

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049674.D
 Acq On : 6 Aug 2021 12:03
 Operator : CG/JU
 Sample : M3124-13
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0094

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-BG080421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049674.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.074	3	6	15	rVB3	72840	141568	21.35%	1.465%
2	3.150	15	19	27	rVB	204577	322676	48.66%	3.339%
3	3.426	63	66	73	rBV7	4970	8821	1.33%	0.091%
4	3.855	135	139	147	rBV2	32472	64291	9.70%	0.665%
5	3.955	152	156	164	rVB3	18176	31004	4.68%	0.321%
6	4.084	176	178	182	rVB4	2570	3005	0.45%	0.031%
7	4.184	192	195	197	rBV3	3217	3395	0.51%	0.035%
8	4.401	228	232	239	rVV2	10679	16658	2.51%	0.172%
9	4.519	250	252	256	rVB3	2535	3169	0.48%	0.033%
10	4.742	287	290	296	rVB4	3064	4867	0.73%	0.050%
11	4.824	300	304	307	rBV2	4922	6085	0.92%	0.063%
12	5.118	351	354	357	rBV2	3882	4368	0.66%	0.045%
13	5.212	366	370	372	rBV3	2719	3467	0.52%	0.036%
14	7.198	701	708	716	rVV	93988	173547	26.17%	1.796%
15	7.362	728	736	747	rBV	61184	108209	16.32%	1.120%
16	7.568	765	771	782	rBV2	93889	166375	25.09%	1.722%
17	8.043	845	852	859	rBV	95991	168262	25.37%	1.741%
18	8.748	965	972	978	rBV	102942	176600	26.63%	1.827%
19	8.872	989	993	998	rVB3	3993	6150	0.93%	0.064%
20	9.212	1043	1051	1059	rBV	96631	179579	27.08%	1.858%
21	9.277	1059	1062	1064	rVB	2706	3137	0.47%	0.032%
22	9.935	1168	1174	1181	rVV3	89118	166449	25.10%	1.722%
23	10.481	1257	1267	1276	rVV	126665	235026	35.44%	2.432%
24	10.628	1285	1292	1296	rBV3	15860	28280	4.26%	0.293%
25	10.863	1320	1332	1338	rBV	132248	246825	37.22%	2.554%
26	10.910	1338	1340	1345	rVV	2710	4165	0.63%	0.043%
27	10.998	1348	1355	1364	rVV2	70939	144584	21.80%	1.496%
28	12.267	1569	1571	1576	rVB4	2695	3842	0.58%	0.040%
29	12.449	1598	1602	1607	rVB2	2318	3686	0.56%	0.038%
30	12.526	1611	1615	1618	rVB2	2960	4258	0.64%	0.044%
31	13.048	1699	1704	1707	rBV2	1611	2786	0.42%	0.029%
32	13.301	1742	1747	1750	rBV3	2487	4076	0.61%	0.042%
33	13.513	1777	1783	1788	rVB5	4613	8739	1.32%	0.090%
34	14.000	1862	1866	1867	rBV2	2593	2734	0.41%	0.028%
35	14.094	1873	1882	1889	rBV	240945	366552	55.28%	3.793%
36	14.165	1891	1894	1897	rVB3	2312	2962	0.45%	0.031%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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Sampling : 1

Min Area: 0.1 % of largest Peak

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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-BG080421.M
 Title : SVOA CALIBRATION

37	14.241	1904	1907	1910	rBV3	1792	2822	0.43%	0.029%
38	14.388	1924	1932	1943	rBV	251247	406113	61.24%	4.202%
39	14.582	1959	1965	1969	rVB4	4561	7240	1.09%	0.075%
40	14.693	1976	1984	1991	rVB	208721	328411	49.53%	3.398%
41	14.758	1991	1995	1998	rBV3	1979	2757	0.42%	0.029%
42	14.893	2009	2018	2030	rBV	87390	166547	25.12%	1.723%
43	15.046	2040	2044	2048	rBV4	4203	6370	0.96%	0.066%
44	15.093	2048	2052	2058	rVB3	8202	10879	1.64%	0.113%
45	15.310	2084	2089	2092	rBV2	2577	3695	0.56%	0.038%
46	15.480	2113	2118	2122	rBV2	6939	10365	1.56%	0.107%
47	15.522	2124	2125	2130	rVB2	5668	5897	0.89%	0.061%
48	15.686	2144	2153	2159	rBV	337546	500500	75.48%	5.179%
49	15.739	2159	2162	2167	rVB3	8365	12789	1.93%	0.132%
50	15.804	2167	2173	2182	rBV	109061	163101	24.60%	1.688%
51	15.927	2190	2194	2197	rBV4	4758	7118	1.07%	0.074%
52	16.039	2208	2213	2216	rVB2	4349	5313	0.80%	0.055%
53	16.185	2235	2238	2244	rVB5	3583	5703	0.86%	0.059%
54	16.268	2250	2252	2257	rBV2	2323	3321	0.50%	0.034%
55	16.356	2263	2267	2270	rVV2	2520	3056	0.46%	0.032%
56	16.391	2270	2273	2283	rVB5	7485	17816	2.69%	0.184%
57	16.638	2312	2315	2318	rBV2	5689	7359	1.11%	0.076%
58	16.755	2333	2335	2339	rVB4	3600	2756	0.42%	0.029%
59	16.796	2339	2342	2346	rBV3	2706	3369	0.51%	0.035%
60	16.978	2367	2373	2376	rBV5	6900	11421	1.72%	0.118%
61	17.020	2377	2380	2381	rVV2	3910	4159	0.63%	0.043%
62	17.102	2388	2394	2398	rBV	9650	15637	2.36%	0.162%
63	17.184	2404	2408	2413	rVB4	8510	13054	1.97%	0.135%
64	17.266	2419	2422	2426	rVB	4847	7081	1.07%	0.073%
65	17.443	2446	2452	2457	rBV	242661	390257	58.85%	4.038%
66	17.484	2457	2459	2464	rVV	103719	145247	21.90%	1.503%
67	17.542	2464	2469	2480	rVB	346206	542266	81.78%	5.611%
68	17.636	2482	2485	2488	rVV5	3745	4722	0.71%	0.049%
69	17.754	2503	2505	2508	rVV3	3994	3569	0.54%	0.037%
70	17.848	2516	2521	2526	rBV2	11173	20429	3.08%	0.211%
71	17.930	2532	2535	2538	rVB2	6589	7444	1.12%	0.077%
72	17.965	2538	2541	2544	rBV4	4217	4604	0.69%	0.048%
73	18.048	2553	2555	2558	rVB3	3356	3307	0.50%	0.034%
74	18.100	2560	2564	2566	rBV2	8144	7860	1.19%	0.081%
75	18.206	2580	2582	2586	rBV5	3580	5196	0.78%	0.054%
76	18.324	2598	2602	2606	rBV3	34491	49615	7.48%	0.513%

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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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77	18.371	2606	2610	2618	rVB3	25188	39959	6.03%	0.413%
78	18.447	2621	2623	2628	rVB5	6826	8277	1.25%	0.086%
79	18.518	2630	2635	2638	rVV3	25342	41384	6.24%	0.428%
80	18.553	2638	2641	2646	rVB3	12352	19956	3.01%	0.207%
81	18.623	2649	2653	2657	rVB6	6833	10840	1.63%	0.112%
82	18.817	2682	2686	2689	rBV2	20157	23937	3.61%	0.248%
83	18.858	2689	2693	2698	rVB3	29022	41216	6.22%	0.427%
84	19.070	2725	2729	2731	rVB5	5142	5641	0.85%	0.058%
85	19.234	2753	2757	2762	rBV6	15477	27171	4.10%	0.281%
86	19.375	2778	2781	2790	rVB8	13293	30776	4.64%	0.318%
87	19.499	2796	2802	2808	rVB	325243	475350	71.68%	4.919%
88	19.598	2815	2819	2824	rBV7	11471	18113	2.73%	0.187%
89	19.833	2853	2859	2862	rBV	395885	663113	100.00%	6.862%
90	20.033	2891	2893	2897	rVB3	14278	19971	3.01%	0.207%
91	20.350	2944	2947	2951	rVB2	43349	56060	8.45%	0.580%
92	20.444	2961	2963	2967	rVB	21436	24225	3.65%	0.251%
93	21.120	3074	3078	3083	rVB2	29040	42266	6.37%	0.437%
94	21.349	3114	3117	3121	rBV5	25380	32358	4.88%	0.335%
95	21.602	3155	3160	3166	rVB	438597	609465	91.91%	6.307%
96	21.725	3173	3181	3186	rBV2	246873	595052	89.74%	6.158%
97	21.772	3186	3189	3197	rVB	123554	203392	30.67%	2.105%
98	23.963	3555	3562	3570	rBV	88993	281733	42.49%	2.915%
99	24.774	3694	3700	3707	rBV2	117584	298637	45.04%	3.090%
100	25.009	3731	3740	3749	rBV2	126016	355431	53.60%	3.678%

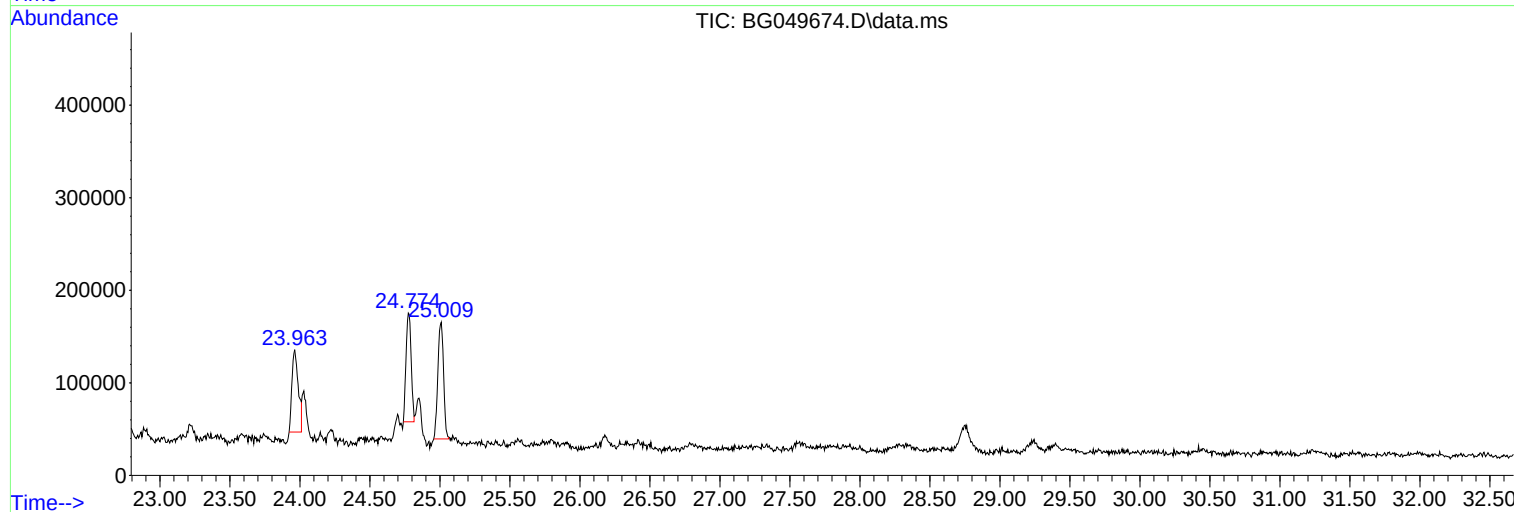
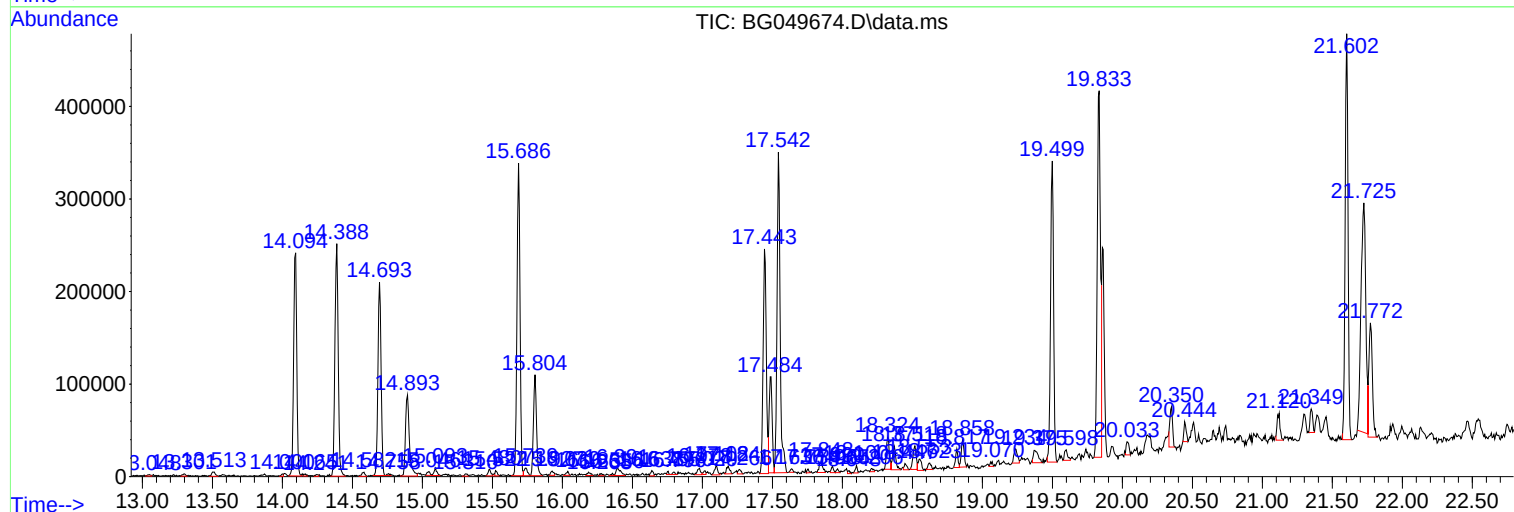
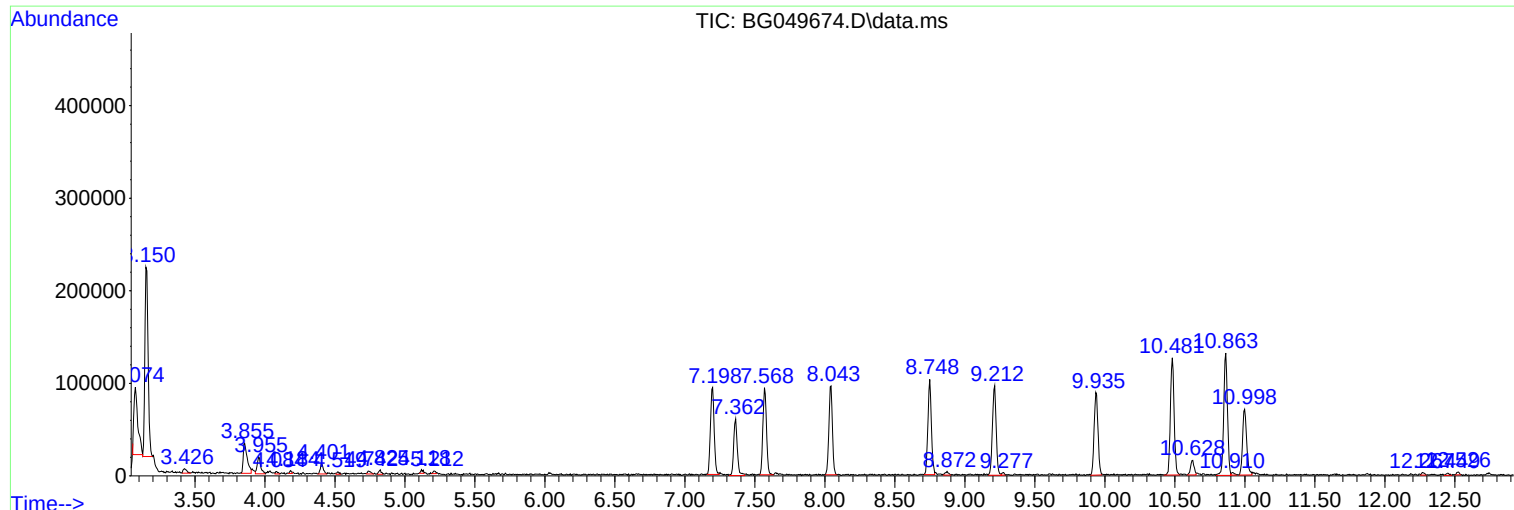
Sum of corrected areas: 9663685

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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
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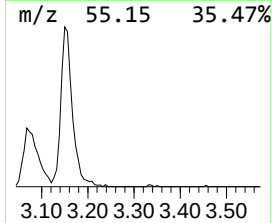
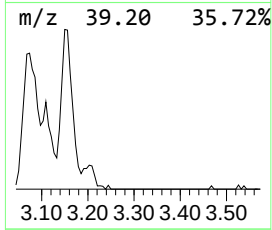
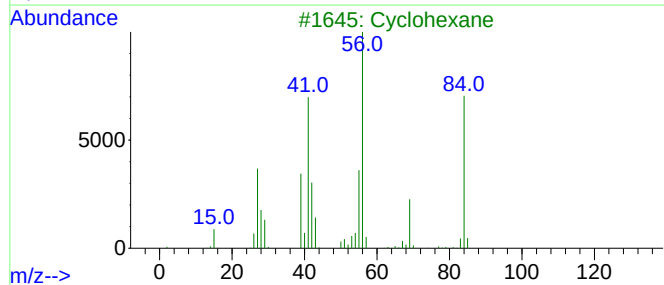
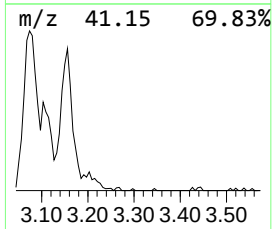
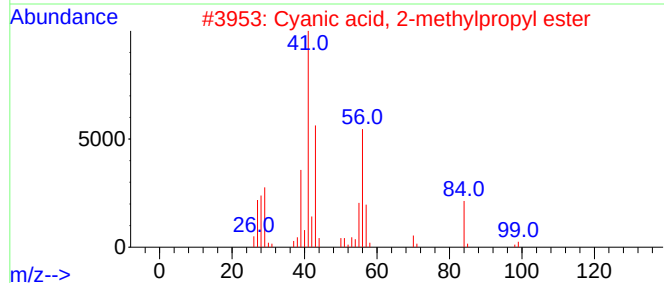
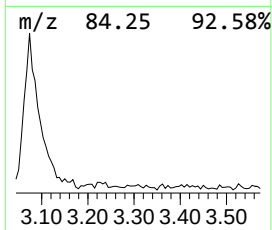
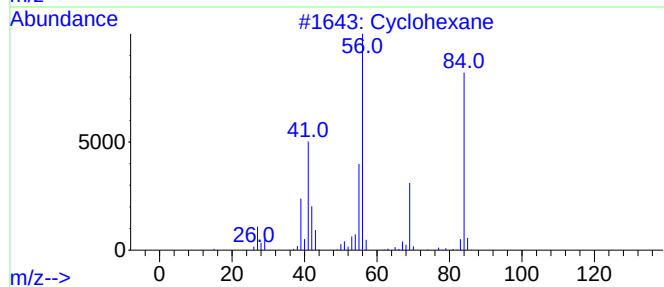
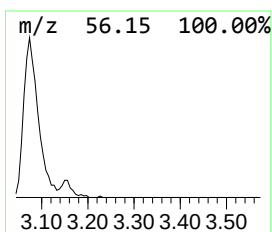
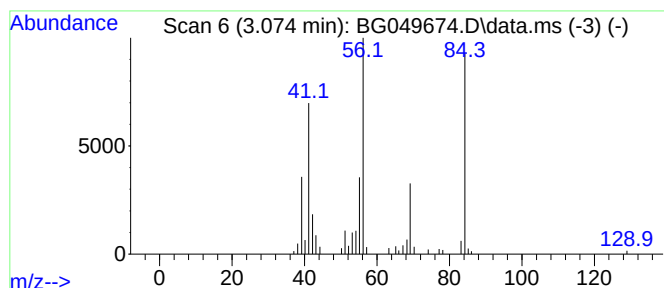
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.074 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD			R.T.
3.074	16.83 ng/ul	141568	1,4-Dichlorobenzene-d4			8.043
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexane		84	C6H12	000110-82-7	91
2	Cyanic acid, 2-methylpropyl ester		99	C5H9NO	001768-25-8	83
3	Cyclohexane		84	C6H12	000110-82-7	78
4	Cyclopentane, methyl-		84	C6H12	000096-37-7	72
5	Cyclohexane		84	C6H12	000110-82-7	72



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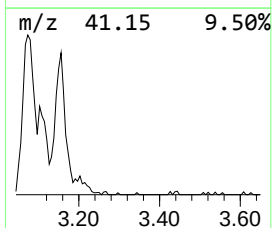
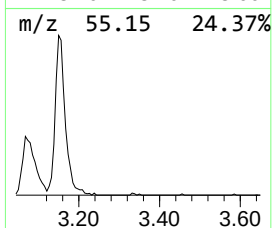
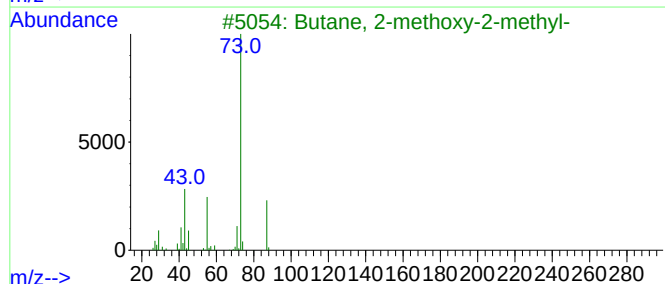
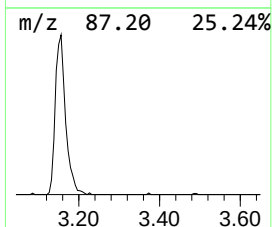
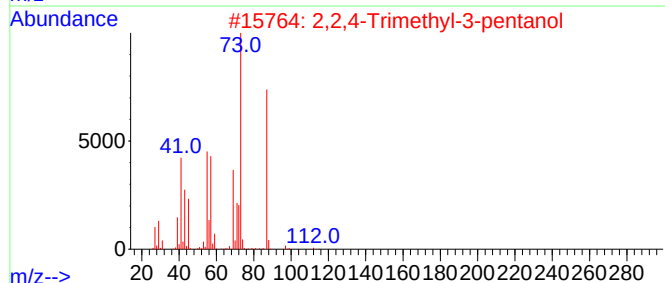
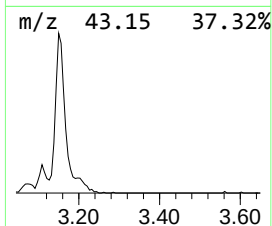
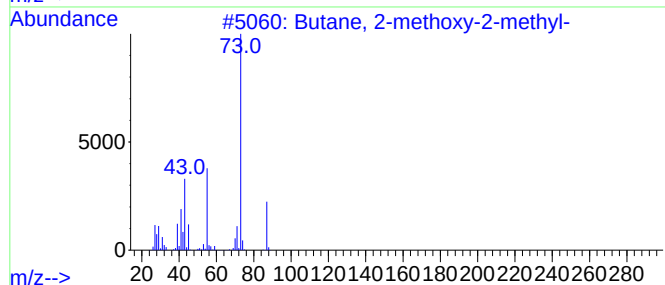
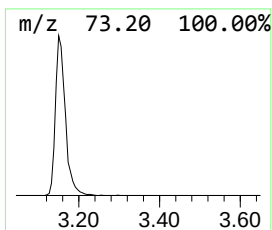
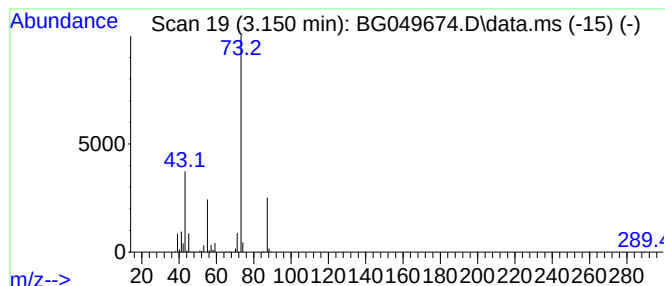
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.150	38.35 ng/ul	322676	1,4-Dichlorobenzene-d4	8.043

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	43		
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
3	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	38		
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9		
5	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	9		



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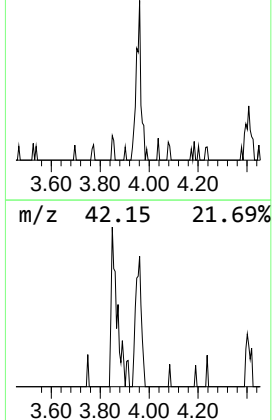
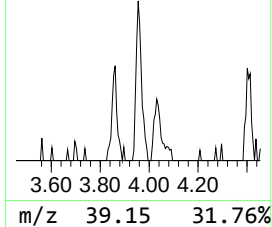
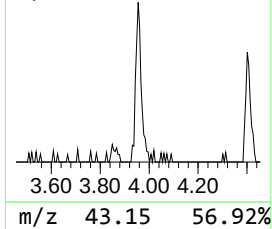
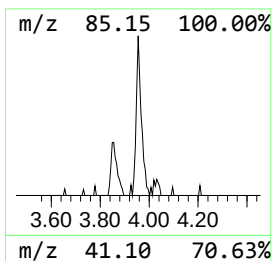
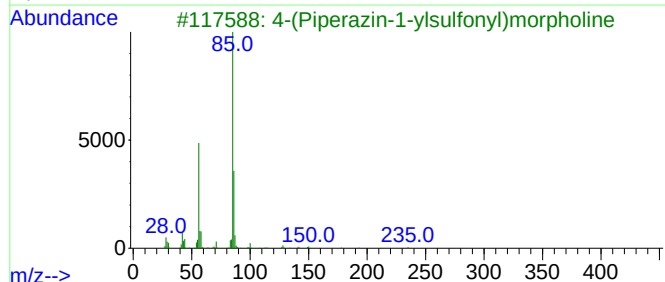
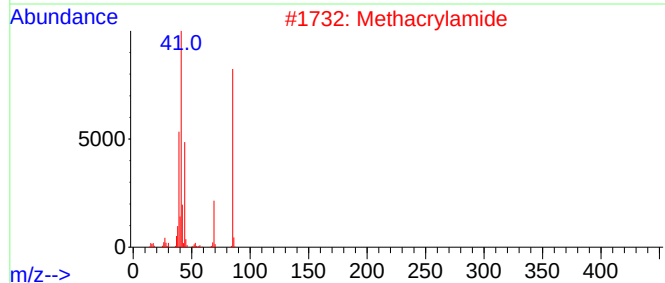
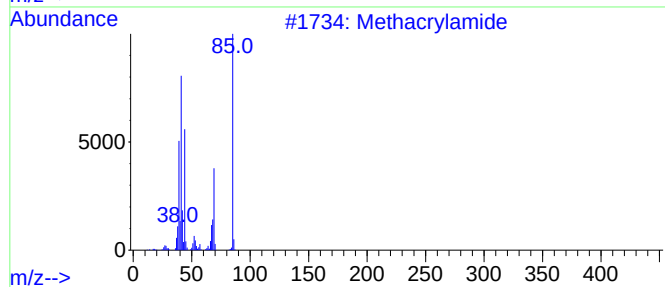
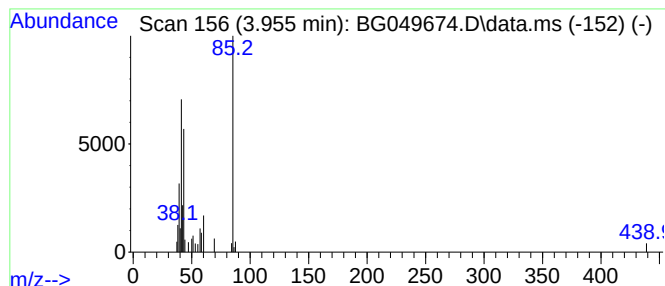
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.955	3.69 ng/ul	31004	1,4-Dichlorobenzene-d4	8.043

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	45
2			Methacrylamide	85	C4H7NO	000079-39-0	40
3			4-(Piperazin-1-ylsulfonyl)morpho...	235	C8H17N3O3S	005625-93-4	40
4			Methacrylamide	85	C4H7NO	000079-39-0	40
5			N,N'-Diacetylenediamine	144	C6H12N2O2	000871-78-3	39



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049674.D
Acq On : 6 Aug 2021 12:03
Operator : CG/JU
Sample : M3124-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0094

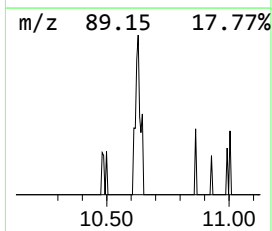
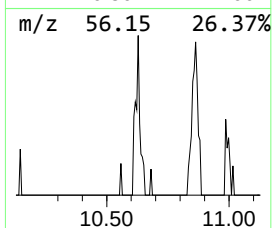
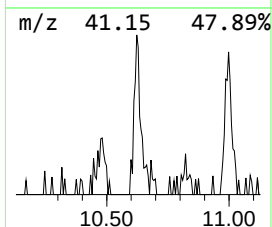
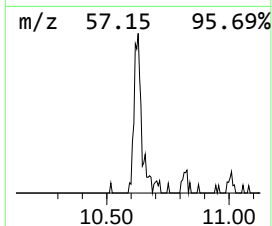
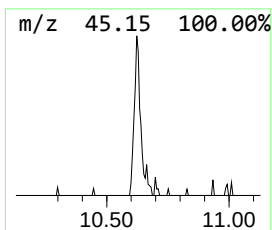
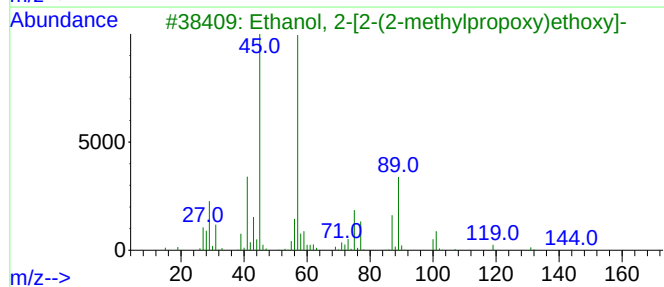
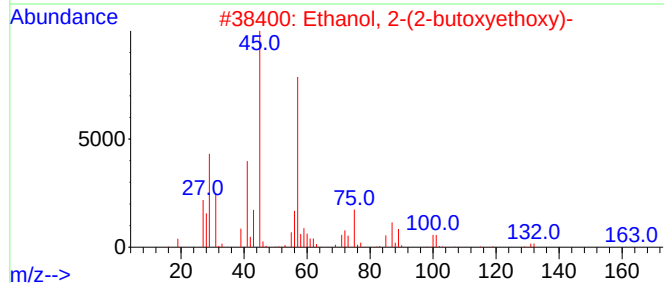
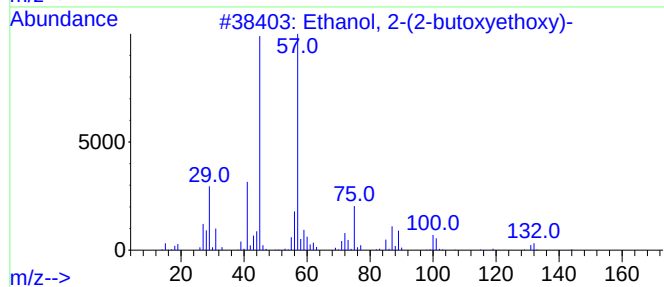
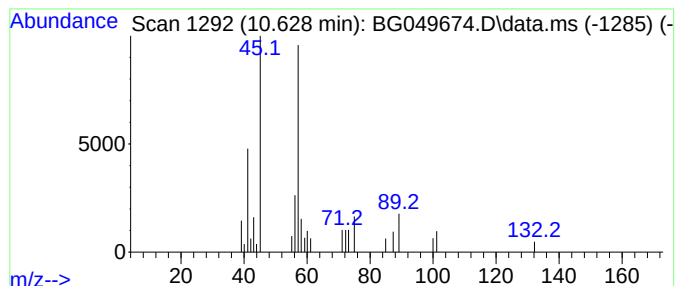
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.628	2.29 ng/ul	28280	Naphthalene-d8	10.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	59
3			Ethanol, 2-[2-(2-methylpropoxy)ethoxy]-	162	C8H18O3	018912-80-6	59
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	59
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049674.D
Acq On : 6 Aug 2021 12:03
Operator : CG/JU
Sample : M3124-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0094

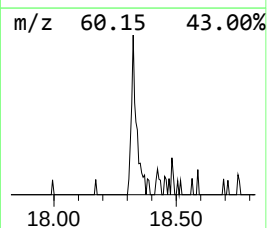
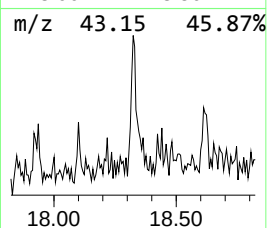
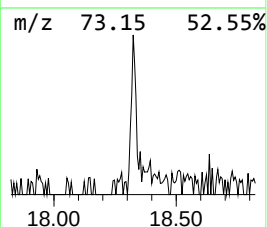
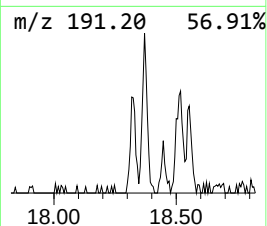
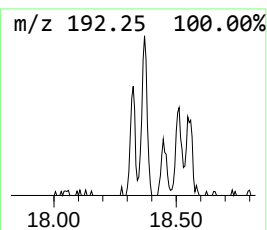
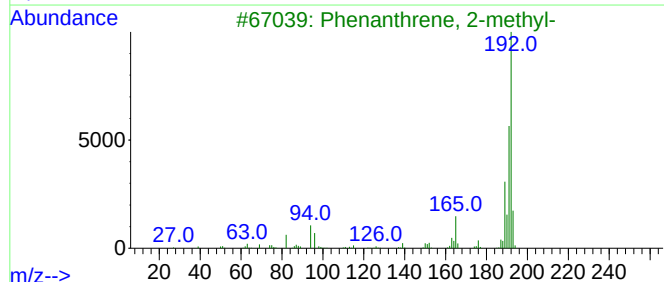
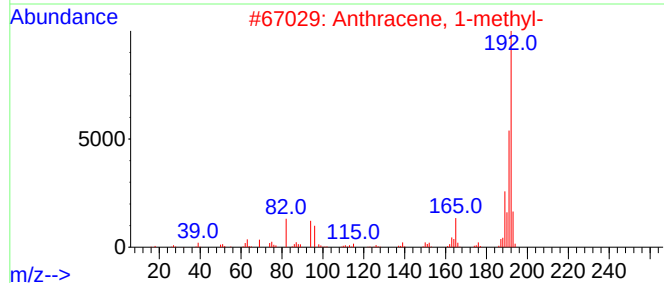
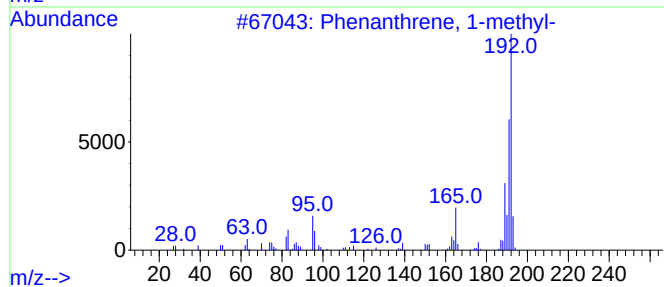
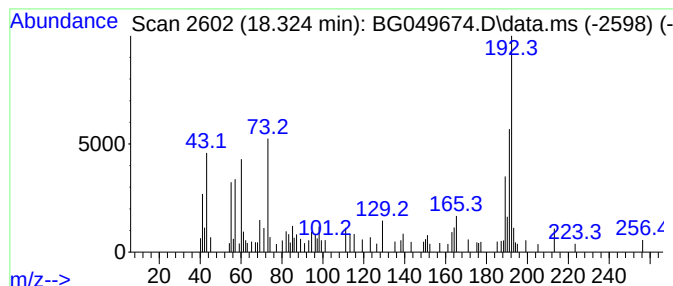
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.324	2.54 ng/ul	49615	Phenanthrene-d10	17.442

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	64
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	64
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049674.D
Acq On : 6 Aug 2021 12:03
Operator : CG/JU
Sample : M3124-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0094

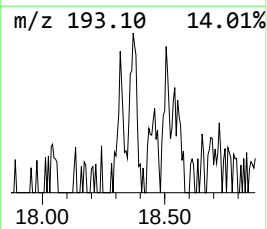
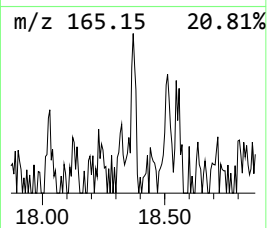
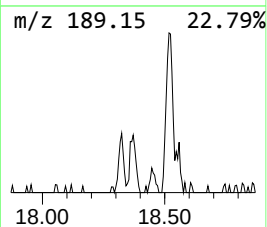
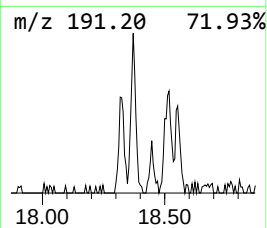
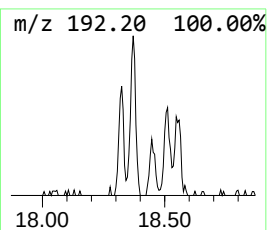
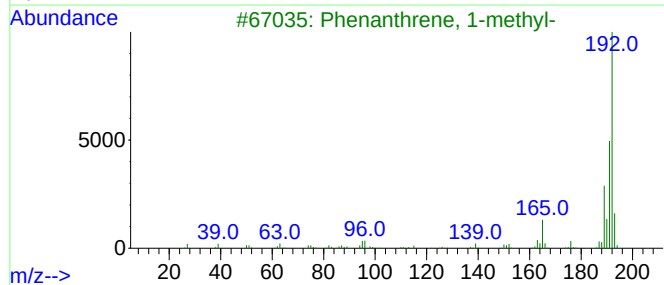
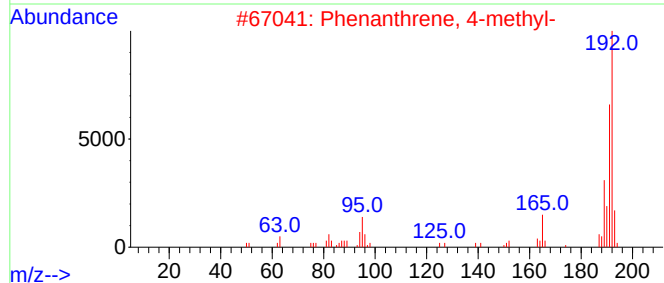
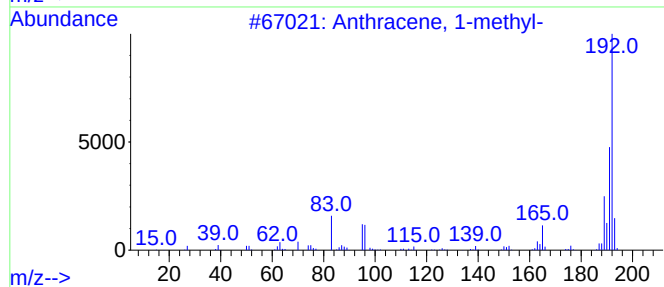
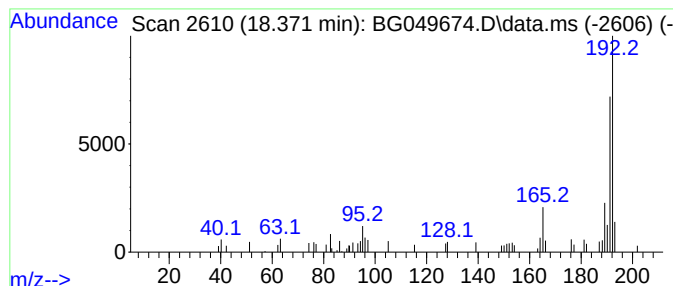
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.371	2.05 ng/ul	39959	Phenanthrene-d10	17.442

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
2			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049674.D
Acq On : 6 Aug 2021 12:03
Operator : CG/JU
Sample : M3124-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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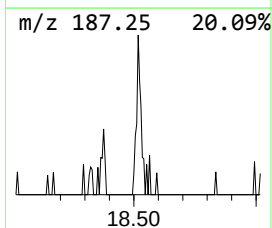
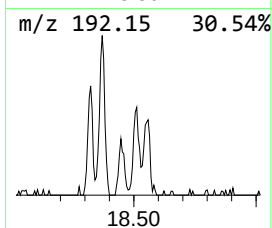
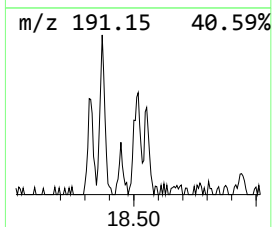
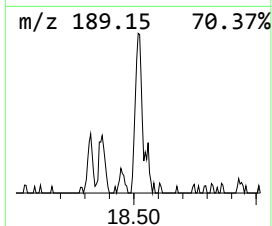
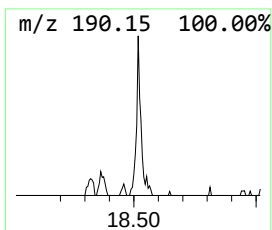
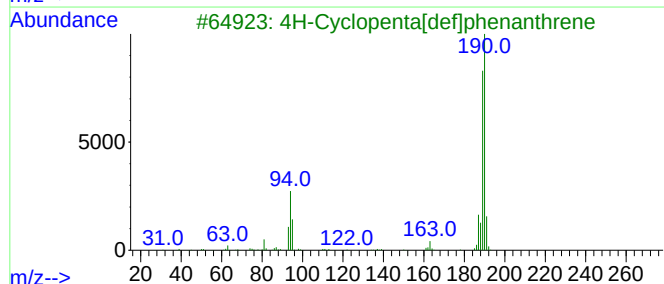
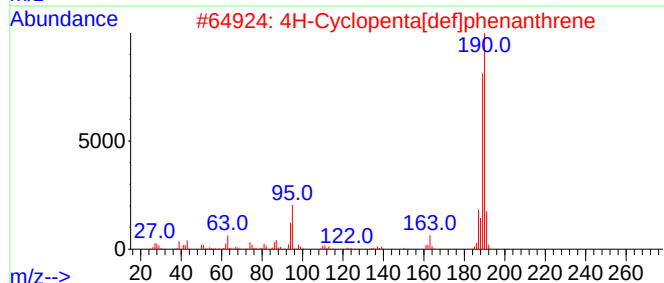
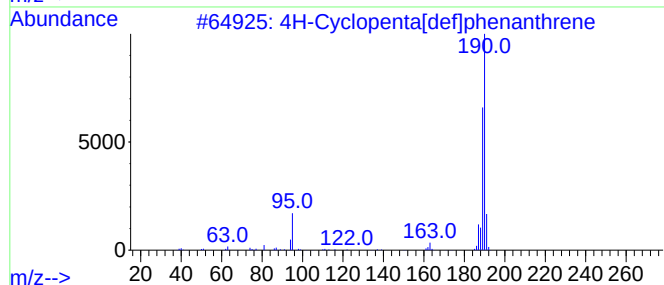
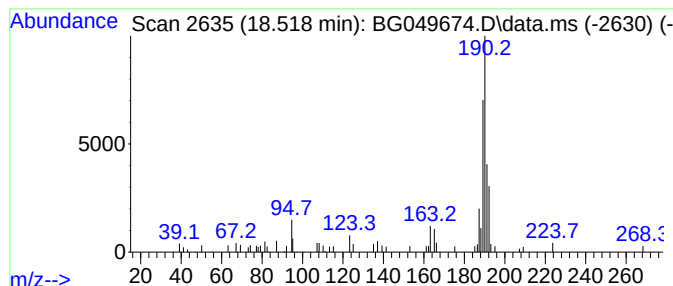
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.517	2.12 ng/ul	41384	Phenanthrene-d10	17.442

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	58
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
4			Pyrimidine, 6-oxo-4-(3-fluorophe...	190	C10H7FN2O	085979-55-1	43
5			Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049674.D
Acq On : 6 Aug 2021 12:03
Operator : CG/JU
Sample : M3124-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
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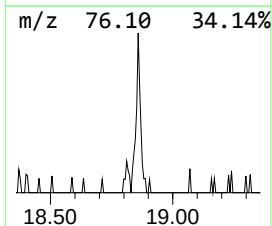
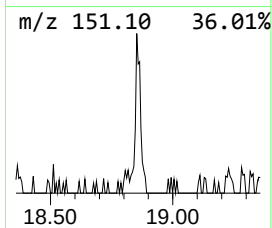
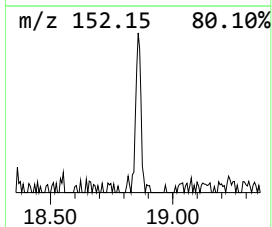
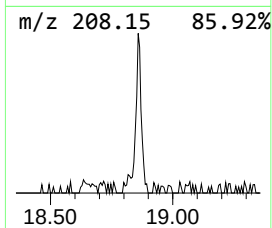
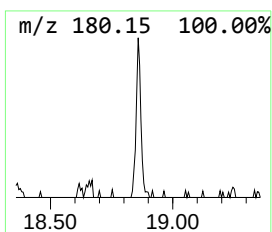
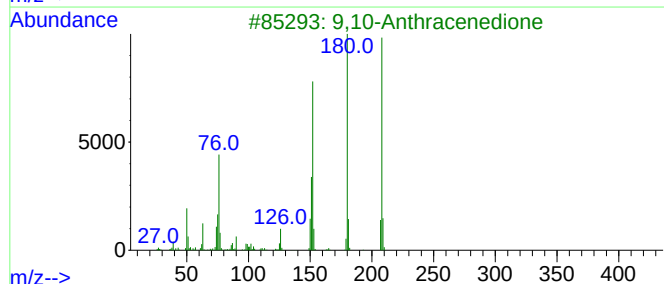
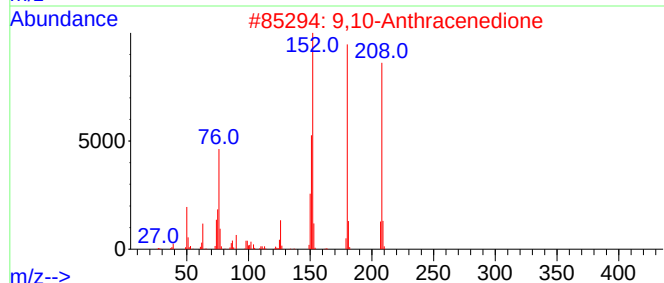
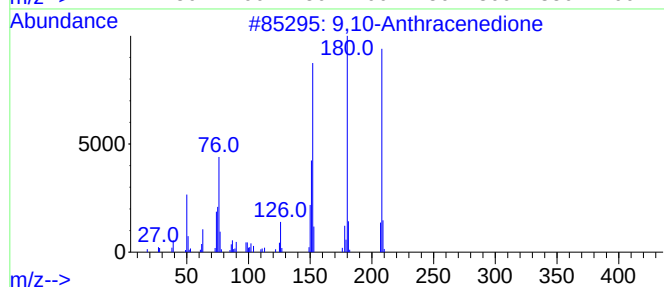
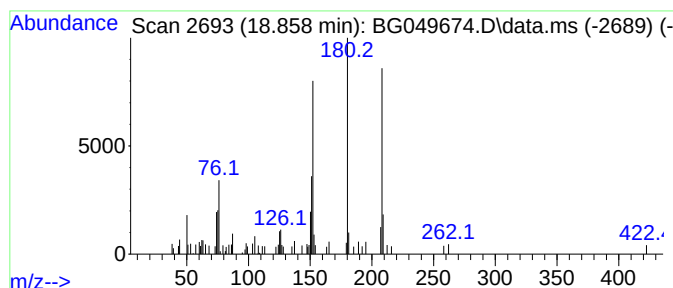
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.858	2.11 ng/ul	41216	Phenanthrene-d10	17.442

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	98
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
4		9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
5		9H-Fluoren-9-one	180	C13H8O	000486-25-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049674.D
Acq On : 6 Aug 2021 12:03
Operator : CG/JU
Sample : M3124-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.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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unknown-01	3.150	38.4	ng/ul	322676	1	8.043	168262	20.0
unknown-02	3.955	3.7	ng/ul	31004	1	8.043	168262	20.0
Ethanol, 2-(2-b...	10.628	2.3	ng/ul	28280	2	10.863	246825	20.0
Phenanthrene, 1...	18.324	2.5	ng/ul	49615	4	17.442	390257	20.0
Anthracene, 1-m...	18.371	2.0	ng/ul	39959	4	17.442	390257	20.0
4H-Cyclopenta[d...	18.517	2.1	ng/ul	41384	4	17.442	390257	20.0
9,10-Anthracene...	18.858	2.1	ng/ul	41216	4	17.442	390257	20.0