

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
 Data File : BN031724.D
 Acq On : 30 May 2024 20:17
 Operator : MA/JU
 Sample : P2569-21DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BH6A7DL

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-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031724.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.599	101	104	116	rBV	245305	404305	2.21%	0.196%
2	6.852	647	657	668	rVB	455732	792198	4.33%	0.385%
3	7.022	680	686	698	rVB	383197	607845	3.32%	0.295%
4	7.216	709	719	727	rBV	459790	728862	3.98%	0.354%
5	7.687	791	799	806	rBV	1682077	2669881	14.58%	1.297%
6	8.387	911	918	926	rBV	555257	890113	4.86%	0.432%
7	8.840	987	995	1004	rBV	453644	749869	4.09%	0.364%
8	9.557	1110	1117	1125	rVV	324914	554194	3.03%	0.269%
9	10.087	1199	1207	1220	rVB	629052	1032894	5.64%	0.502%
10	10.469	1263	1272	1279	rBV	2658268	4483320	24.48%	2.178%
11	10.610	1287	1296	1306	rBV	427270	775488	4.23%	0.377%
12	12.134	1549	1555	1559	rVV	102270	177753	0.97%	0.086%
13	12.357	1583	1593	1600	rVB	119860	208156	1.14%	0.101%
14	13.645	1802	1812	1817	rBV2	241136	432884	2.36%	0.210%
15	13.698	1817	1821	1824	rBV	123869	175483	0.96%	0.085%
16	13.740	1824	1828	1833	rVB	1274912	1708988	9.33%	0.830%
17	13.881	1844	1852	1856	rBV	115628	189754	1.04%	0.092%
18	14.016	1865	1875	1879	rBV	1427231	2310081	12.61%	1.122%
19	14.328	1916	1928	1934	rBV	4340182	6597489	36.02%	3.205%
20	14.392	1934	1939	1947	rVV2	234254	426766	2.33%	0.207%
21	14.528	1956	1962	1972	rBV2	473407	875601	4.78%	0.425%
22	14.728	1991	1996	2002	rVV	418472	641814	3.50%	0.312%
23	14.804	2002	2009	2013	rVB	135762	212565	1.16%	0.103%
24	14.945	2025	2033	2038	rBV3	108971	213200	1.16%	0.104%
25	15.116	2054	2062	2065	rBV	89560	182526	1.00%	0.089%
26	15.322	2090	2097	2102	rVV	1982828	2949190	16.10%	1.433%
27	15.375	2102	2106	2111	rVV	548871	836428	4.57%	0.406%
28	15.445	2113	2118	2124	rVV	425685	679080	3.71%	0.330%
29	15.669	2151	2156	2163	rVB	332058	494331	2.70%	0.240%
30	15.816	2174	2181	2190	rVB	333016	683839	3.73%	0.332%
31	16.051	2216	2221	2228	rVB3	146049	234338	1.28%	0.114%
32	16.381	2267	2277	2282	rBV4	196922	439774	2.40%	0.214%
33	16.610	2308	2316	2319	rBV3	190742	330738	1.81%	0.161%
34	16.651	2319	2323	2326	rVV2	199868	303333	1.66%	0.147%
35	16.728	2331	2336	2344	rVB	531319	801454	4.38%	0.389%
36	16.804	2344	2349	2352	rBV	209145	365792	2.00%	0.178%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
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37	16.839	2352	2355	2359	rVV	440913	594785	3.25%	0.289%
38	16.892	2359	2364	2370	rVV	678873	1022196	5.58%	0.497%
39	17.075	2388	2395	2398	rBV	5277136	7748119	42.31%	3.764%
40	17.116	2398	2402	2407	rVV	9498425	13880887	75.79%	6.743%
41	17.175	2407	2412	2415	rVV	2129784	3253623	17.77%	1.581%
42	17.204	2415	2417	2423	rVV	1881007	2368359	12.93%	1.151%
43	17.257	2423	2426	2432	rVB4	105393	186207	1.02%	0.090%
44	17.481	2457	2464	2467	rBV	218753	361084	1.97%	0.175%
45	17.598	2479	2484	2489	rVV	290495	438920	2.40%	0.213%
46	17.651	2489	2493	2502	rVB2	341829	582827	3.18%	0.283%
47	17.798	2513	2518	2525	rVB	179251	336849	1.84%	0.164%
48	17.869	2525	2530	2534	rBV	173763	254188	1.39%	0.123%
49	17.951	2538	2544	2547	rVV	1355320	2124365	11.60%	1.032%
50	17.992	2547	2551	2559	rVV	2116893	3803912	20.77%	1.848%
51	18.075	2561	2565	2570	rVV	522481	778611	4.25%	0.378%
52	18.139	2570	2576	2580	rVV2	1600783	2934137	16.02%	1.425%
53	18.180	2580	2583	2591	rVB	933795	1279524	6.99%	0.622%
54	18.451	2624	2629	2633	rVV	1341290	1827832	9.98%	0.888%
55	18.492	2633	2636	2642	rVB	739752	1028247	5.61%	0.500%
56	18.698	2668	2671	2675	rVV2	282894	412021	2.25%	0.200%
57	18.751	2676	2680	2683	rVV	314154	403781	2.20%	0.196%
58	18.786	2683	2686	2691	rVB	275312	365149	1.99%	0.177%
59	18.875	2695	2701	2705	rVV	677537	1126484	6.15%	0.547%
60	18.933	2705	2711	2714	rVV4	379287	848110	4.63%	0.412%
61	18.963	2714	2716	2720	rVV	302154	338506	1.85%	0.164%
62	19.010	2720	2724	2732	rVV	871139	1429325	7.80%	0.694%
63	19.139	2740	2746	2753	rVB	12785743	18313894	100.00%	8.897%
64	19.204	2753	2757	2760	rBV	214278	269847	1.47%	0.131%
65	19.233	2760	2762	2765	rBV2	172026	196376	1.07%	0.095%
66	19.286	2765	2771	2775	rVV2	682678	1343769	7.34%	0.653%
67	19.333	2776	2779	2783	rVB	194810	253499	1.38%	0.123%
68	19.380	2783	2787	2790	rBV	473344	599205	3.27%	0.291%
69	19.469	2797	2802	2804	rBV2	4339487	6053306	33.05%	2.941%
70	19.504	2804	2808	2815	rVV	11601420	16470291	89.93%	8.001%
71	19.569	2815	2819	2826	rVB2	562284	912982	4.99%	0.444%
72	19.680	2833	2838	2843	rVB	698636	951521	5.20%	0.462%
73	19.739	2844	2848	2854	rVB2	265565	444335	2.43%	0.216%
74	19.833	2857	2864	2870	rBV3	825828	2104125	11.49%	1.022%
75	19.957	2880	2885	2888	rBV3	641057	1148110	6.27%	0.558%
76	19.998	2888	2892	2896	rVB	1162900	1669929	9.12%	0.811%

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Integration Parameters: LSCINT.P

Integrator: RTE

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77	20.098	2905	2909	2912	rBV	432694	502823	2.75%	0.244%
78	20.157	2912	2919	2923	rVV2	1137544	1908410	10.42%	0.927%
79	20.239	2929	2933	2938	rBV3	194152	339735	1.86%	0.165%
80	20.292	2938	2942	2946	rVV	580262	769451	4.20%	0.374%
81	20.339	2946	2950	2955	rVV	700165	999245	5.46%	0.485%
82	20.745	3015	3019	3025	rVB	1454295	1887858	10.31%	0.917%
83	20.916	3042	3048	3052	rBV2	994271	1842310	10.06%	0.895%
84	20.957	3052	3055	3059	rVV	1046057	1415151	7.73%	0.687%
85	20.998	3059	3062	3068	rVV2	1469904	2434497	13.29%	1.183%
86	21.057	3068	3072	3078	rVB2	1264797	1920680	10.49%	0.933%
87	21.222	3097	3100	3104	rVB	624636	699310	3.82%	0.340%
88	21.298	3107	3113	3119	rVV2	7113352	17263488	94.26%	8.386%
89	21.351	3119	3122	3134	rVB	4656013	7254670	39.61%	3.524%
90	21.492	3142	3146	3152	rBV3	394010	857124	4.68%	0.416%
91	21.686	3174	3179	3184	rBV2	387397	795147	4.34%	0.386%
92	21.980	3226	3229	3235	rVB	713411	1098391	6.00%	0.534%
93	22.069	3237	3244	3250	rVB2	704104	1515349	8.27%	0.736%
94	23.327	3450	3458	3464	rBV2	2286377	6959893	38.00%	3.381%
95	23.433	3472	3476	3483	rVB2	990048	1843995	10.07%	0.896%
96	23.974	3563	3568	3575	rBV	1081447	2477511	13.53%	1.204%
97	24.045	3576	3580	3585	rVV2	978215	2133906	11.65%	1.037%
98	24.110	3586	3591	3602	rVB	1905485	4578065	25.00%	2.224%
99	24.251	3607	3615	3623	rBV	2393109	6426639	35.09%	3.122%
100	25.304	3789	3794	3804	rVB4	491742	1308903	7.15%	0.636%

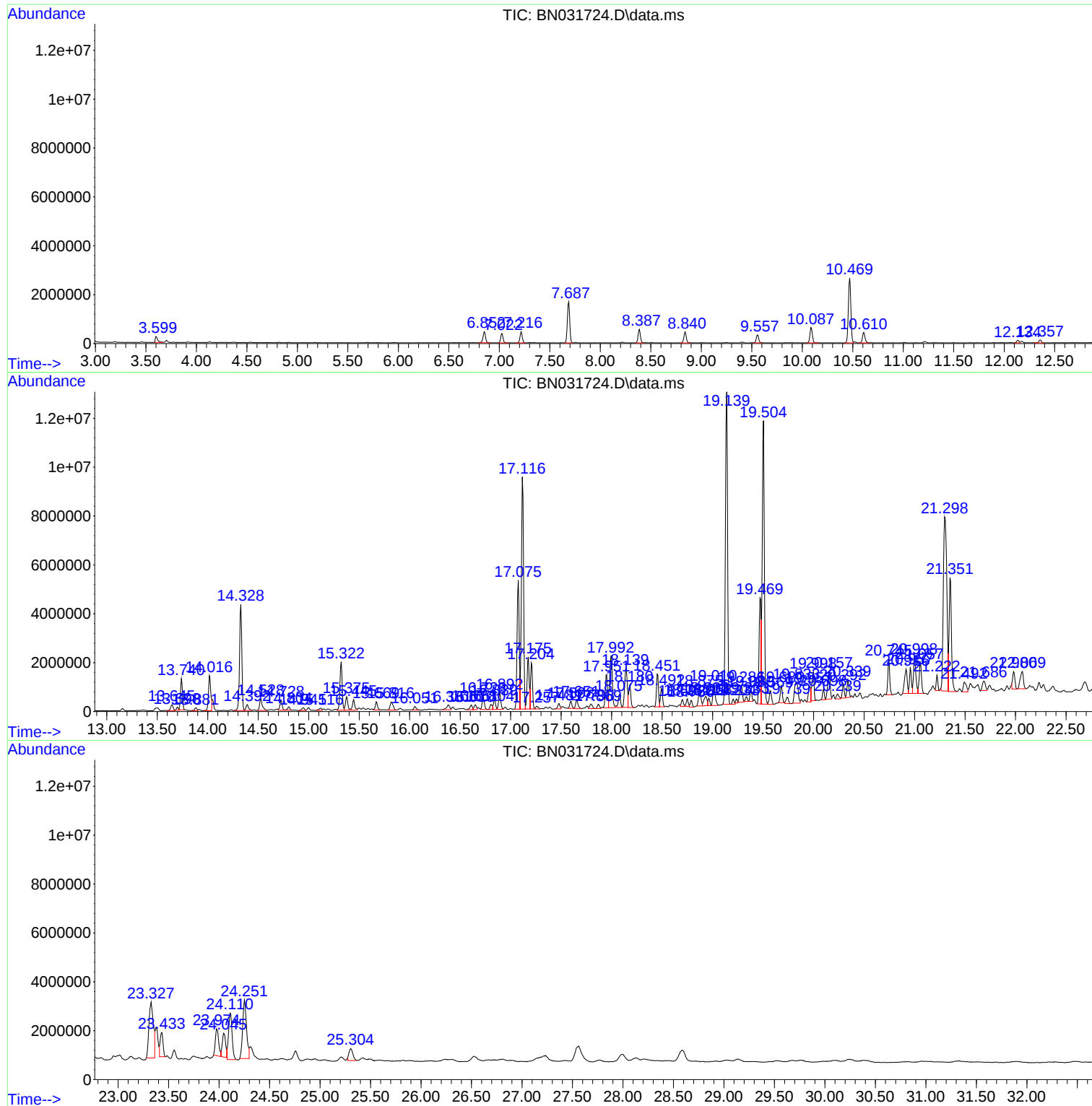
Sum of corrected areas: 205852114

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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH6A7DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
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Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
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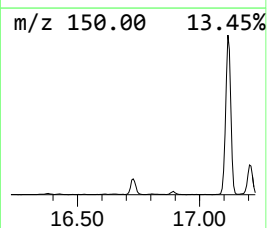
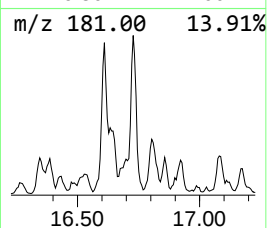
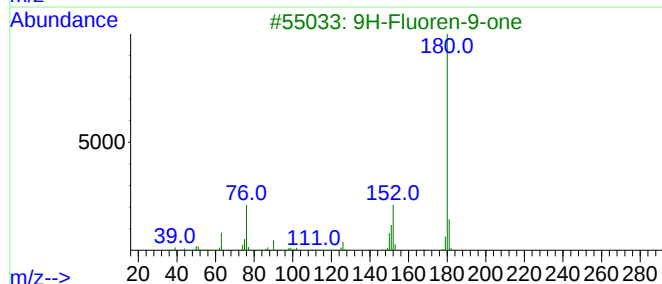
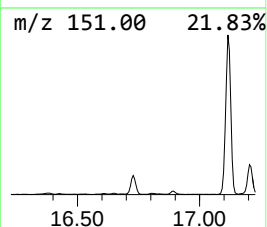
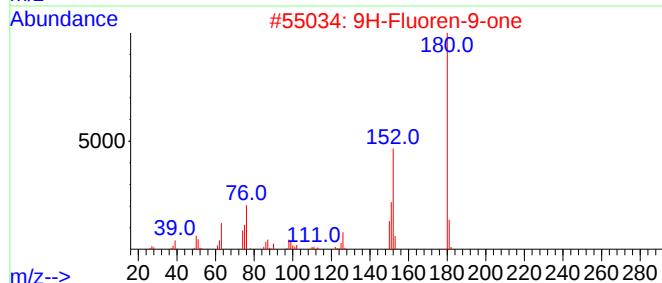
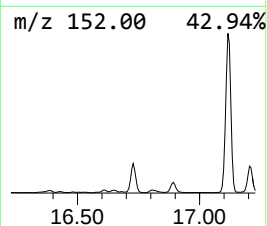
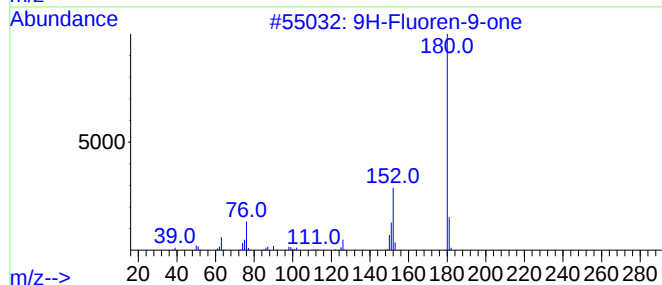
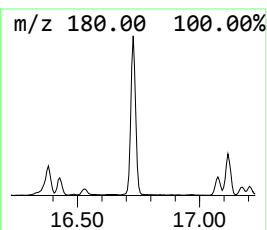
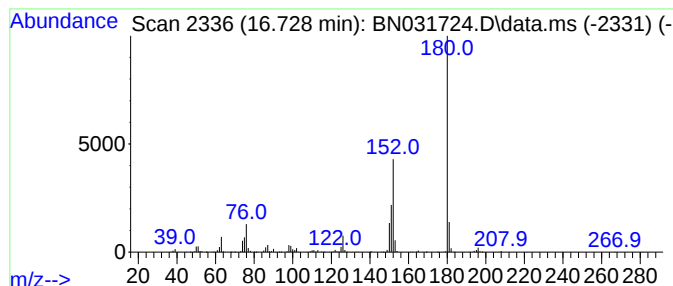
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 9H-Fluoren-9-one Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.728	2.07 ng/ul	801454	Phenanthrene-d10	17.075

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	96
3	9H-Fluoren-9-one			180	C13H8O	000486-25-9	95
4	9H-Fluoren-9-one			180	C13H8O	000486-25-9	95
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	95



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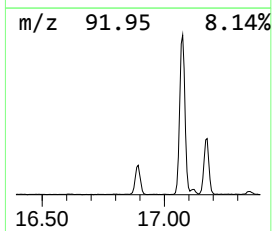
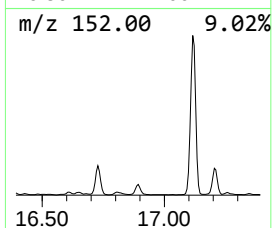
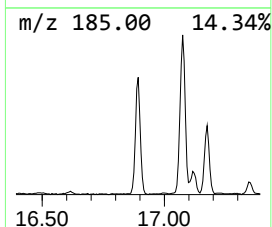
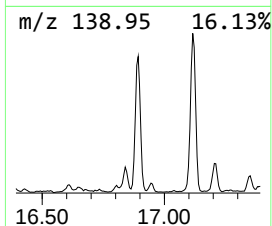
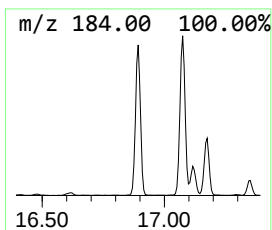
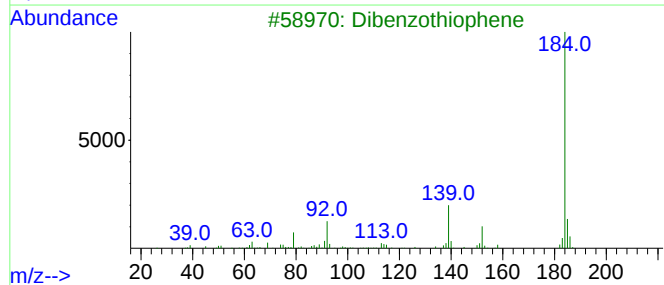
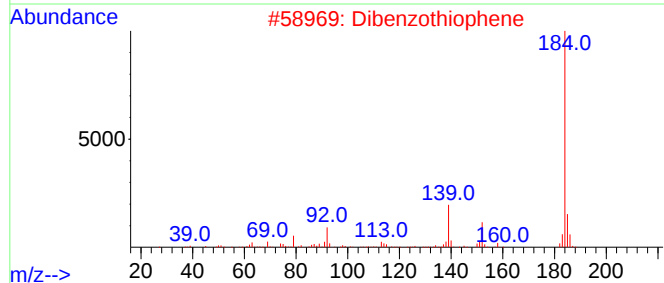
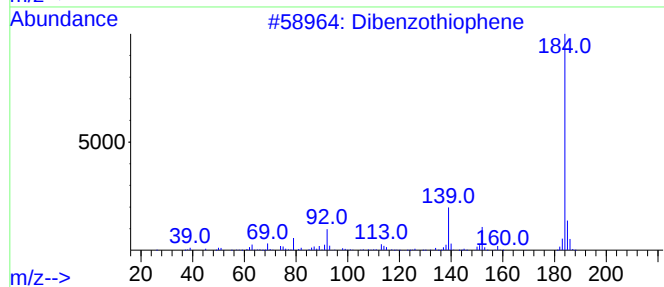
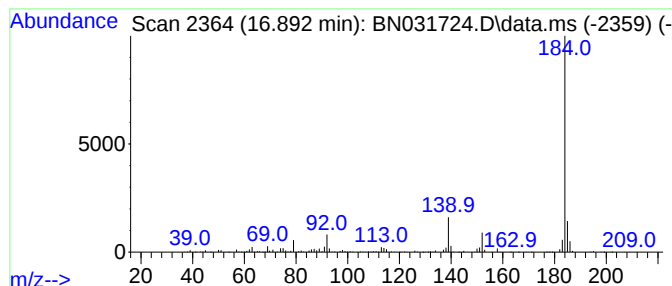
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Dibenzothiophene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.892	2.64 ng/ul	1022200	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Dibenzothiophene	184	C12H8S	000132-65-0	97
3			Dibenzothiophene	184	C12H8S	000132-65-0	97
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	96
5			Dibenzothiophene	184	C12H8S	000132-65-0	96



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ClientSampleId :
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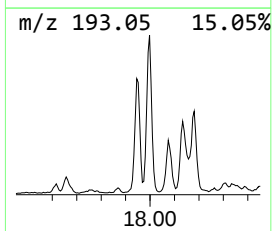
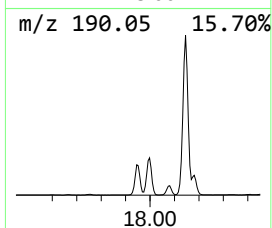
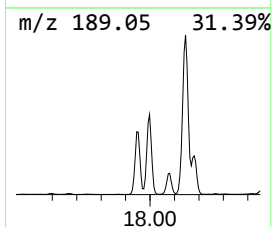
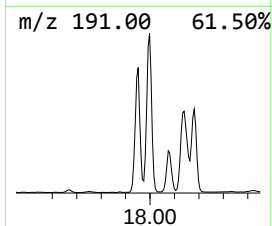
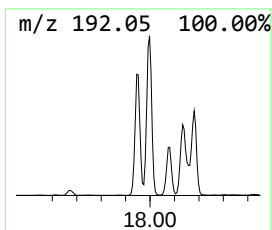
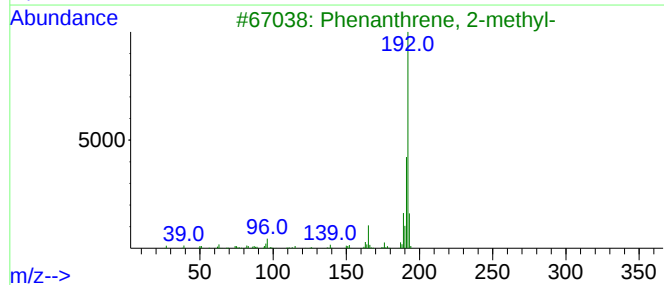
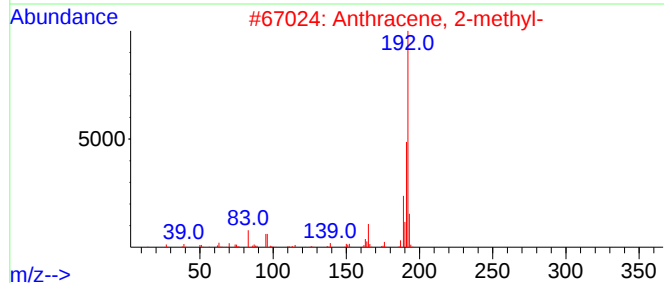
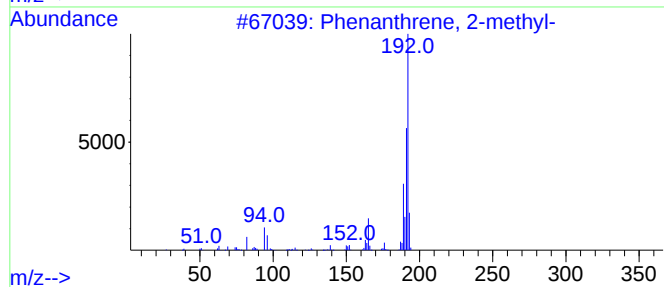
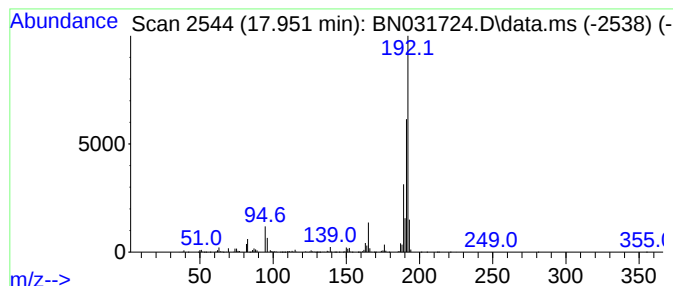
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.951	5.48 ng/ul	2124370	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



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Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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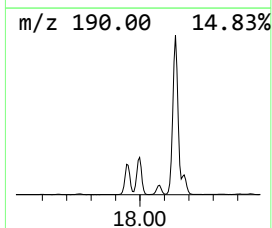
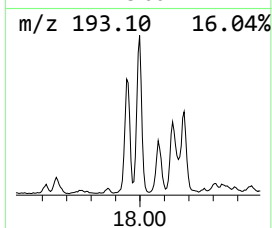
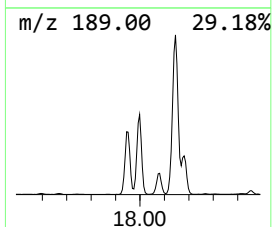
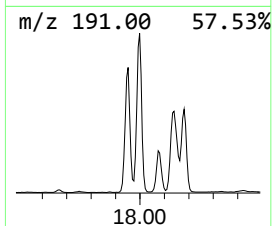
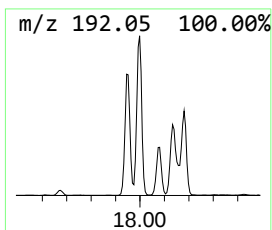
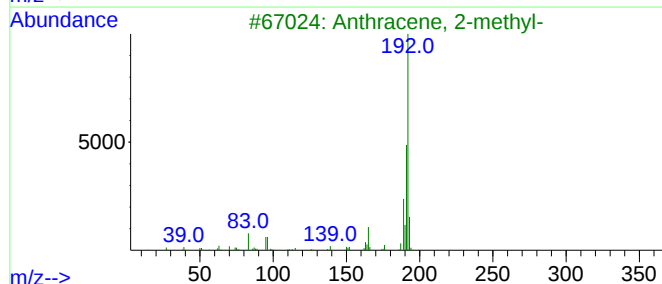
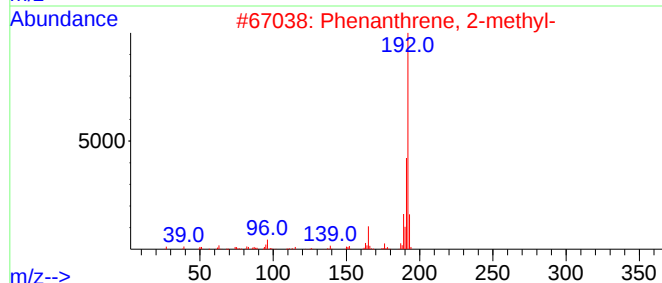
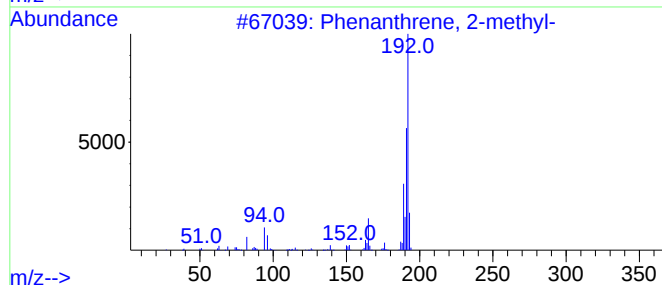
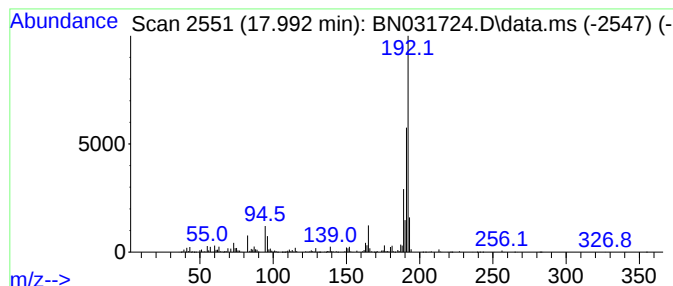
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.992	9.82 ng/ul	3803910	Phenanthrene-d10	17.075

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	96
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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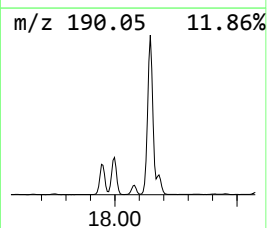
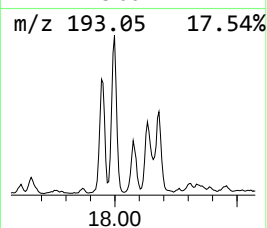
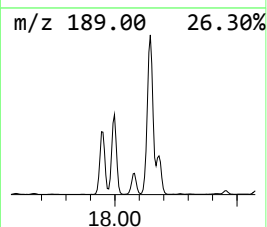
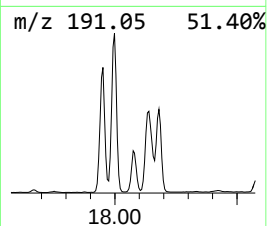
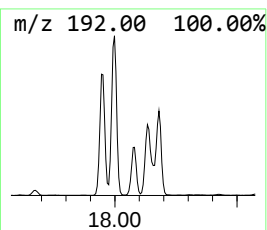
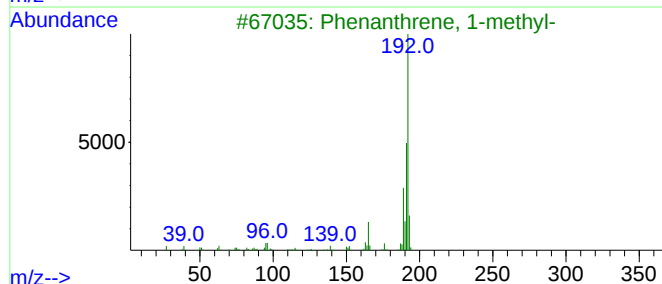
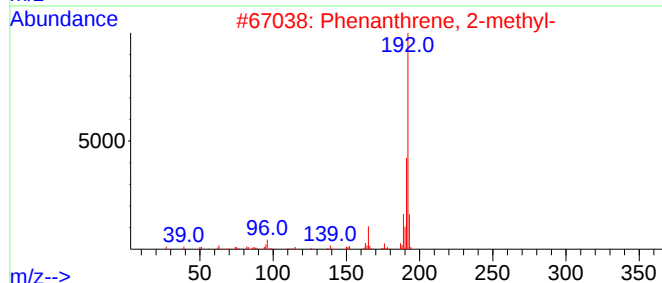
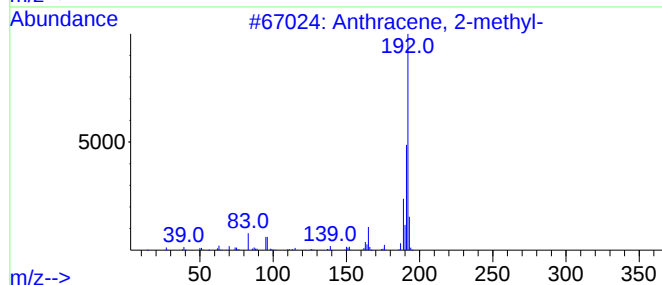
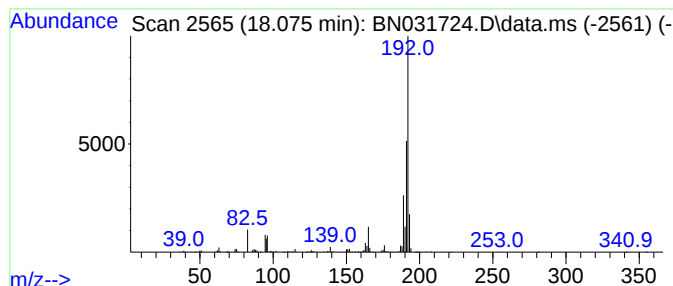
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.075	2.01 ng/ul	778611	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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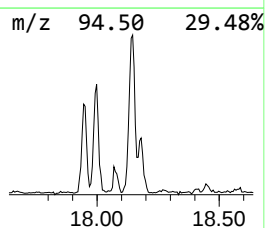
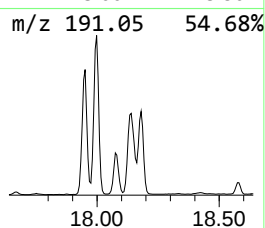
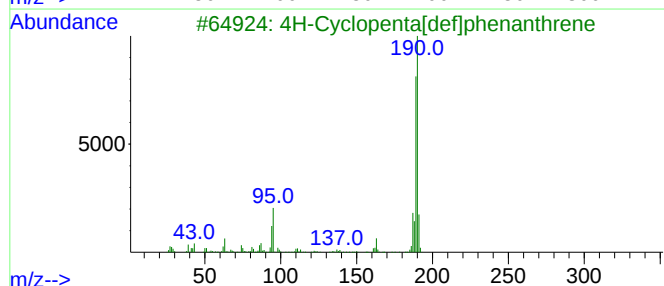
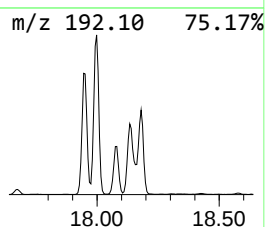
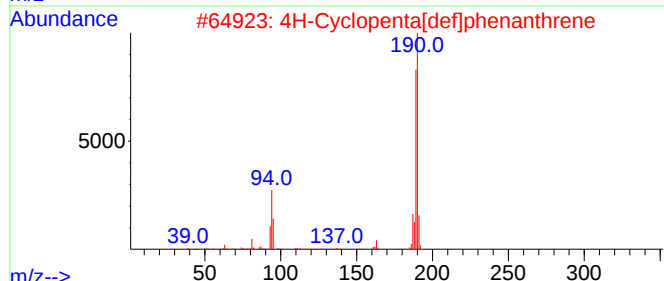
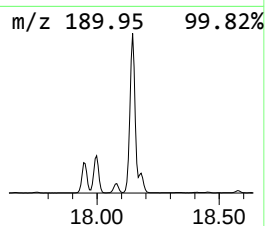
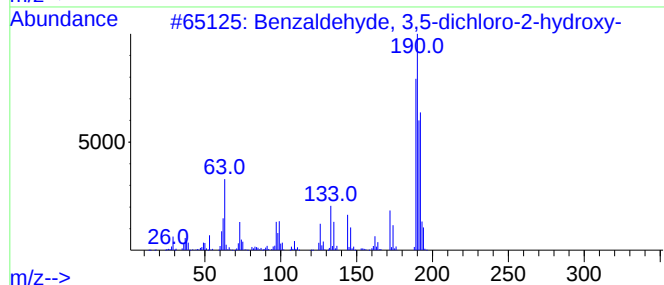
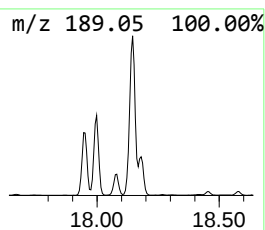
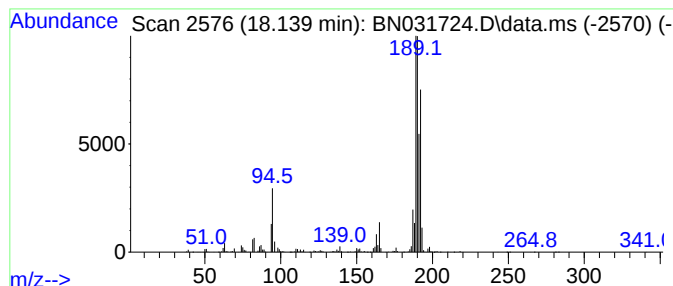
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzaldehyde, 3,5-dichloro-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	7.57 ng/ul	2934140	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
4			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
5			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
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Instrument :
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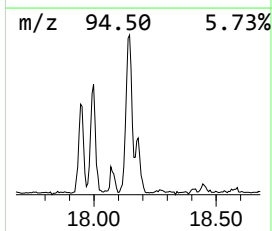
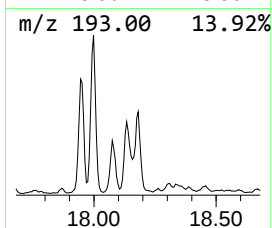
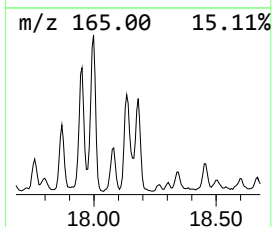
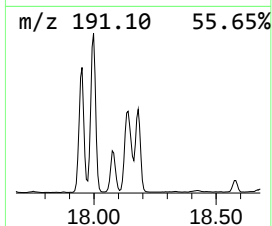
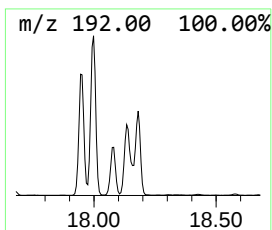
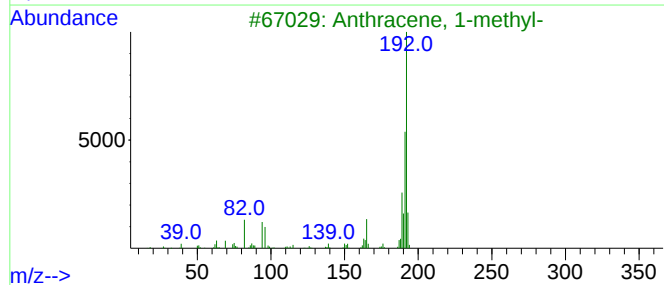
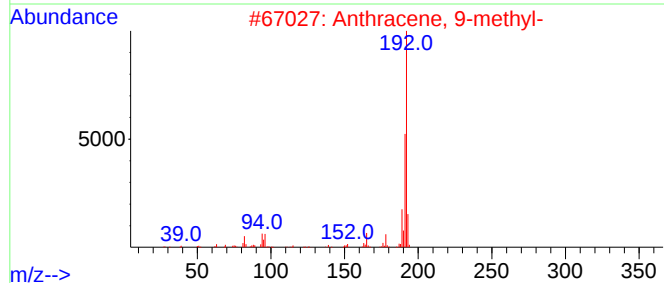
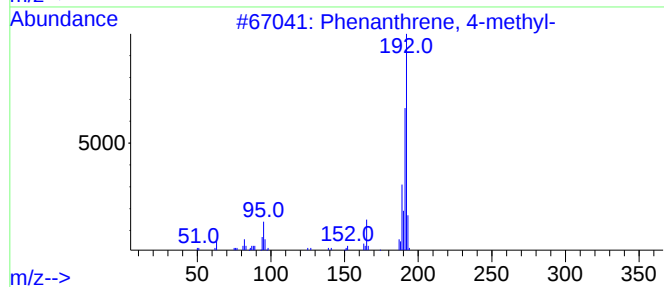
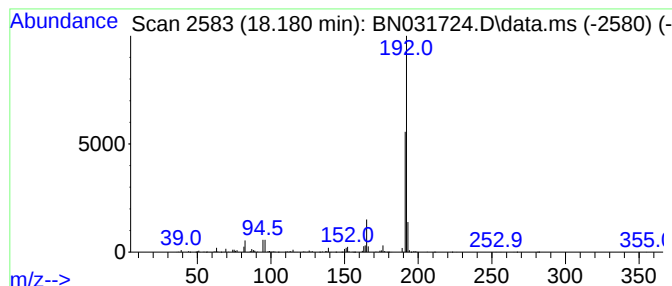
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenanthrene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	3.30 ng/ul	1279520	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
5			1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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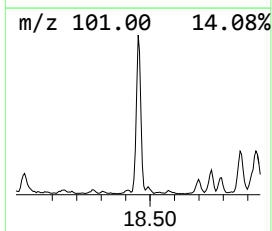
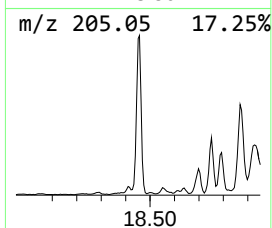
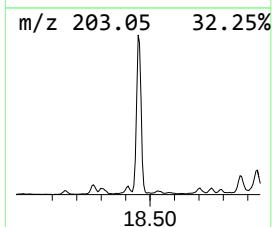
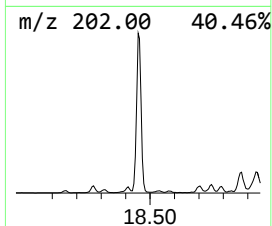
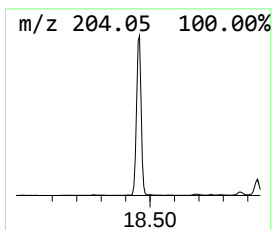
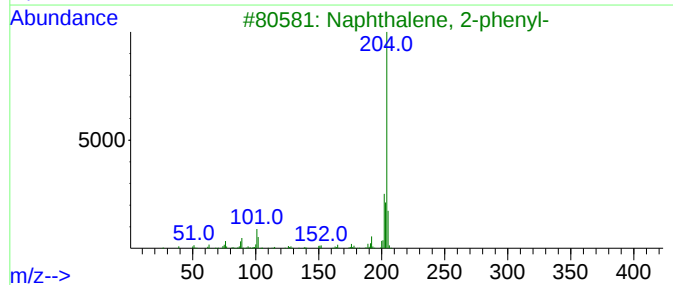
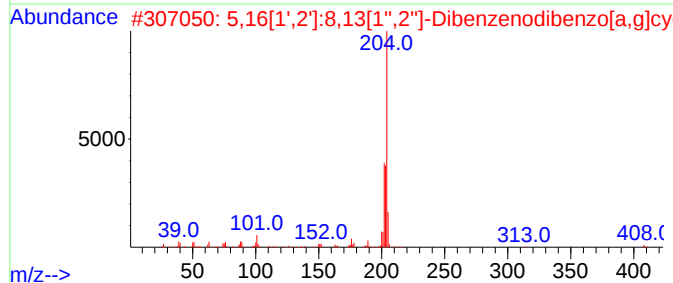
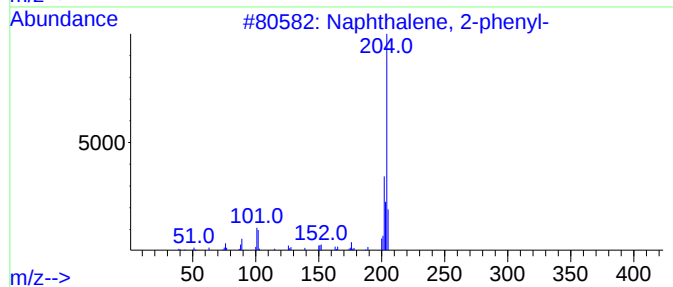
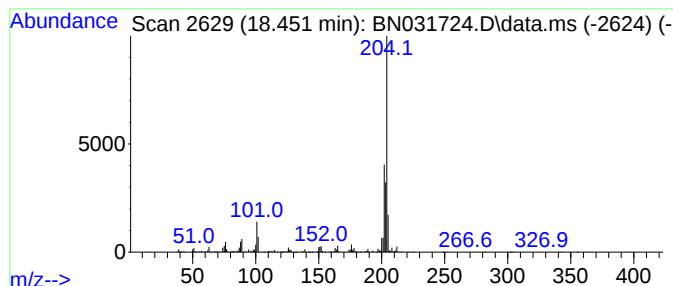
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.451	4.72 ng/ul	1827830	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	96
2			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91
3			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	89
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
5			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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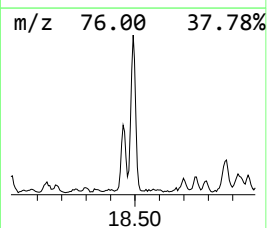
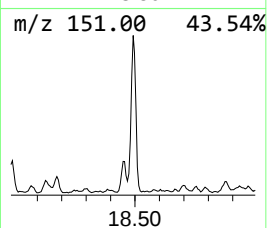
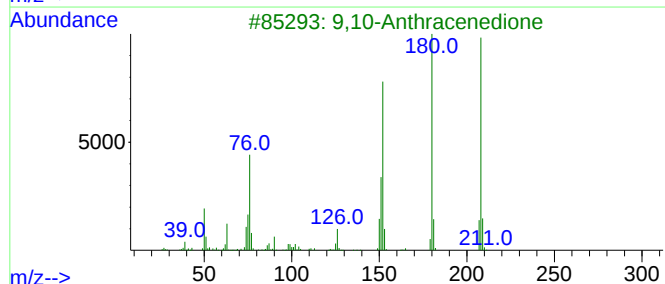
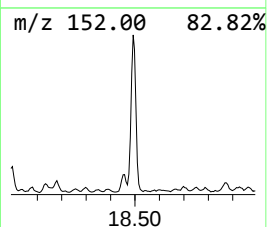
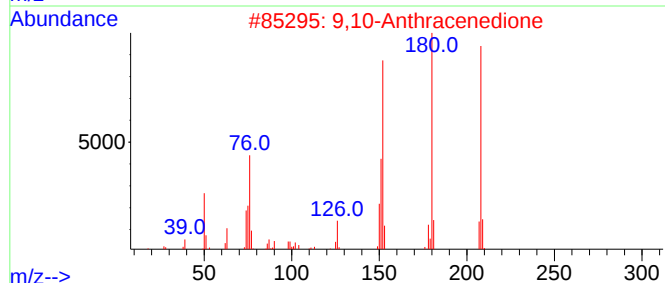
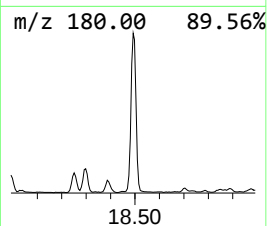
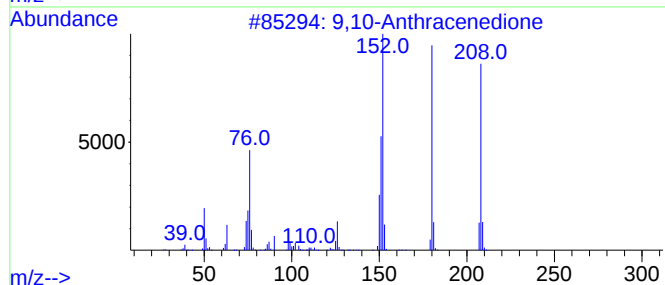
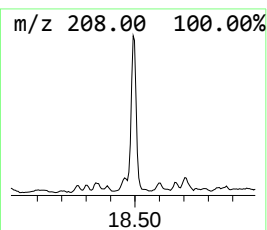
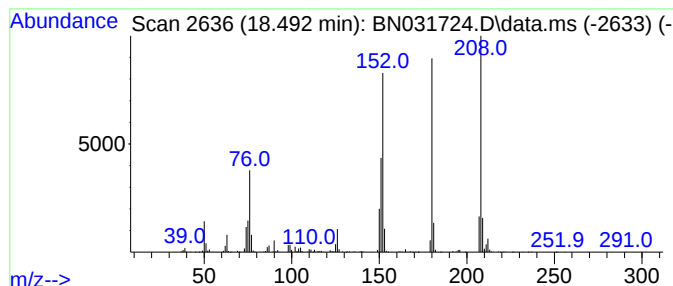
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.492	2.65 ng/ul	1028250	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
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Sample : P2569-21DL 2X
Misc :
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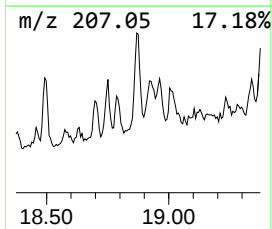
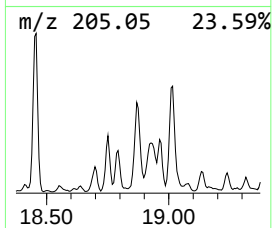
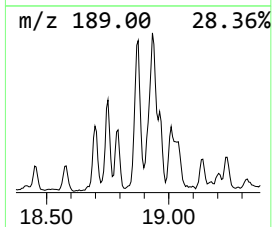
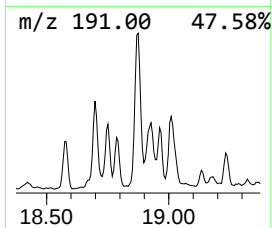
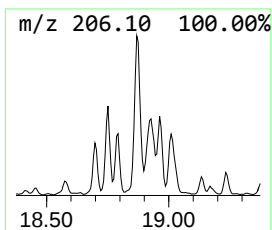
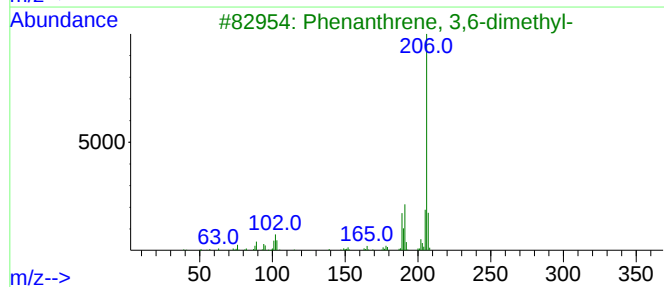
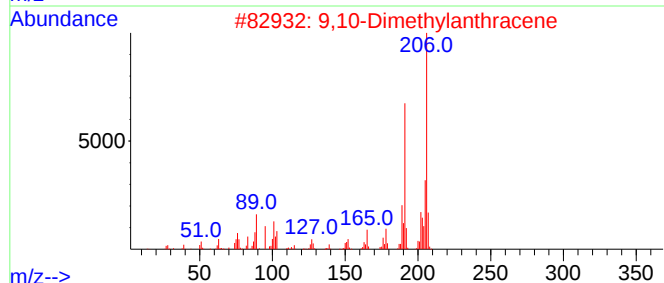
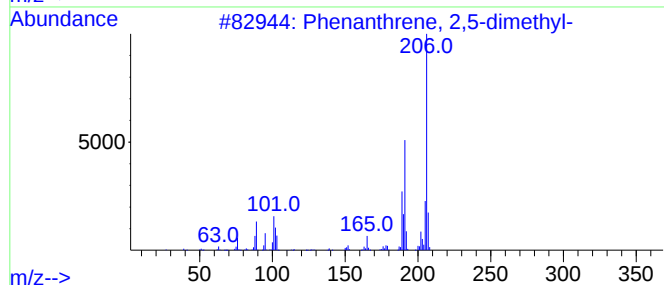
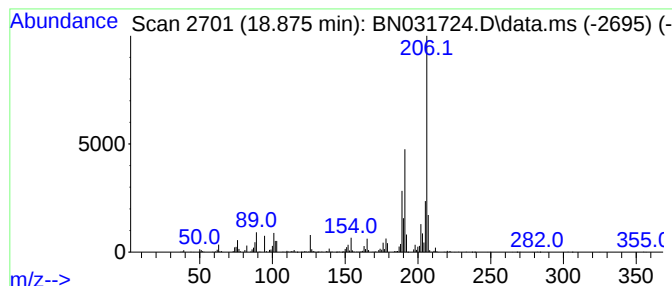
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ClientSampleId :
BH6A7DL

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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.875	2.91 ng/ul	1126480	Phenanthrene-d10		17.075	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	93
2	9,10-Dimethylanthracene		206	C16H14	000781-43-1	91
3	Phenanthrene, 3,6-dimethyl-		206	C16H14	001576-67-6	91
4	Phenanthrene, 3,6-dimethyl-		206	C16H14	001576-67-6	90
5	3,4-Dimethylphenanthrene		206	C16H14	1000452-24-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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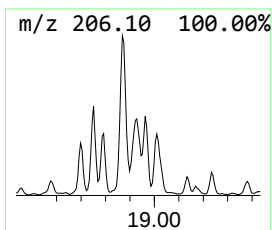
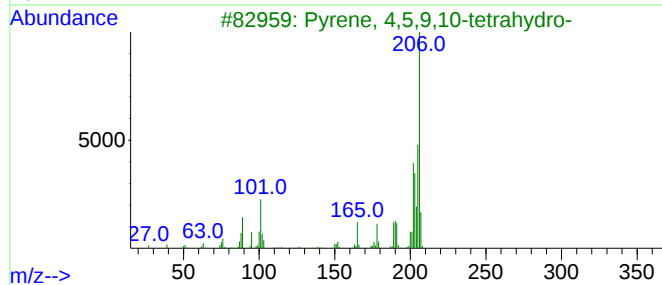
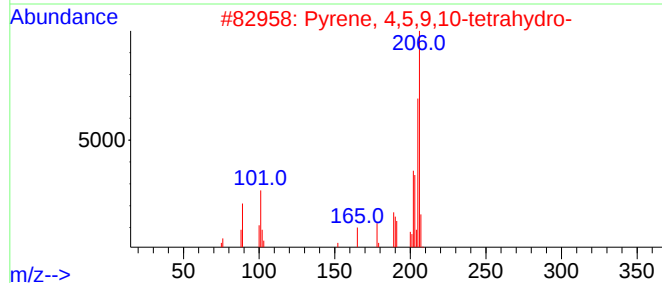
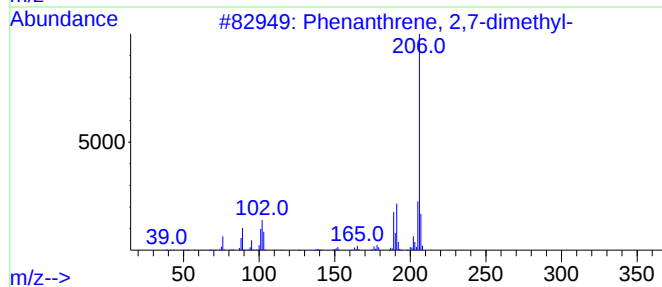
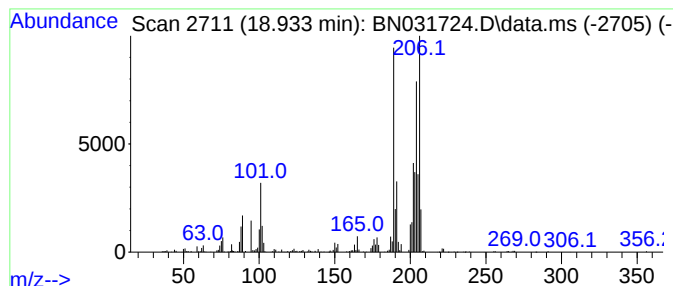
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.933	2.19 ng/ul	848110	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	43
2			Pyrene, 4,5,9,10-tetrahydro-	206	C16H14	000781-17-9	42
3			Pyrene, 4,5,9,10-tetrahydro-	206	C16H14	000781-17-9	38
4			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	38
5			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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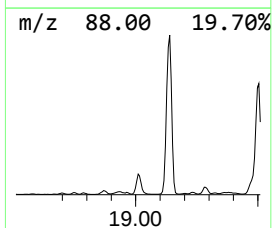
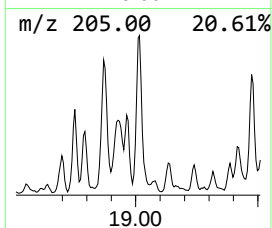
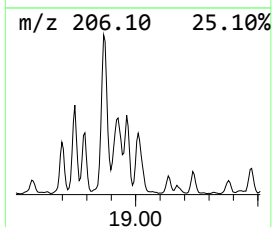
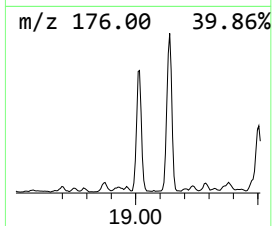
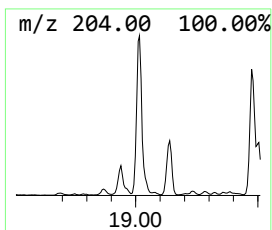
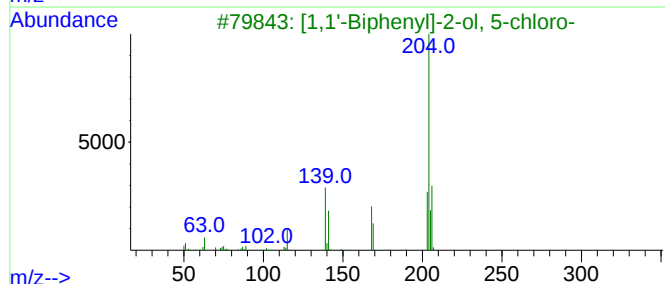
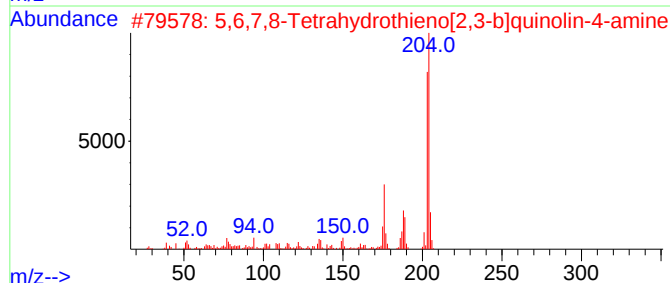
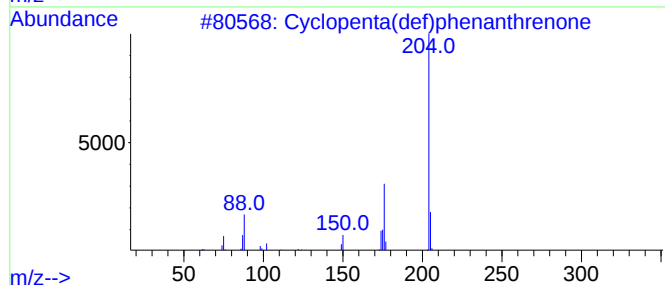
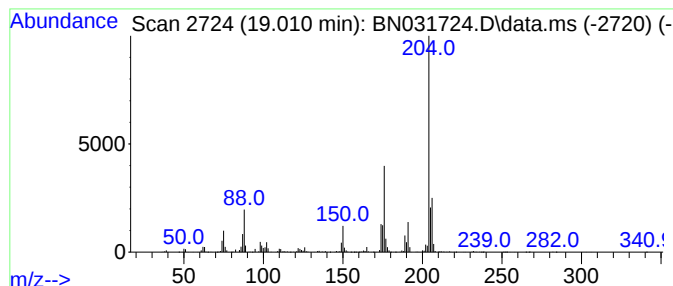
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.010	3.69 ng/ul	1429330	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	97
2			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	47
3			[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	43
4			2-Chloro-4-phenylphenol	204	C12H9ClO	000092-04-6	38
5			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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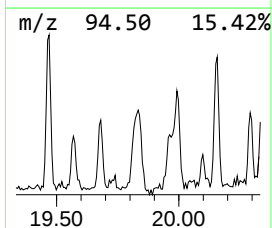
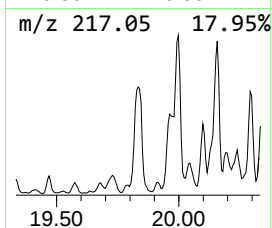
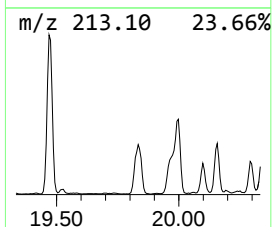
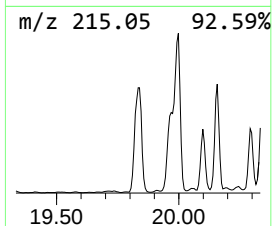
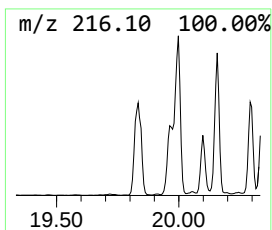
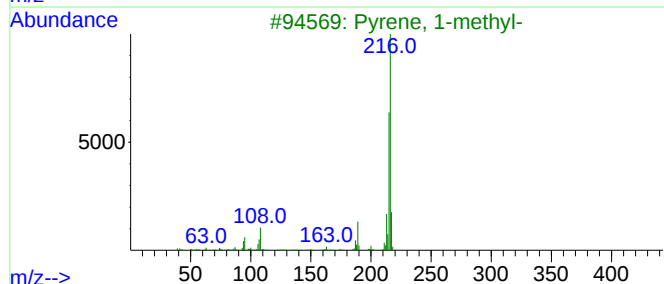
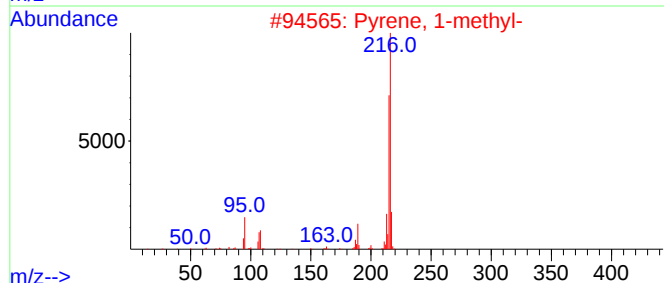
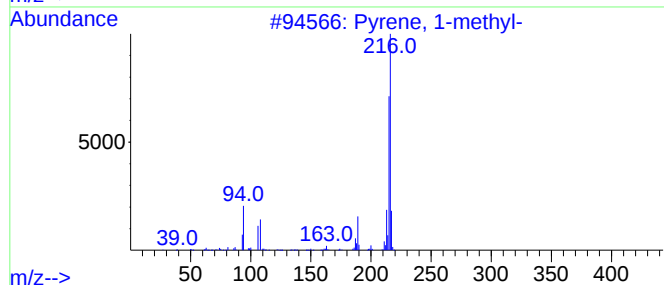
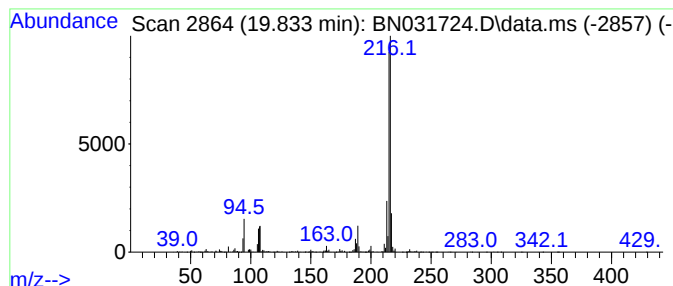
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Pyrene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.833	2.44 ng/ul	2104130	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-			216	C17H12	002381-21-7	98
2	Pyrene, 1-methyl-			216	C17H12	002381-21-7	96
3	Pyrene, 1-methyl-			216	C17H12	002381-21-7	95
4	Fluoranthene, 2-methyl-			216	C17H12	033543-31-6	92
5	Fluoranthene, 2-methyl-			216	C17H12	033543-31-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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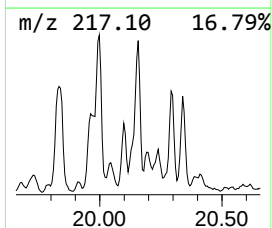
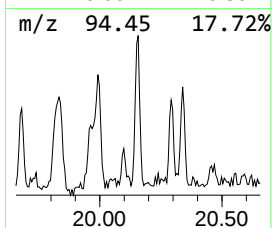
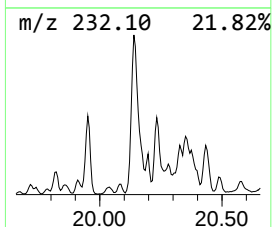
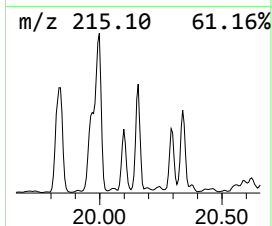
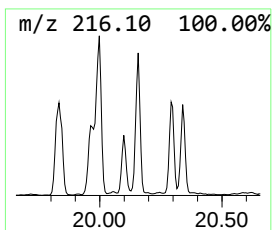
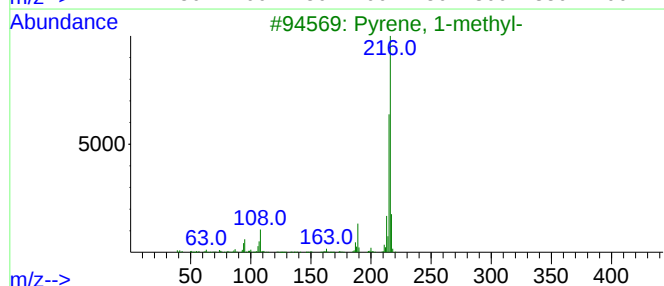
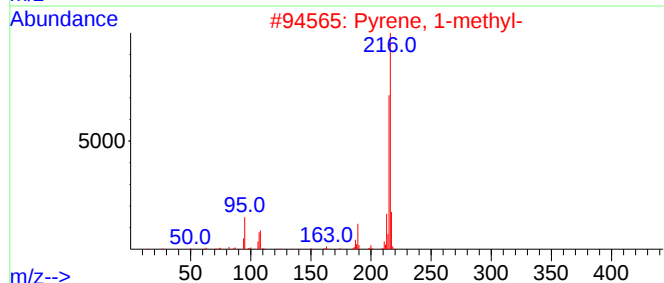
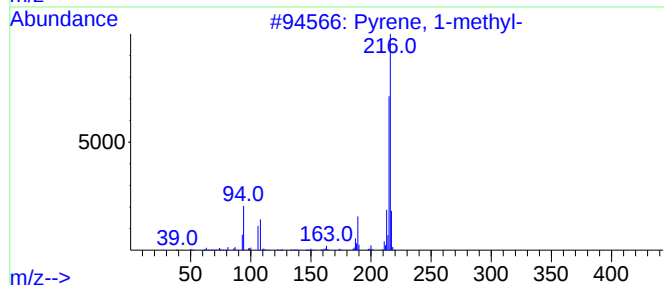
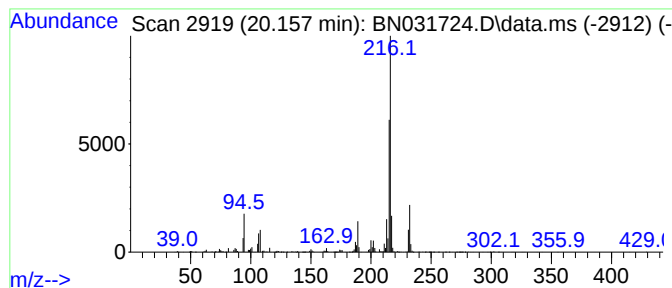
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Fluoranthene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.157	2.21 ng/ul	1908410	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-			216	C17H12	002381-21-7	98
2	Pyrene, 1-methyl-			216	C17H12	002381-21-7	96
3	Pyrene, 1-methyl-			216	C17H12	002381-21-7	96
4	Fluoranthene, 2-methyl-			216	C17H12	033543-31-6	95
5	Pyrene, 2-methyl-			216	C17H12	003442-78-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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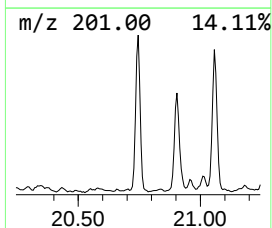
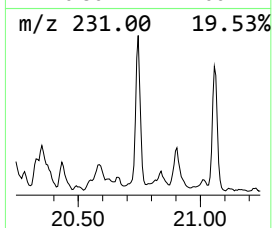
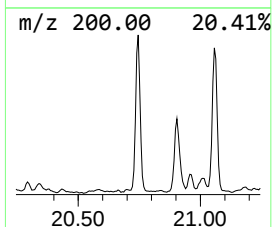
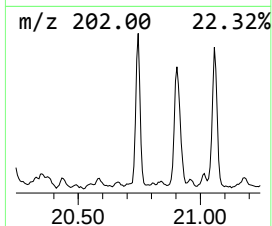
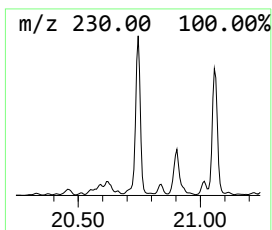
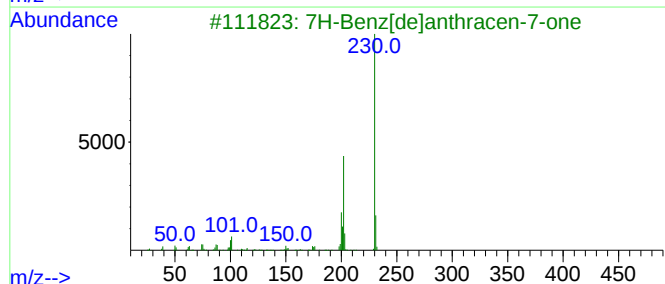
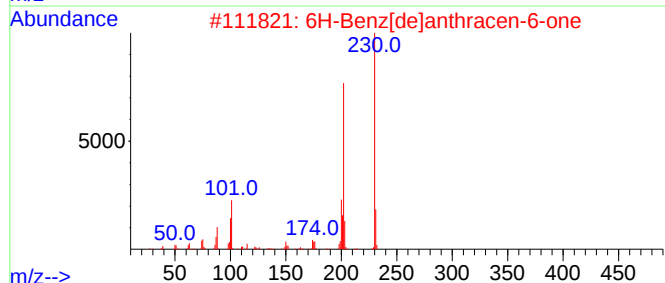
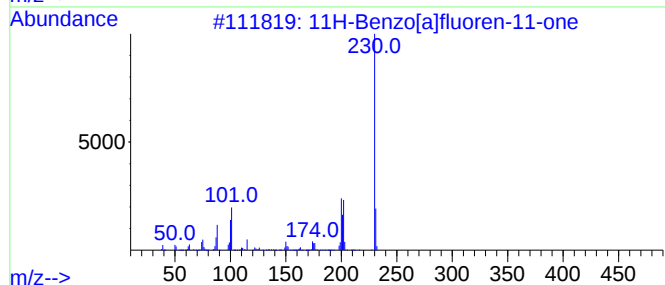
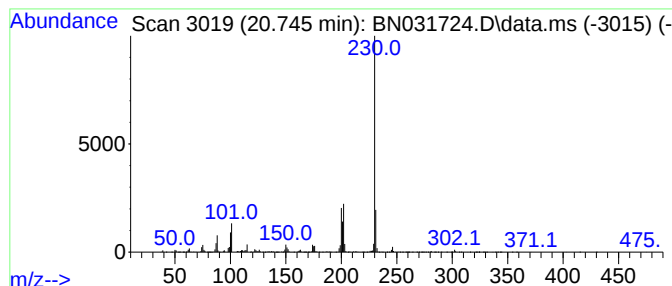
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 11H-Benzo[a]fluoren-11-one Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.745	2.19 ng/ul	1887860	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	98
2			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	91
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	91
4			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
5			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
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Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

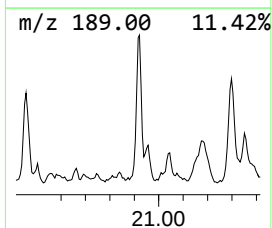
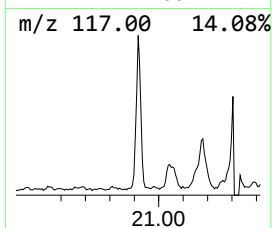
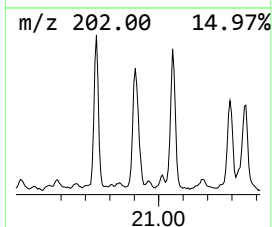
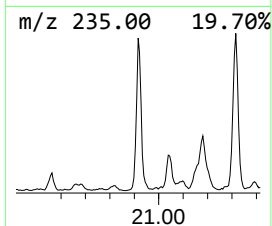
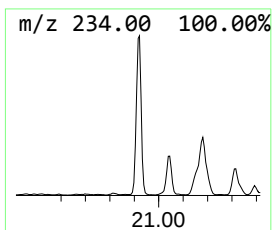
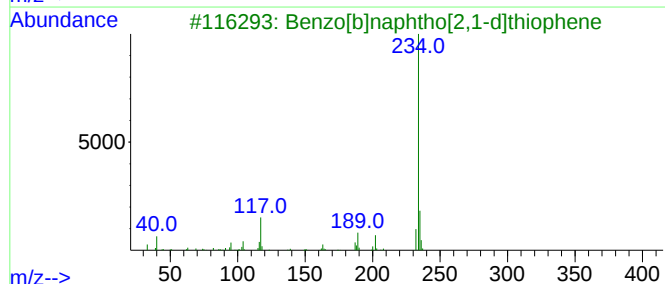
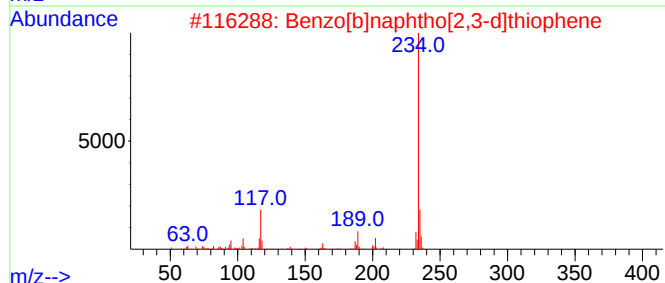
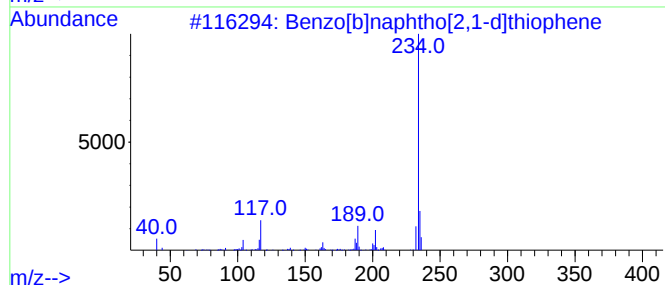
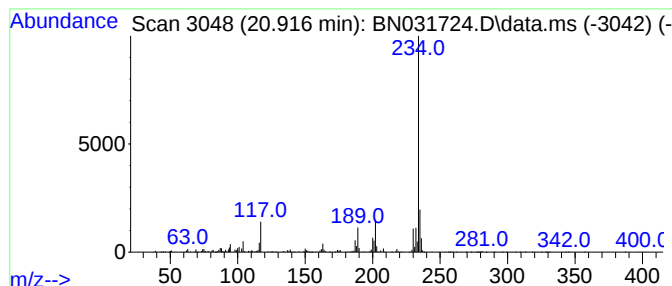
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.916	2.13 ng/ul	1842310	Chrysene-d12		21.310	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene		234	C16H10S	000239-35-0	96
2	Benzo[b]naphtho[2,3-d]thiophene		234	C16H10S	000243-46-9	95
3	Benzo[b]naphtho[2,1-d]thiophene		234	C16H10S	000239-35-0	94
4	Benzo[b]naphtho[2,1-d]thiophene		234	C16H10S	000239-35-0	94
5	Benzo[b]naphtho[2,3-d]thiophene		234	C16H10S	000243-46-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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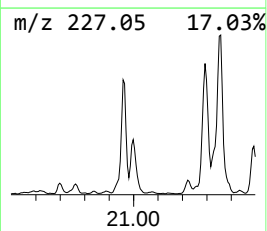
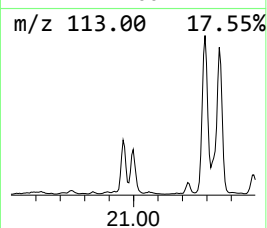
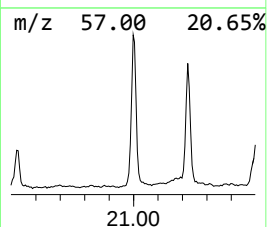
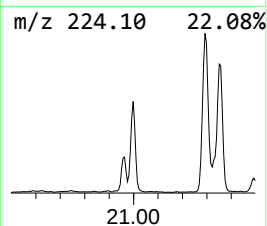
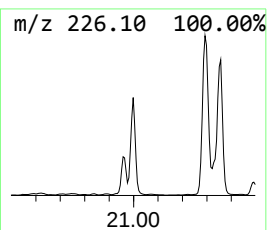
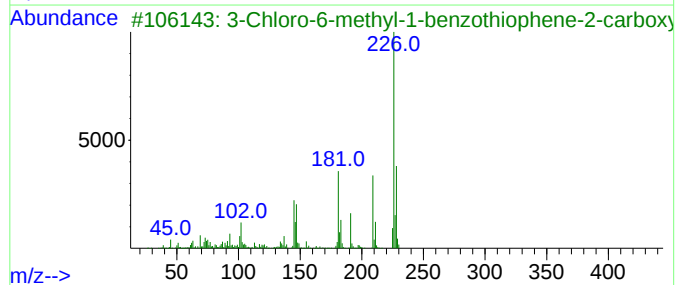
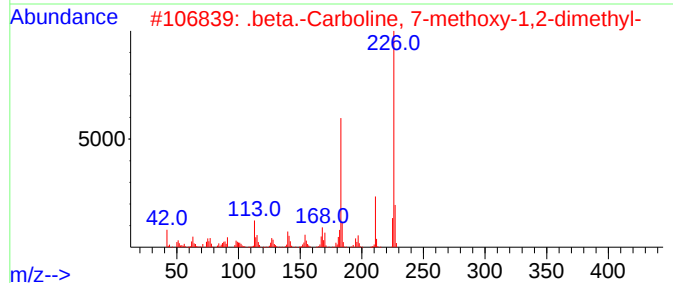
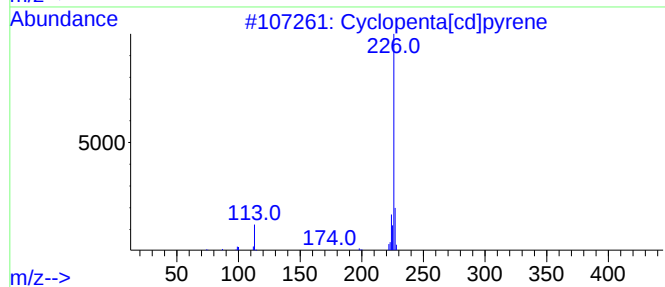
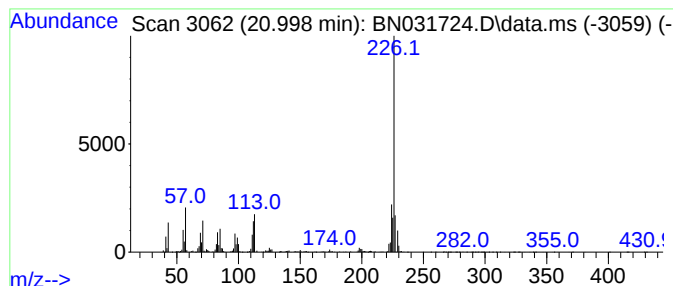
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Cyclopenta[cd]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.998	2.82 ng/ul	2434500	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	70
2			.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	50
3			3-Chloro-6-methyl-1-benzothiophe...	226	C10H7ClO2S	1010484-31-0	40
4			4-Naphthalen-2-yl-thiazol-2-ylamine	226	C13H10N2S	1000300-53-4	40
5			9H-Thioxanthen-9-one, 2-methyl-	226	C14H10O5	015774-82-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

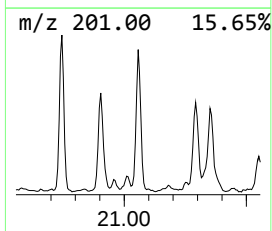
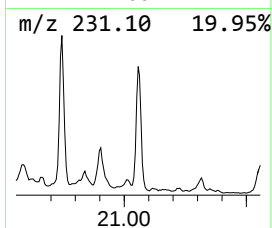
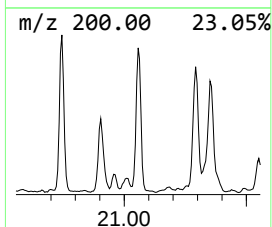
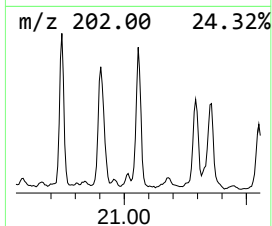
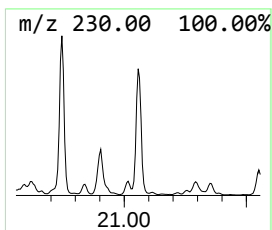
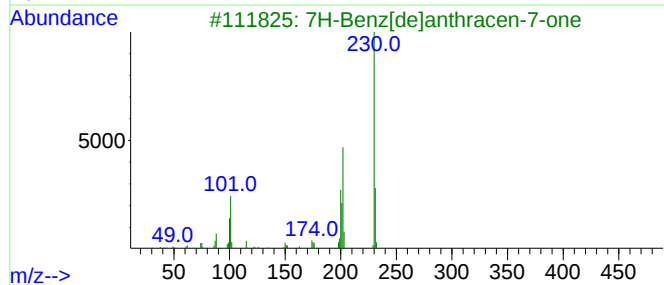
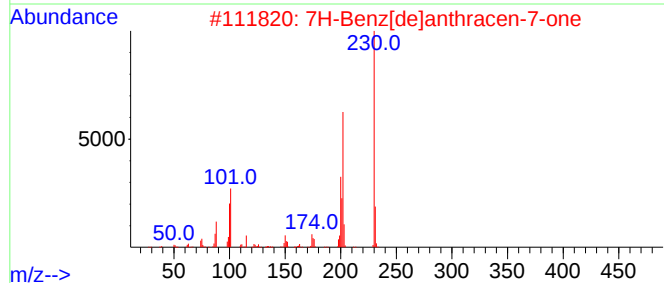
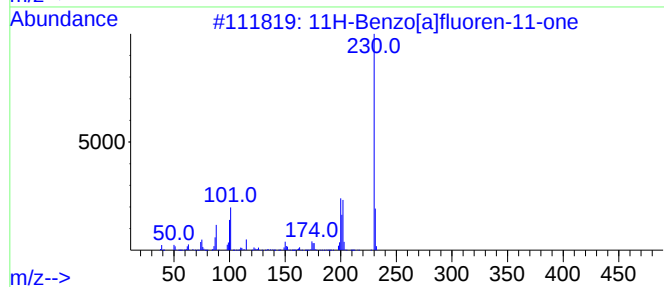
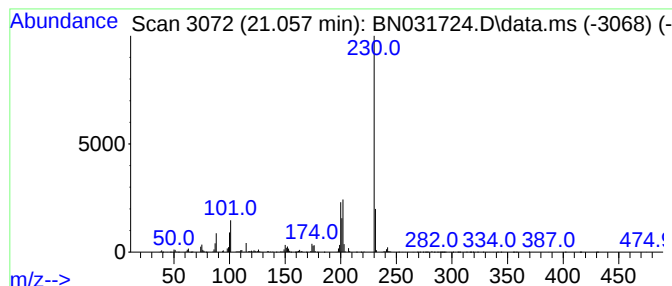
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 7H-Benz[de]anthracen-7-one Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.057	2.23 ng/ul	1920680	Chrysene-d12	21.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	95
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
3	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	74
5	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH6A7DL

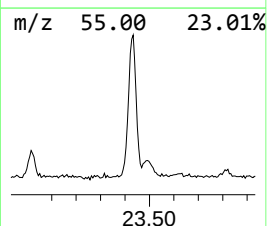
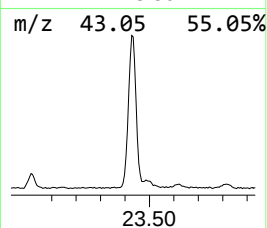
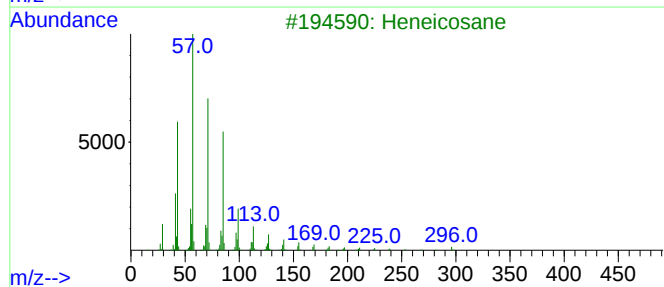
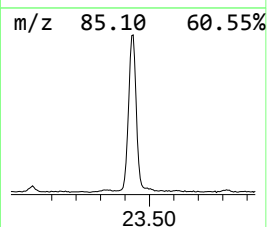
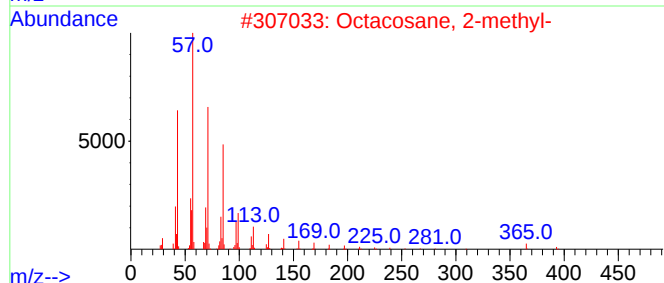
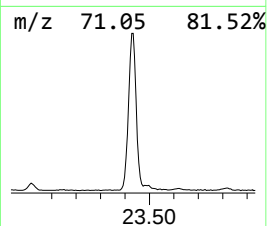
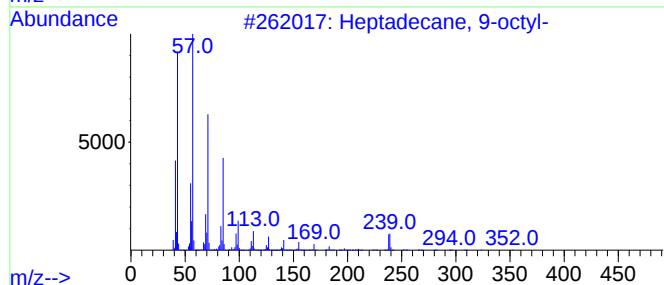
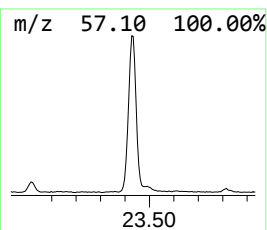
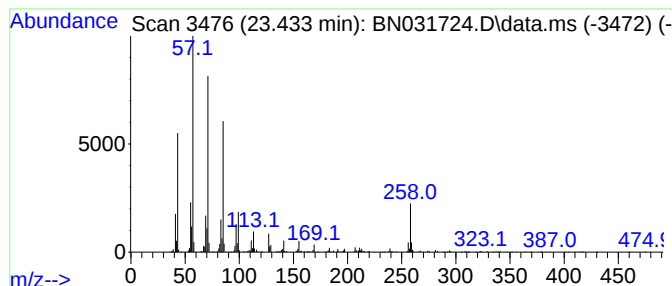
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.433	5.74 ng/ul	1844000	Perylene-d12	24.251

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	81
3		Heneicosane	296	C21H44	000629-94-7	81
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	74
5		Octadecane	254	C18H38	000593-45-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH6A7DL

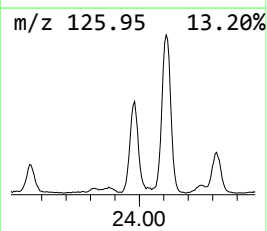
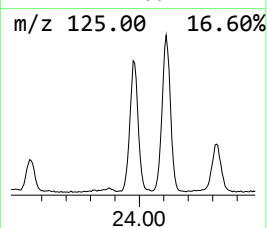
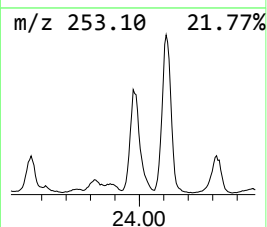
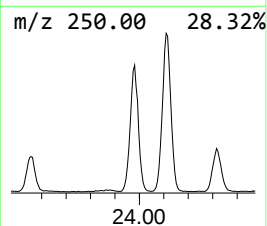
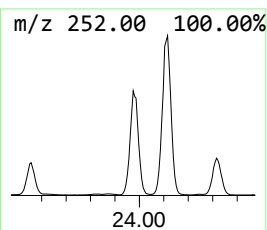
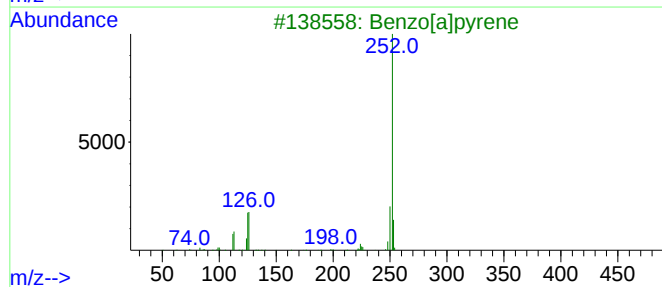
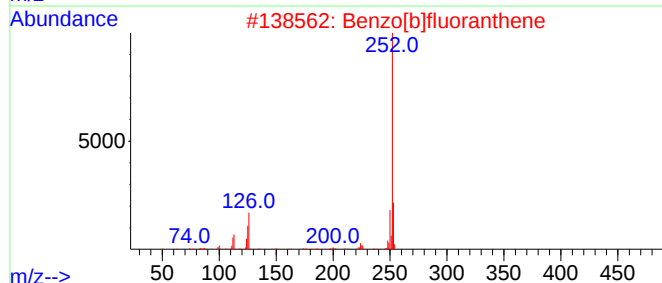
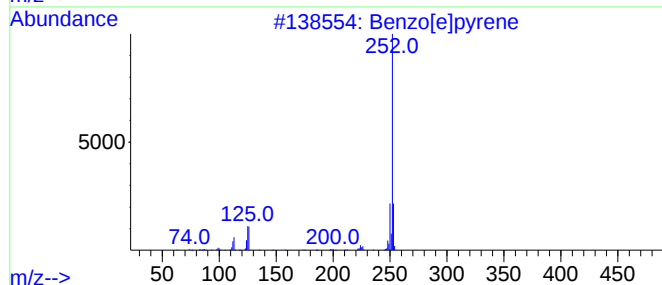
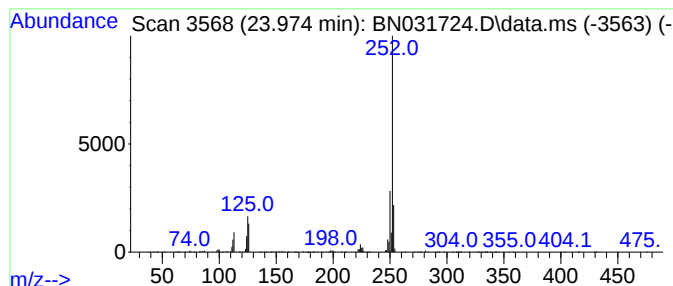
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.974	7.71 ng/ul	2477510	Perylene-d12	24.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	96
3			Benzo[a]pyrene	252	C20H12	000050-32-8	95
4			Benzo[a]pyrene	252	C20H12	000050-32-8	93
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
Data File : BN031724.D
Acq On : 30 May 2024 20:17
Operator : MA/JU
Sample : P2569-21DL 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH6A7DL

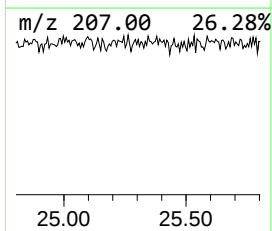
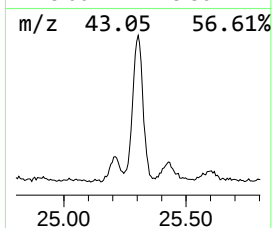
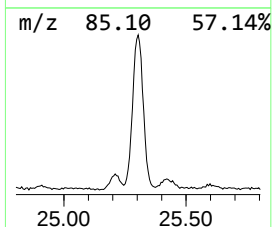
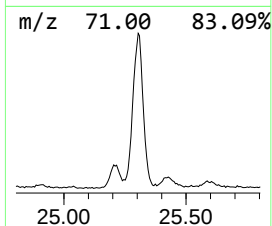
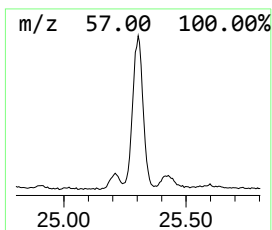
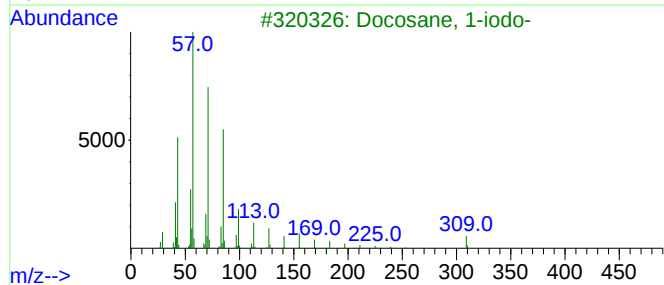
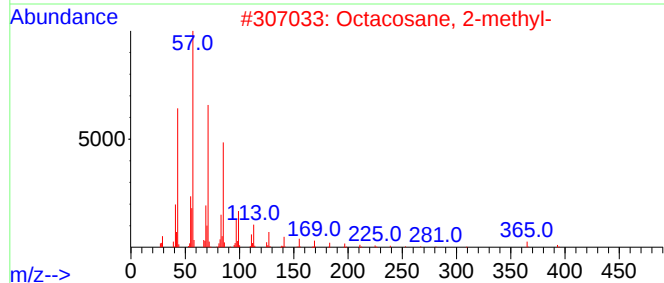
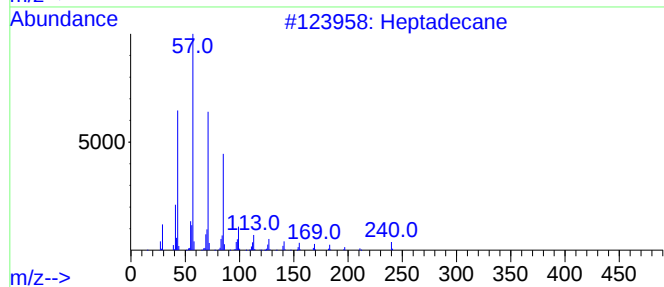
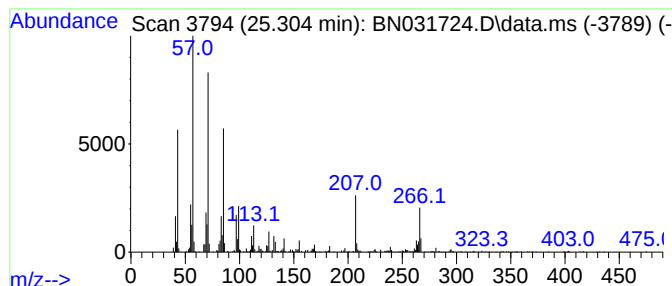
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.304	4.07 ng/ul	1308900	Perylene-d12	24.251

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	83
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	70
3		Docosane, 1-iodo-	436	C22H45I	1000406-31-9	70
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	64
5		Tetratriacontane, 2-methyl-	493	C35H72	014167-65-8	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN053024\
 Data File : BN031724.D
 Acq On : 30 May 2024 20:17
 Operator : MA/JU
 Sample : P2569-21DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BH6A7DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
9H-Fluoren-9-one	16.728	2.1	ng/ul	801454	4	17.075	7748120	20.0
Dibenzothiophene	16.892	2.6	ng/ul	1022200	4	17.075	7748120	20.0
Phenanthrene, 2...	17.951	5.5	ng/ul	2124370	4	17.075	7748120	20.0
Anthracene, 2-m...	17.992	9.8	ng/ul	3803910	4	17.075	7748120	20.0
Phenanthrene, 1...	18.075	2.0	ng/ul	778611	4	17.075	7748120	20.0
Benzaldehyde, 3...	18.139	7.6	ng/ul	2934140	4	17.075	7748120	20.0
Phenanthrene, 4...	18.180	3.3	ng/ul	1279520	4	17.075	7748120	20.0
Naphthalene, 2-...	18.451	4.7	ng/ul	1827830	4	17.075	7748120	20.0
9,10-Anthracene...	18.492	2.6	ng/ul	1028250	4	17.075	7748120	20.0
Phenanthrene, 2...	18.875	2.9	ng/ul	1126480	4	17.075	7748120	20.0
unknown-01	18.933	2.2	ng/ul	848110	4	17.075	7748120	20.0
Cyclopenta(def)...	19.010	3.7	ng/ul	1429330	4	17.075	7748120	20.0
Pyrene, 1-methyl-	19.833	2.4	ng/ul	2104130	5	21.310	17263500	20.0
Fluoranthene, 2...	20.157	2.2	ng/ul	1908410	5	21.310	17263500	20.0
11H-Benzo[a]flu...	20.745	2.2	ng/ul	1887860	5	21.310	17263500	20.0
Benzo[b]naphtho...	20.916	2.1	ng/ul	1842310	5	21.310	17263500	20.0
Cyclopenta[cd]p...	20.998	2.8	ng/ul	2434500	5	21.310	17263500	20.0
7H-Benz[de]anth...	21.057	2.2	ng/ul	1920680	5	21.310	17263500	20.0
(DEL) Alkane: S...	23.433	5.7	ng/ul	1844000	6	24.251	6426640	20.0
Benzo[e]pyrene	23.974	7.7	ng/ul	2477510	6	24.251	6426640	20.0
(DEL) Alkane: S...	25.304	4.1	ng/ul	1308900	6	24.251	6426640	20.0