

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
 Data File : BN031853.D
 Acq On : 08 Jun 2024 11:16
 Operator : MA/JU
 Sample : P2634-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBW8

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: BN031853.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.446	74	78	93	rVB	558598	811936	7.10%	0.371%
2	3.587	98	102	116	rBV	1150148	1802784	15.77%	0.823%
3	6.846	645	656	669	rBV	1870001	3063485	26.80%	1.398%
4	7.011	677	684	695	rVB	1511906	2344583	20.51%	1.070%
5	7.205	710	717	724	rBV	1899212	3070661	26.87%	1.402%
6	7.275	724	729	734	rVV	217126	347465	3.04%	0.159%
7	7.669	787	796	803	rVV	1658941	2653357	23.22%	1.211%
8	8.375	909	916	926	rBV	2175957	3673613	32.14%	1.677%
9	8.828	985	993	1000	rVV	1785992	2957478	25.88%	1.350%
10	9.393	1078	1089	1100	rBV	255306	415681	3.64%	0.190%
11	9.540	1107	1114	1125	rBV	1439379	2488613	21.77%	1.136%
12	10.075	1198	1205	1217	rVV	2190455	3821351	33.44%	1.744%
13	10.452	1260	1269	1275	rBV	2396454	4031166	35.27%	1.840%
14	10.593	1286	1293	1307	rVV	1693845	3035717	26.56%	1.386%
15	11.199	1389	1396	1407	rBV	504579	954864	8.35%	0.436%
16	13.728	1821	1826	1831	rBV	3541157	4976041	43.54%	2.271%
17	14.004	1867	1873	1890	rVV2	4255827	7582645	66.34%	3.461%
18	14.316	1915	1926	1932	rBV	3387390	5220790	45.68%	2.383%
19	14.375	1932	1936	1945	rVB2	97675	180281	1.58%	0.082%
20	14.534	1955	1963	1972	rBV2	3415416	7630407	66.76%	3.483%
21	14.716	1990	1994	2002	rVB2	119628	223774	1.96%	0.102%
22	14.934	2023	2031	2035	rBV3	94141	221979	1.94%	0.101%
23	15.104	2052	2060	2064	rBV6	80833	182682	1.60%	0.083%
24	15.310	2088	2095	2101	rBV	5526175	7887416	69.01%	3.600%
25	15.363	2101	2104	2112	rVB2	147094	231459	2.03%	0.106%
26	15.439	2112	2117	2122	rBV	1166730	1576380	13.79%	0.720%
27	16.039	2213	2219	2225	rVB5	195763	359119	3.14%	0.164%
28	16.369	2267	2275	2280	rBV4	115881	280120	2.45%	0.128%
29	16.598	2306	2314	2317	rBV4	98138	206661	1.81%	0.094%
30	16.722	2330	2335	2341	rVB2	222331	355968	3.11%	0.162%
31	16.881	2358	2362	2370	rVB	225532	366586	3.21%	0.167%
32	17.063	2385	2393	2397	rBV	4186794	6357421	55.62%	2.902%
33	17.104	2397	2400	2405	rVV	3446009	4722912	41.32%	2.156%
34	17.163	2405	2410	2414	rVV	5748105	8424999	73.72%	3.846%
35	17.198	2414	2416	2420	rVV	651516	716615	6.27%	0.327%
36	17.251	2420	2425	2431	rVB3	140152	264228	2.31%	0.121%

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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 >

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

37	17.469	2455	2462	2467	rBV	295809	504176	4.41%	0.230%
38	17.581	2475	2481	2485	rVV2	127013	222678	1.95%	0.102%
39	17.645	2487	2492	2499	rVB2	246957	405409	3.55%	0.185%
40	17.939	2537	2542	2544	rBV	972097	1436539	12.57%	0.656%
41	17.969	2544	2547	2555	rVV2	2303271	4561381	39.91%	2.082%
42	18.069	2560	2564	2568	rVV	294775	445173	3.90%	0.203%
43	18.128	2569	2574	2578	rVV2	1240452	2222454	19.45%	1.015%
44	18.169	2578	2581	2588	rVB	762936	1051347	9.20%	0.480%
45	18.257	2592	2596	2599	rBV2	178391	249407	2.18%	0.114%
46	18.292	2599	2602	2605	rVV4	152986	229377	2.01%	0.105%
47	18.333	2606	2609	2614	rVB2	165240	239477	2.10%	0.109%
48	18.445	2623	2628	2631	rBV	578136	810149	7.09%	0.370%
49	18.486	2631	2635	2645	rVB2	710122	1084708	9.49%	0.495%
50	18.686	2666	2669	2674	rVV	306403	432623	3.79%	0.197%
51	18.739	2674	2678	2681	rVV	413297	508263	4.45%	0.232%
52	18.780	2681	2685	2689	rVB	330224	431211	3.77%	0.197%
53	18.863	2689	2699	2703	rBV	1038015	1947222	17.04%	0.889%
54	18.922	2704	2709	2712	rVV2	517584	973859	8.52%	0.445%
55	18.957	2712	2715	2718	rVB	360837	416841	3.65%	0.190%
56	19.004	2718	2723	2730	rBV3	612337	1272495	11.13%	0.581%
57	19.128	2736	2744	2751	rBV	7463270	10702924	93.65%	4.886%
58	19.192	2752	2755	2758	rBV	204584	239835	2.10%	0.109%
59	19.228	2758	2761	2762	rBV	231851	225014	1.97%	0.103%
60	19.251	2762	2765	2768	rBV	723151	798523	6.99%	0.365%
61	19.333	2775	2779	2782	rBV2	178928	237722	2.08%	0.109%
62	19.375	2782	2786	2788	rVV	284453	378200	3.31%	0.173%
63	19.404	2788	2791	2795	rVB3	243095	333899	2.92%	0.152%
64	19.463	2795	2801	2804	rBV3	7057996	11082359	96.97%	5.059%
65	19.492	2804	2806	2814	rVV	6909873	9282220	81.22%	4.237%
66	19.569	2814	2819	2824	rVB2	576324	900033	7.87%	0.411%
67	19.675	2834	2837	2842	rVB	318043	469109	4.10%	0.214%
68	19.733	2843	2847	2852	rVB4	356511	603869	5.28%	0.276%
69	19.816	2856	2861	2868	rBV4	719150	1715468	15.01%	0.783%
70	19.957	2879	2885	2886	rBV2	586672	913496	7.99%	0.417%
71	19.986	2887	2890	2898	rVB2	1218282	1932716	16.91%	0.882%
72	20.092	2904	2908	2911	rBV2	695736	833383	7.29%	0.380%
73	20.151	2912	2918	2922	rVV2	974212	1560114	13.65%	0.712%
74	20.186	2922	2924	2928	rVB2	176945	212078	1.86%	0.097%
75	20.286	2937	2941	2945	rBV	686581	869428	7.61%	0.397%
76	20.333	2945	2949	2953	rVV2	649035	885440	7.75%	0.404%

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77	20.398	2958	2960	2964	rVB2	250960	280038	2.45%	0.128%
78	20.451	2966	2969	2974	rVB4	241937	332444	2.91%	0.152%
79	20.510	2976	2979	2982	rVB2	242447	281937	2.47%	0.129%
80	20.739	3014	3018	3025	rVB	806068	1208592	10.57%	0.552%
81	20.910	3044	3047	3051	rVB2	598695	836699	7.32%	0.382%
82	20.951	3051	3054	3057	rVV	596906	704396	6.16%	0.322%
83	20.986	3057	3060	3067	rVV3	1868770	2794739	24.45%	1.276%
84	21.051	3068	3071	3079	rVB	668342	1052079	9.21%	0.480%
85	21.210	3095	3098	3102	rVB	872727	1010114	8.84%	0.461%
86	21.251	3102	3105	3107	rVV	695308	887400	7.76%	0.405%
87	21.298	3107	3113	3118	rVV3	4665216	11429127	100.00%	5.217%
88	21.345	3118	3121	3133	rVB	3288003	5179141	45.32%	2.364%
89	21.492	3141	3146	3152	rBV2	581492	1123574	9.83%	0.513%
90	21.974	3225	3228	3235	rVB	687089	1094745	9.58%	0.500%
91	22.051	3236	3241	3249	rVB3	636528	1250171	10.94%	0.571%
92	22.227	3267	3271	3274	rBV	352633	559364	4.89%	0.255%
93	22.669	3342	3346	3355	rVB6	684185	1485077	12.99%	0.678%
94	23.316	3449	3456	3463	rBV2	2145563	6277197	54.92%	2.865%
95	23.545	3492	3495	3503	rVB	384930	717874	6.28%	0.328%
96	23.974	3561	3568	3573	rBV	1203731	2835468	24.81%	1.294%
97	24.045	3573	3580	3586	rVV2	2748480	6760906	59.16%	3.086%
98	24.110	3586	3591	3602	rVB	1917133	4496307	39.34%	2.052%
99	24.245	3606	3614	3622	rBV	2184273	5774802	50.53%	2.636%
100	27.556	4170	4177	4199	rVB3	878235	3633811	31.79%	1.659%

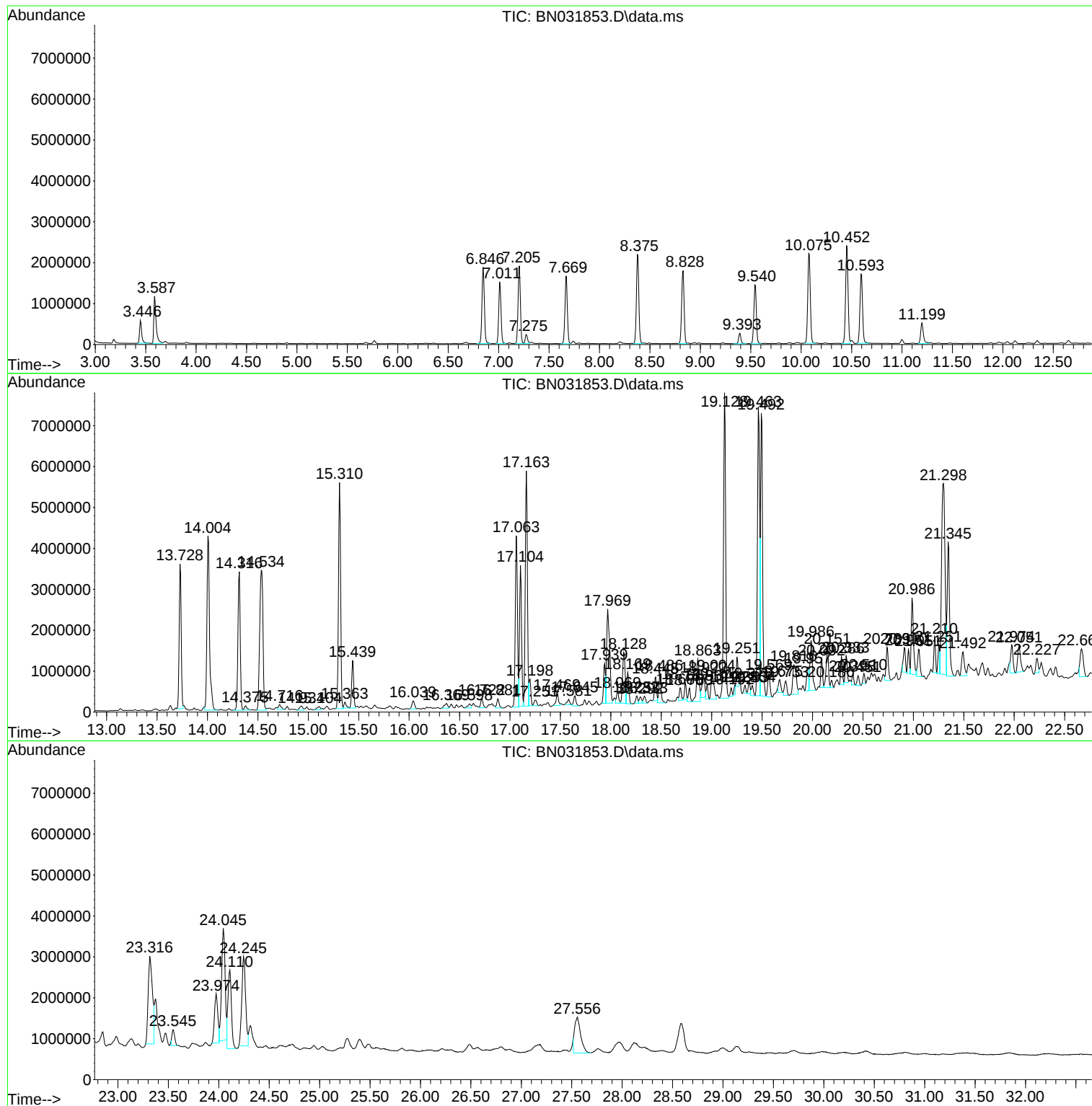
Sum of corrected areas: 219067858

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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



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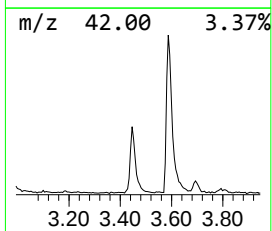
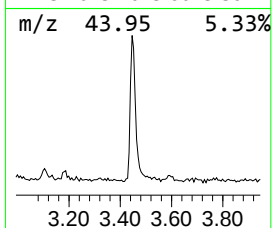
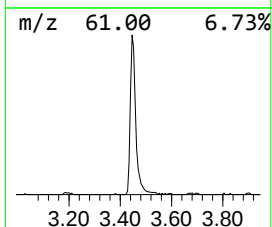
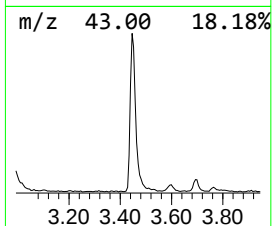
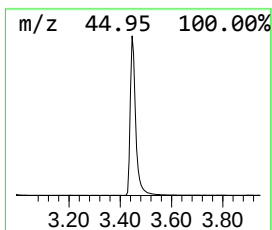
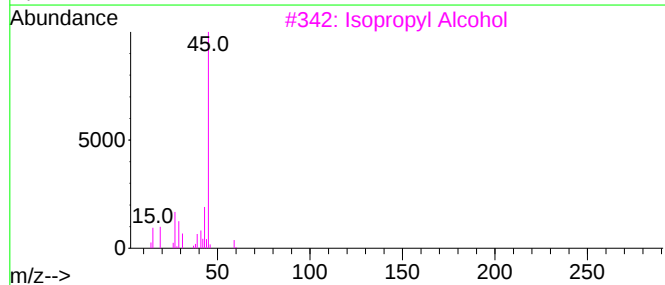
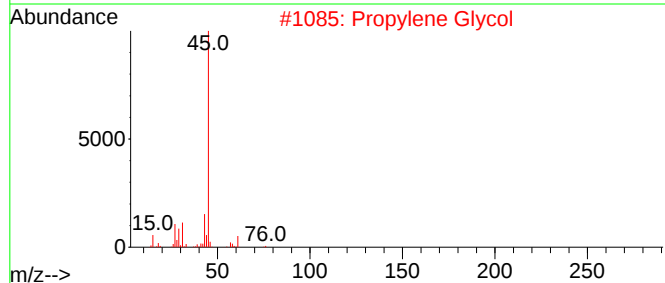
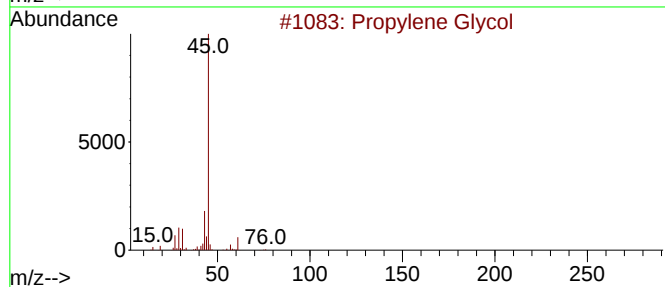
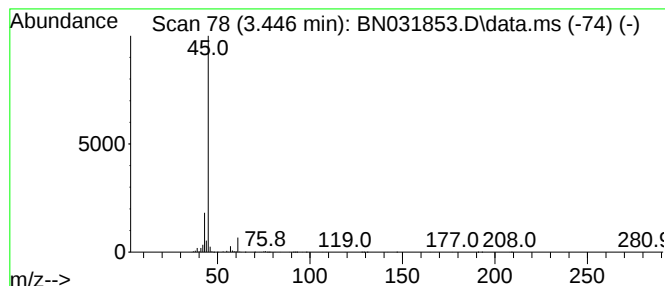
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.446	6.12 ng/ul	811936	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			Propylene Glycol	76	C3H8O2	000057-55-6	83
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	37



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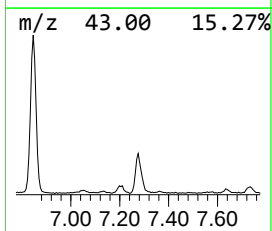
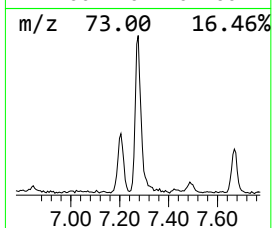
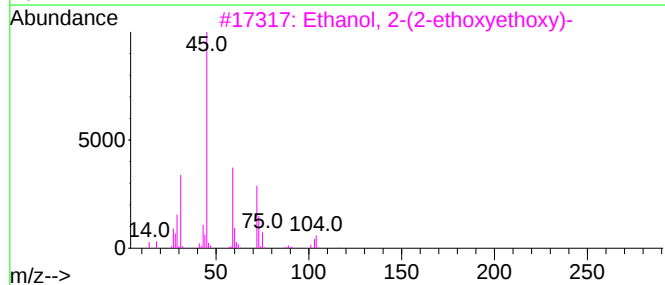
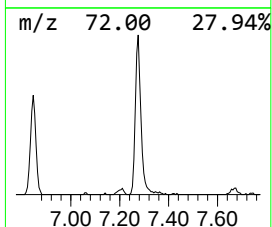
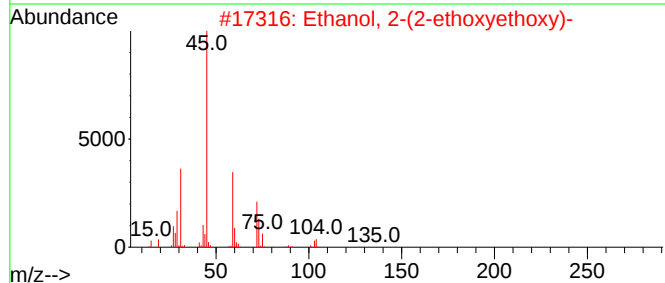
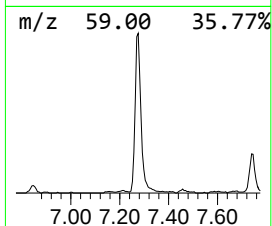
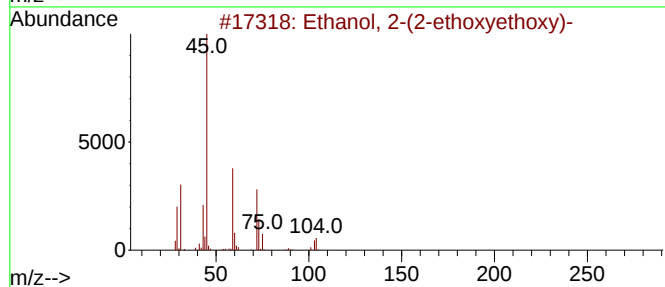
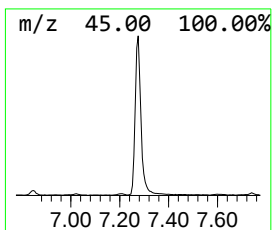
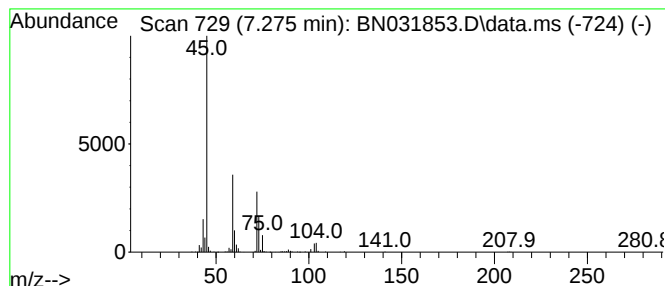
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.275	2.62 ng/ul	347465	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
3			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	50
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	50
5			Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	45



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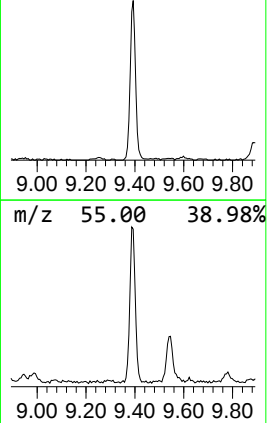
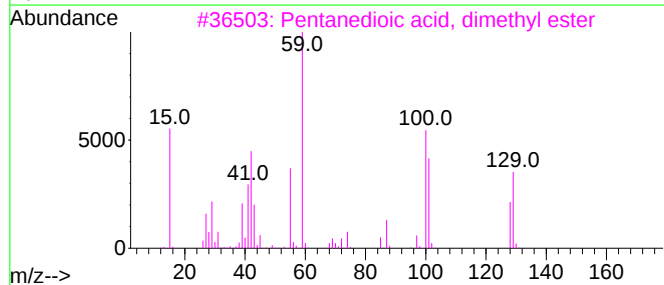
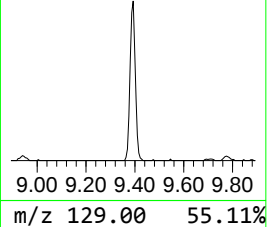
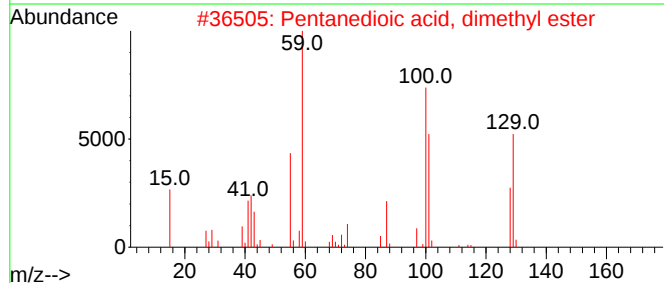
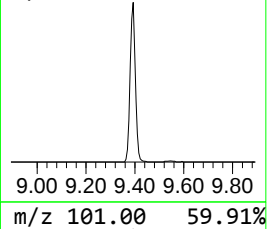
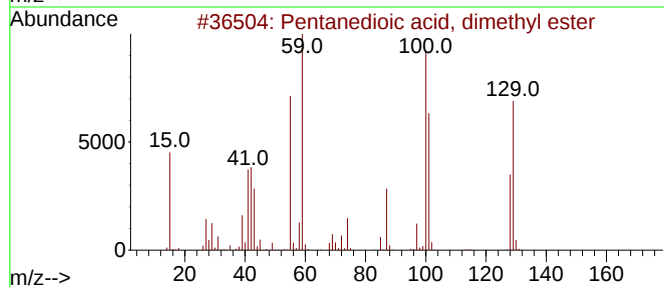
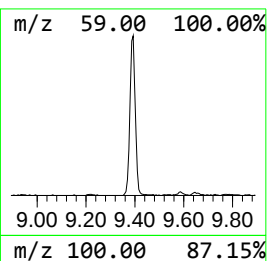
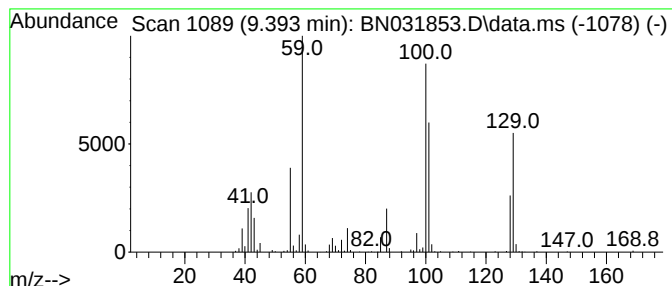
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanedioic acid, dimethyl... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.393	2.06 ng/ul	415681	Naphthalene-d8	10.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	91
2			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	91
3			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	83
4			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	83
5			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
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Instrument :
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ClientSampleId :
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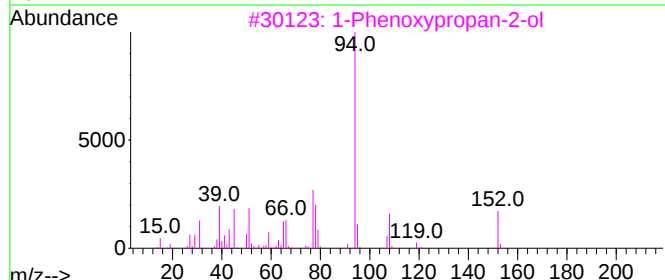
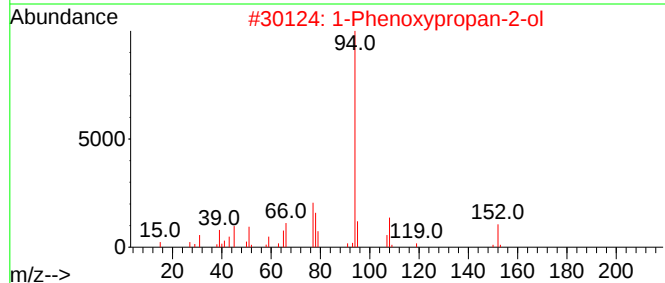
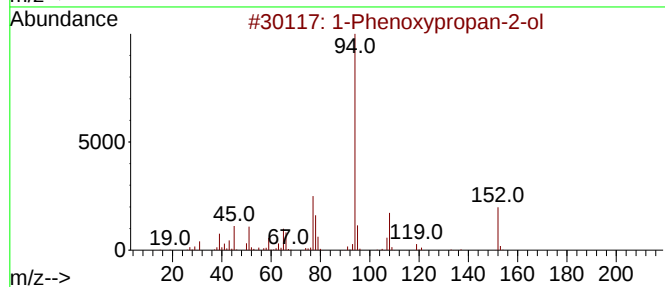
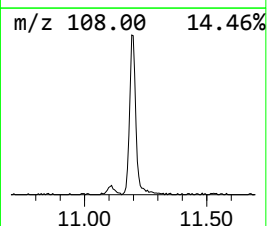
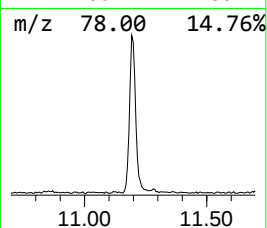
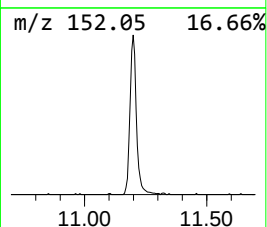
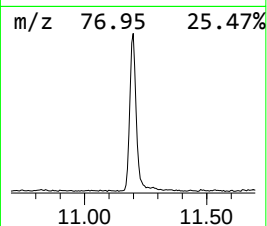
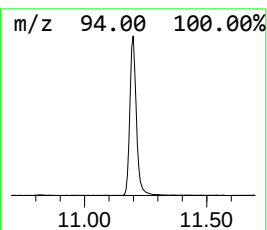
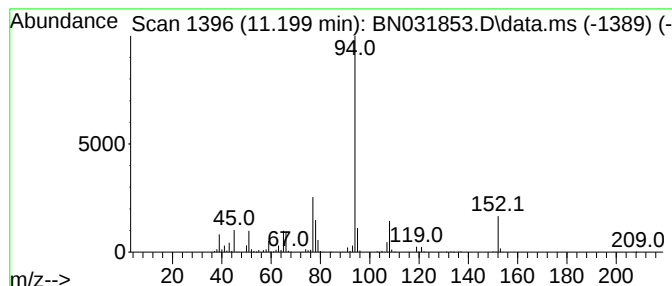
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.198	4.74 ng/ul	954864	Naphthalene-d8	10.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
5			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
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Instrument :
BNA_N
ClientSampleId :
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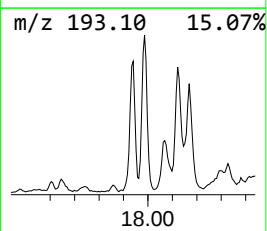
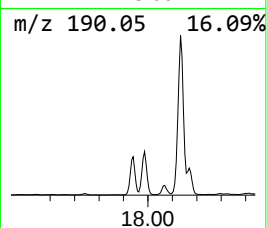
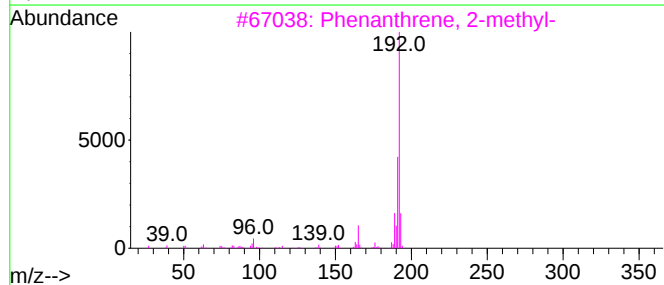
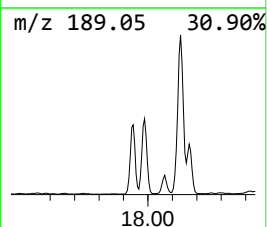
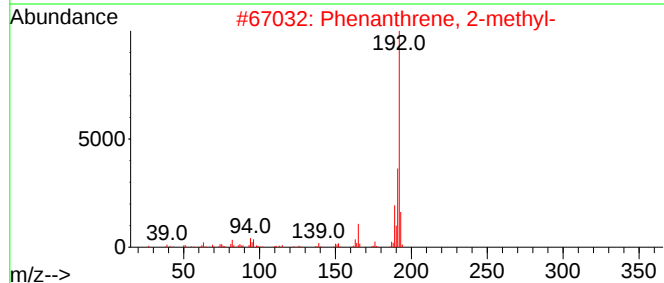
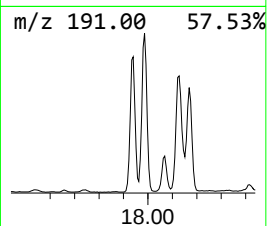
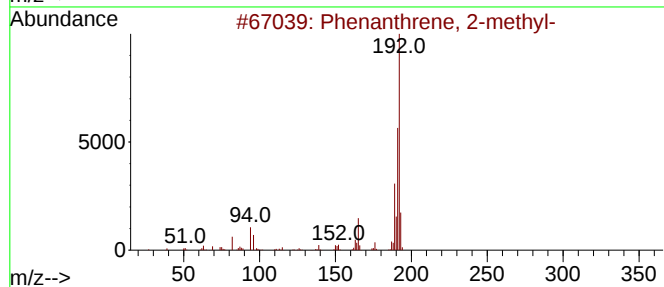
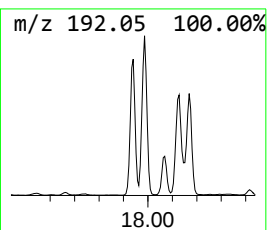
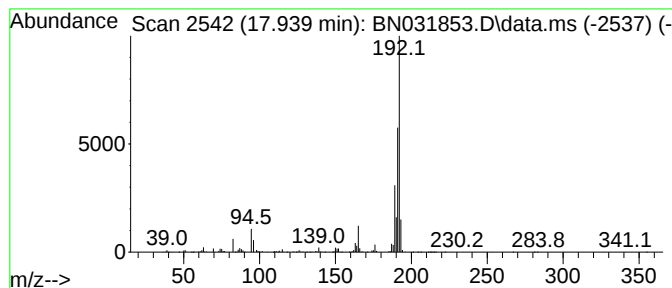
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.939	4.52 ng/ul	1436540	Phenanthrene-d10	17.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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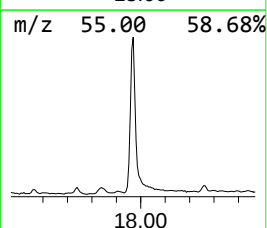
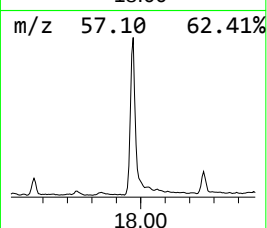
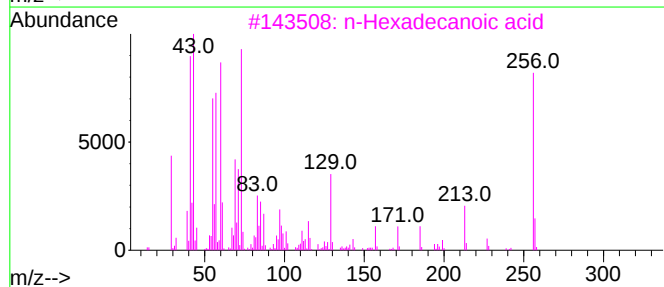
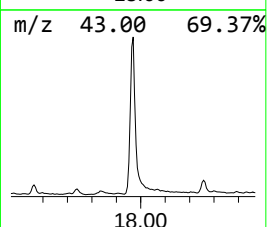
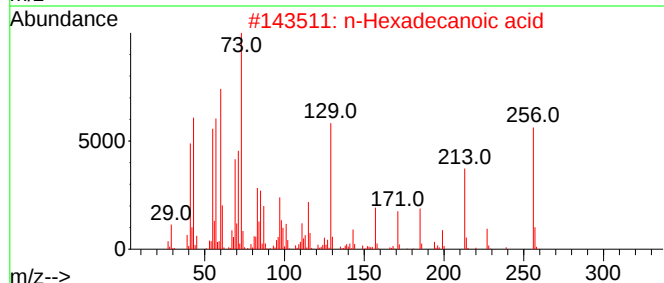
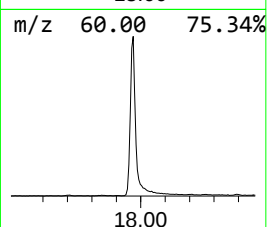
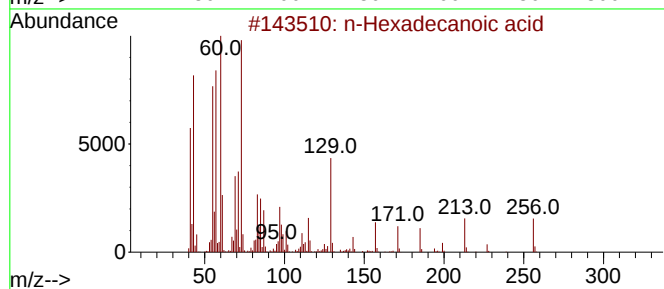
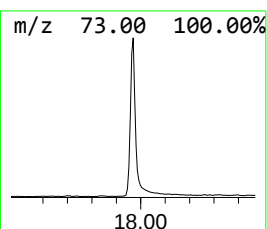
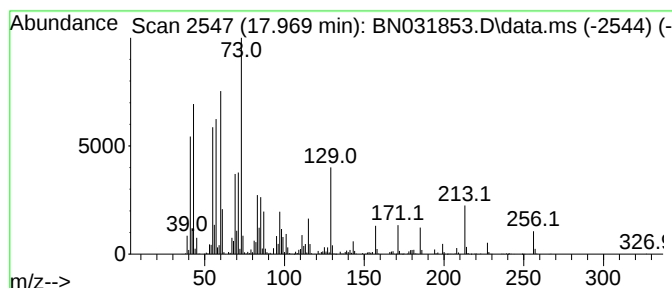
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.969	14.35 ng/ul	4561380	Phenanthrene-d10	17.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98		
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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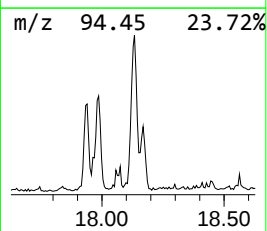
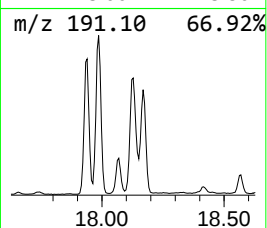
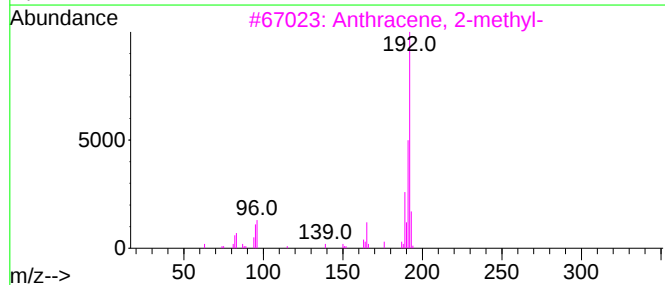
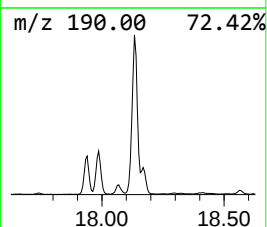
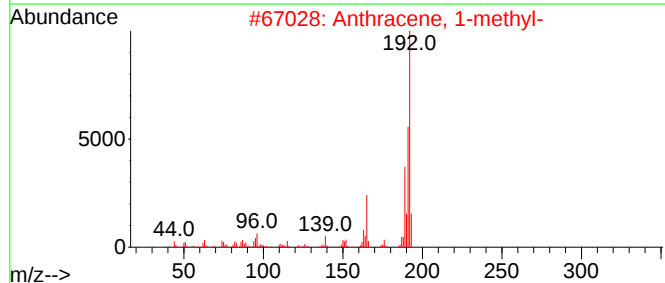
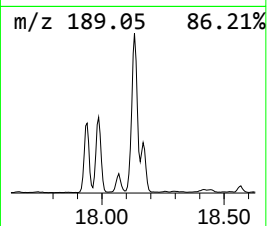
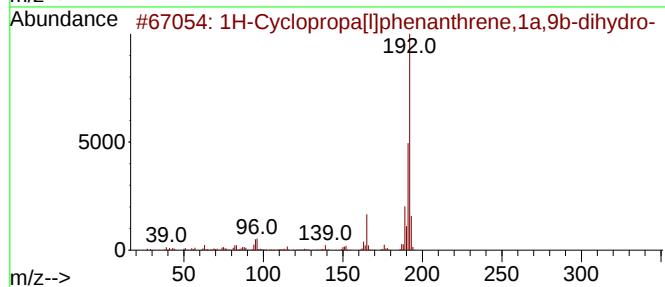
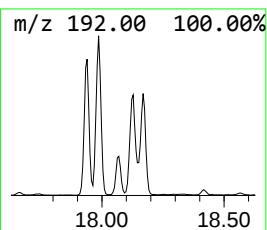
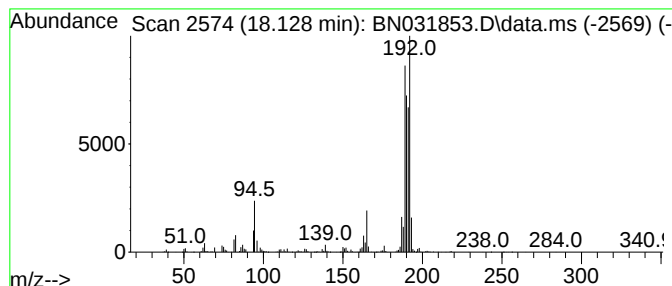
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1H-Cyclopropa[l]phenanthren... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	6.99 ng/ul	2222450	Phenanthrene-d10	17.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	89
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	50
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	50
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	50
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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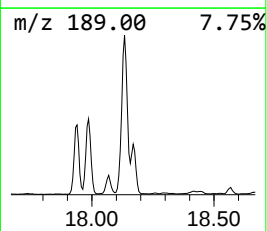
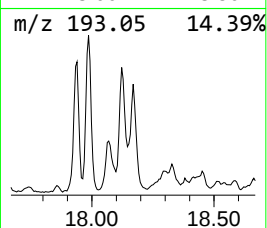
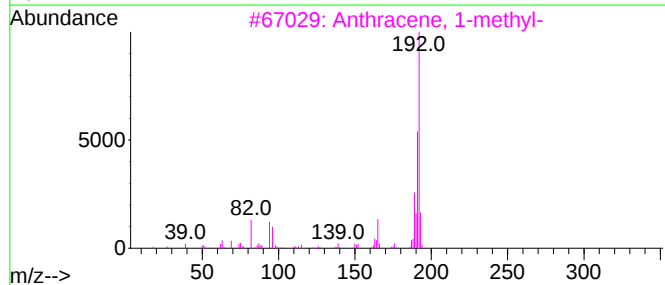
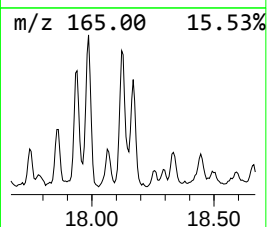
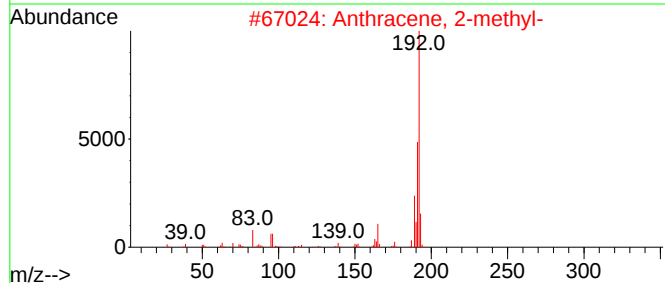
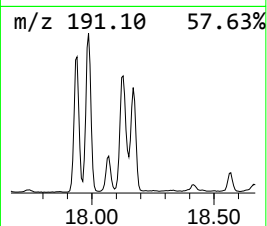
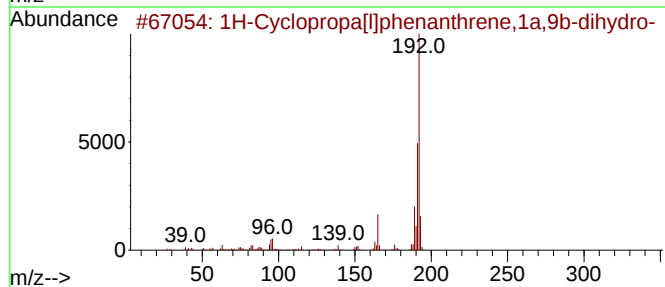
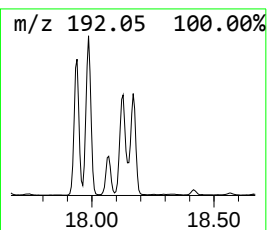
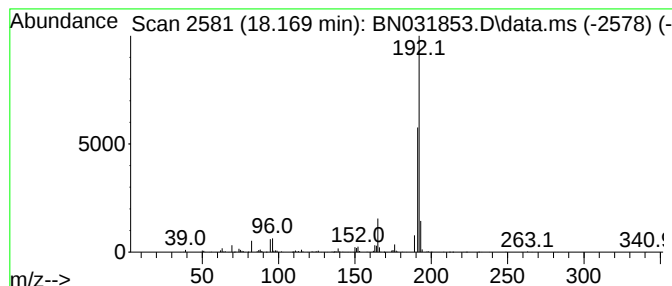
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Anthracene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.169	3.31 ng/ul	1051350	Phenanthrene-d10	17.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
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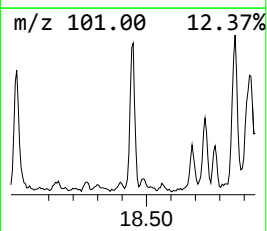
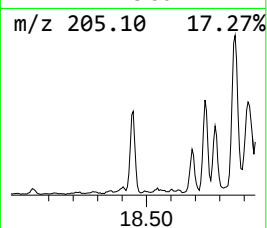
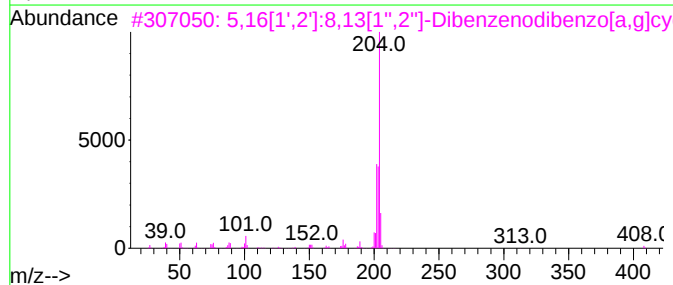
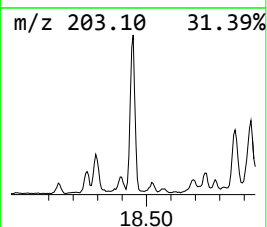
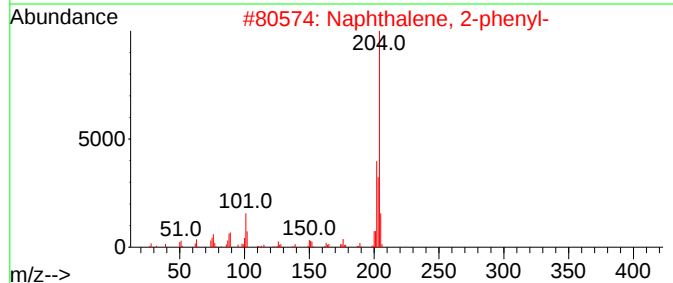
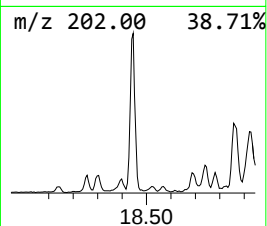
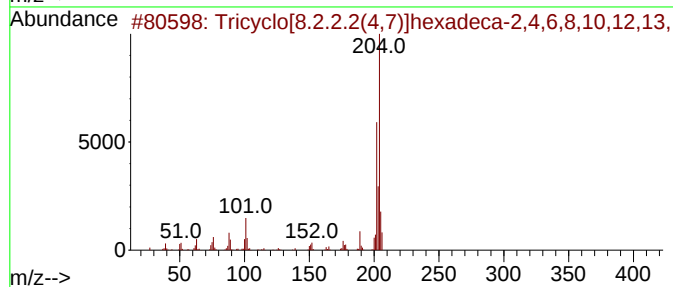
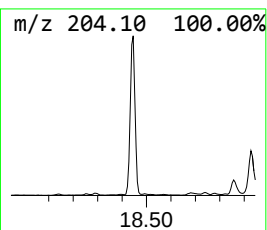
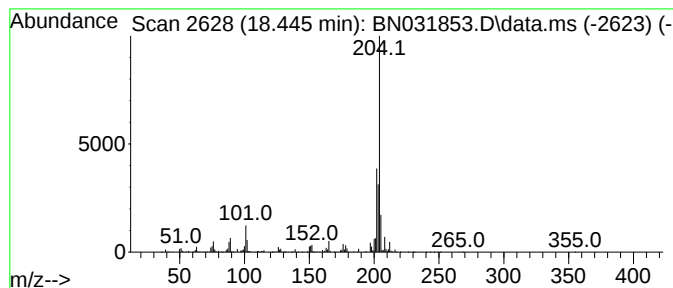
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.445	2.55 ng/ul	810149	Phenanthrene-d10	17.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	86
2			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
3			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
5			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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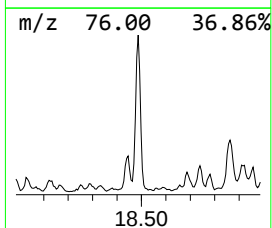
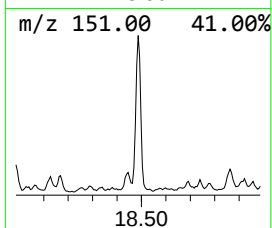
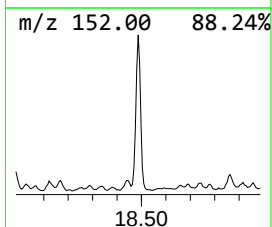
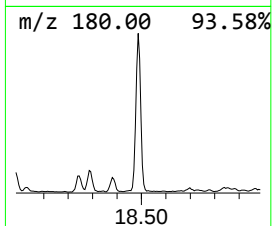
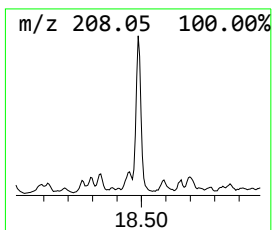
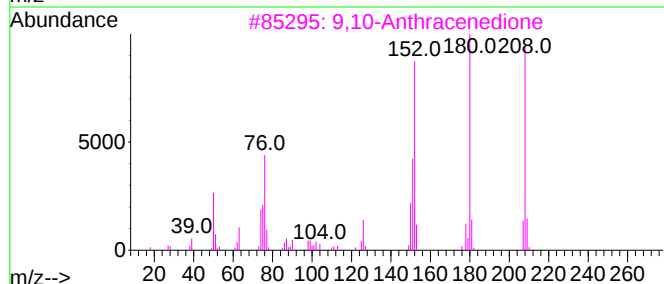
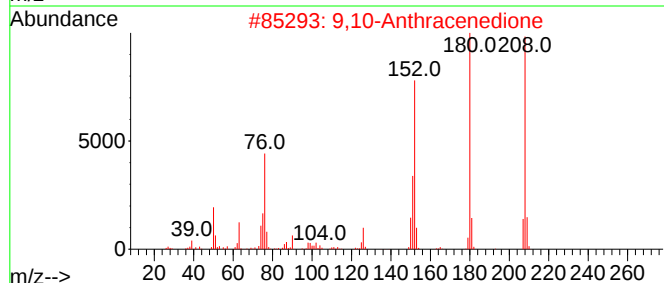
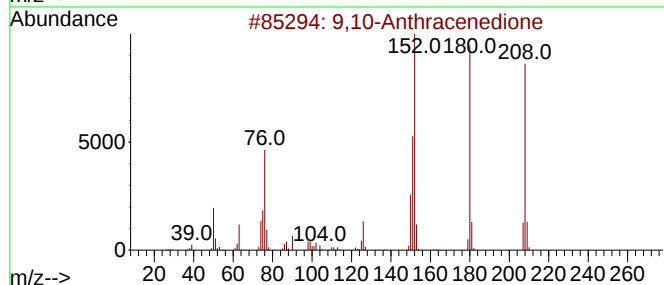
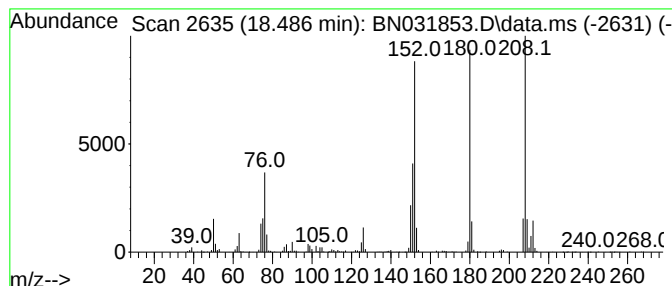
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
18.486	3.41 ng/ul	1084710	Phenanthrene-d10		17.063

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	99	
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98	
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98	
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95	
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	83	



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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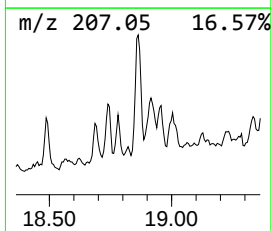
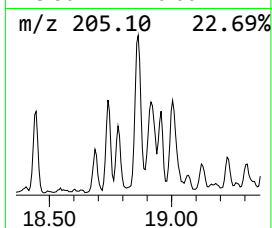
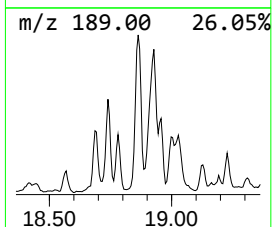
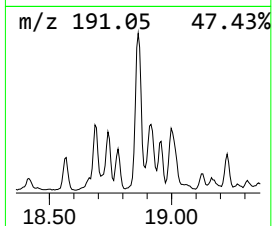
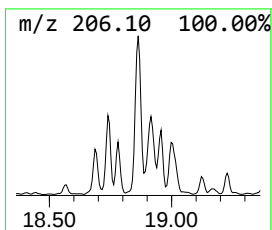
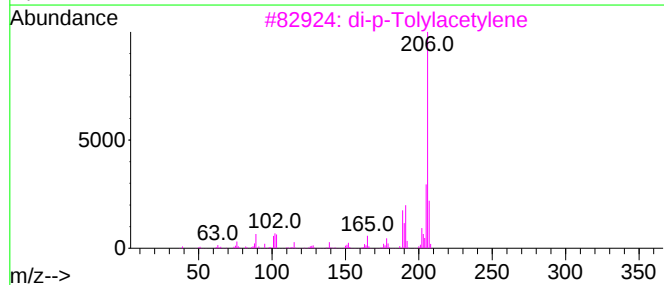
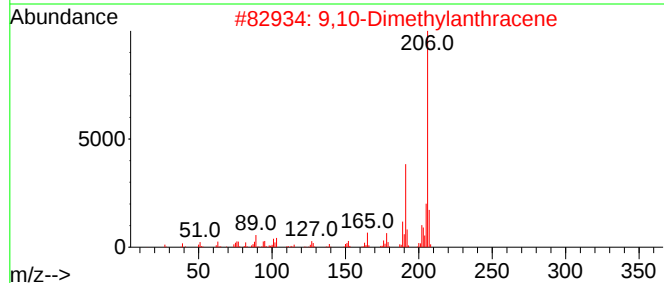
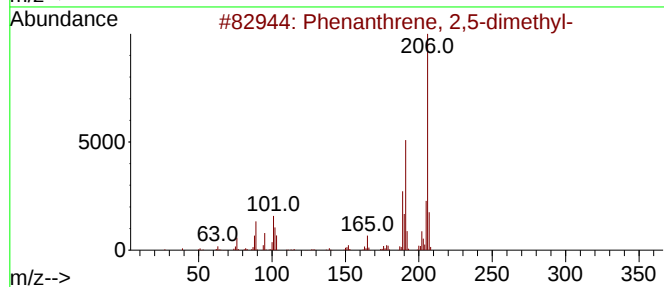
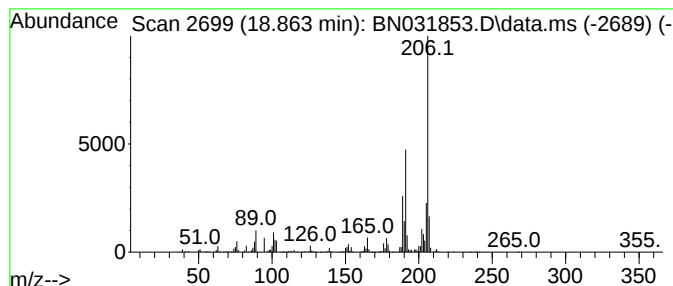
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenanthrene, 2,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.863	6.13 ng/ul	1947220	Phenanthrene-d10	17.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	98
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	96
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	96
4			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	94
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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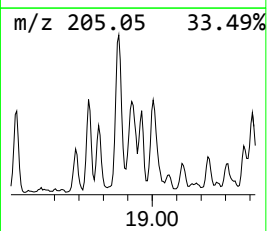
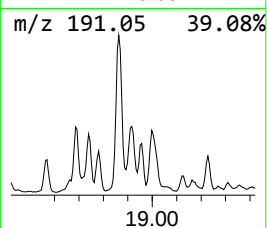
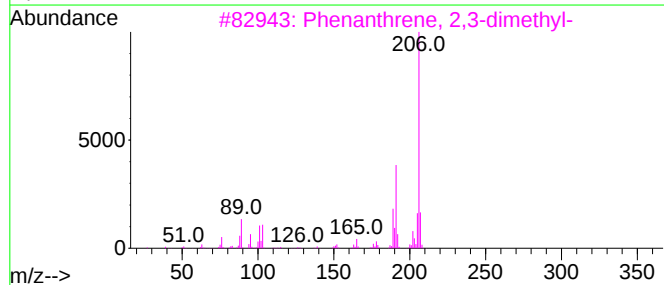
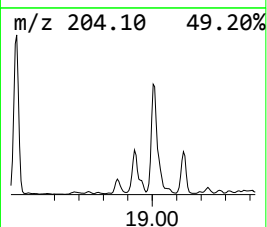
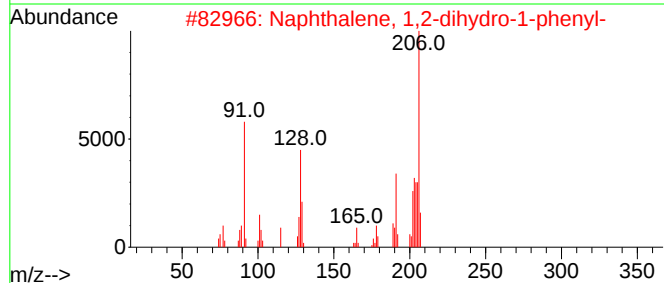
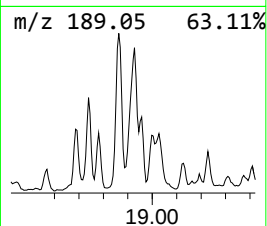
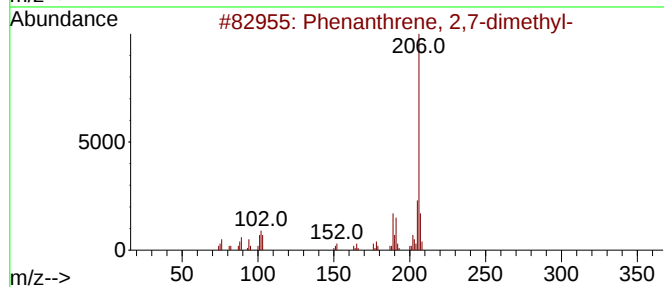
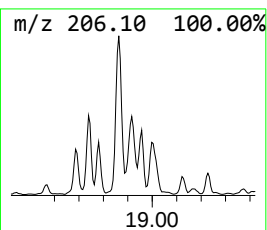
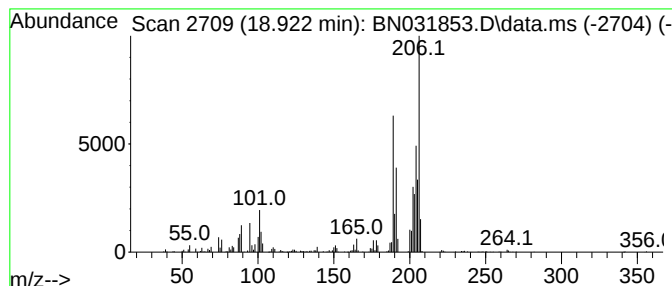
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenanthrene, 2,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.922	3.06 ng/ul	973859	Phenanthrene-d10	17.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	53
2			Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	53
3			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	49
4			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	49
5			1,4-Diphenyl-1,3-butadiene	206	C16H14	000886-65-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

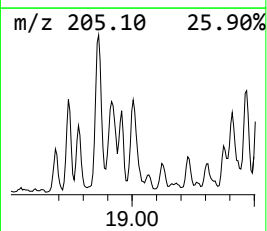
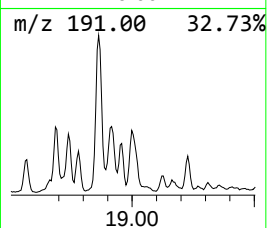
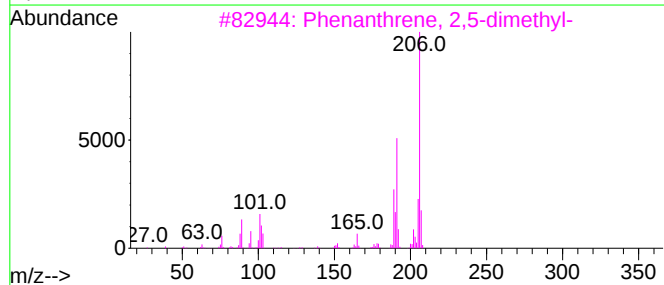
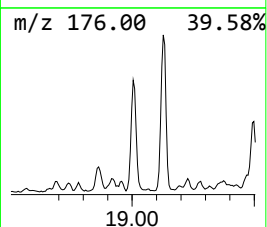
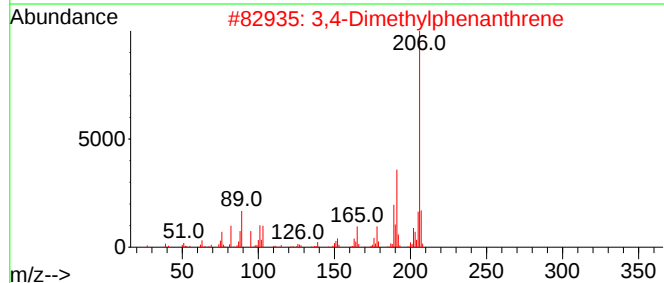
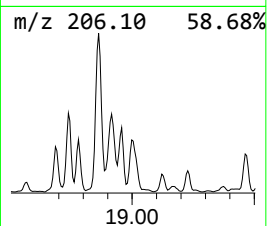
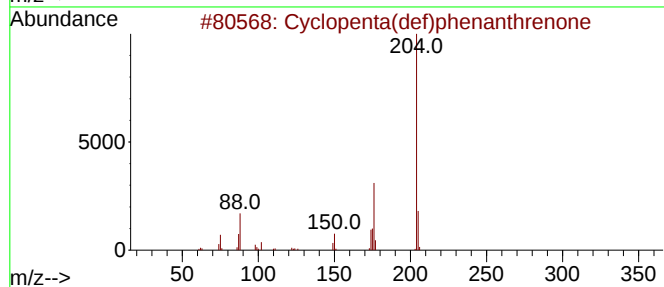
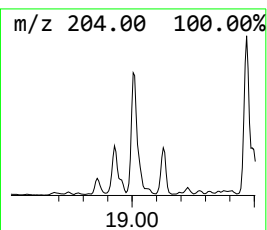
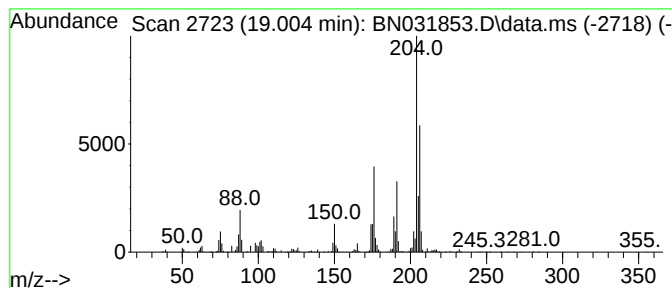
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Cyclopenta(def)phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD			R.T.
19.004	4.00 ng/ul	1272500	Phenanthrene-d10			17.063
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone		204	C15H8O	005737-13-3	95
2	3,4-Dimethylphenanthrene		206	C16H14	1000452-24-5	42
3	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	42
4	9,10-Dimethylanthracene		206	C16H14	000781-43-1	42
5	[14]Annulene, 1,6:8,13-bis(metha...		206	C16H14	085385-68-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
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Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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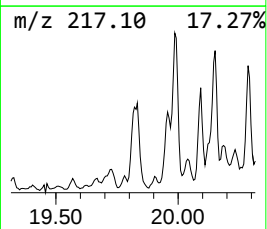
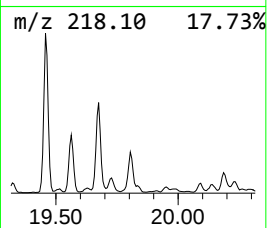
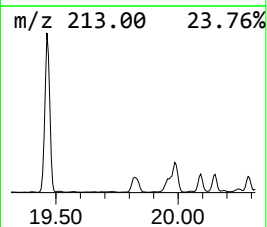
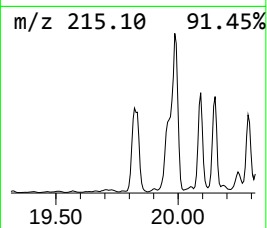
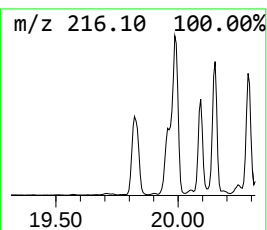
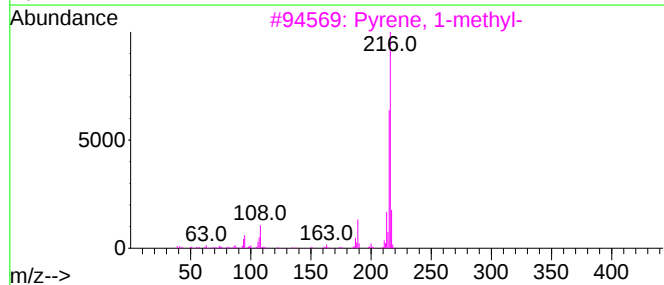
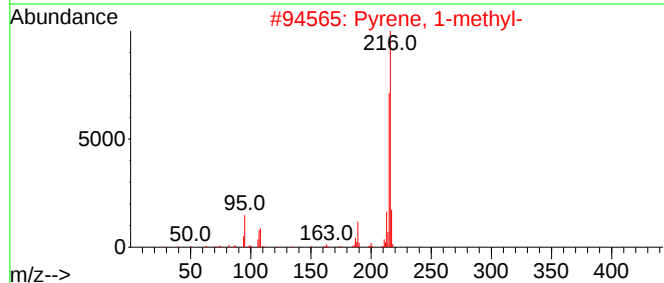
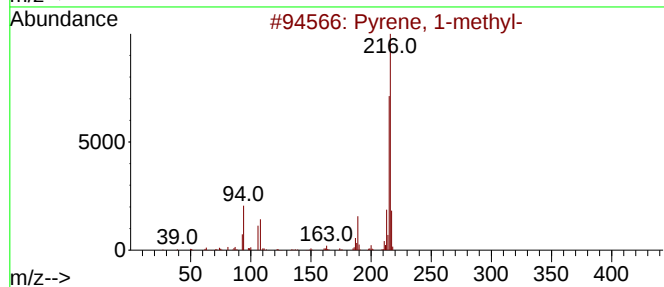
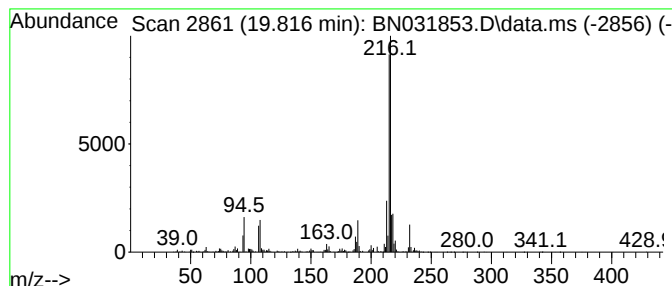
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.816	3.00 ng/ul	1715470	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-			216	C17H12	002381-21-7	97
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	93
5	11H-Benzo[b]fluorene			216	C17H12	000243-17-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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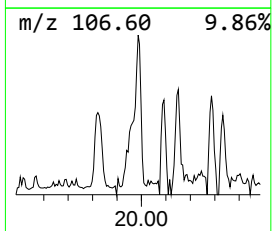
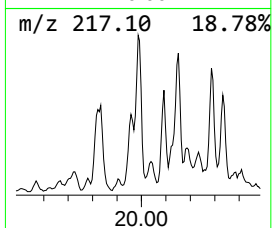
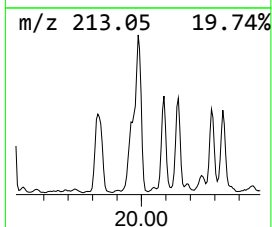
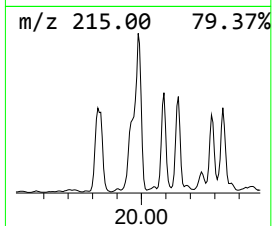
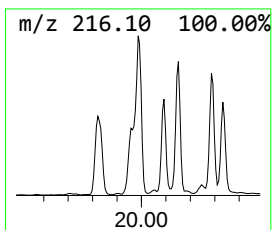
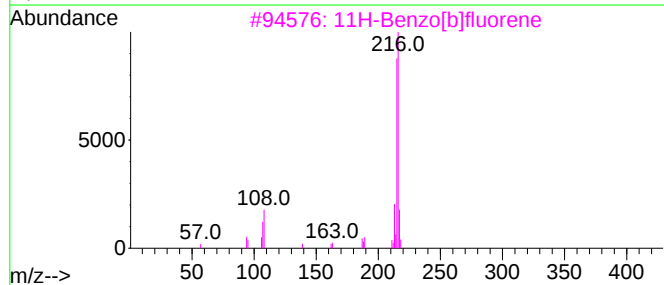
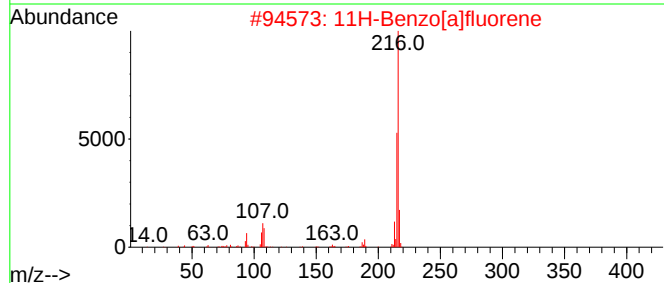
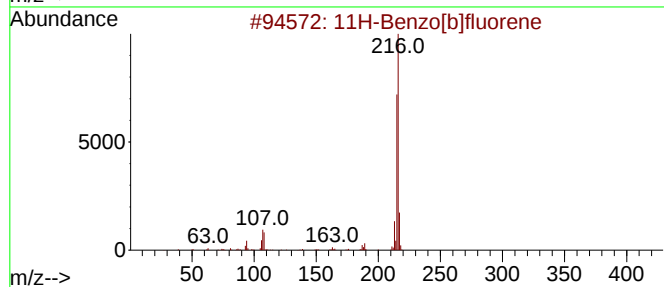
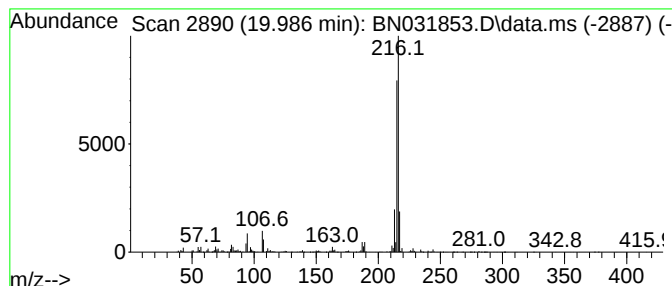
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.986	3.38 ng/ul	1932720	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
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Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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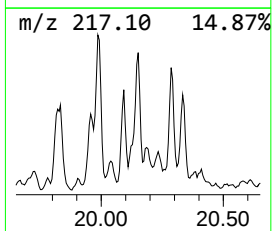
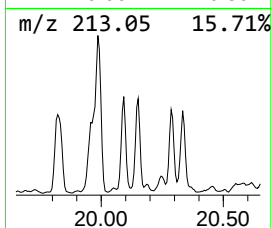
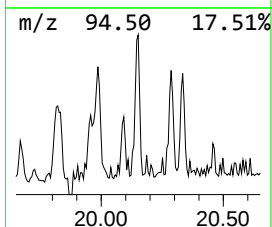
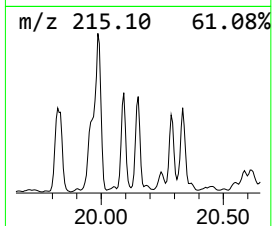
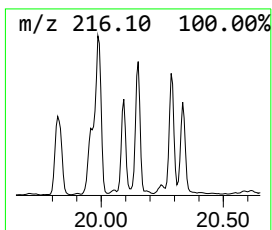
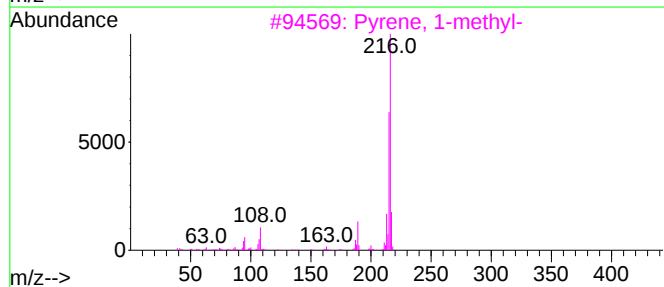
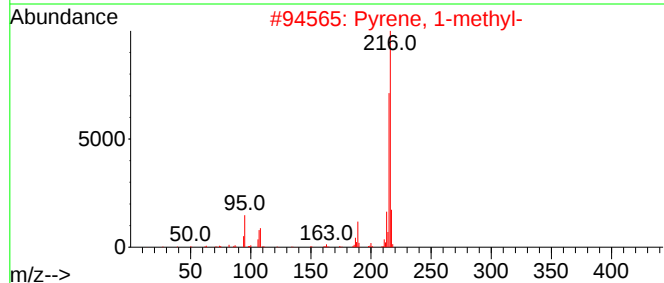
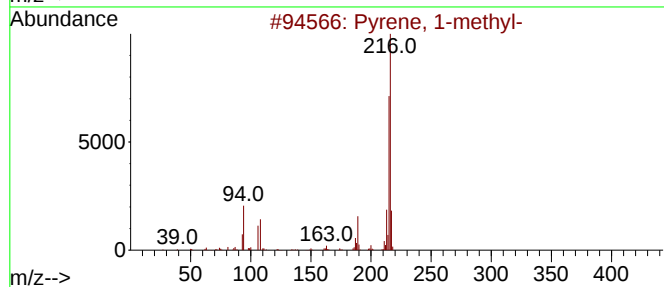
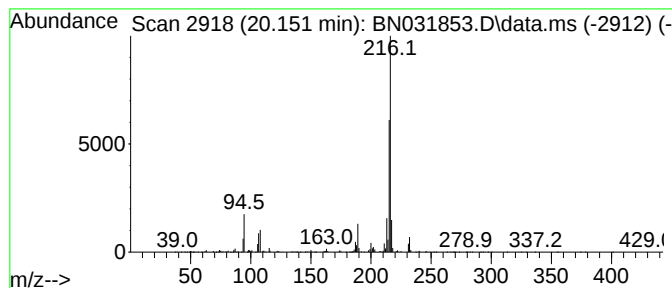
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Fluoranthene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.151	2.73 ng/ul	1560110	Chrysene-d12	21.304

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	94
5	Pyrene, 4-methyl-		216	C17H12		003353-12-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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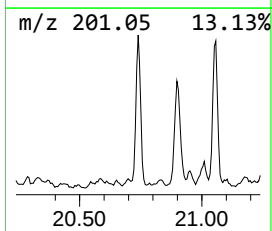
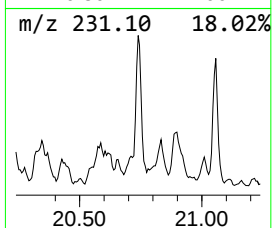
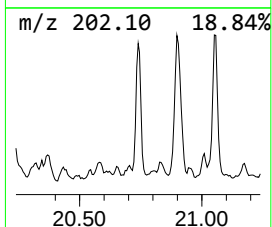
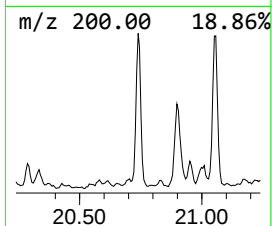
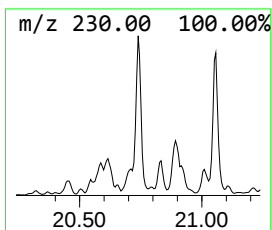
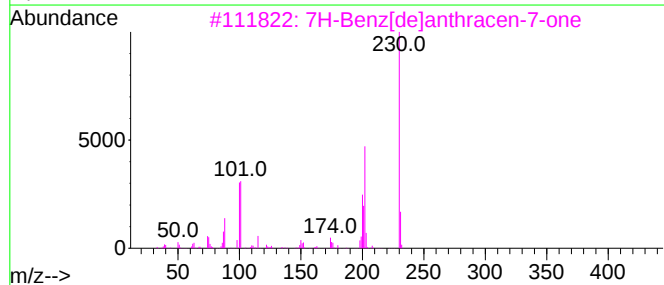
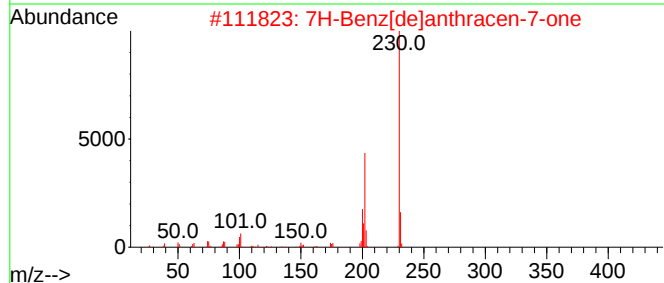
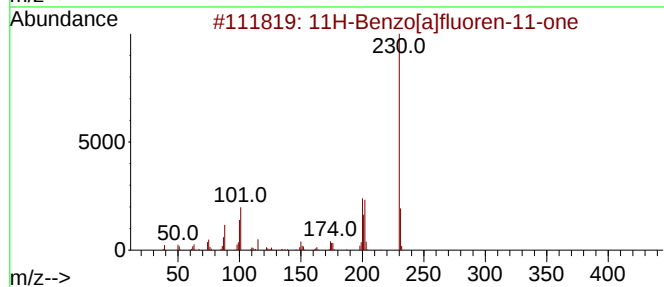
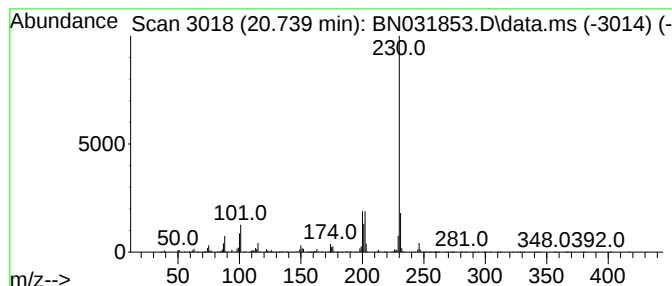
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 11H-Benzo[a]fluoren-11-one Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.739	2.11 ng/ul	1208590	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			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
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	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBW8

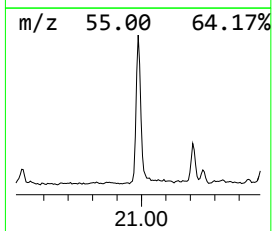
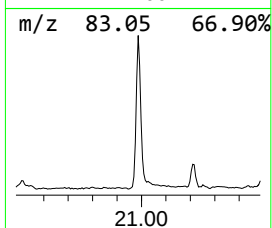
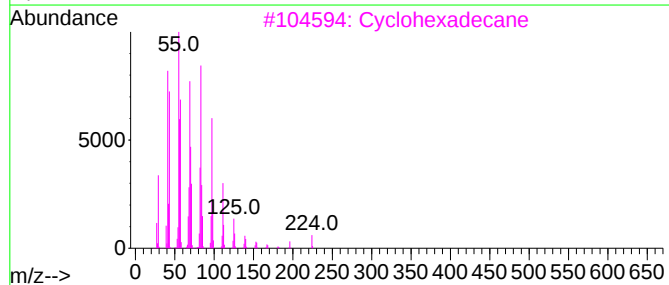
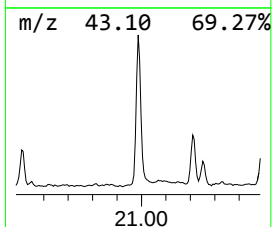
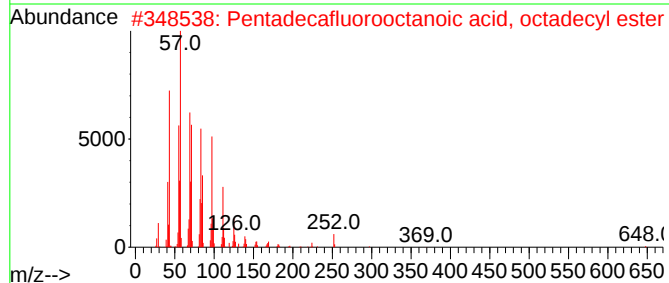
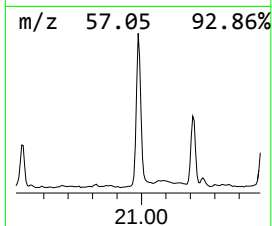
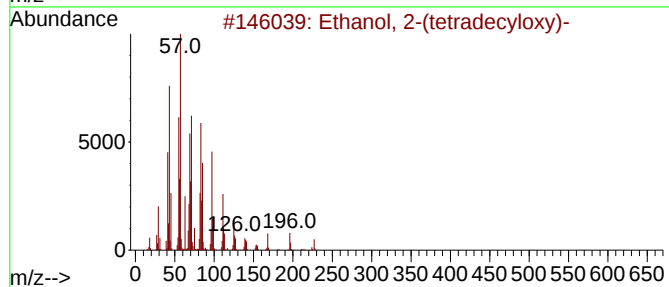
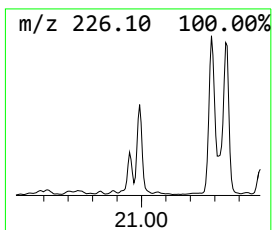
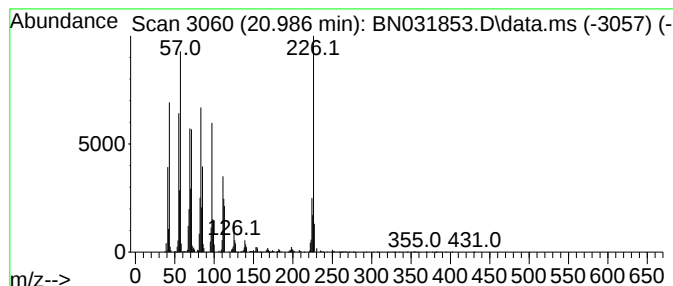
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Ethanol, 2-(tetradecyloxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	4.89 ng/ul	2794740	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	96
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	90
3			Cyclohexadecane	224	C16H32	000295-65-8	89
4			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	60
5			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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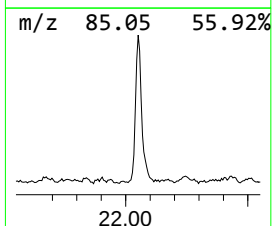
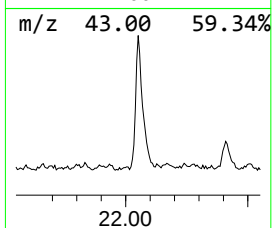
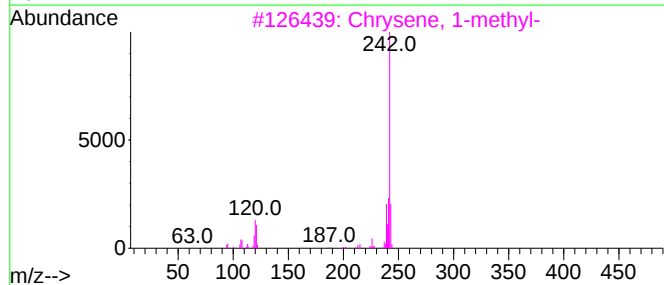
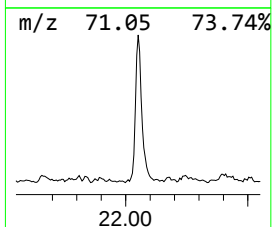
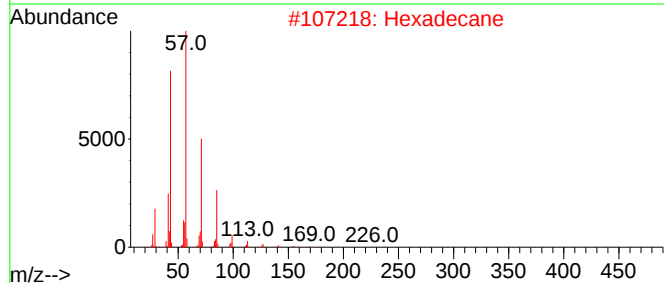
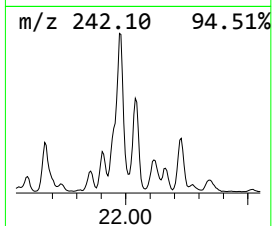
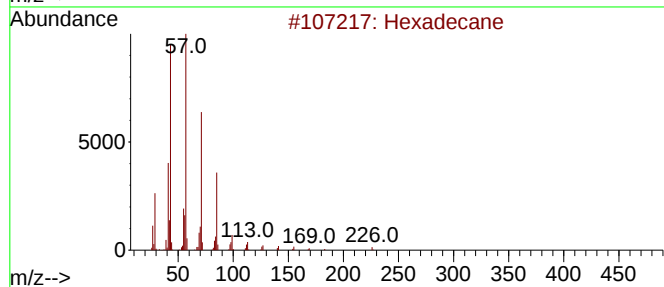
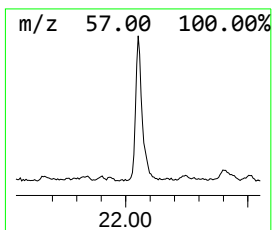
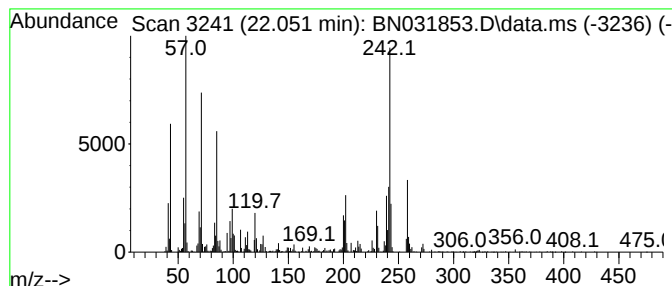
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.051	2.19 ng/ul	1250170	Chrysene-d12	21.304

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	74
2		Hexadecane	226	C16H34	000544-76-3	72
3		Chrysene, 1-methyl-	242	C19H14	003351-28-8	60
4		Chrysene, 5-methyl-	242	C19H14	003697-24-3	55
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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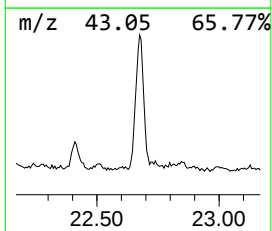
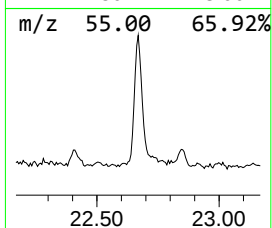
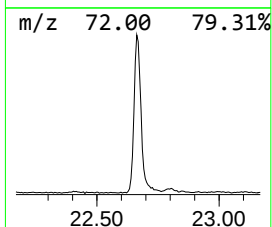
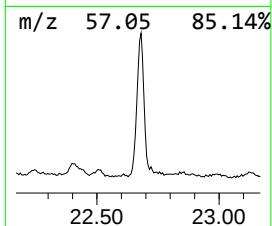
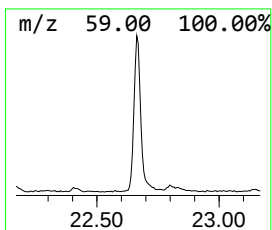
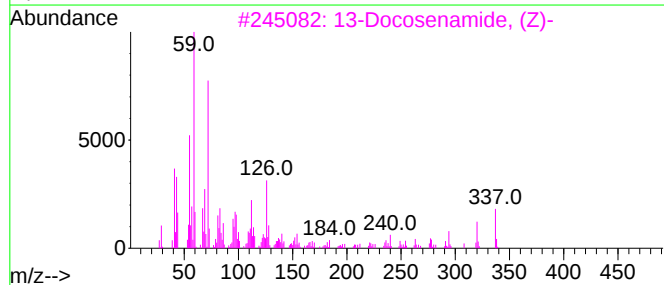
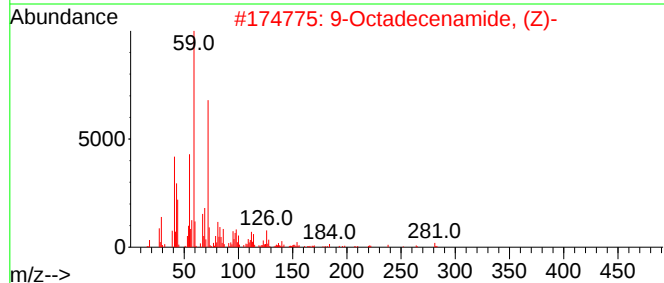
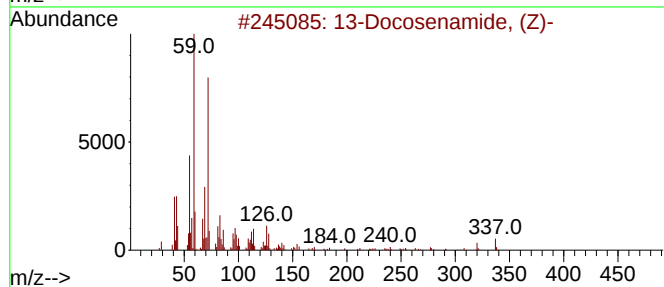
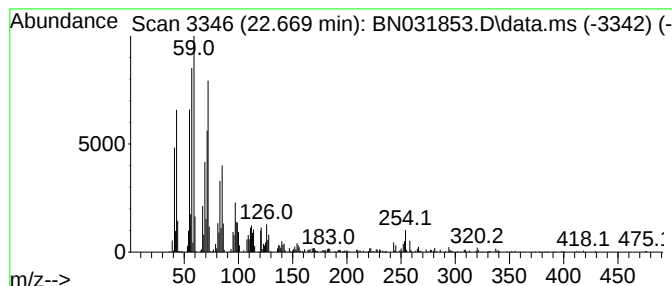
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 13-Docosenamide, (Z)- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.669	2.60 ng/ul	1485080	Chrysene-d12	21.304

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	96
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
3		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	78
4		Palmitoleamide	253	C16H31NO	106010-22-4	78
5		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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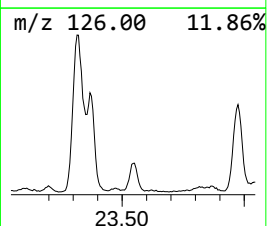
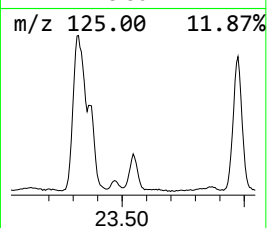
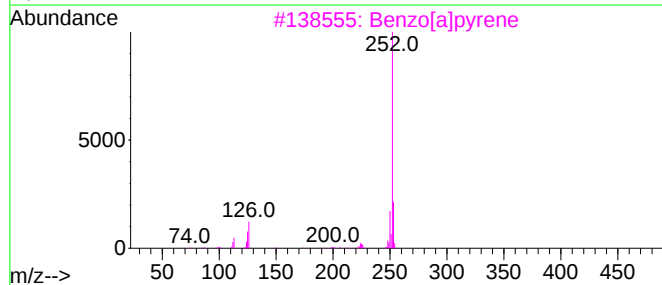
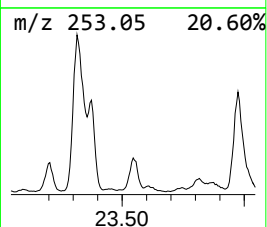
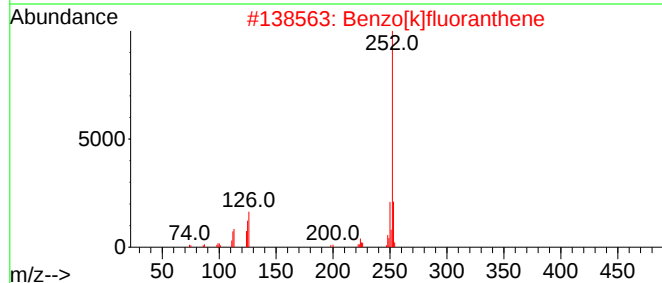
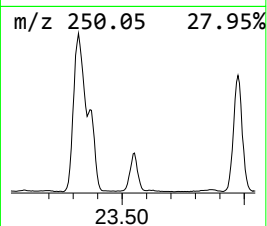
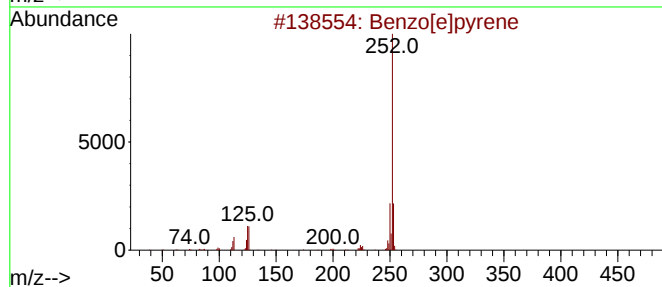
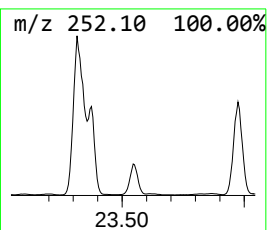
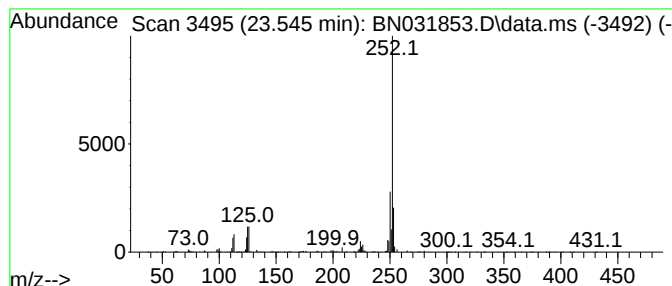
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzo[e]pyrene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.545	2.49 ng/ul	717874	Perylene-d12	24.245

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBW8

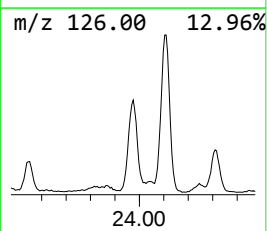
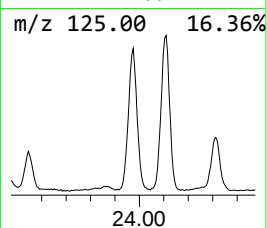
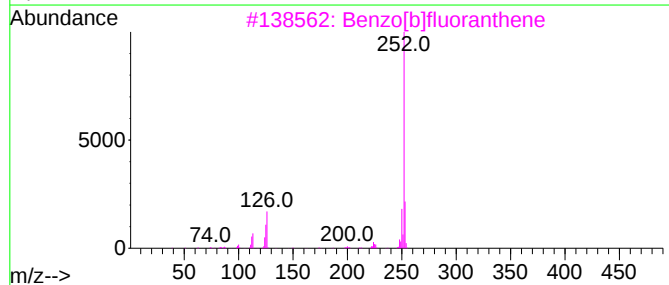
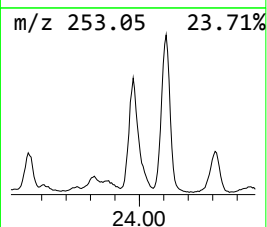
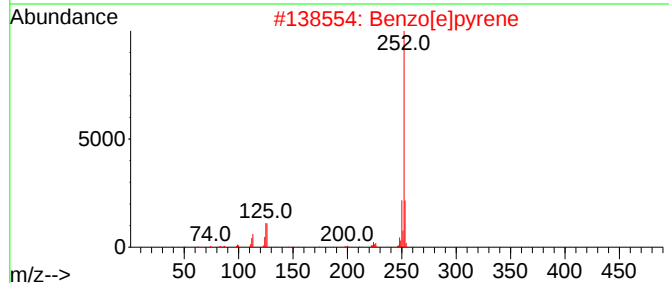
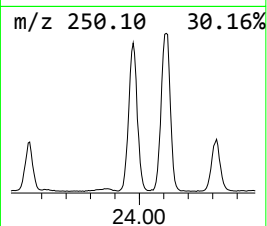
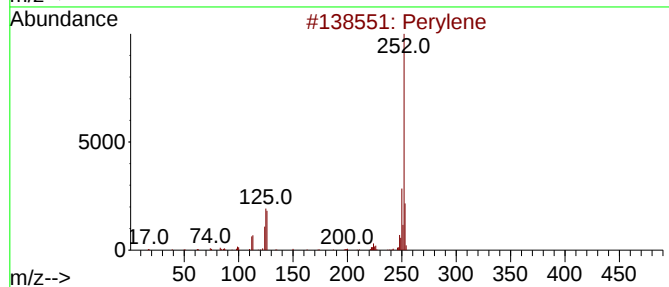
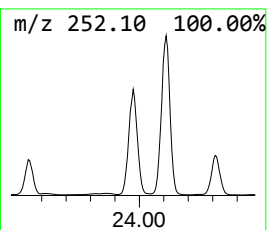
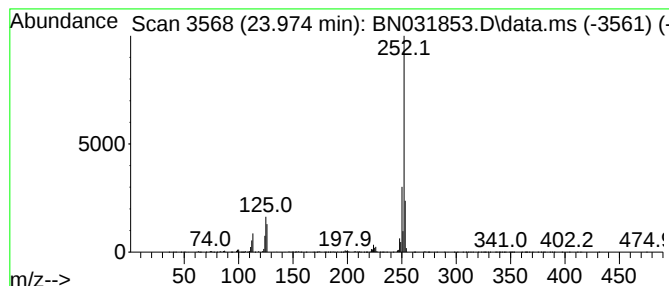
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Perylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.974	9.82 ng/ul	2835470	Perylene-d12	24.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	98
2			Benzo[e]pyrene	252	C20H12	000192-97-2	98
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	96
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
5			Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060824\
Data File : BN031853.D
Acq On : 08 Jun 2024 11:16
Operator : MA/JU
Sample : P2634-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBW8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.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
Propylene Glycol	3.446	6.1	ng/ul	811936	1	7.669	2653360	20.0
Ethanol, 2-(2-e...	7.275	2.6	ng/ul	347465	1	7.669	2653360	20.0
Pentanedioic ac...	9.393	2.1	ng/ul	415681	2	10.451	4031170	20.0
1-Phenoxypropan...	11.198	4.7	ng/ul	954864	2	10.451	4031170	20.0
Phenanthrene, 2...	17.939	4.5	ng/ul	1436540	4	17.063	6357420	20.0
n-Hexadecanoic ...	17.969	14.4	ng/ul	4561380	4	17.063	6357420	20.0
1H-Cyclopropa[1...	18.128	7.0	ng/ul	2222450	4	17.063	6357420	20.0
Anthracene, 2-m...	18.169	3.3	ng/ul	1051350	4	17.063	6357420	20.0
Tricyclo[8.2.2....	18.445	2.5	ng/ul	810149	4	17.063	6357420	20.0
9,10-Anthracene...	18.486	3.4	ng/ul	1084710	4	17.063	6357420	20.0
Phenanthrene, 2...	18.863	6.1	ng/ul	1947220	4	17.063	6357420	20.0
Phenanthrene, 2...	18.922	3.1	ng/ul	973859	4	17.063	6357420	20.0
Cyclopenta(def)...	19.004	4.0	ng/ul	1272500	4	17.063	6357420	20.0
Pyrene, 1-methyl-	19.816	3.0	ng/ul	1715470	5	21.304	11429100	20.0
11H-Benzo[b]flu...	19.986	3.4	ng/ul	1932720	5	21.304	11429100	20.0
Fluoranthene, 2...	20.151	2.7	ng/ul	1560110	5	21.304	11429100	20.0
11H-Benzo[a]flu...	20.739	2.1	ng/ul	1208590	5	21.304	11429100	20.0
Ethanol, 2-(tet...	20.986	4.9	ng/ul	2794740	5	21.304	11429100	20.0
(DEL) Alkane: S...	22.051	2.2	ng/ul	1250170	5	21.304	11429100	20.0
13-Docosenamide...	22.669	2.6	ng/ul	1485080	5	21.304	11429100	20.0
Benzo[e]pyrene	23.545	2.5	ng/ul	717874	6	24.245	5774800	20.0
Perylene	23.974	9.8	ng/ul	2835470	6	24.245	5774800	20.0