

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
 Data File : BN031806.D
 Acq On : 05 Jun 2024 19:05
 Operator : MA/JU
 Sample : P2591-15DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBQ6DL

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: BN031806.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	98916	178301	1.52%	0.166%
2	3.693	116	120	127	rVV	27876	48848	0.42%	0.046%
3	6.846	643	656	669	rVB	306497	530421	4.53%	0.494%
4	7.011	677	684	698	rVB	226364	374287	3.20%	0.349%
5	7.205	710	717	727	rVB	294346	473724	4.05%	0.441%
6	7.669	789	796	804	rBV	1130584	1873023	16.00%	1.745%
7	8.375	907	916	924	rBV	363141	618036	5.28%	0.576%
8	8.828	983	993	1000	rBV	272864	466196	3.98%	0.434%
9	9.399	1085	1090	1097	rVB	33005	56350	0.48%	0.052%
10	9.546	1108	1115	1128	rVB	218321	378200	3.23%	0.352%
11	10.081	1199	1206	1216	rBV	371558	668068	5.71%	0.622%
12	10.458	1261	1270	1276	rBV	1789225	3081667	26.32%	2.871%
13	10.505	1276	1278	1284	rVV	54926	76769	0.66%	0.072%
14	10.599	1287	1294	1310	rVV	237406	461530	3.94%	0.430%
15	11.205	1390	1397	1401	rBV	40787	76341	0.65%	0.071%
16	12.122	1548	1553	1564	rVB	37567	67567	0.58%	0.063%
17	12.346	1586	1591	1598	rVB	36240	60034	0.51%	0.056%
18	13.634	1803	1810	1815	rBV2	37536	69272	0.59%	0.065%
19	13.728	1821	1826	1838	rVB	722592	1026223	8.77%	0.956%
20	14.010	1861	1874	1884	rBV	829556	1370698	11.71%	1.277%
21	14.316	1917	1926	1932	rBV	2919312	4430489	37.85%	4.128%
22	14.381	1932	1937	1945	rVB	162456	258859	2.21%	0.241%
23	14.522	1954	1961	1971	rBV	314773	560315	4.79%	0.522%
24	14.681	1983	1988	1990	rBV2	32914	52422	0.45%	0.049%
25	14.716	1990	1994	2003	rVV	156803	349730	2.99%	0.326%
26	15.110	2054	2061	2065	rBV4	17556	42681	0.36%	0.040%
27	15.310	2089	2095	2101	rVV	1118857	1680692	14.36%	1.566%
28	15.363	2101	2104	2111	rVV	192432	284912	2.43%	0.265%
29	15.440	2111	2117	2123	rVV	238469	347285	2.97%	0.324%
30	15.663	2150	2155	2164	rVB	88144	145544	1.24%	0.136%
31	15.810	2175	2180	2187	rVB	87476	166058	1.42%	0.155%
32	16.040	2213	2219	2225	rVB4	78115	126565	1.08%	0.118%
33	16.198	2238	2246	2252	rVV4	25525	59600	0.51%	0.056%
34	16.345	2264	2271	2272	rBV2	33719	55589	0.47%	0.052%
35	16.375	2273	2276	2280	rVV2	44838	66830	0.57%	0.062%
36	16.475	2288	2293	2298	rBV	79710	123457	1.05%	0.115%

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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

37	16.598	2306	2314	2317	rBV2	49094	86276	0.74%	0.080%
38	16.640	2318	2321	2324	rVB3	34979	44841	0.38%	0.042%
39	16.722	2330	2335	2342	rVB	318059	462236	3.95%	0.431%
40	16.798	2343	2348	2355	rVV2	46700	126145	1.08%	0.118%
41	16.881	2358	2362	2370	rVB	182291	274494	2.34%	0.256%
42	16.998	2376	2382	2386	rBV	134580	182858	1.56%	0.170%
43	17.063	2387	2393	2397	rVV	3694961	5504412	47.02%	5.128%
44	17.110	2397	2401	2406	rVV	3468086	5044660	43.09%	4.700%
45	17.163	2406	2410	2413	rVV	1296070	1825221	15.59%	1.700%
46	17.198	2413	2416	2421	rVV	677287	916717	7.83%	0.854%
47	17.251	2421	2425	2431	rVB3	72386	134845	1.15%	0.126%
48	17.375	2444	2446	2452	rVB2	32479	49946	0.43%	0.047%
49	17.469	2452	2462	2474	rBV2	463660	937875	8.01%	0.874%
50	17.587	2478	2482	2487	rVV3	56650	88990	0.76%	0.083%
51	17.645	2488	2492	2501	rVB3	87645	150431	1.29%	0.140%
52	17.740	2505	2508	2512	rBV4	36756	47929	0.41%	0.045%
53	17.781	2512	2515	2522	rVB3	31554	49131	0.42%	0.046%
54	17.857	2525	2528	2533	rVB	45582	63202	0.54%	0.059%
55	17.939	2533	2542	2544	rBV	396530	591296	5.05%	0.551%
56	17.987	2544	2550	2555	rVV2	630557	1264293	10.80%	1.178%
57	18.039	2555	2559	2561	rVV	81466	105116	0.90%	0.098%
58	18.069	2561	2564	2568	rVB	146791	191650	1.64%	0.179%
59	18.134	2569	2575	2579	rBV2	472942	863073	7.37%	0.804%
60	18.169	2579	2581	2588	rVB	248232	310133	2.65%	0.289%
61	18.245	2591	2594	2598	rBV4	45525	74974	0.64%	0.070%
62	18.292	2598	2602	2606	rVV	65901	85839	0.73%	0.080%
63	18.439	2623	2627	2631	rVV2	485754	678696	5.80%	0.632%
64	18.487	2631	2635	2643	rVV2	425807	591196	5.05%	0.551%
65	18.692	2666	2670	2674	rBV2	87049	110325	0.94%	0.103%
66	18.739	2675	2678	2682	rVB	78020	84735	0.72%	0.079%
67	18.781	2682	2685	2689	rVB2	87320	106106	0.91%	0.099%
68	18.834	2690	2694	2695	rBV3	64475	83137	0.71%	0.077%
69	18.863	2695	2699	2704	rVB2	189044	310885	2.66%	0.290%
70	19.004	2718	2723	2729	rBV2	396338	617485	5.27%	0.575%
71	19.128	2737	2744	2752	rBV	6271817	8408104	71.82%	7.833%
72	19.192	2752	2755	2758	rBV	69732	94424	0.81%	0.088%
73	19.228	2758	2761	2763	rVV	125469	170501	1.46%	0.159%
74	19.251	2763	2765	2773	rVB3	175329	297976	2.55%	0.278%
75	19.369	2783	2785	2789	rVB	107127	120942	1.03%	0.113%
76	19.463	2795	2801	2803	rBV2	2341013	3447591	29.45%	3.212%

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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

77	19.492	2803	2806	2813	rVV	4152630	5356292	45.75%	4.990%
78	19.563	2814	2818	2824	rVB2	252494	369060	3.15%	0.344%
79	19.675	2832	2837	2841	rVB	273144	381374	3.26%	0.355%
80	19.734	2843	2847	2852	rVB	285023	420216	3.59%	0.391%
81	19.816	2855	2861	2862	rBV2	316469	508637	4.34%	0.474%
82	19.951	2879	2884	2887	rBV5	259093	501500	4.28%	0.467%
83	19.986	2887	2890	2894	rVB	330079	459974	3.93%	0.429%
84	20.086	2904	2907	2911	rBV2	125248	157876	1.35%	0.147%
85	20.145	2911	2917	2921	rBV2	434098	771918	6.59%	0.719%
86	20.286	2939	2941	2944	rVB	180718	178497	1.52%	0.166%
87	20.333	2946	2949	2953	rVB3	197521	264541	2.26%	0.246%
88	20.739	3013	3018	3023	rBV	984347	1427838	12.20%	1.330%
89	20.904	3041	3046	3050	rBV2	514212	966570	8.26%	0.900%
90	20.951	3050	3054	3057	rVV	419624	602577	5.15%	0.561%
91	20.986	3057	3060	3066	rVV2	638632	1232227	10.53%	1.148%
92	21.051	3067	3071	3083	rVB	874653	1349804	11.53%	1.258%
93	21.210	3095	3098	3101	rBV2	229310	240083	2.05%	0.224%
94	21.298	3105	3113	3117	rVV3	5303789	11706675	100.00%	10.906%
95	21.345	3117	3121	3132	rVB	2572037	4022590	34.36%	3.748%
96	23.310	3448	3455	3462	rBV	1923740	5711332	48.79%	5.321%
97	24.033	3574	3578	3583	rVB2	469383	834336	7.13%	0.777%
98	24.098	3584	3589	3602	rVB	903765	2188091	18.69%	2.038%
99	24.239	3605	3613	3623	rBV	2742876	7010689	59.89%	6.531%
100	27.492	4158	4166	4197	rVB4	944056	5371774	45.89%	5.004%

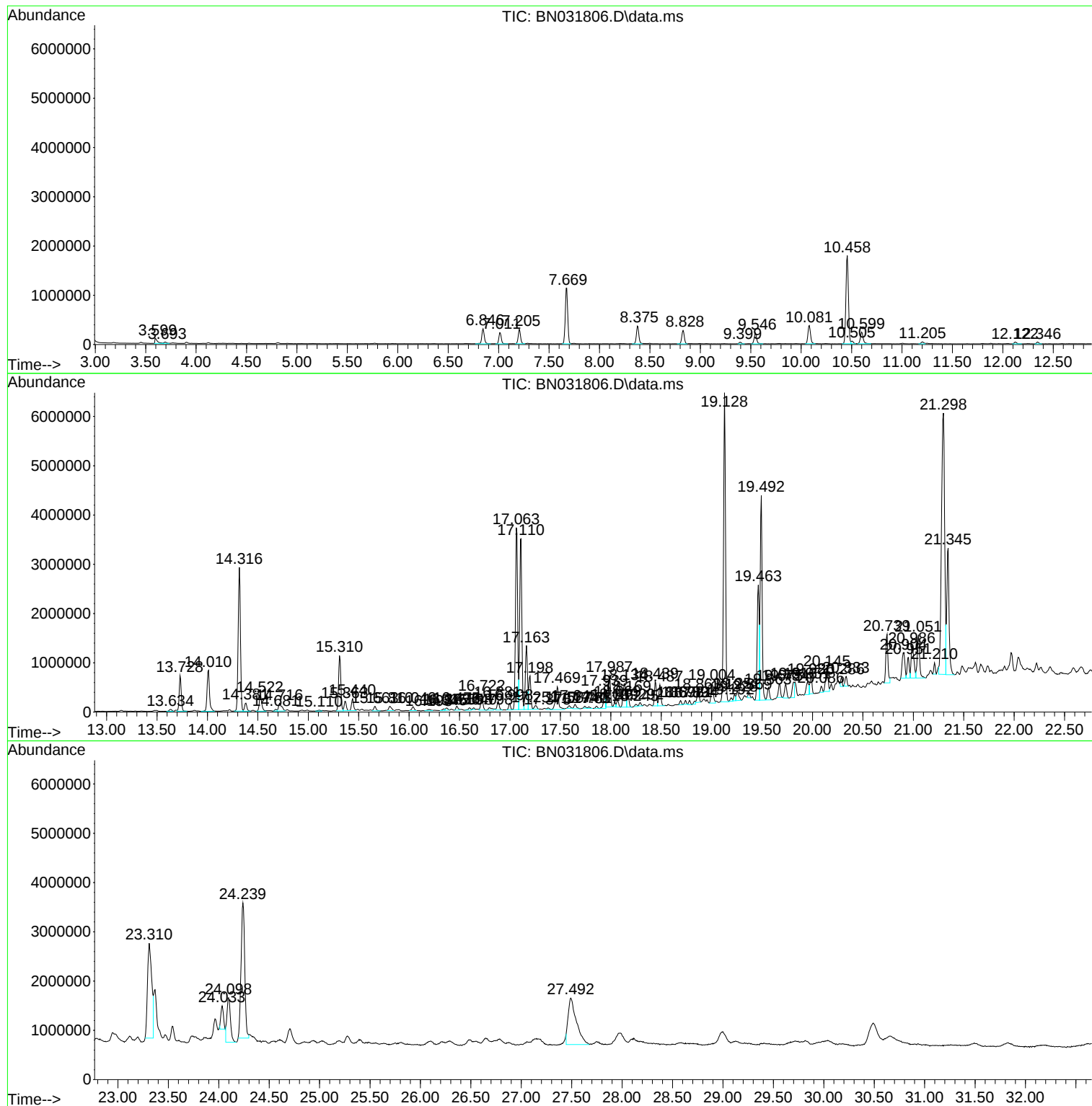
Sum of corrected areas: 107339740

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Instrument :
BNA_N
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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
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Sample : P2591-15DL 5X
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
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BHBQ6DL

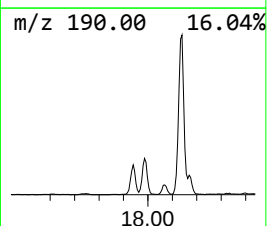
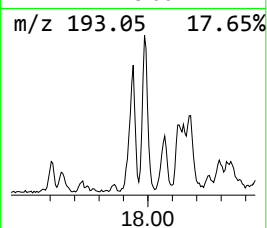
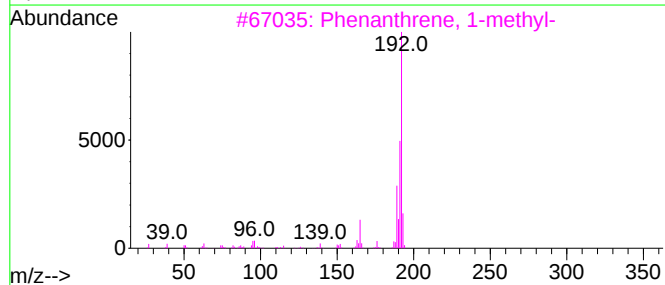
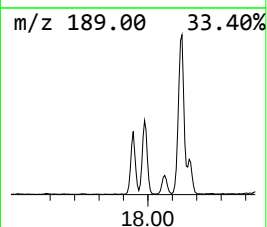
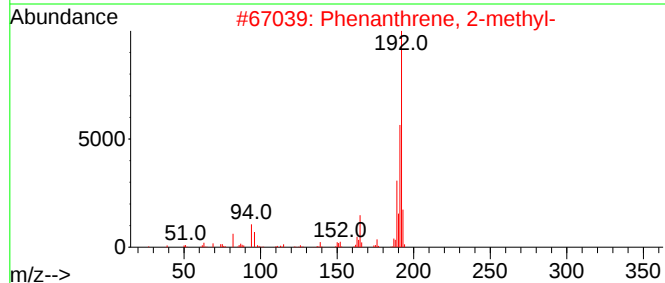
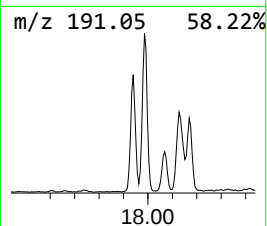
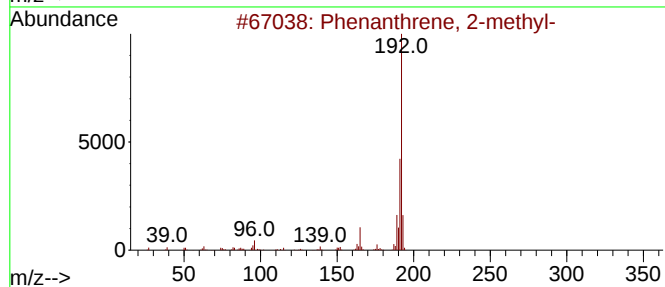
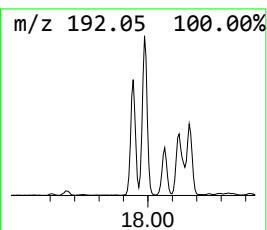
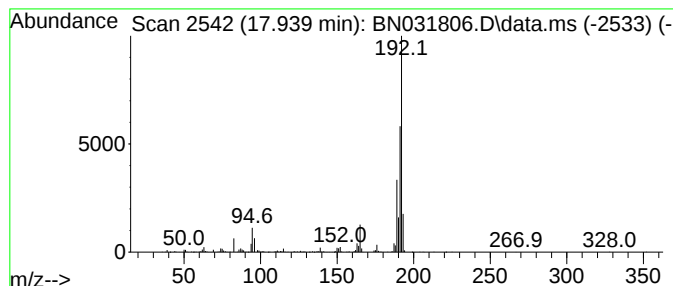
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.939	2.15 ng/ul	591296	Phenanthrene-d10	17.069

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91



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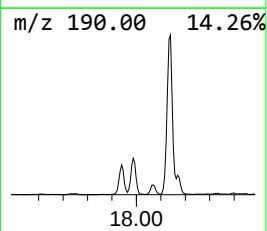
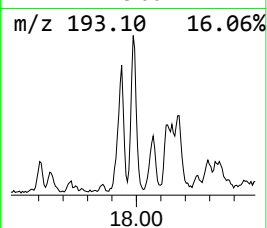
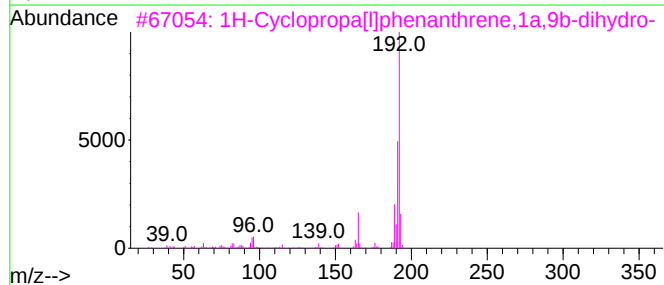
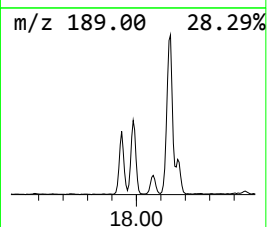
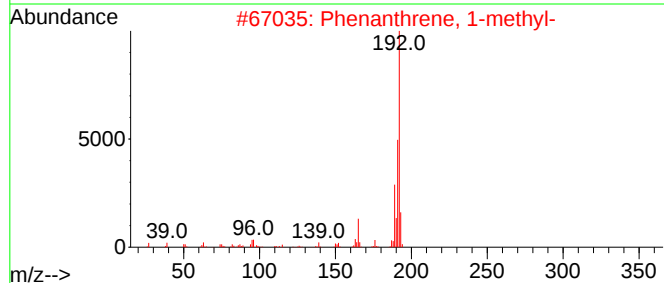
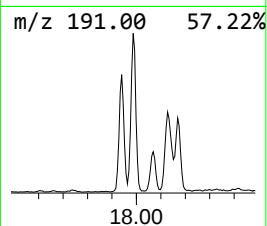
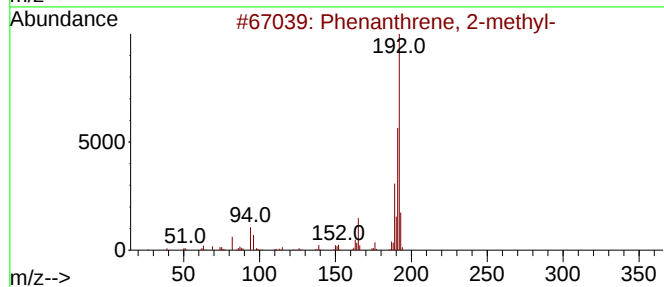
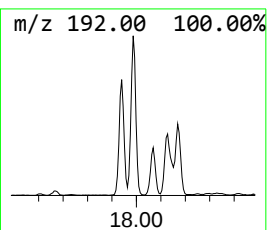
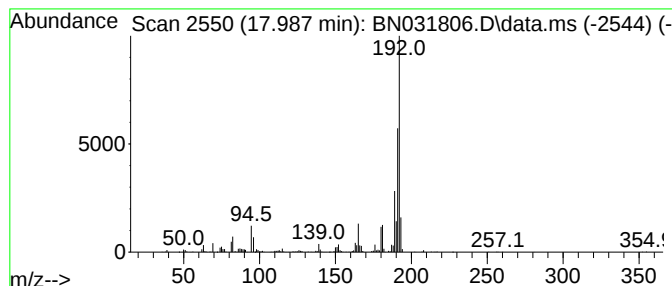
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.987	4.59 ng/ul	1264290	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, 1-methyl-	192	C15H12	000832-69-9	97
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96



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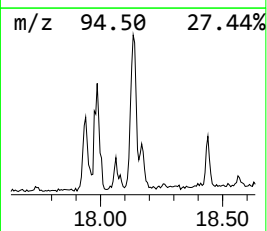
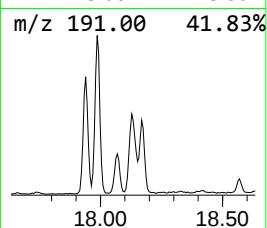
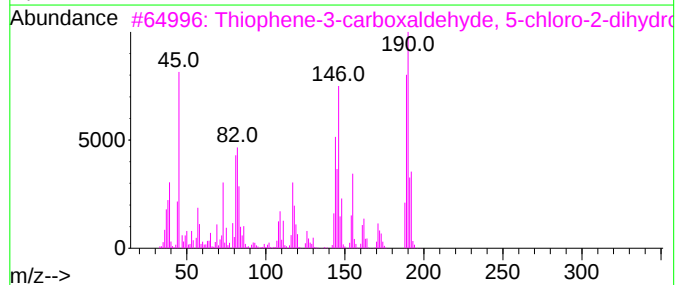
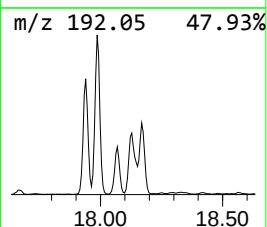
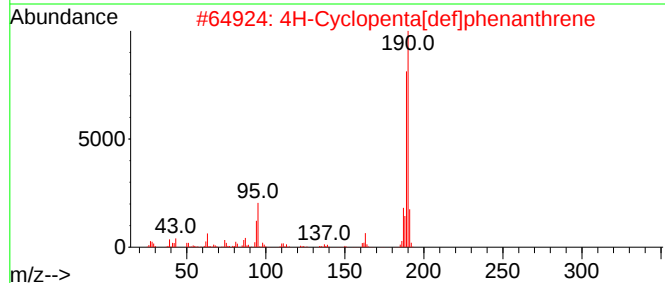
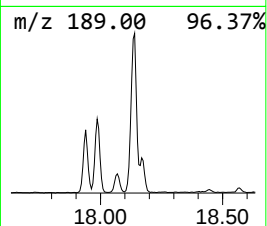
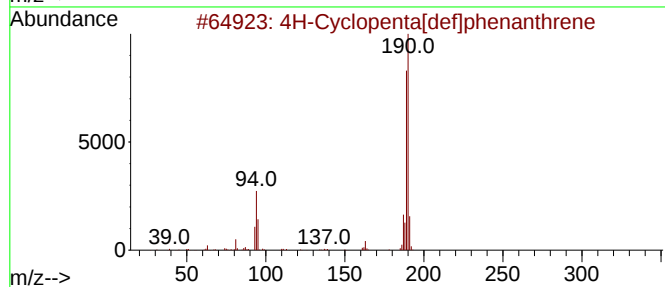
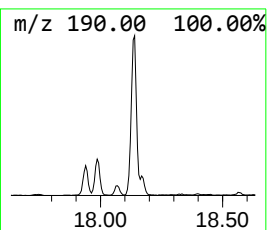
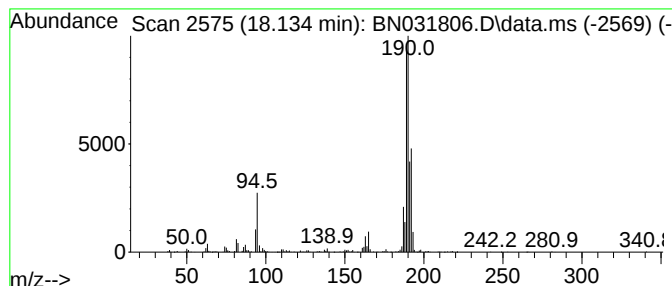
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.134	3.14 ng/ul	863073	Phenanthrene-d10	17.069

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	60
3			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
5			4,6-Dichloro-2-ethylpyrimidin-5-...	191	C6H7Cl2N3	006237-96-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031806.D
Acq On : 05 Jun 2024 19:05
Operator : MA/JU
Sample : P2591-15DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ6DL

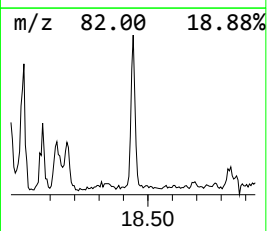
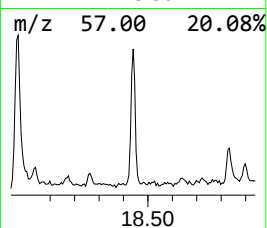
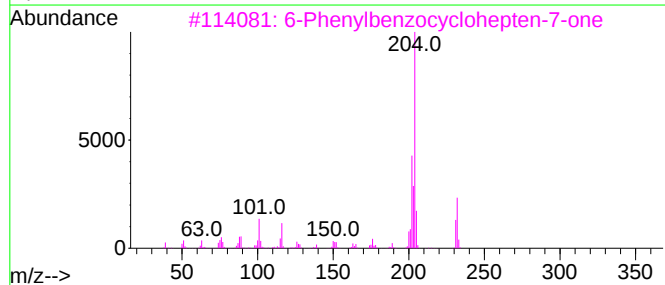
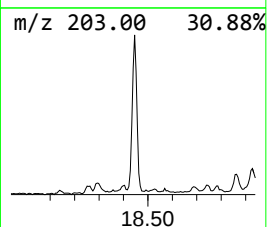
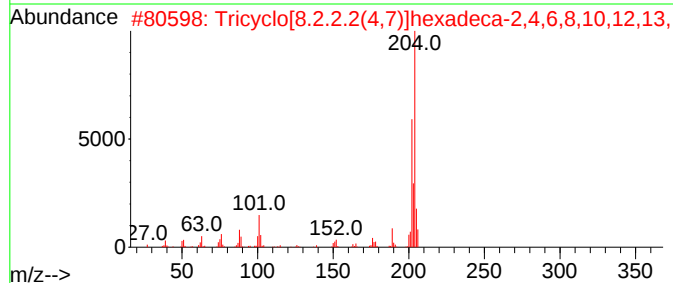
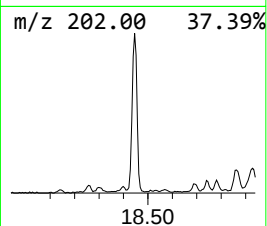
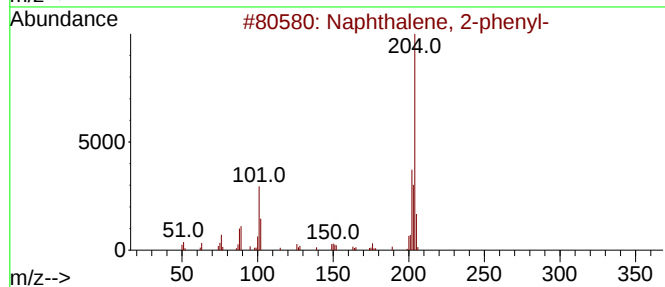
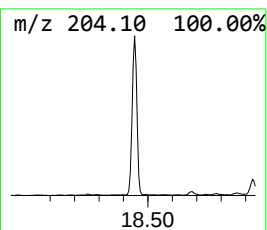
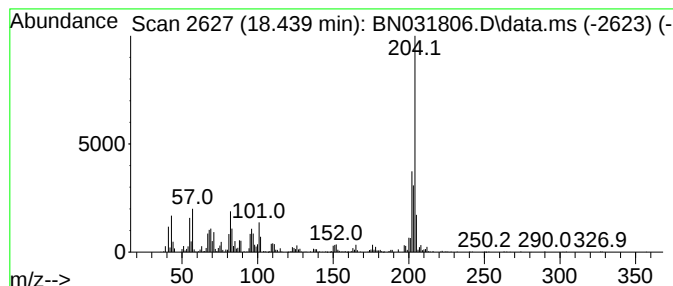
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 2-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.439	2.47 ng/ul	678696	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	89
3			6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	86
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
5			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031806.D
Acq On : 05 Jun 2024 19:05
Operator : MA/JU
Sample : P2591-15DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ6DL

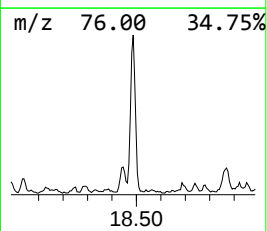
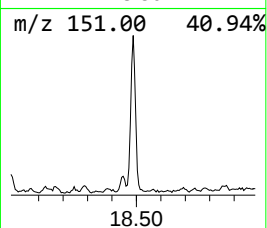
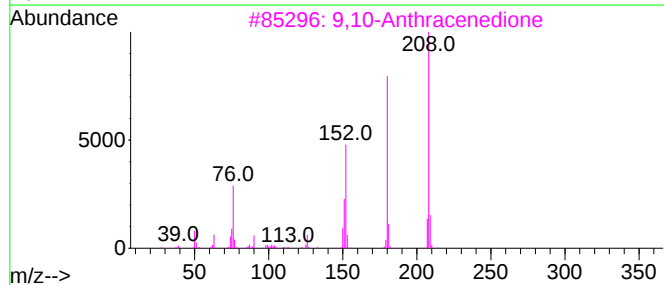
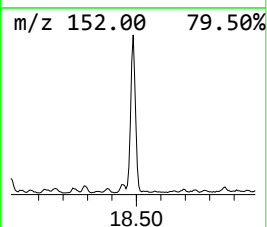
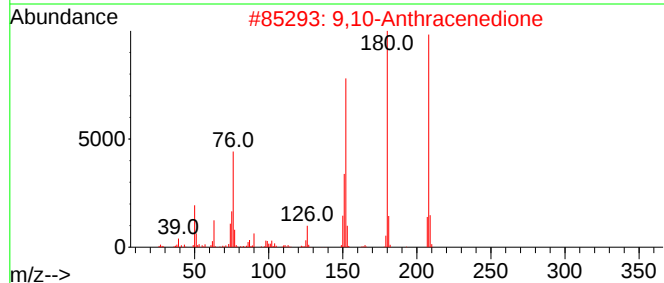
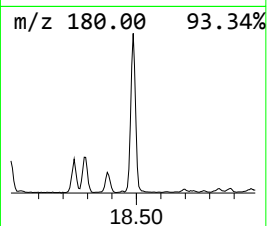
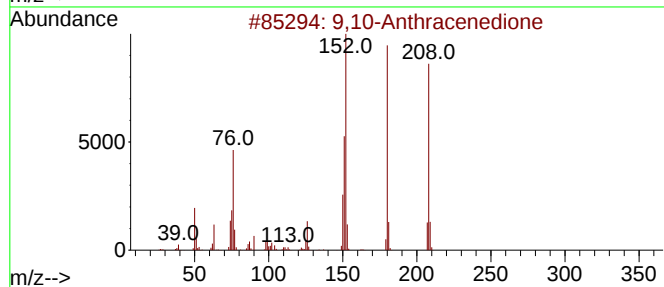
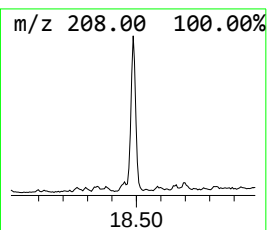
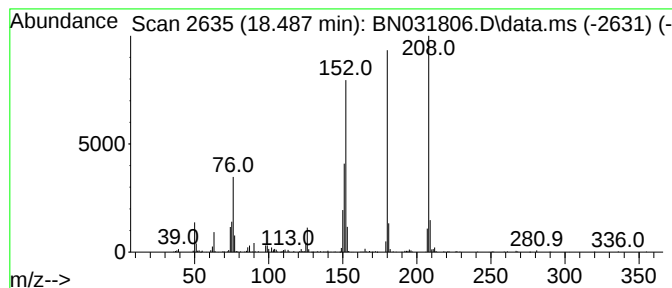
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 9,10-Anthracenedione Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.486	2.15 ng/ul	591196	Phenanthrene-d10	17.069

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	94
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
5			1H-Phenalen-1-one	180	C13H8O	000548-39-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031806.D
Acq On : 05 Jun 2024 19:05
Operator : MA/JU
Sample : P2591-15DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

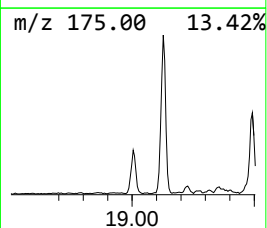
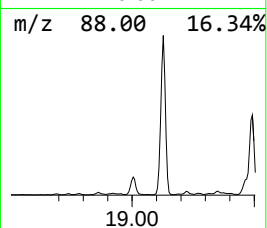
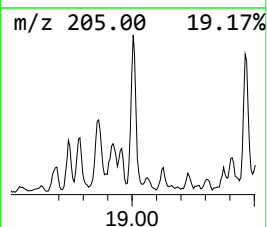
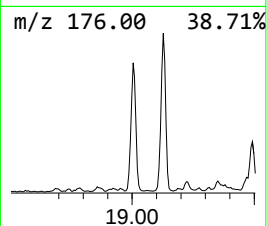
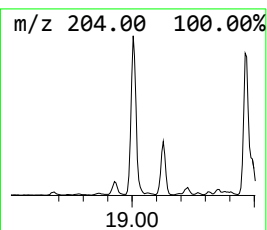
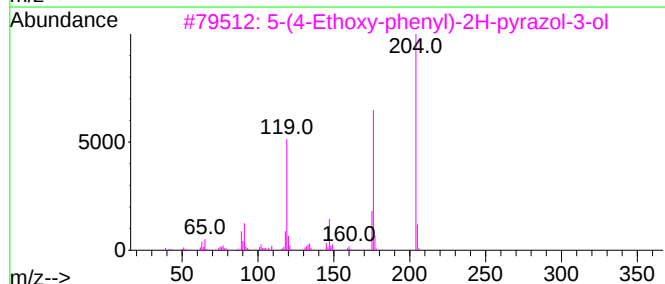
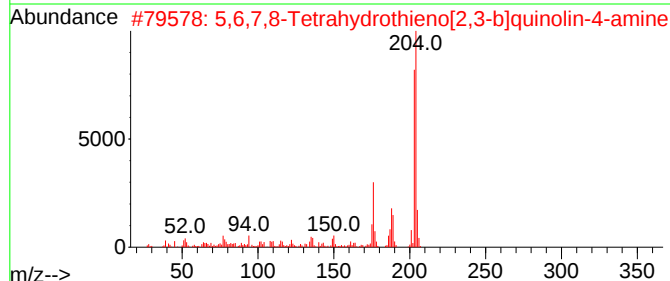
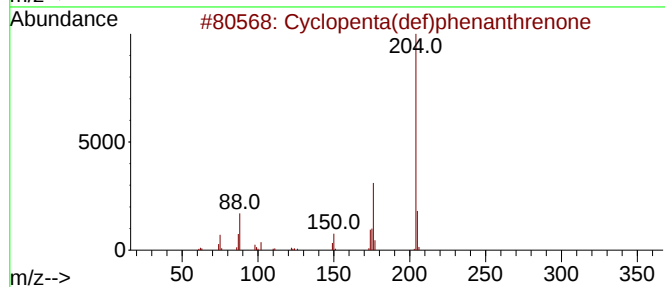
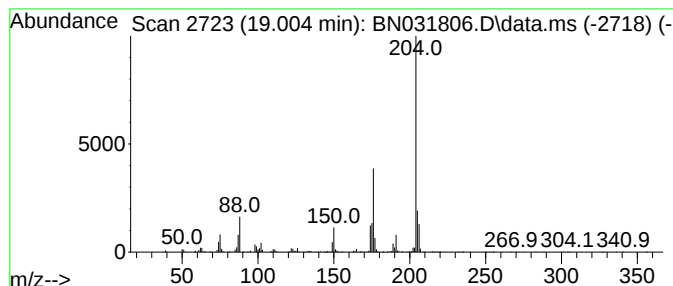
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cyclopenta(def)phenanthrenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.004	2.24 ng/ul	617485	Phenanthrene-d10	17.069	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	98
2	5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	64
3	5-(4-Ethoxy-phenyl)-2H-pyrazol-3-ol	204	C11H12N2O2	1000278-26-8	53
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	49
5	Quinoline, 8-methoxy-5-nitro-	204	C10H8N2O3	1000298-88-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031806.D
Acq On : 05 Jun 2024 19:05
Operator : MA/JU
Sample : P2591-15DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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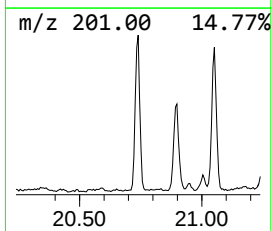
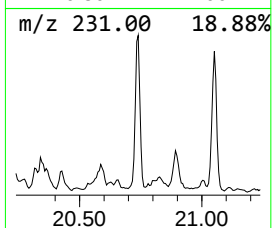
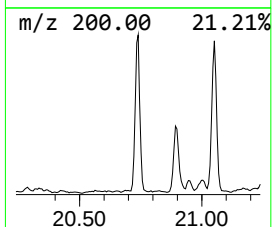
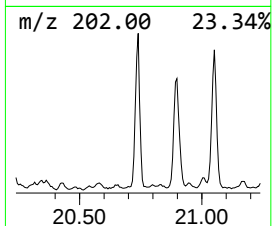
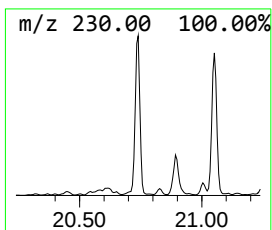
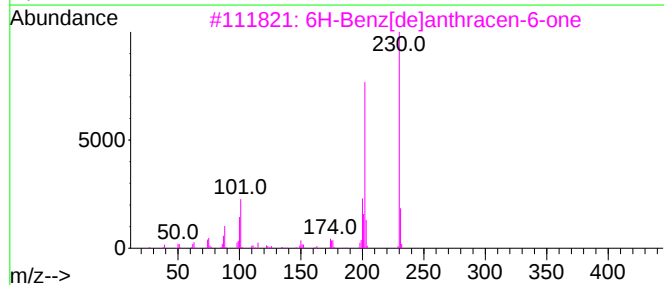
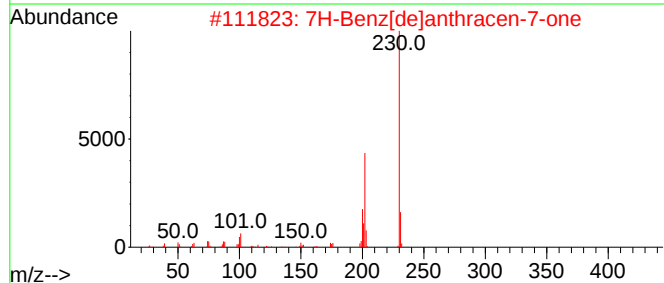
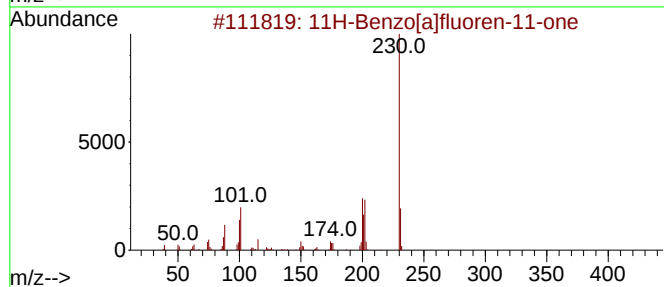
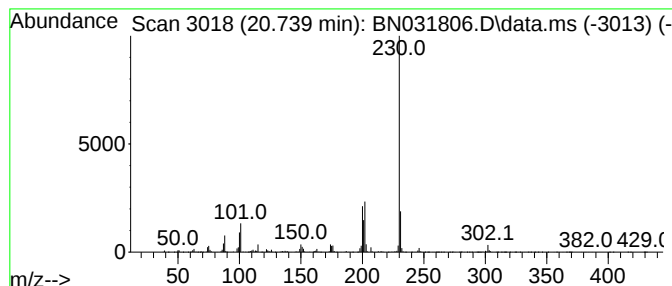
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 11H-Benzo[a]fluoren-11-one Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.739	2.44 ng/ul	1427840	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	98
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	91
3			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	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\BN060524\
Data File : BN031806.D
Acq On : 05 Jun 2024 19:05
Operator : MA/JU
Sample : P2591-15DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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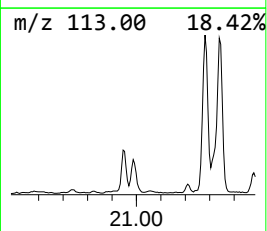
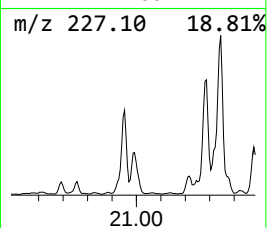
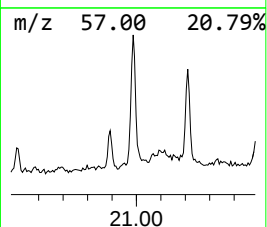
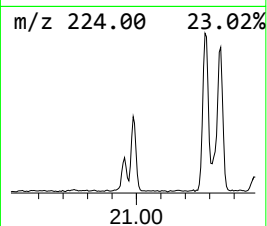
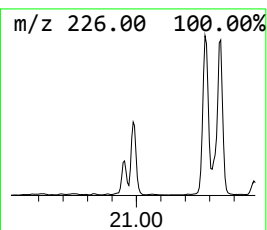
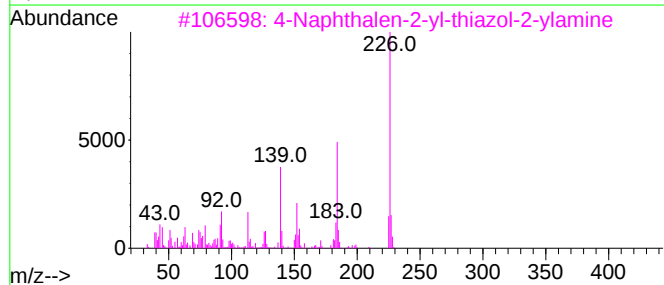
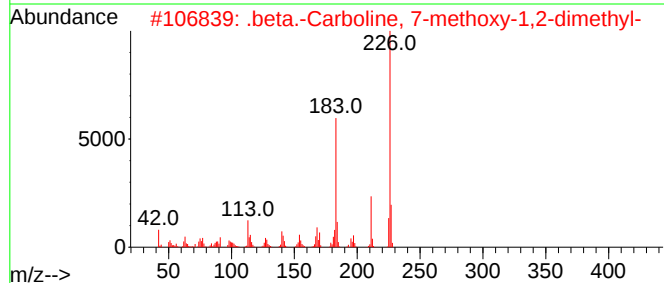
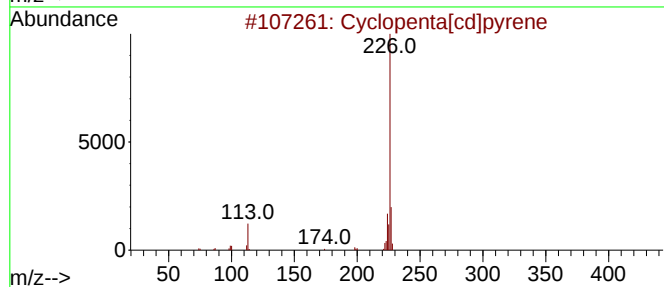
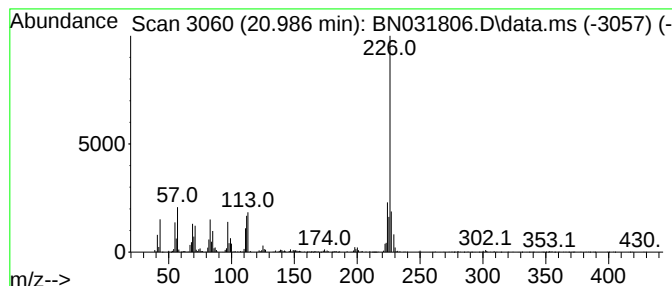
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cyclopenta[cd]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	2.11 ng/ul	1232230	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	92
2			.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	49
3			4-Naphthalen-2-yl-thiazol-2-ylamine	226	C13H10N2S	1000300-53-4	47
4			Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	46
5			Ethyl 2-(allylamino)-4-methyl-1,...	226	C10H14N2O2S	006721-80-8	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031806.D
Acq On : 05 Jun 2024 19:05
Operator : MA/JU
Sample : P2591-15DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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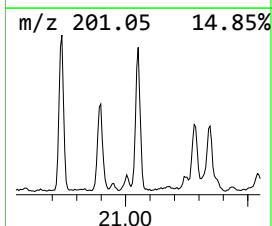
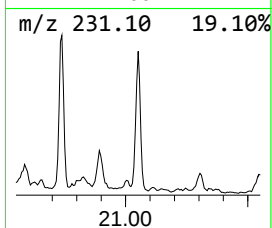
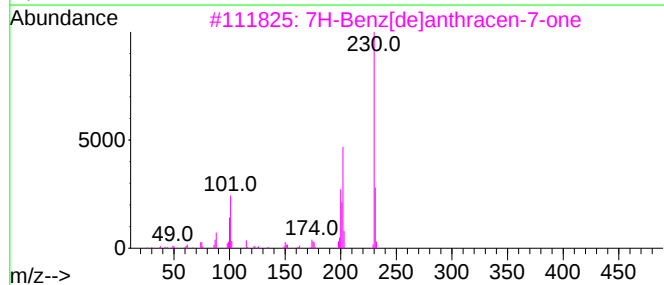
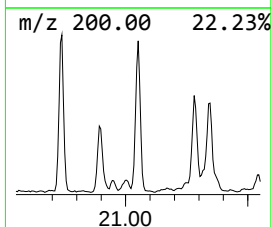
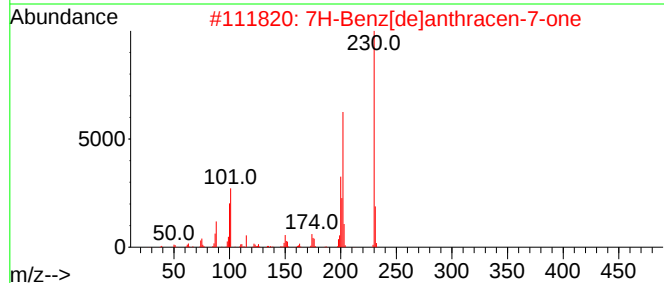
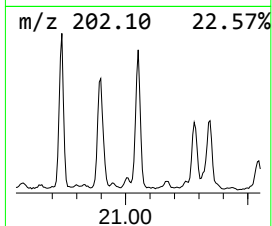
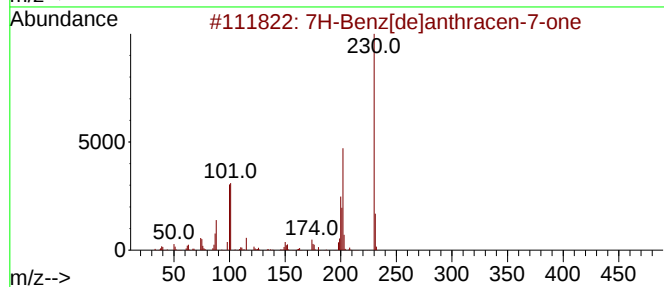
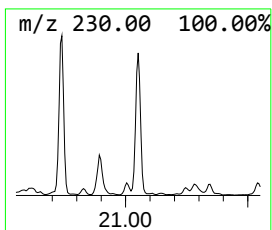
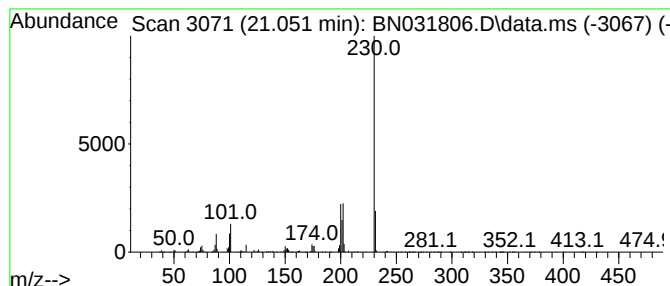
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 7H-Benz[de]anthracen-7-one Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.051	2.31 ng/ul	1349800	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
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	87
5			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031806.D
Acq On : 05 Jun 2024 19:05
Operator : MA/JU
Sample : P2591-15DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ6DL

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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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Phenanthrene, 1...	17.987	4.6	ng/u1	1264290	4	17.069	5504410	20.0
4H-Cyclopenta[d...	18.134	3.1	ng/u1	863073	4	17.069	5504410	20.0
Naphthalene, 2-...	18.439	2.5	ng/u1	678696	4	17.069	5504410	20.0
9,10-Anthracene...	18.486	2.1	ng/u1	591196	4	17.069	5504410	20.0
Cyclopenta(def)...	19.004	2.2	ng/u1	617485	4	17.069	5504410	20.0
11H-Benzo[a]flu...	20.739	2.4	ng/u1	1427840	5	21.304	11706700	20.0
Cyclopenta[cd]p...	20.986	2.1	ng/u1	1232230	5	21.304	11706700	20.0
7H-Benz[de]anth...	21.051	2.3	ng/u1	1349800	5	21.304	11706700	20.0