

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
 Data File : BG061649.D
 Acq On : 22 Jun 2024 15:03
 Operator : MA/JU
 Sample : P2775-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4D46

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_G\Methods\SFAM-EPA-BG062024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061649.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.469	27	31	36	rBV4	14218	22730	1.07%	0.081%
2	3.740	73	77	83	rBV	18570	29523	1.38%	0.106%
3	3.898	101	104	109	rBV2	22911	32030	1.50%	0.115%
4	4.004	117	122	128	rVB4	18385	30739	1.44%	0.110%
5	4.603	222	224	228	rBV3	4102	6894	0.32%	0.025%
6	4.738	244	247	250	rVV4	8393	11564	0.54%	0.041%
7	4.791	252	256	264	rVB2	17707	30199	1.42%	0.108%
8	5.150	312	317	325	rVB	41436	74345	3.49%	0.266%
9	5.238	328	332	338	rVB7	8864	14934	0.70%	0.053%
10	5.761	419	421	425	rVB3	3995	5778	0.27%	0.021%
11	6.137	480	485	493	rVB3	13247	23797	1.12%	0.085%
12	6.930	618	620	627	rBV5	4579	6829	0.32%	0.024%
13	7.024	631	636	638	rBV3	4536	7346	0.34%	0.026%
14	7.218	662	669	680	rVB	308585	531116	24.90%	1.899%
15	7.400	694	700	707	rBV	195523	317502	14.89%	1.135%
16	7.600	727	734	742	rBV	259214	451710	21.18%	1.615%
17	7.653	742	743	748	rBV2	5780	7906	0.37%	0.028%
18	7.764	758	762	765	rBV2	3615	5813	0.27%	0.021%
19	8.076	808	815	822	rBV	259904	432175	20.26%	1.545%
20	8.123	822	823	830	rVB5	9027	14170	0.66%	0.051%
21	8.769	925	933	940	rVV	334394	572100	26.83%	2.046%
22	8.898	949	955	962	rBV4	21459	49123	2.30%	0.176%
23	9.239	1006	1013	1020	rBV	268711	475381	22.29%	1.700%
24	9.797	1103	1108	1114	rVB3	27504	41813	1.96%	0.150%
25	9.962	1128	1136	1143	rBV2	230877	399364	18.73%	1.428%
26	10.067	1151	1154	1158	rBV3	6864	9224	0.43%	0.033%
27	10.156	1165	1169	1173	rVB4	6955	10427	0.49%	0.037%
28	10.326	1195	1198	1200	rVB4	5993	5869	0.28%	0.021%
29	10.496	1218	1227	1237	rBV	363851	645032	30.24%	2.306%
30	10.890	1284	1294	1300	rBV	389882	718828	33.71%	2.570%
31	10.943	1300	1303	1307	rVV4	12710	14824	0.70%	0.053%
32	11.019	1309	1316	1325	rVB	215505	406819	19.08%	1.455%
33	11.413	1379	1383	1389	rVB6	13912	22865	1.07%	0.082%
34	11.624	1411	1419	1427	rBV3	41366	97240	4.56%	0.348%
35	12.294	1529	1533	1536	rBV4	8187	10853	0.51%	0.039%
36	12.394	1547	1550	1555	rVB4	9871	13834	0.65%	0.049%

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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_G\Methods\SFAM-EPA-BG062024.MA.M
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37	12.465	1558	1562	1564	rBV3	7267	11716	0.55%	0.042%
38	12.541	1569	1575	1579	rVV2	12884	19721	0.92%	0.071%
39	12.606	1583	1586	1590	rBV4	5115	5711	0.27%	0.020%
40	12.758	1605	1612	1618	rBV4	9873	20155	0.95%	0.072%
41	13.534	1738	1744	1751	rBV5	15214	31978	1.50%	0.114%
42	13.740	1775	1779	1780	rVV	5270	6510	0.31%	0.023%
43	13.875	1798	1802	1808	rVB5	10810	20580	0.96%	0.074%
44	14.027	1823	1828	1834	rBV4	16155	28663	1.34%	0.102%
45	14.110	1834	1842	1848	rVV	600986	905369	42.45%	3.237%
46	14.180	1851	1854	1857	rVB3	4564	6257	0.29%	0.022%
47	14.221	1857	1861	1863	rBV4	5586	7310	0.34%	0.026%
48	14.339	1878	1881	1884	rBV5	9343	11074	0.52%	0.040%
49	14.398	1884	1891	1905	rVB	696034	1123343	52.67%	4.017%
50	14.597	1921	1925	1929	rVB3	12540	15671	0.73%	0.056%
51	14.639	1929	1932	1935	rBV2	9571	12887	0.60%	0.046%
52	14.703	1937	1943	1949	rVB	652840	991605	46.50%	3.546%
53	14.762	1949	1953	1960	rVB	88330	137273	6.44%	0.491%
54	14.874	1966	1972	1975	rBV	322463	549760	25.78%	1.966%
55	15.103	2005	2011	2019	rVB2	60994	105320	4.94%	0.377%
56	15.167	2019	2022	2028	rBV6	8529	14349	0.67%	0.051%
57	15.291	2039	2043	2044	rBV3	5079	6071	0.28%	0.022%
58	15.314	2045	2047	2052	rVB3	9061	8652	0.41%	0.031%
59	15.361	2052	2055	2057	rBV4	6987	8596	0.40%	0.031%
60	15.490	2072	2077	2082	rBV7	11187	25752	1.21%	0.092%
61	15.690	2104	2111	2117	rVV	893662	1351136	63.35%	4.831%
62	15.749	2117	2121	2124	rVV2	94053	139183	6.53%	0.498%
63	15.796	2124	2129	2133	rVB2	189277	278249	13.05%	0.995%
64	15.955	2152	2156	2160	rVB6	36445	51937	2.44%	0.186%
65	16.037	2166	2170	2175	rVB2	26592	39025	1.83%	0.140%
66	16.119	2181	2184	2185	rBV3	5184	6322	0.30%	0.023%
67	16.184	2191	2195	2201	rBV4	30646	57844	2.71%	0.207%
68	16.407	2230	2233	2240	rVB5	19355	28563	1.34%	0.102%
69	16.554	2255	2258	2259	rBV3	8180	8808	0.41%	0.031%
70	16.671	2275	2278	2283	rBV2	17974	25006	1.17%	0.089%
71	16.977	2327	2330	2333	rBV4	21359	20685	0.97%	0.074%
72	17.094	2347	2350	2358	rVB4	27823	42735	2.00%	0.153%
73	17.200	2361	2368	2372	rBV8	29897	66463	3.12%	0.238%
74	17.265	2375	2379	2382	rVB	51364	67502	3.17%	0.241%
75	17.324	2385	2389	2392	rBV6	19255	29507	1.38%	0.106%
76	17.359	2392	2395	2399	rVB5	22751	29349	1.38%	0.105%

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77	17.447	2402	2410	2413	rBV	770317	1260716	59.11%	4.508%
78	17.488	2413	2417	2421	rVB	953783	1325382	62.15%	4.739%
79	17.541	2421	2426	2431	rBV	857541	1359586	63.75%	4.862%
80	17.623	2437	2440	2445	rVB5	42474	69439	3.26%	0.248%
81	17.841	2474	2477	2481	rVB2	65821	88002	4.13%	0.315%
82	17.964	2496	2498	2504	rVB5	25402	35305	1.66%	0.126%
83	18.040	2505	2511	2518	rVB6	64012	154660	7.25%	0.553%
84	18.328	2554	2560	2563	rBV3	260729	418829	19.64%	1.498%
85	18.364	2563	2566	2572	rVB	157232	237248	11.12%	0.848%
86	18.446	2577	2580	2585	rVB3	62638	84997	3.99%	0.304%
87	18.510	2585	2591	2595	rBV3	195315	355215	16.66%	1.270%
88	18.610	2606	2608	2610	rBV3	27864	25813	1.21%	0.092%
89	18.816	2639	2643	2646	rBV	71381	95817	4.49%	0.343%
90	19.057	2682	2684	2689	rVB3	47552	54666	2.56%	0.195%
91	19.362	2733	2736	2746	rVB8	61416	133949	6.28%	0.479%
92	19.492	2752	2758	2764	rVB	1373557	1954881	91.66%	6.990%
93	19.821	2809	2814	2817	rBV2	1171308	1879717	88.14%	6.721%
94	19.850	2817	2819	2827	rVB	967991	1214506	56.95%	4.343%
95	20.438	2916	2919	2923	rVB	121002	147178	6.90%	0.526%
96	21.366	3073	3077	3084	rVB5	230834	399422	18.73%	1.428%
97	21.701	3127	3134	3139	rBV2	858313	2132703	100.00%	7.626%
98	21.754	3140	3143	3155	rVB	496132	817157	38.32%	2.922%
99	23.910	3505	3510	3519	rBV3	235833	687472	32.23%	2.458%
100	24.932	3677	3684	3693	rVB2	439875	1165954	54.67%	4.169%

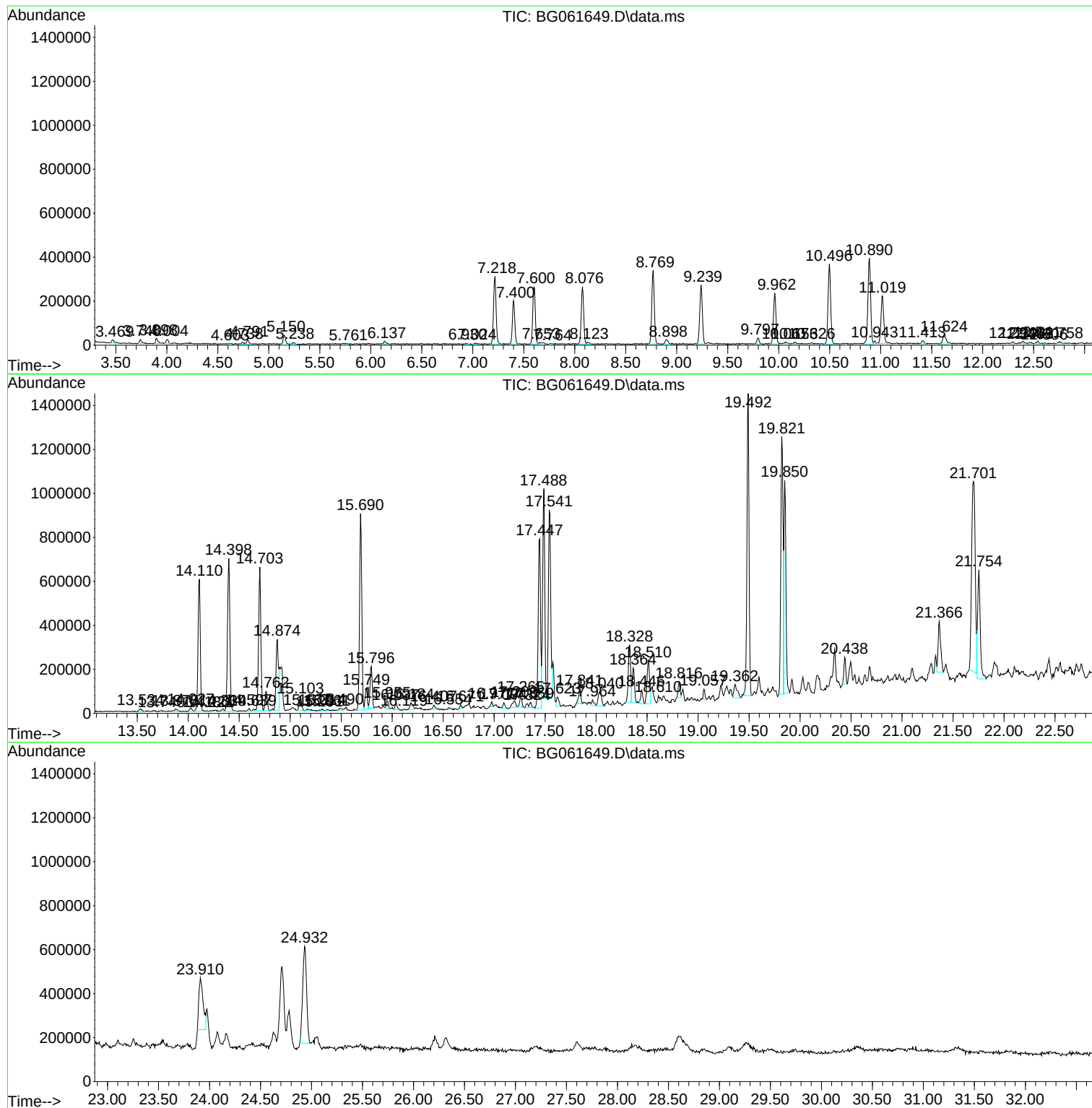
Sum of corrected areas: 27965947

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
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Instrument :
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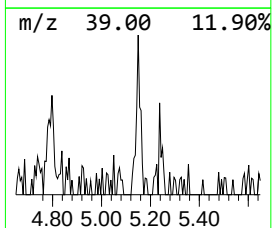
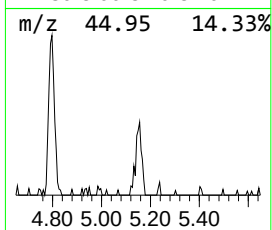
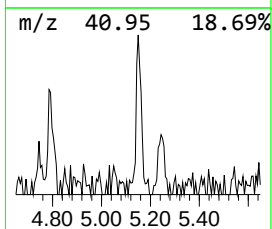
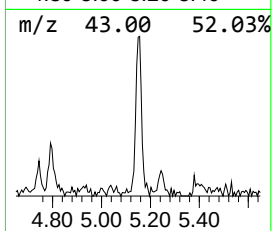
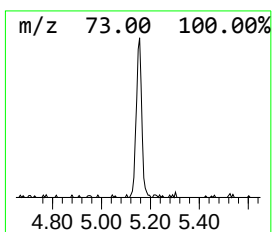
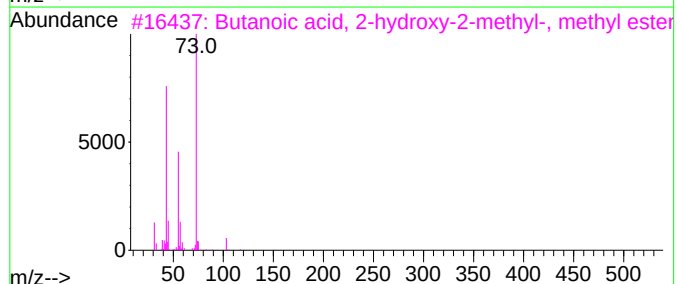
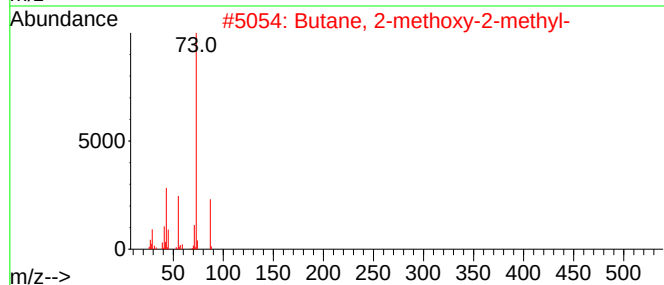
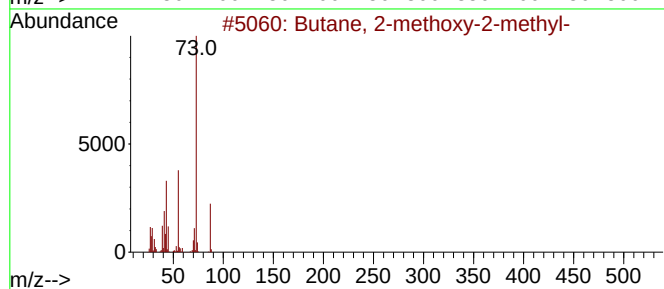
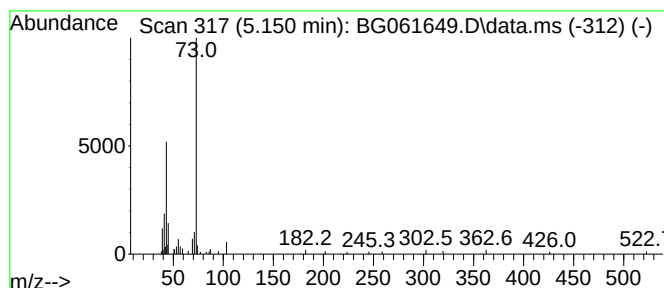
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.150	3.44 ng/ul	74345	1,4-Dichlorobenzene-d4	8.076

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	25
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	25
3			Butanoic acid, 2-hydroxy-2-methyl-	132	C6H12O3	032793-34-3	17
4			1,4-Diacetyl-3-acetoxymethyl-2,5-dimethyl-2,5-dimethyl-	318	C14H22O8	1000128-41-9	9
5			5-(2-Methoxypropan-2-yl)-2-methyl-2,5-dimethyl-	184	C11H20O2	126657-18-9	9



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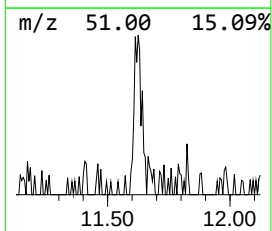
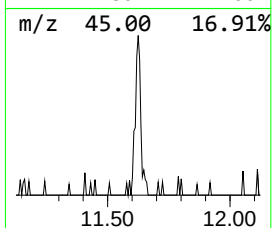
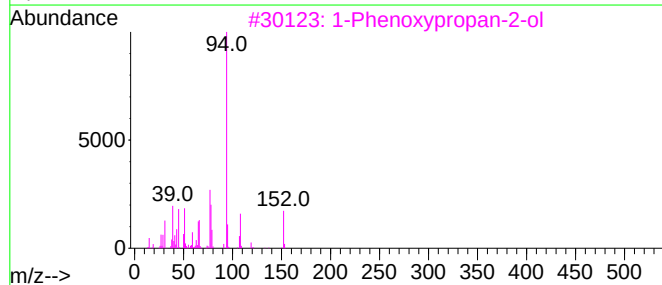
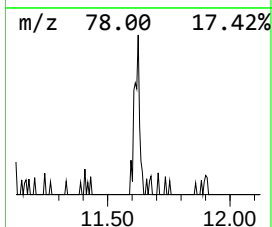
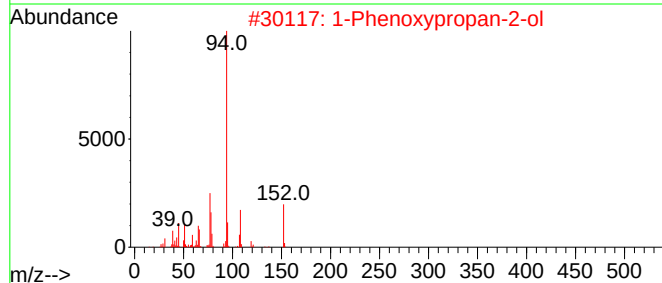
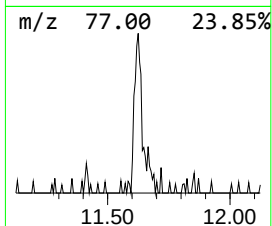
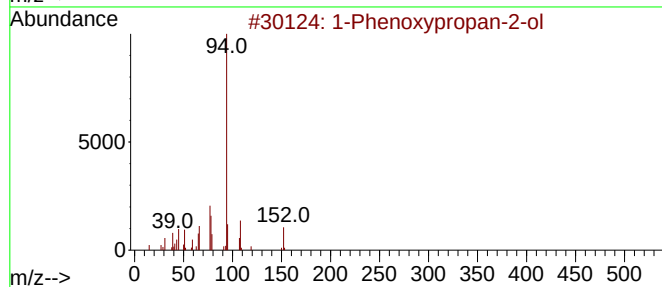
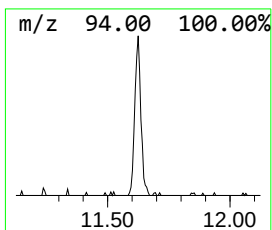
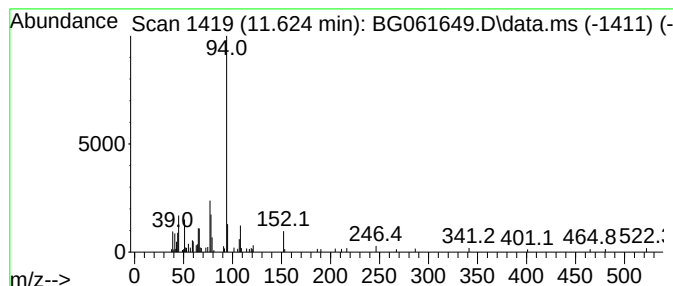
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.624	2.71 ng/ul	97240	Naphthalene-d8	10.890

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	95	
3	1-Phenoxypropan-2-ol		152	C9H12O2	000770-35-4	93	
4	1-Phenoxypropan-2-ol		152	C9H12O2	000770-35-4	87	
5	1-Propanol, 3-phenoxy-		152	C9H12O2	006180-61-6	83	



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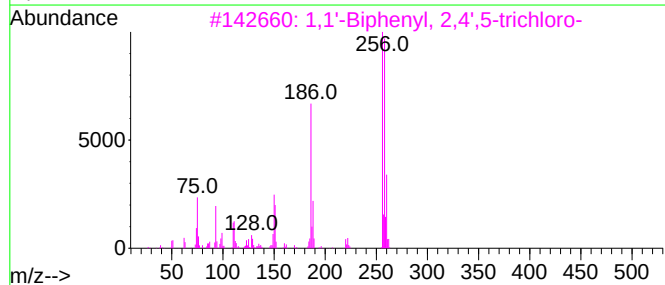
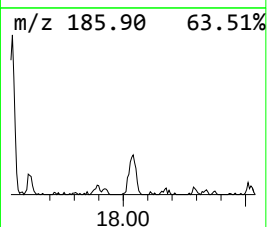
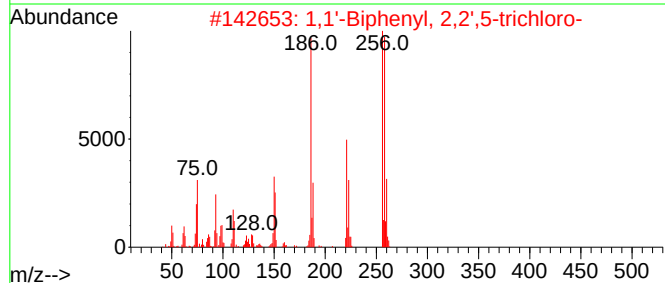
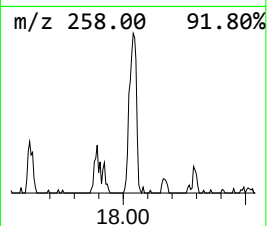
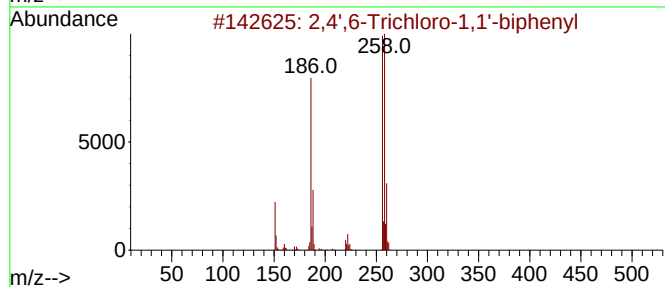
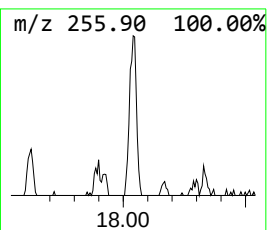
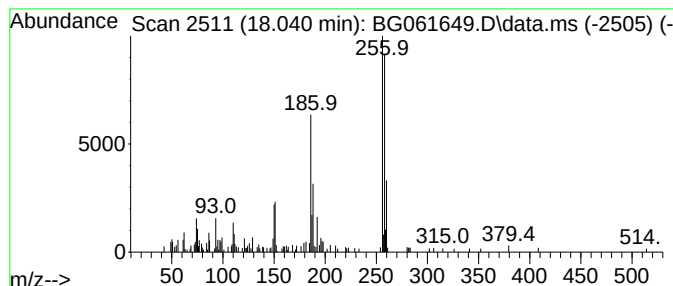
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,4',6-Trichloro-1,1'-biphenyl Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.040	2.45 ng/ul	154660	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	91		
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	91		
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	91		
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	91		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	90		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061649.D
Acq On : 22 Jun 2024 15:03
Operator : MA/JU
Sample : P2775-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D46

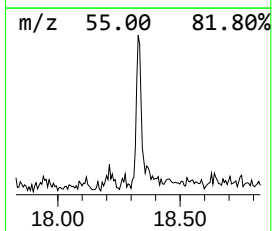
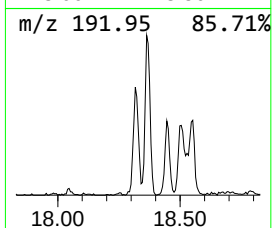
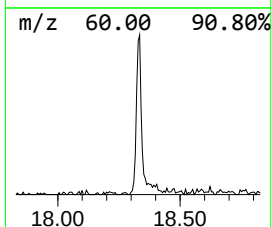
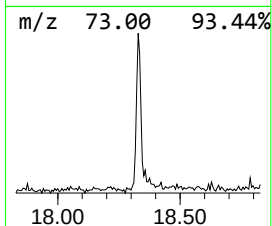
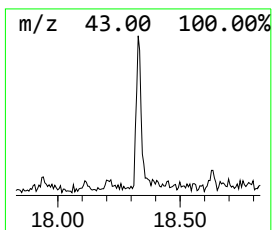
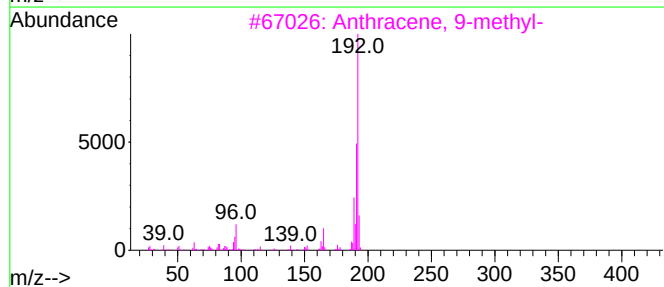
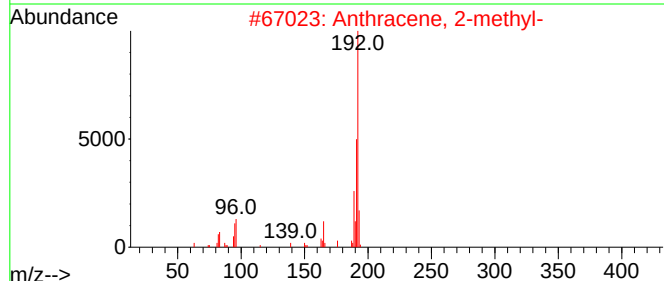
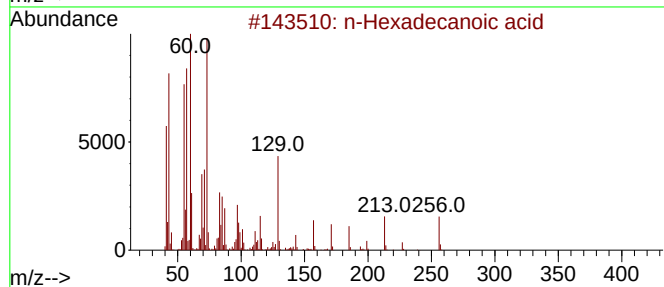
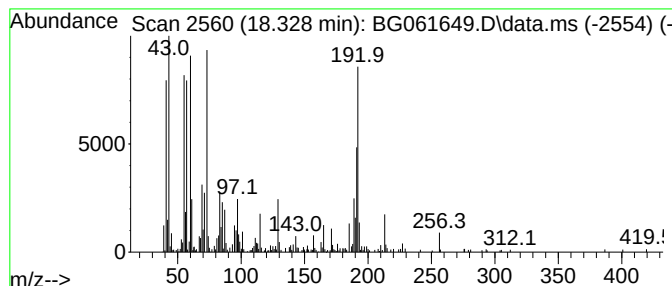
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.328	6.64 ng/ul	418829	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	78
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	70
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061649.D
Acq On : 22 Jun 2024 15:03
Operator : MA/JU
Sample : P2775-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D46

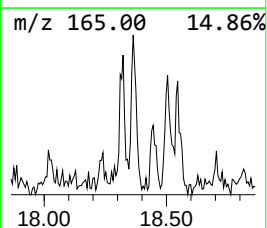
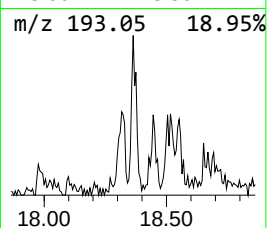
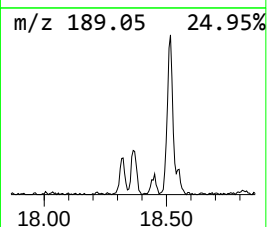
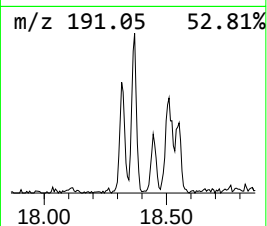
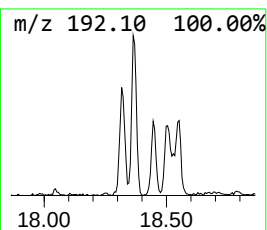
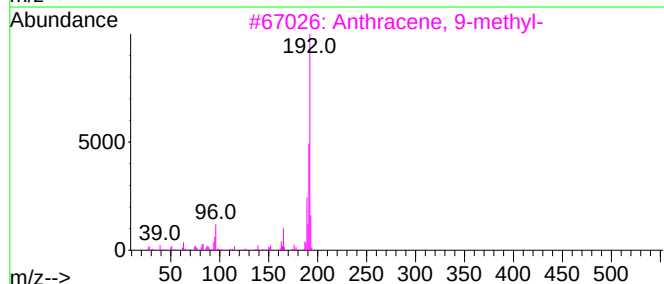
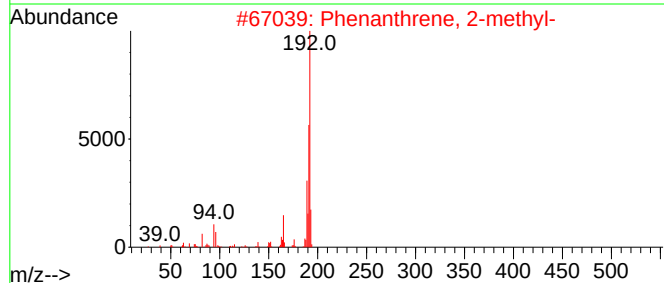
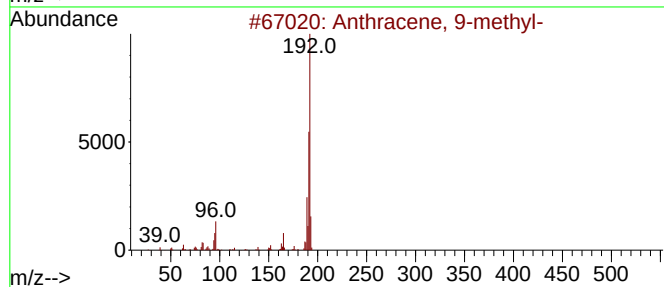
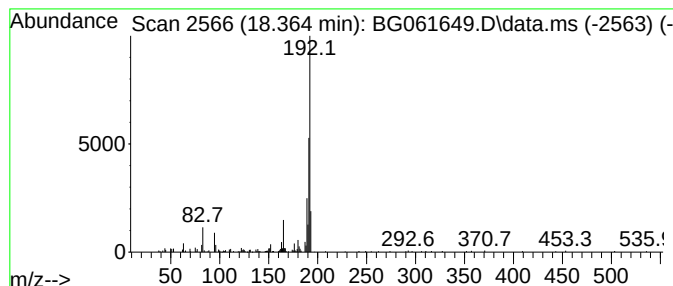
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Anthracene, 9-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.364	3.76 ng/ul	237248	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061649.D
Acq On : 22 Jun 2024 15:03
Operator : MA/JU
Sample : P2775-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

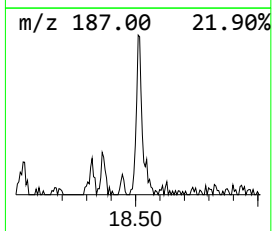
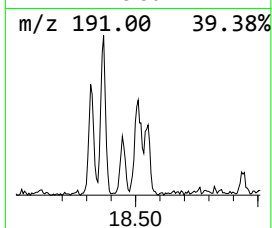
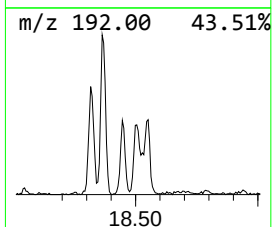
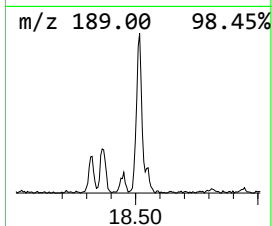
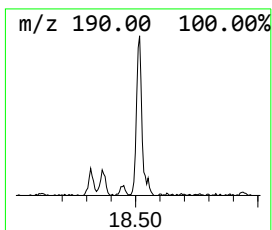
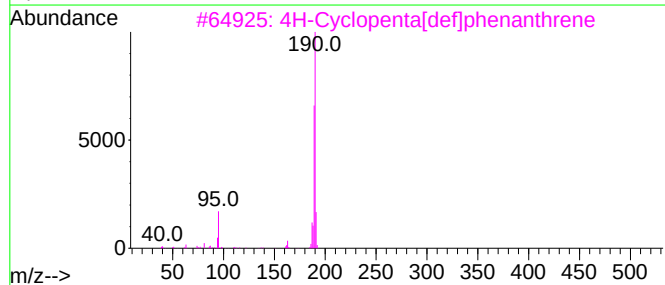
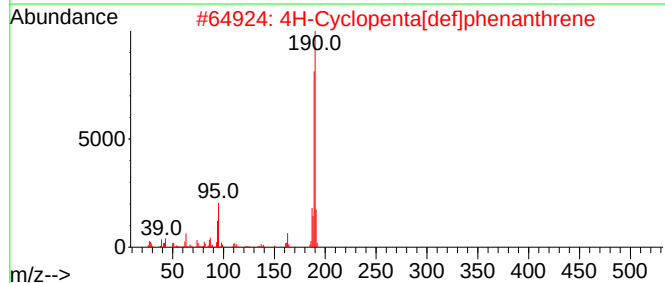
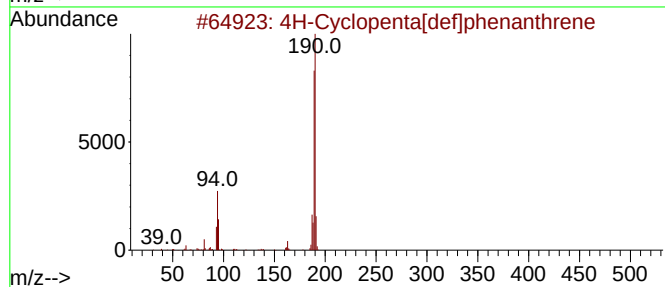
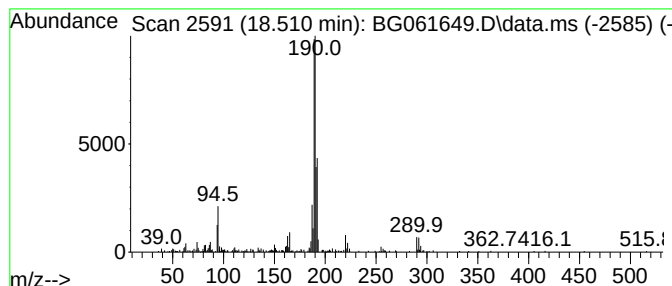
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.510	5.64 ng/ul	355215	Phenanthrene-d10	17.447	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
5	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061649.D
Acq On : 22 Jun 2024 15:03
Operator : MA/JU
Sample : P2775-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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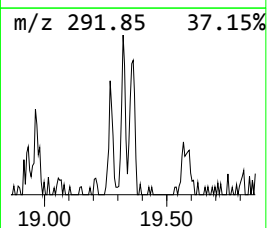
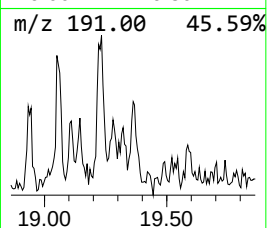
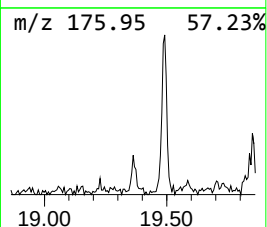
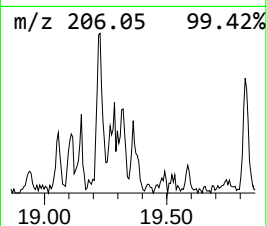
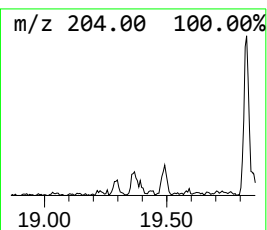
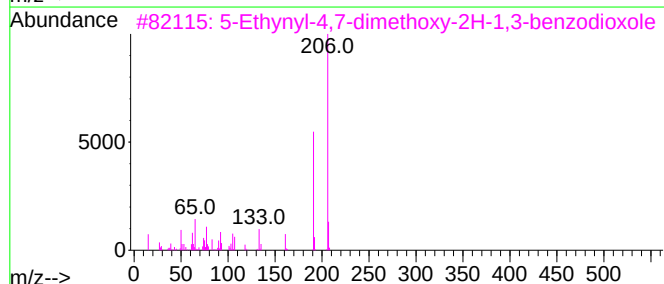
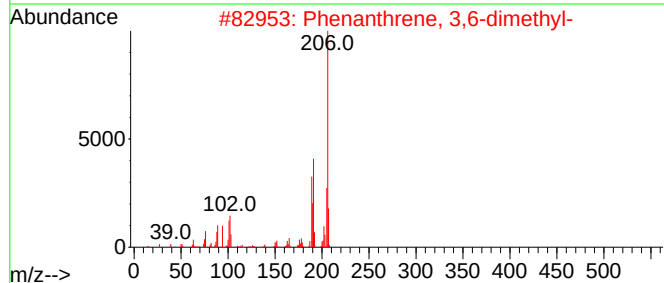
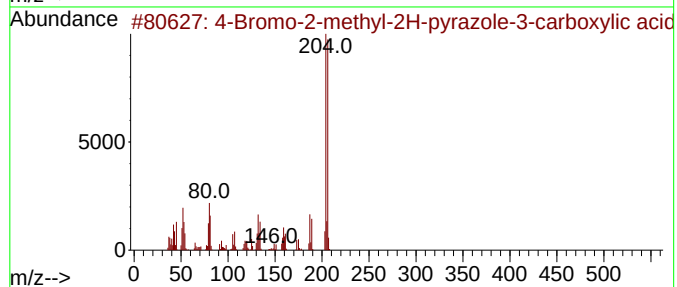
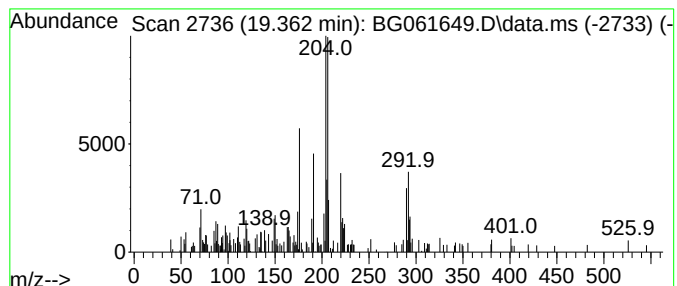
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.362	2.12 ng/ul	133949	Phenanthrene-d10	17.447

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Bromo-2-methyl-2H-pyrazole-3-c...	204	C5H5BrN2O2	1000300-50-8	35
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	30
3			5-Ethynyl-4,7-dimethoxy-2H-1,3-b...	206	C11H10O4	1000411-41-8	25
4			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	25
5			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061649.D
Acq On : 22 Jun 2024 15:03
Operator : MA/JU
Sample : P2775-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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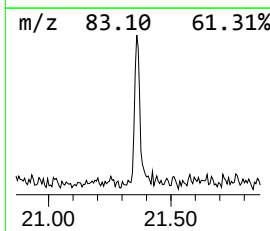
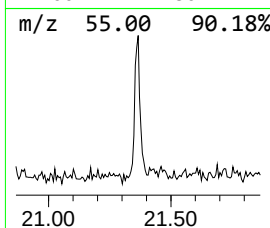
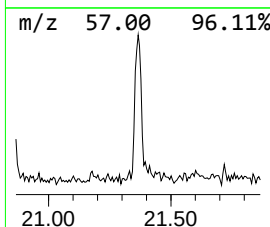
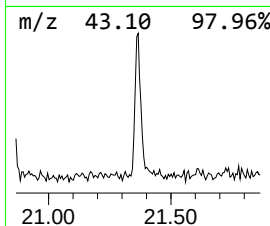
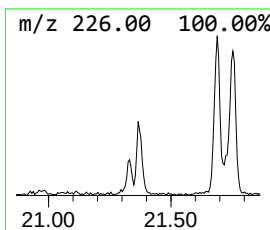
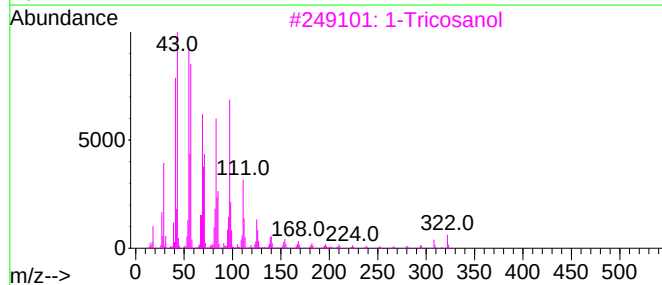
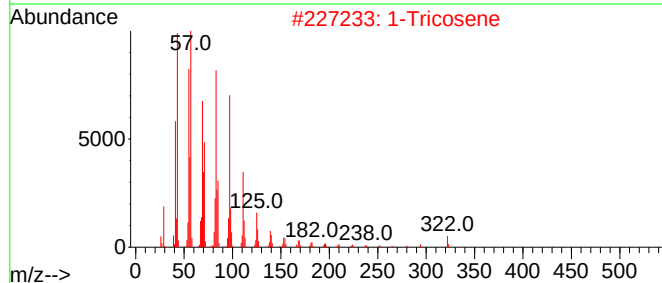
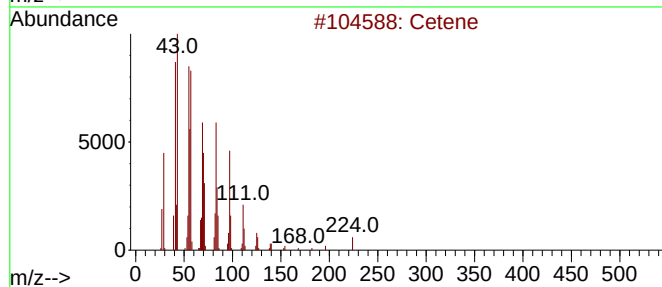
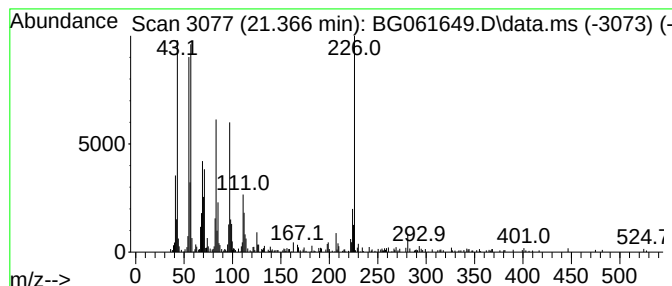
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.366	3.75 ng/ul	399422	Chrysene-d12	21.707

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cetene	224	C16H32	000629-73-2	60
2			1-Tricosene	322	C23H46	018835-32-0	53
3			1-Tricosanol	340	C23H48O	003133-01-5	53
4			1-Pentadecene	210	C15H30	013360-61-7	49
5			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061649.D
Acq On : 22 Jun 2024 15:03
Operator : MA/JU
Sample : P2775-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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1-Phenoxypropan...	11.624	2.7	ng/ul	97240	2	10.890	718828	20.0
2,4',6-Trichlor...	18.040	2.5	ng/ul	154660	4	17.447	1260720	20.0
n-Hexadecanoic ...	18.328	6.6	ng/ul	418829	4	17.447	1260720	20.0
Anthracene, 9-m...	18.364	3.8	ng/ul	237248	4	17.447	1260720	20.0
4H-Cyclopenta[d...	18.510	5.6	ng/ul	355215	4	17.447	1260720	20.0
unknown-02	19.362	2.1	ng/ul	133949	4	17.447	1260720	20.0
(DEL) Alkane: S...	21.366	3.8	ng/ul	399422	5	21.707	2132700	20.0