

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052324\
 Data File : BN031642.D
 Acq On : 23 May 2024 21:32
 Operator : MA/JU
 Sample : P2547-18
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BH5X5

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_N\Methods\SFAM-EPA-BN052124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031642.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.205	33	37	43	rVB	78575	99009	1.44%	0.092%
2	3.463	77	81	90	rBV	153158	206620	3.01%	0.192%
3	3.605	101	105	118	rBV	923669	1312422	19.13%	1.221%
4	6.857	650	658	675	rVB	1089947	1798273	26.21%	1.673%
5	7.028	681	687	697	rBV	982405	1554913	22.67%	1.447%
6	7.222	713	720	729	rBV	1209034	1873019	27.30%	1.743%
7	7.693	792	800	807	rBV	1530676	2444854	35.64%	2.275%
8	8.387	908	918	928	rBV	1167599	1857938	27.08%	1.729%
9	8.845	988	996	1005	rBV	1051011	1736941	25.32%	1.616%
10	9.410	1085	1092	1101	rVB	71188	110360	1.61%	0.103%
11	9.557	1109	1117	1128	rBV	747874	1270759	18.52%	1.183%
12	10.092	1200	1208	1218	rBV	1237807	2086208	30.41%	1.941%
13	10.257	1229	1236	1242	rBV	51017	83544	1.22%	0.078%
14	10.475	1264	1273	1279	rBV	1893821	3261282	47.54%	3.035%
15	10.522	1279	1281	1287	rVV	31587	44813	0.65%	0.042%
16	10.610	1287	1296	1310	rVB	1155818	1972188	28.75%	1.835%
17	11.216	1392	1399	1409	rBV	97122	180480	2.63%	0.168%
18	12.063	1536	1543	1551	rVV2	26024	49072	0.72%	0.046%
19	12.139	1551	1556	1562	rVB2	24197	41905	0.61%	0.039%
20	12.363	1588	1594	1599	rVB	27494	46553	0.68%	0.043%
21	13.645	1806	1812	1817	rBV2	34922	58220	0.85%	0.054%
22	13.739	1823	1828	1841	rVB	1755664	2498088	36.42%	2.325%
23	14.022	1869	1876	1886	rVB	2178401	3488821	50.86%	3.247%
24	14.328	1918	1928	1934	rBV	2331516	3507337	51.13%	3.264%
25	14.392	1934	1939	1946	rVB	201556	302081	4.40%	0.281%
26	14.522	1954	1961	1966	rBV	622392	1006148	14.67%	0.936%
27	14.639	1972	1981	1985	rVB4	18994	42025	0.61%	0.039%
28	14.728	1991	1996	2004	rVB	102463	171095	2.49%	0.159%
29	14.945	2026	2033	2038	rBV4	22678	49060	0.72%	0.046%
30	15.122	2054	2063	2065	rBV5	22354	48233	0.70%	0.045%
31	15.322	2090	2097	2103	rVV	2606115	3753733	54.72%	3.493%
32	15.375	2103	2106	2111	rVV	170503	247178	3.60%	0.230%
33	15.439	2111	2117	2124	rVV	379099	571609	8.33%	0.532%
34	15.498	2124	2127	2130	rVV4	39452	60159	0.88%	0.056%
35	15.539	2130	2134	2140	rVV3	37642	82654	1.20%	0.077%
36	15.669	2152	2156	2167	rVB	63834	111327	1.62%	0.104%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052124.MA.M
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37	15.816	2177	2181	2189	rVB	57355	116461	1.70%	0.108%
38	16.051	2216	2221	2226	rBV3	48013	73123	1.07%	0.068%
39	16.357	2266	2273	2274	rBV5	34700	62298	0.91%	0.058%
40	16.380	2275	2277	2282	rVV2	45962	67761	0.99%	0.063%
41	16.422	2282	2284	2289	rVV2	28070	42214	0.62%	0.039%
42	16.610	2308	2316	2319	rBV3	38504	70998	1.03%	0.066%
43	16.645	2319	2322	2326	rVB3	33771	45602	0.66%	0.042%
44	16.727	2331	2336	2343	rVB	101169	168204	2.45%	0.157%
45	16.827	2345	2353	2359	rVV5	61883	176551	2.57%	0.164%
46	16.892	2359	2364	2373	rVB	120887	194924	2.84%	0.181%
47	17.069	2388	2394	2398	rBV	2321784	3489144	50.86%	3.247%
48	17.116	2398	2402	2406	rVV	2402217	3505711	51.10%	3.262%
49	17.169	2406	2411	2415	rVV2	2484484	3664212	53.41%	3.410%
50	17.204	2415	2417	2422	rVV	553815	661287	9.64%	0.615%
51	17.263	2422	2427	2432	rVB3	72550	123090	1.79%	0.115%
52	17.474	2453	2463	2467	rBV2	222842	386138	5.63%	0.359%
53	17.592	2479	2483	2489	rBV3	49552	77396	1.13%	0.072%
54	17.651	2489	2493	2500	rVB4	45487	72231	1.05%	0.067%
55	17.945	2534	2543	2546	rBV	305958	523628	7.63%	0.487%
56	17.992	2546	2551	2559	rVV2	471434	943771	13.76%	0.878%
57	18.074	2559	2565	2570	rVV	146671	218742	3.19%	0.204%
58	18.139	2570	2576	2580	rBV2	434274	798317	11.64%	0.743%
59	18.174	2580	2582	2590	rVB	206263	275352	4.01%	0.256%
60	18.298	2600	2603	2606	rVB4	41760	53979	0.79%	0.050%
61	18.451	2625	2629	2632	rBV	223331	302306	4.41%	0.281%
62	18.486	2632	2635	2642	rVB	184921	264622	3.86%	0.246%
63	18.574	2648	2650	2656	rVB3	32809	44617	0.65%	0.042%
64	18.698	2667	2671	2675	rBV	80929	105517	1.54%	0.098%
65	18.745	2676	2679	2682	rVV	70537	85589	1.25%	0.080%
66	18.786	2682	2686	2690	rVB2	72292	98174	1.43%	0.091%
67	18.868	2695	2700	2705	rBV	192981	313081	4.56%	0.291%
68	18.933	2705	2711	2714	rBV4	96267	190961	2.78%	0.178%
69	19.004	2719	2723	2731	rBV2	177318	342394	4.99%	0.319%
70	19.133	2739	2745	2750	rVB	4099880	5493325	80.08%	5.112%
71	19.233	2759	2762	2764	rBV3	55514	66325	0.97%	0.062%
72	19.257	2764	2766	2773	rVB3	104862	174727	2.55%	0.163%
73	19.374	2783	2786	2790	rVB	104474	122417	1.78%	0.114%
74	19.468	2796	2802	2804	rBV	3404097	5136961	74.88%	4.780%
75	19.492	2804	2806	2815	rVV	3375042	4569916	66.62%	4.253%
76	19.568	2815	2819	2823	rVV2	146516	208586	3.04%	0.194%

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 Title : SVOA CALIBRATION

77	19.610	2823	2826	2831	rVB	119180	135743	1.98%	0.126%
78	19.674	2833	2837	2842	rBV	183037	254260	3.71%	0.237%
79	19.733	2843	2847	2853	rVB	192230	291403	4.25%	0.271%
80	19.821	2856	2862	2867	rBV4	226817	565366	8.24%	0.526%
81	19.957	2881	2885	2887	rBV3	142328	227947	3.32%	0.212%
82	19.992	2887	2891	2895	rVB	447926	626017	9.13%	0.583%
83	20.092	2904	2908	2911	rBV2	287229	373340	5.44%	0.347%
84	20.145	2912	2917	2922	rVV2	349321	650379	9.48%	0.605%
85	20.186	2922	2924	2927	rVV2	92852	112558	1.64%	0.105%
86	20.286	2938	2941	2945	rBV2	184673	249167	3.63%	0.232%
87	20.739	3013	3018	3028	rBV4	527325	1328600	19.37%	1.236%
88	20.951	3051	3054	3057	rBV	272544	380035	5.54%	0.354%
89	20.992	3057	3061	3066	rVV3	545521	1043474	15.21%	0.971%
90	21.045	3068	3070	3083	rVB	447861	895856	13.06%	0.834%
91	21.298	3106	3113	3117	rBV3	3089200	6860037	100.00%	6.384%
92	21.345	3117	3121	3130	rVB	1702553	2675293	39.00%	2.490%
93	21.486	3141	3145	3152	rVB2	223267	394771	5.75%	0.367%
94	21.974	3225	3228	3234	rVB	293086	431978	6.30%	0.402%
95	23.309	3448	3455	3462	rBV	1276474	3890048	56.71%	3.620%
96	23.968	3561	3567	3572	rBV	659721	1449068	21.12%	1.348%
97	24.039	3572	3579	3584	rVV	1738049	4052413	59.07%	3.771%
98	24.098	3585	3589	3603	rVB2	1223431	2853579	41.60%	2.655%
99	24.233	3605	3612	3620	rBV2	1725008	4137881	60.32%	3.851%
100	27.533	4167	4173	4192	rVB3	533927	2222195	32.39%	2.068%

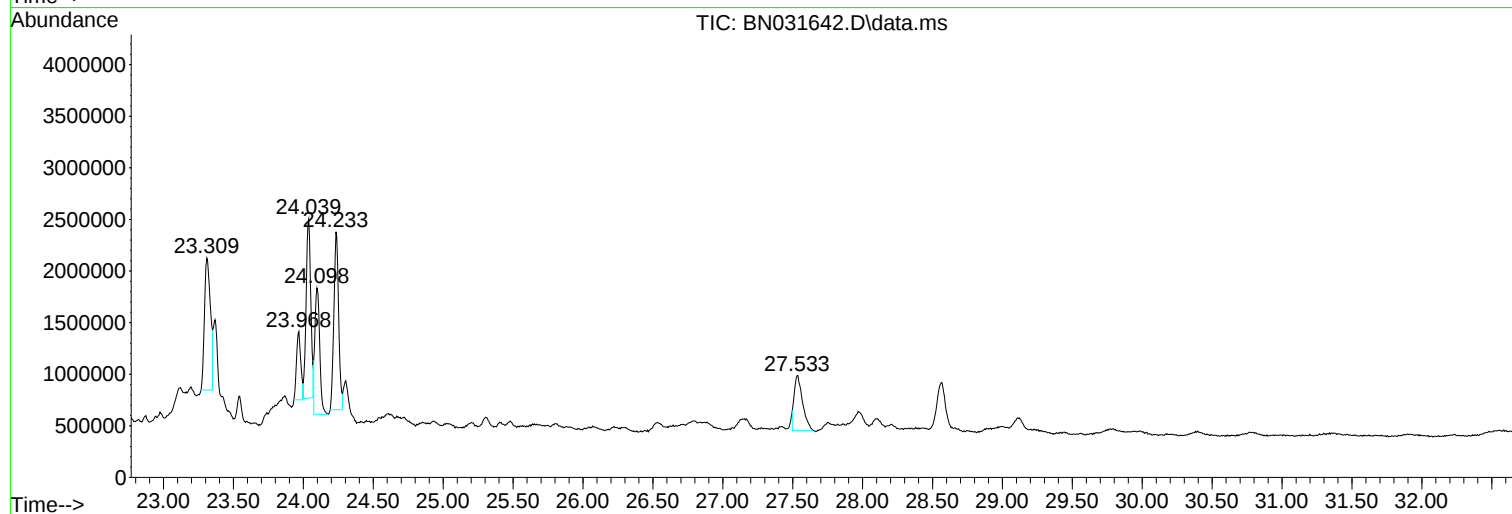
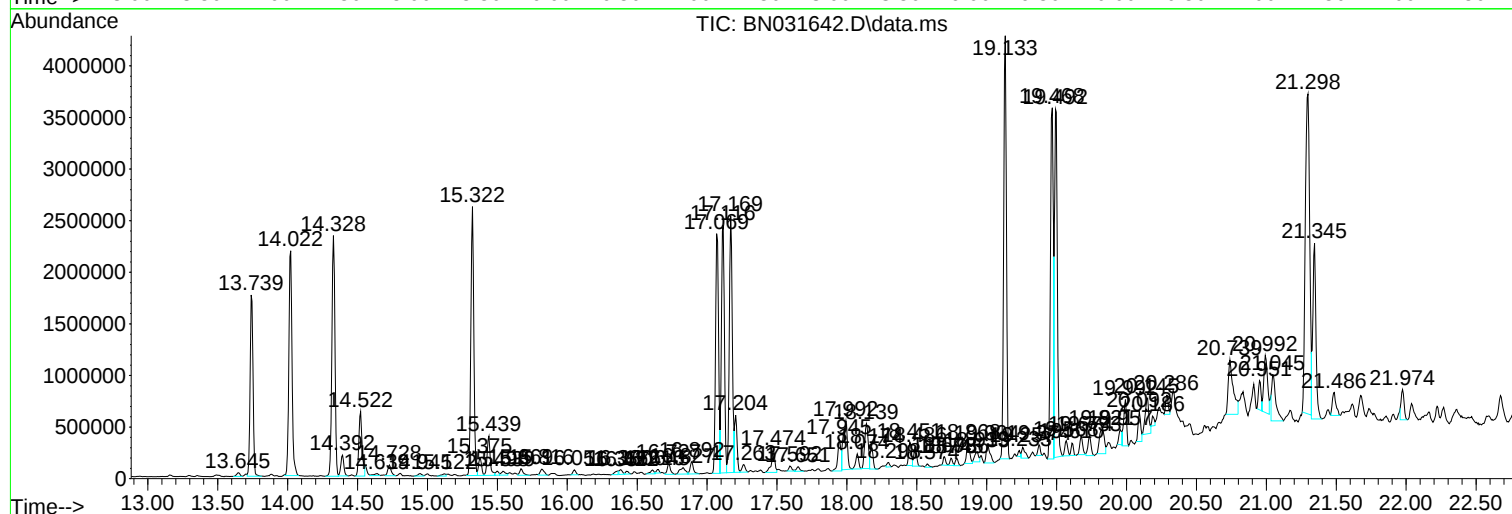
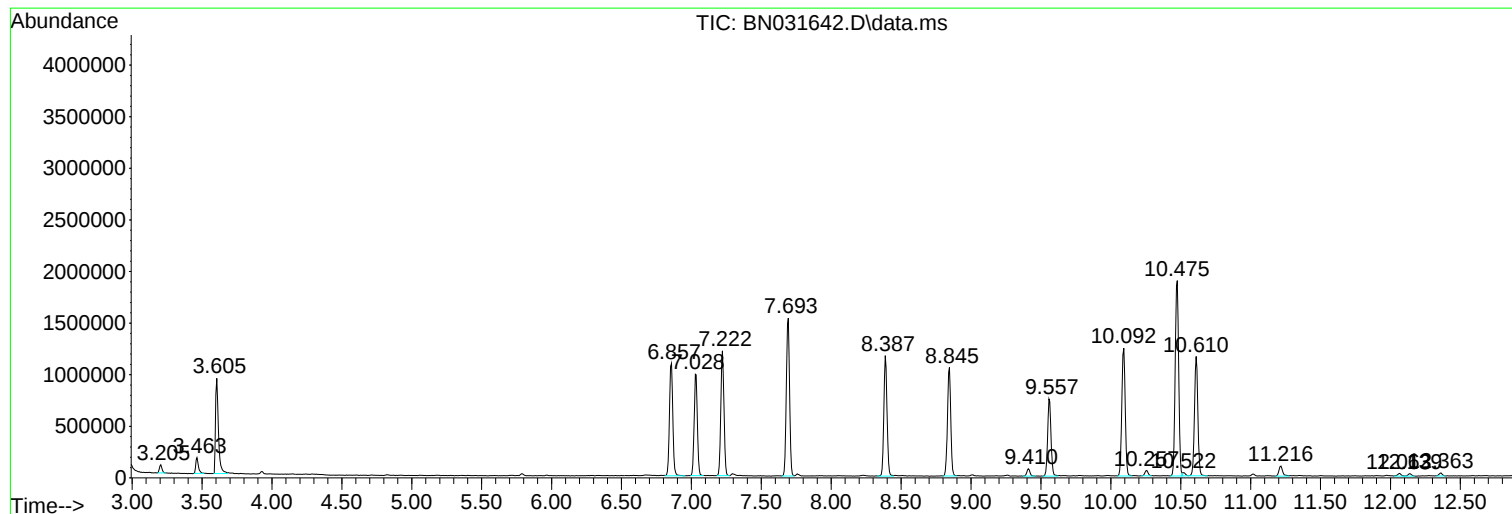
Sum of corrected areas: 107460981

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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052124.MA.M
Quant Title : SVOA CALIBRATION

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



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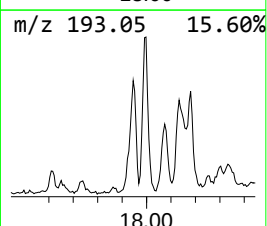
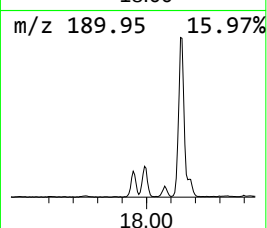
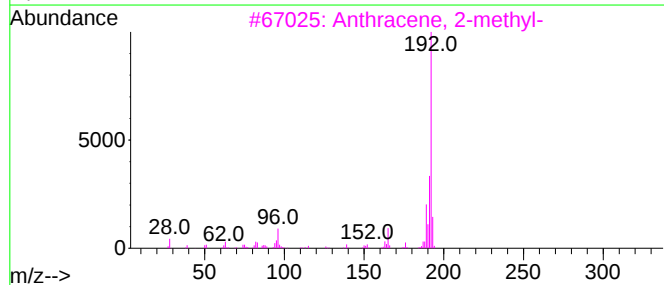
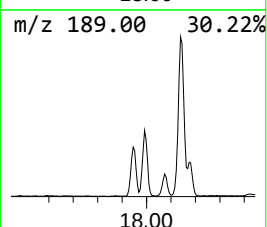
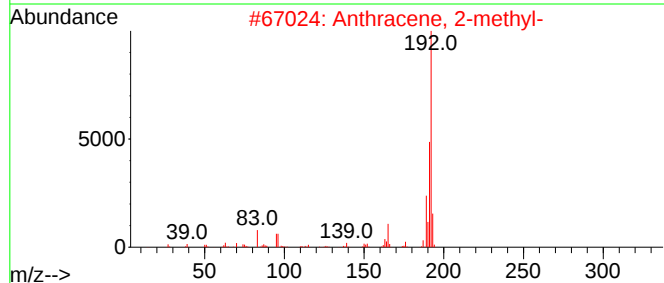
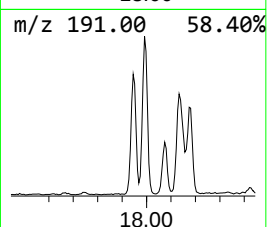
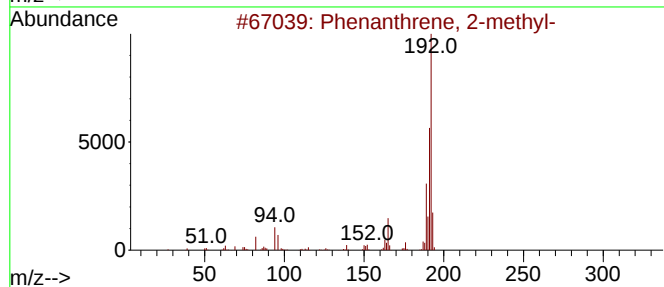
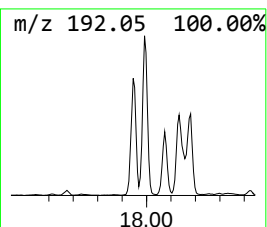
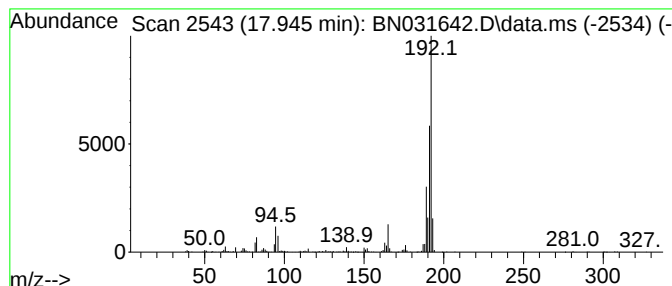
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.945	3.00 ng/ul	523628	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



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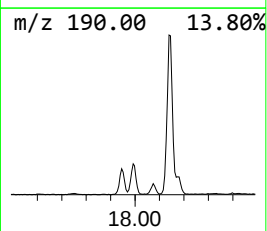
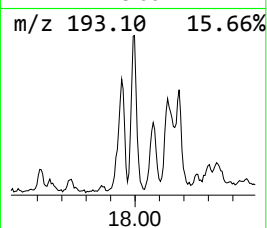
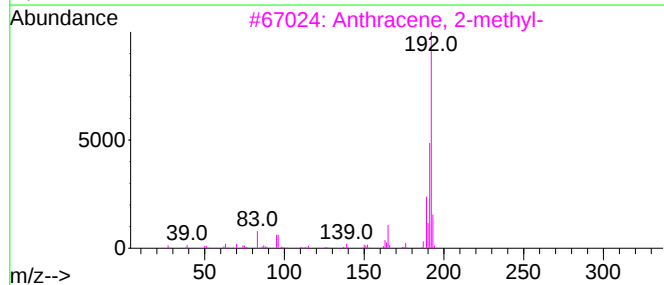
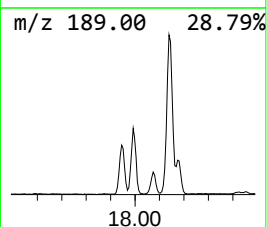
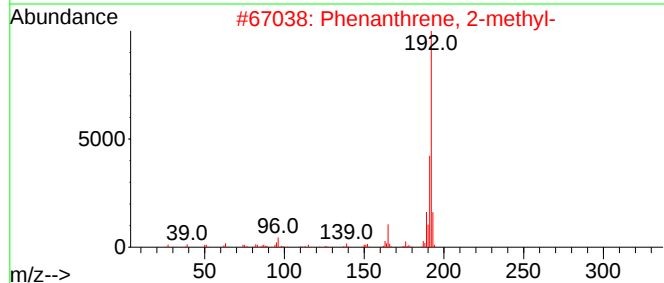
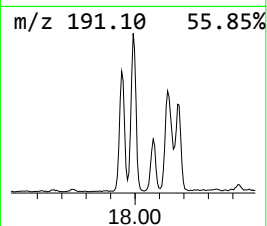
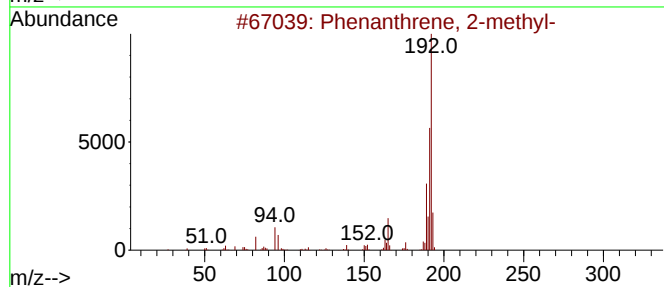
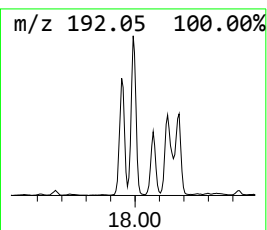
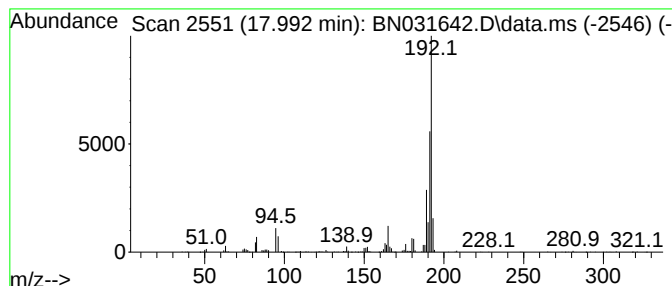
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052124.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.992	5.41 ng/ul	943771	Phenanthrene-d10	17.069

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	97
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



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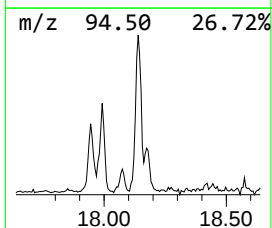
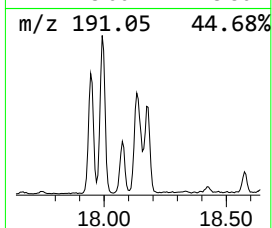
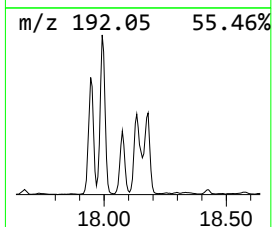
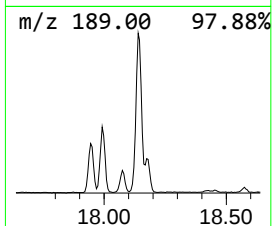
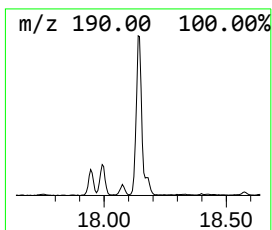
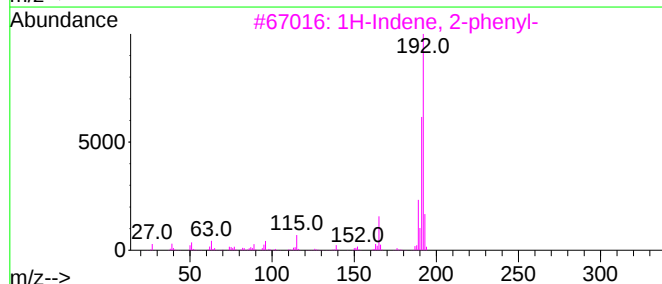
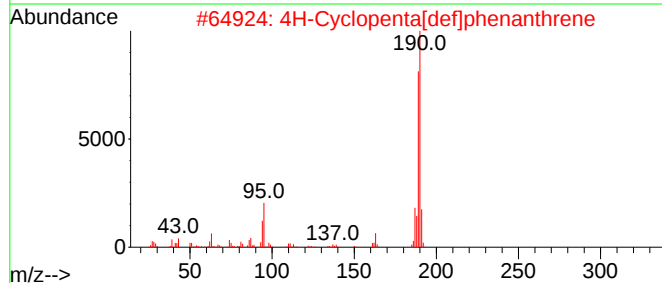
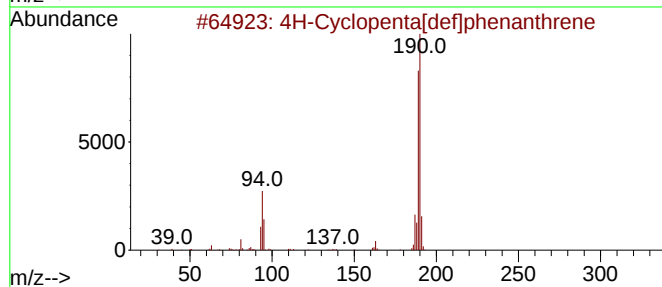
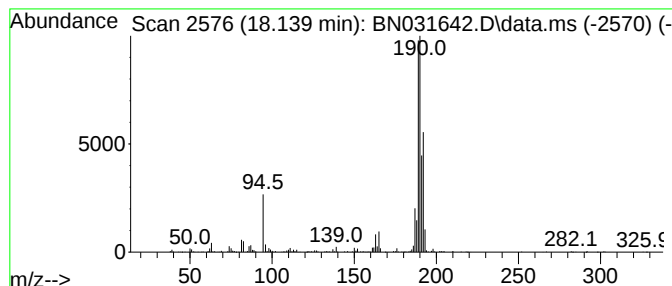
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	4.58 ng/ul	798317	Phenanthrene-d10	17.069

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	60
3			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30
4			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	25
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052324\
Data File : BN031642.D
Acq On : 23 May 2024 21:32
Operator : MA/JU
Sample : P2547-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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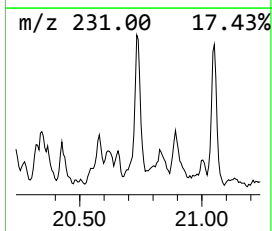
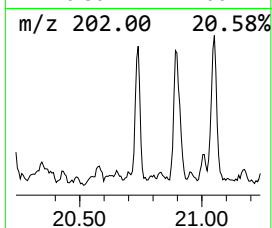
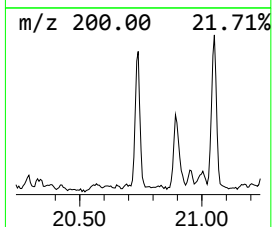
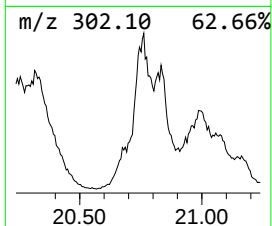
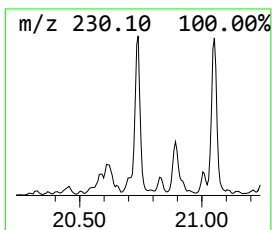
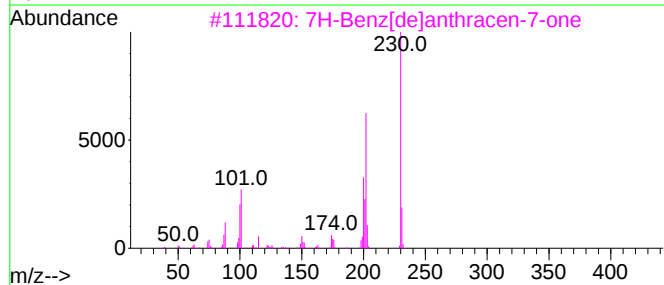
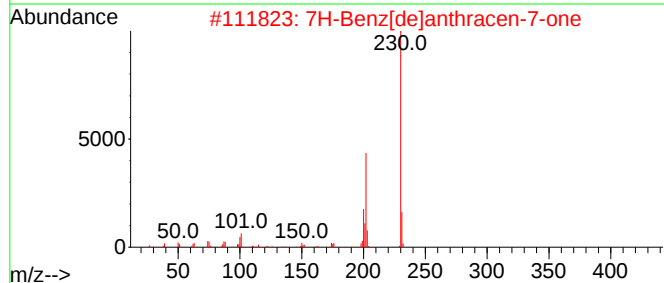
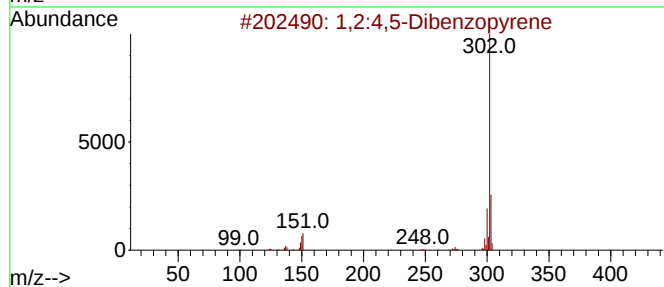
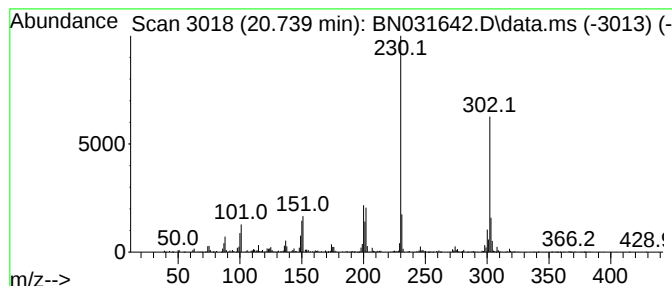
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,2:4,5-Dibenzopyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.739	3.87 ng/ul	1328600	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	91
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	70
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	70
4			Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	59
5			1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052324\
Data File : BN031642.D
Acq On : 23 May 2024 21:32
Operator : MA/JU
Sample : P2547-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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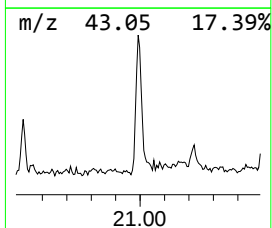
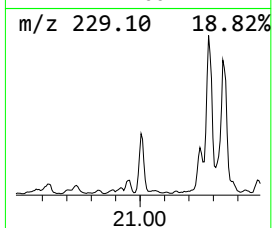
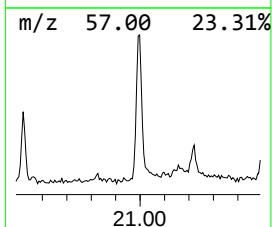
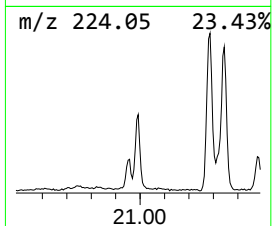
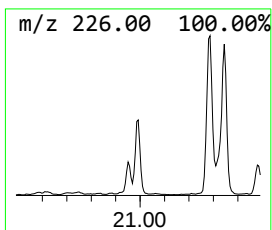
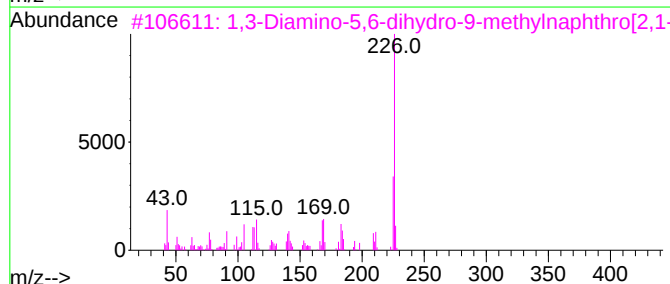
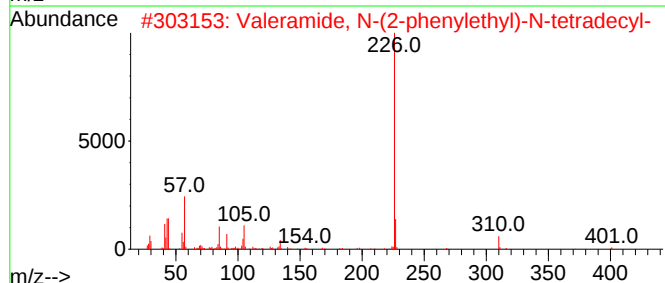
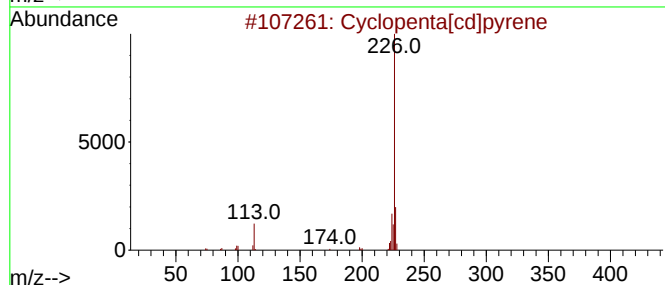
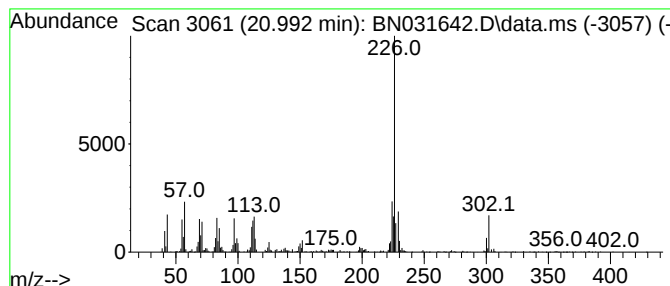
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclopenta[cd]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.992	3.04 ng/ul	1043470	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	53
2			Valeramide, N-(2-phenylethyl)-N...	401	C27H47NO	1000451-53-7	47
3			1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	43
4			.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	40
5			benzenamine, 4-(2-benzothiazolyl)-	226	C13H10N2S	1000400-93-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052324\
Data File : BN031642.D
Acq On : 23 May 2024 21:32
Operator : MA/JU
Sample : P2547-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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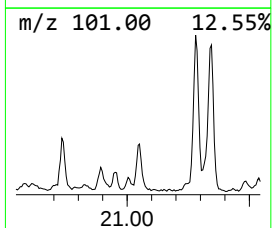
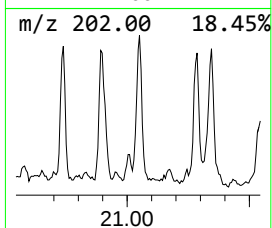
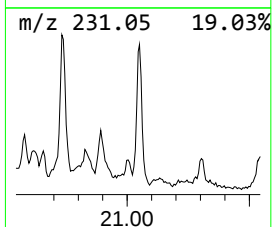
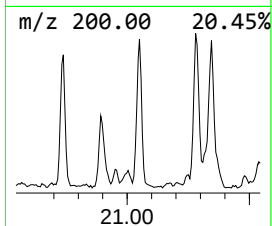
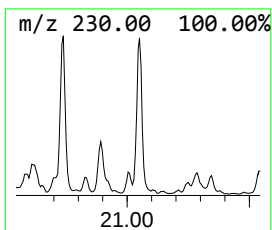
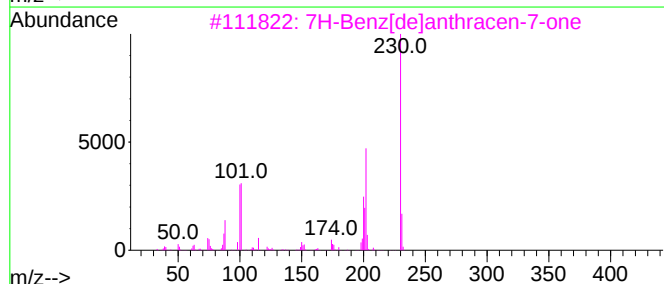
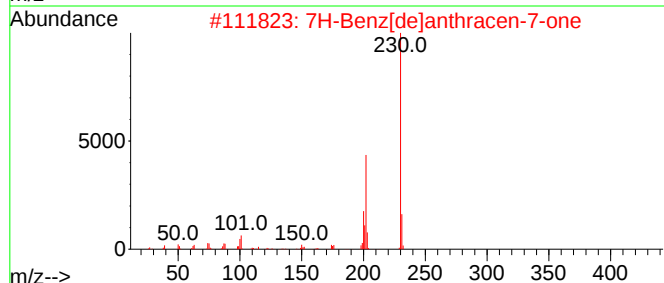
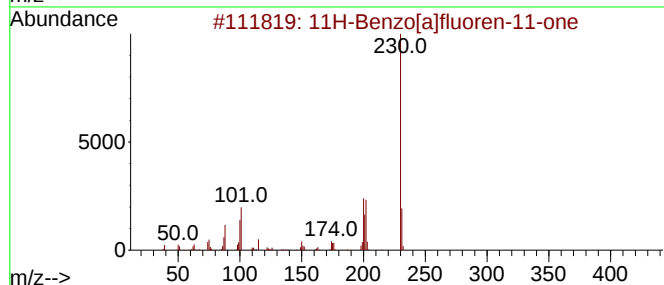
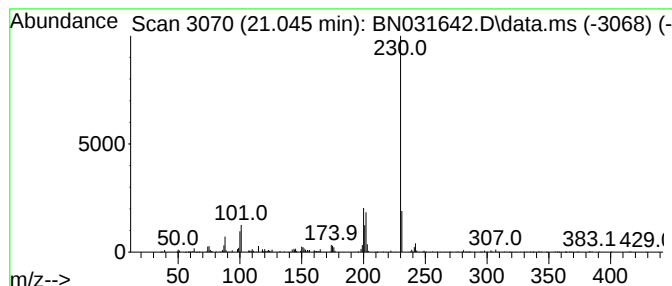
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 11H-Benzo[a]fluoren-11-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.045	2.61 ng/ul	895856	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	97
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	93
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
4			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
5			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052324\
Data File : BN031642.D
Acq On : 23 May 2024 21:32
Operator : MA/JU
Sample : P2547-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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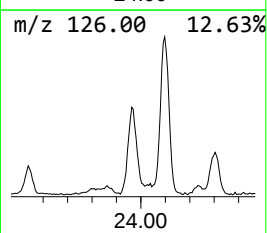
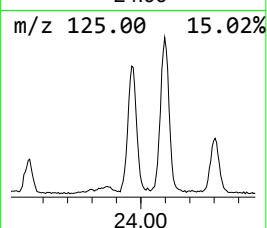
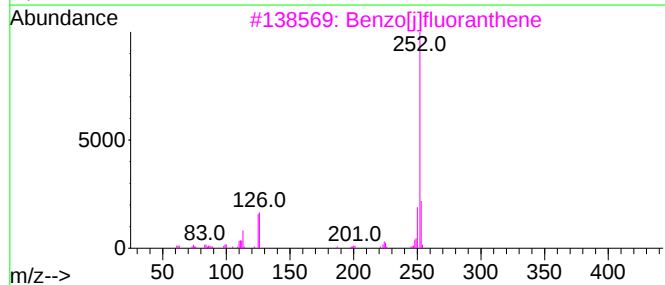
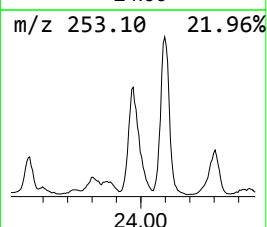
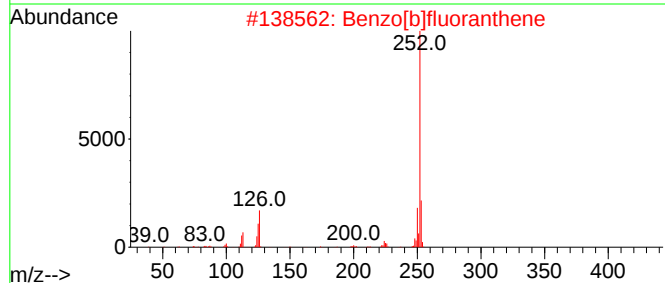
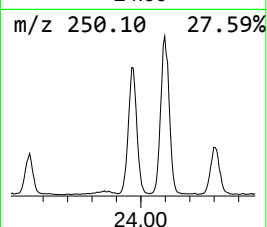
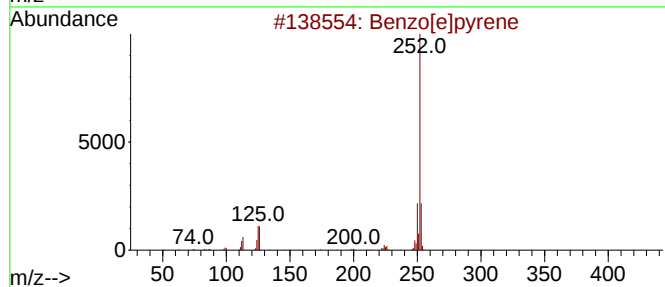
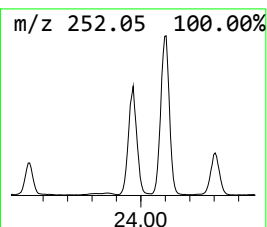
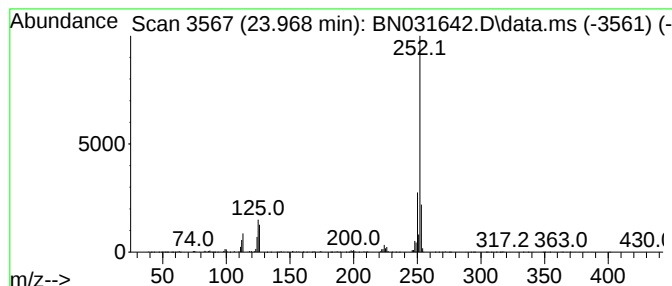
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.968	7.00 ng/ul	1449070	Perylene-d12	24.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	98
3			Benzo[j]fluoranthene	252	C20H12	000205-82-3	97
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5			Perylene	252	C20H12	000198-55-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052324\
Data File : BN031642.D
Acq On : 23 May 2024 21:32
Operator : MA/JU
Sample : P2547-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH5X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052124.MA.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
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Anthracene, 2-m...	17.992	5.4	ng/u1	943771	4	17.069	3489140	20.0
4H-Cyclopenta[d...	18.139	4.6	ng/u1	798317	4	17.069	3489140	20.0
1,2:4,5-Dibenzo...	20.739	3.9	ng/u1	1328600	5	21.304	6860040	20.0
Cyclopenta[cd]p...	20.992	3.0	ng/u1	1043470	5	21.304	6860040	20.0
11H-Benzo[a]flu...	21.045	2.6	ng/u1	895856	5	21.304	6860040	20.0
Benzo[e]pyrene	23.968	7.0	ng/u1	1449070	6	24.233	4137880	20.0