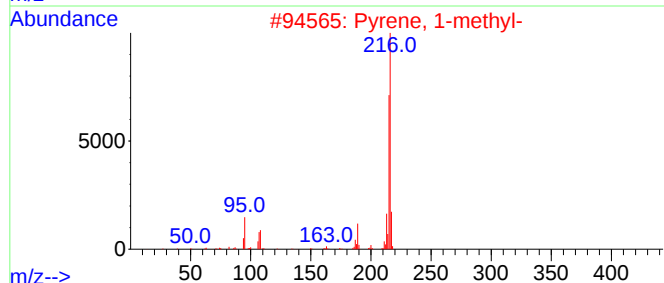
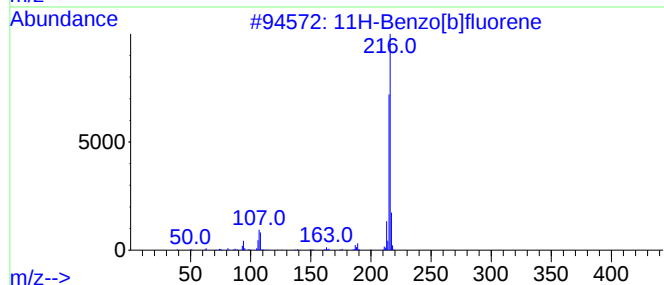
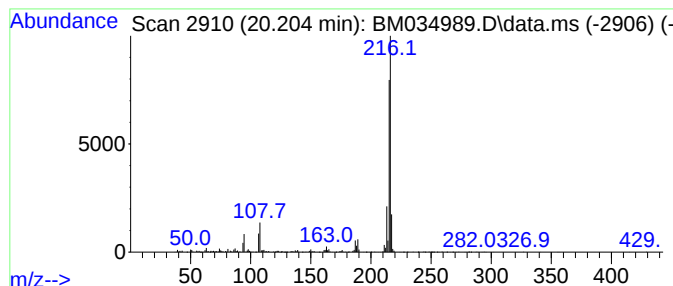
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 11H-Benzo[b]fluorene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.204	2.64 ng/ul	1044090	Chrysene-d12	21.474

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	97
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
5			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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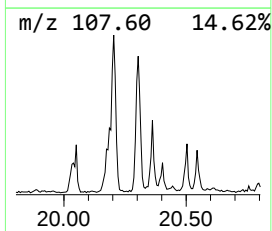
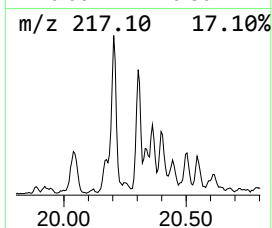
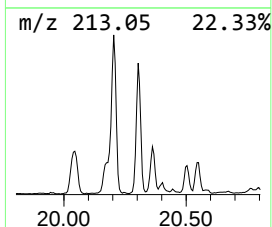
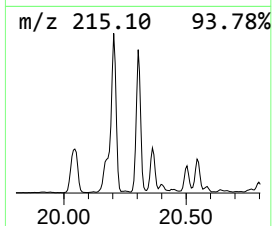
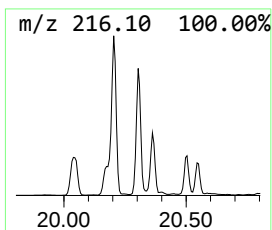
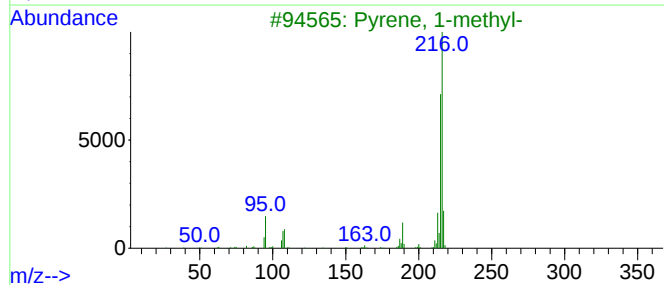
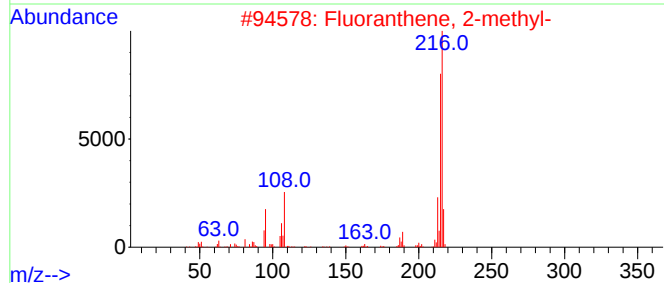
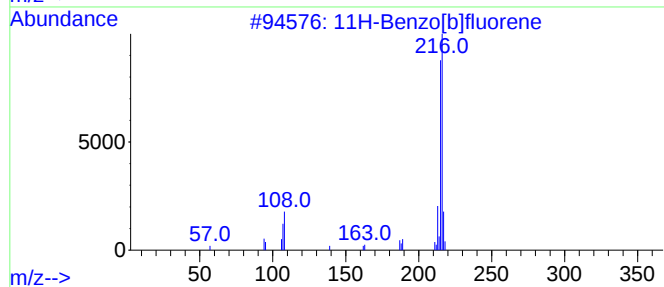
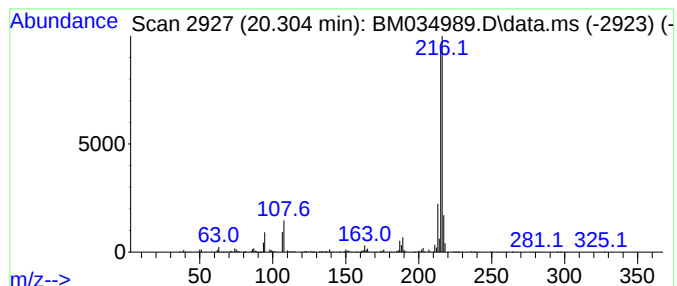
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Fluoranthene, 2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.304	2.17 ng/ul	856881	Chrysene-d12	21.474

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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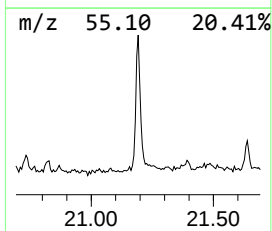
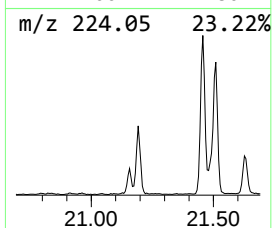
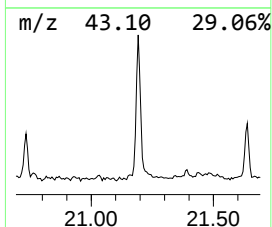
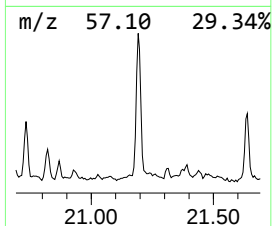
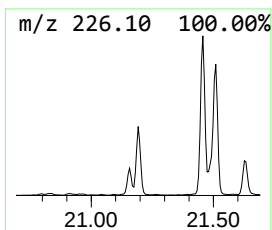
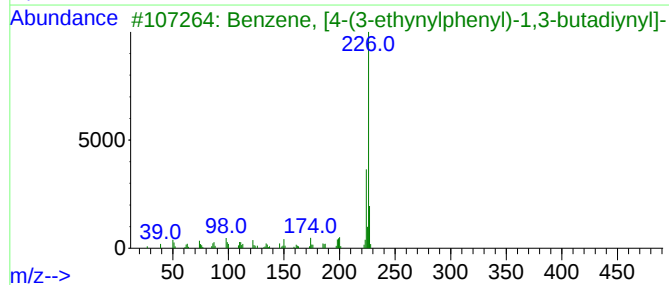
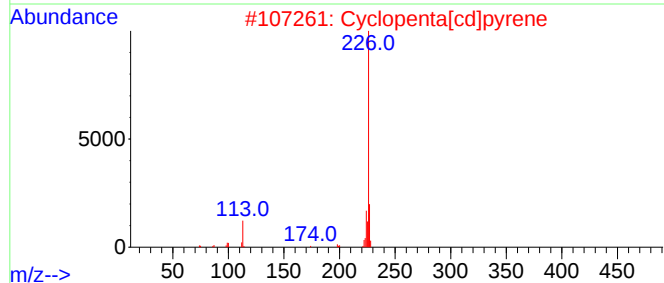
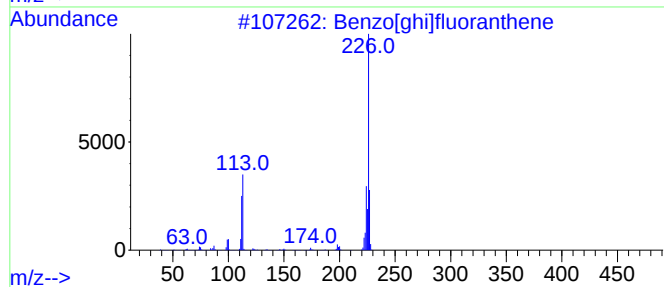
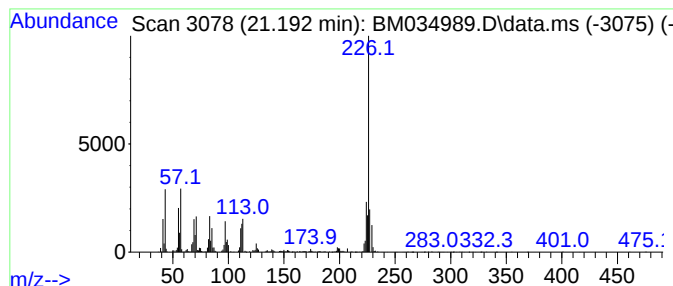
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzo[ghi]fluoranthene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.192	3.38 ng/ul	1334080	Chrysene-d12	21.474

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	76
2			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	64
3			Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	53
4			benzenamine, 4-(2-benzothiazolyl)-	226	C13H10N2S	1000400-93-4	50
5			.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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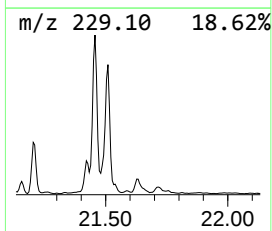
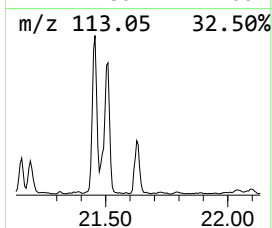
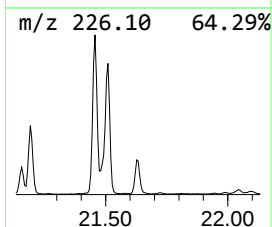
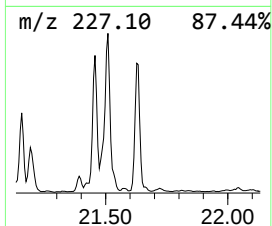
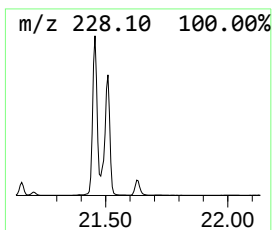
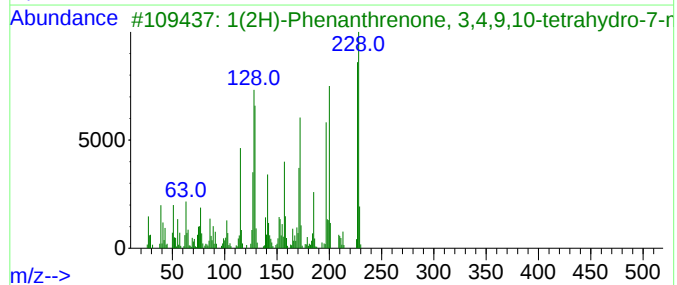
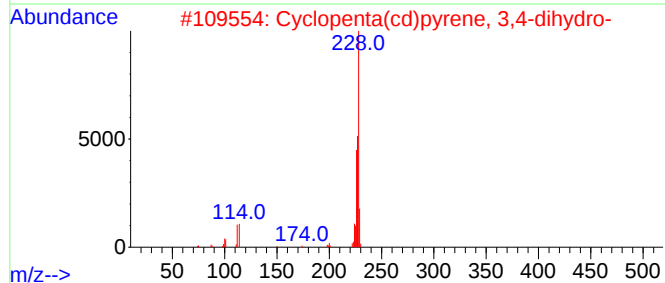
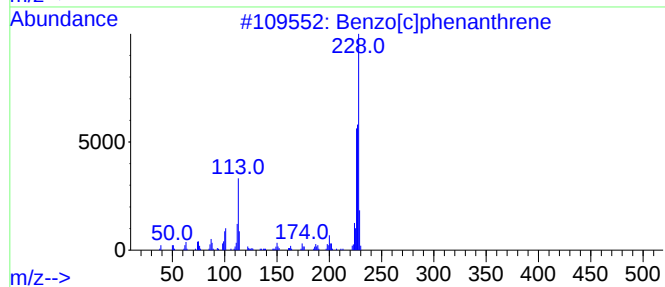
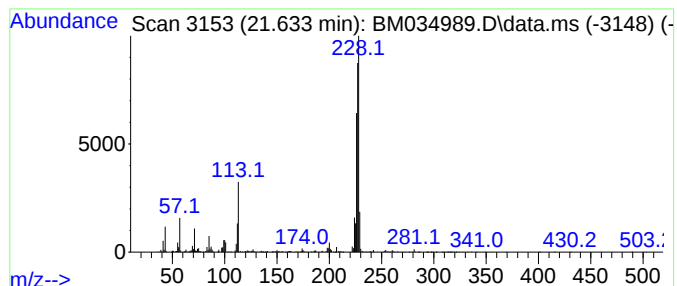
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Benzo[c]phenanthrene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.633	2.29 ng/ul	903362	Chrysene-d12	21.474

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]phenanthrene	228	C18H12	000195-19-7	74
2			Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	55
3			1(2H)-Phenanthrenone, 3,4,9,10-t...	228	C15H16O2	014427-61-3	50
4			Benzo[c]phenanthrene	228	C18H12	000195-19-7	50
5			Benz[a]anthracene	228	C18H12	000056-55-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034989.D
Acq On : 11 May 2022 17:31
Operator : CG/JU
Sample : N2603-19
Misc :
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Instrument :
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ClientSampleId :
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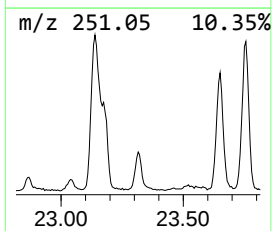
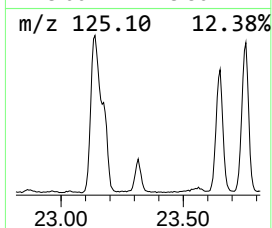
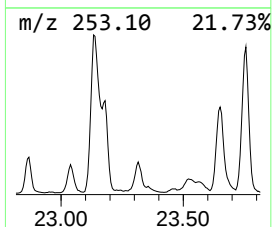
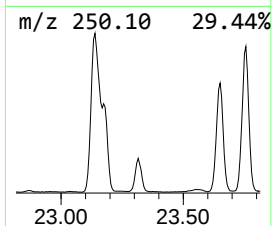
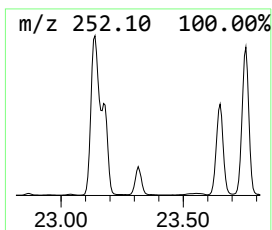
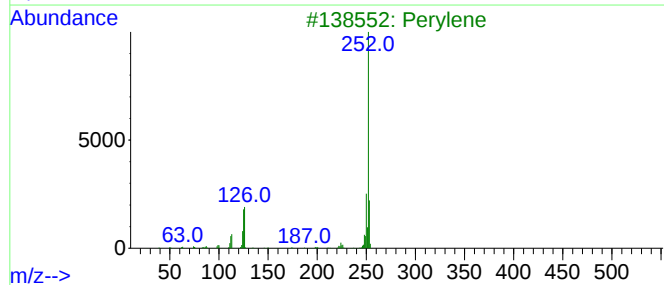
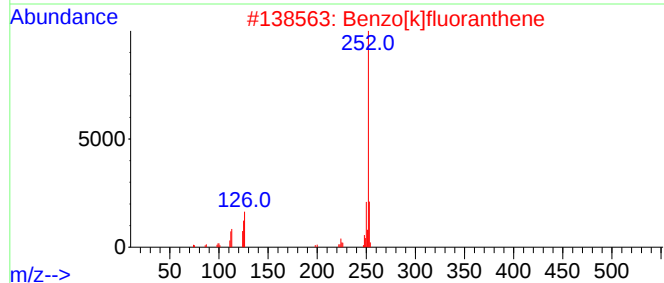
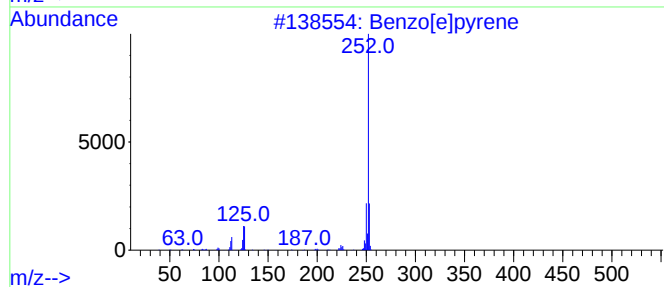
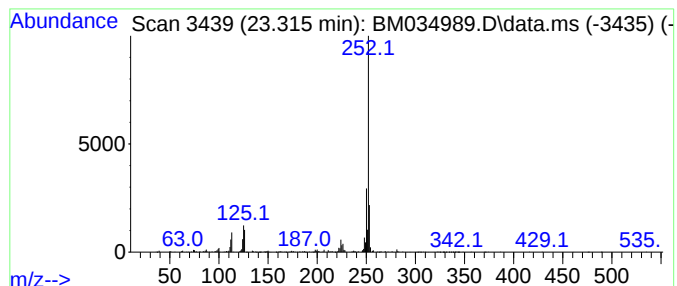
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.315	7.19 ng/ul	782036	Perylene-d12	23.856

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Perylene	252	C20H12	000198-55-0	95
4			Perylene	252	C20H12	000198-55-0	93
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
 Data File : BM034989.D
 Acq On : 11 May 2022 17:31
 Operator : CG/JU
 Sample : N2603-19
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4A0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
9H-Fluoren-9-ol	16.033	2.5	ng/ul	256375	4	17.292	2043040	20.0
(DEL) Alkane: S...	16.245	12.5	ng/ul	1276520	4	17.292	2043040	20.0
9H-Fluorene, 2-...	16.604	2.6	ng/ul	263764	4	17.292	2043040	20.0
9H-Fluoren-9-one	16.945	2.2	ng/ul	226312	4	17.292	2043040	20.0
(DEL) Alkane: S...	17.063	2.9	ng/ul	295323	4	17.292	2043040	20.0
Dibenzothiophene	17.110	4.7	ng/ul	484191	4	17.292	2043040	20.0
1,4-Ethenoanthr...	17.810	2.5	ng/ul	255820	4	17.292	2043040	20.0
Phenanthrene, 1...	18.168	5.2	ng/ul	528487	4	17.292	2043040	20.0
n-Hexadecanoic ...	18.198	5.5	ng/ul	561657	4	17.292	2043040	20.0
Phenanthrene, 4...	18.215	7.4	ng/ul	752876	4	17.292	2043040	20.0
4H-Cyclopenta[d...	18.368	12.8	ng/ul	1308930	4	17.292	2043040	20.0
Anthracene, 9-m...	18.398	2.9	ng/ul	298713	4	17.292	2043040	20.0
5,16[1',2']:8,1...	18.668	4.4	ng/ul	450880	4	17.292	2043040	20.0
9,10-Anthracene...	18.704	3.0	ng/ul	310288	4	17.292	2043040	20.0
Phenanthrene, 2...	19.086	2.1	ng/ul	217675	4	17.292	2043040	20.0
Cyclopenta(def)...	19.221	2.6	ng/ul	261870	4	17.292	2043040	20.0
11H-Benzo[b]flu...	20.204	2.6	ng/ul	1044090	5	21.474	7899270	20.0
Fluoranthene, 2...	20.304	2.2	ng/ul	856881	5	21.474	7899270	20.0
Benzo[ghi]fluor...	21.192	3.4	ng/ul	1334080	5	21.474	7899270	20.0
Benzo[c]phenant...	21.633	2.3	ng/ul	903362	5	21.474	7899270	20.0
Benzo[e]pyrene	23.315	7.2	ng/ul	782036	6	23.856	2174510	20.0