

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
 Data File : BG062162.D
 Acq On : 22 Jul 2024 15:09
 Operator : MA/JU
 Sample : P3263-24
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4DK8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BG071924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062162.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.265	662	677	686	rBV2	677480	1322855	19.15%	1.180%
2	7.389	691	698	704	rVV	75356	122577	1.77%	0.109%
3	7.483	708	714	722	rVB	146456	227577	3.29%	0.203%
4	7.629	727	739	751	rBV3	366312	798698	11.56%	0.712%
5	7.953	787	794	802	rBV	214653	354592	5.13%	0.316%
6	8.099	814	819	831	rVB	825297	1440641	20.85%	1.285%
7	8.417	866	873	881	rBV	571677	998328	14.45%	0.890%
8	8.787	928	936	942	rBV	144624	247560	3.58%	0.221%
9	8.869	942	950	957	rBV	588017	976401	14.13%	0.871%
10	9.145	990	997	1006	rVV	541454	939323	13.60%	0.838%
11	9.239	1006	1013	1016	rVV2	106623	212174	3.07%	0.189%
12	9.280	1016	1020	1030	rVB	194147	335003	4.85%	0.299%
13	9.779	1099	1105	1109	rBV4	7044	10887	0.16%	0.010%
14	9.991	1128	1141	1151	rVB2	253763	622200	9.01%	0.555%
15	10.279	1182	1190	1198	rVB	316275	519950	7.53%	0.464%
16	10.531	1220	1233	1243	rBV4	363649	882599	12.77%	0.787%
17	10.878	1285	1292	1296	rBV	189589	364956	5.28%	0.325%
18	10.931	1296	1301	1309	rVV	1201777	2120617	30.69%	1.891%
19	11.019	1309	1316	1325	rVB	105603	197567	2.86%	0.176%
20	11.612	1408	1417	1425	rBV	38323	72871	1.05%	0.065%
21	12.164	1504	1511	1524	rBV	875309	1434592	20.76%	1.279%
22	13.134	1669	1676	1683	rBV	640087	1025078	14.84%	0.914%
23	13.216	1683	1690	1698	rVV	1479425	2355680	34.10%	2.100%
24	13.281	1698	1701	1709	rVB8	7463	17288	0.25%	0.015%
25	13.492	1732	1737	1741	rVV5	6607	11099	0.16%	0.010%
26	13.574	1741	1751	1764	rVB	767829	1269736	18.38%	1.132%
27	13.751	1776	1781	1785	rVB4	12134	19597	0.28%	0.017%
28	14.097	1830	1840	1847	rBV	412581	621132	8.99%	0.554%
29	14.285	1863	1872	1878	rBV	2045140	3226842	46.70%	2.877%
30	14.426	1884	1896	1904	rVB2	2224191	4073694	58.96%	3.632%
31	14.696	1935	1942	1947	rBV	356820	562659	8.14%	0.502%
32	14.761	1947	1953	1962	rVB	1845106	2957691	42.81%	2.637%
33	14.878	1967	1973	1974	rBV	156199	196398	2.84%	0.175%
34	14.943	1978	1984	1991	rVV	1172615	2139266	30.96%	1.907%
35	15.008	1991	1995	1998	rVV5	43525	78453	1.14%	0.070%
36	15.066	1998	2005	2018	rVB	1213565	1879151	27.20%	1.676%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
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37	15.331	2043	2050	2055	rBV5	21905	38136	0.55%	0.034%
38	15.454	2067	2071	2073	rBV4	7271	11217	0.16%	0.010%
39	15.507	2073	2080	2086	rVV	3120898	4372106	63.28%	3.898%
40	15.554	2086	2088	2092	rVB5	11065	14171	0.21%	0.013%

41	15.689	2102	2111	2114	rBV	571196	932145	13.49%	0.831%
42	15.736	2114	2119	2126	rVV2	3053323	4941857	71.53%	4.406%
43	15.807	2126	2131	2139	rVV	195133	308939	4.47%	0.275%
44	15.989	2157	2162	2169	rVV4	15729	28378	0.41%	0.025%
45	16.224	2196	2202	2205	rVV6	14154	20964	0.30%	0.019%

46	16.282	2206	2212	2217	rVV2	22506	41397	0.60%	0.037%
47	16.335	2217	2221	2225	rVB5	7914	12683	0.18%	0.011%
48	16.376	2225	2228	2230	rBV3	10880	11745	0.17%	0.010%
49	16.412	2230	2234	2242	rVB5	21477	38929	0.56%	0.035%
50	16.564	2256	2260	2265	rVB2	29549	39747	0.58%	0.035%

51	16.747	2283	2291	2298	rBV	4302118	6210169	89.88%	5.537%
52	16.946	2321	2325	2332	rVB7	10680	21746	0.31%	0.019%
53	17.093	2340	2350	2361	rBV	1938025	3093442	44.77%	2.758%
54	17.216	2368	2371	2372	rBV2	13141	14192	0.21%	0.013%
55	17.352	2390	2394	2396	rVB4	8425	11227	0.16%	0.010%

56	17.440	2403	2409	2412	rBV	532443	764752	11.07%	0.682%
57	17.487	2412	2417	2422	rVV	1294195	1969040	28.50%	1.756%
58	17.540	2422	2426	2429	rVV	691325	1148602	16.62%	1.024%
59	17.575	2429	2432	2439	rVB	1771065	2536920	36.72%	2.262%
60	17.698	2448	2453	2456	rBV6	12131	20922	0.30%	0.019%

61	17.822	2470	2474	2476	rBV4	13344	17083	0.25%	0.015%
62	18.033	2505	2510	2514	rVV3	49229	88018	1.27%	0.078%
63	18.080	2514	2518	2527	rVB	128538	149248	2.16%	0.133%
64	18.291	2551	2554	2557	rVB3	9949	13192	0.19%	0.012%
65	18.391	2564	2571	2576	rBV	3686311	4644690	67.23%	4.141%

66	18.497	2586	2589	2598	rBV	17738	38580	0.56%	0.034%
67	18.585	2600	2604	2605	rBV4	9425	12017	0.17%	0.011%
68	18.679	2616	2620	2626	rVV7	18484	43388	0.63%	0.039%
69	18.732	2626	2629	2633	rVV5	19398	30242	0.44%	0.027%
70	18.791	2635	2639	2644	rVV	28565	49367	0.71%	0.044%

71	18.855	2644	2650	2657	rVB	223866	335693	4.86%	0.299%
72	19.161	2697	2702	2705	rBV7	9401	19062	0.28%	0.017%
73	19.214	2708	2711	2714	rVV5	20260	28786	0.42%	0.026%
74	19.243	2714	2716	2720	rVB2	64148	72214	1.05%	0.064%
75	19.378	2735	2739	2745	rBV2	46749	62322	0.90%	0.056%

76	19.496	2751	2759	2765	rBV	4486486	6746527	97.65%	6.015%
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77	19.731	2795	2799	2803	rVB	46767	67283	0.97%	0.060%
78	19.825	2807	2815	2816	rBV	985240	1343464	19.44%	1.198%
79	19.854	2816	2820	2826	rVB	2292385	3285012	47.55%	2.929%
80	19.942	2834	2835	2839	rVB4	17229	15445	0.22%	0.014%
81	20.089	2854	2860	2864	rBV9	10233	25930	0.38%	0.023%
82	20.248	2883	2887	2891	rBV6	16712	24575	0.36%	0.022%
83	20.324	2896	2900	2904	rVB7	17370	29726	0.43%	0.027%
84	20.465	2920	2924	2927	rBV2	50160	64219	0.93%	0.057%
85	20.823	2981	2985	2988	rBV6	19472	24880	0.36%	0.022%
86	21.323	3066	3070	3081	rBV5	77434	174495	2.53%	0.156%
87	21.481	3093	3097	3101	rVB3	72508	98830	1.43%	0.088%
88	21.575	3106	3113	3119	rBV	4574651	6412468	92.81%	5.718%
89	21.693	3124	3133	3139	rBV2	2987790	6128174	88.70%	5.464%
90	21.763	3139	3145	3152	rVB	4113444	6906771	99.97%	6.158%
91	22.468	3263	3265	3270	rBV4	25707	29974	0.43%	0.027%
92	22.674	3296	3300	3307	rVB2	65814	116656	1.69%	0.104%
93	23.990	3511	3524	3535	rBV	2674984	6909130	100.00%	6.160%
94	24.389	3590	3592	3594	rBV3	11723	10881	0.16%	0.010%
95	24.712	3638	3647	3659	rVB	442387	1212112	17.54%	1.081%
96	24.935	3674	3685	3696	rBV	308688	923509	13.37%	0.823%
97	25.940	3854	3856	3859	rBV4	10957	14399	0.21%	0.013%
98	26.034	3866	3872	3883	rBV4	49602	144710	2.09%	0.129%
99	28.713	4308	4328	4346	rVB	1013946	4590068	66.43%	4.093%
100	30.957	4707	4710	4711	rBV3	11425	11326	0.16%	0.010%

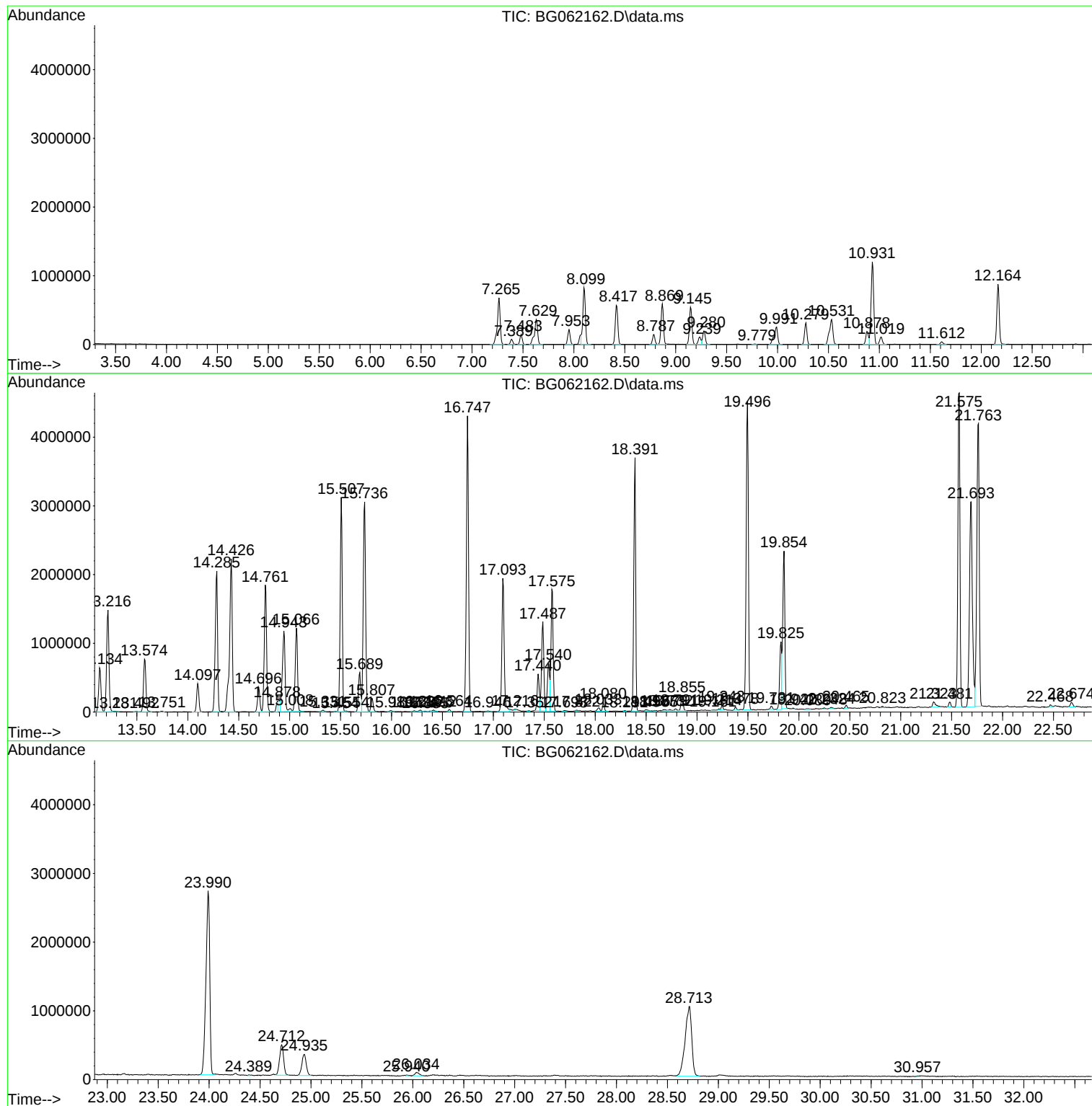
Sum of corrected areas: 112153624

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Instrument :
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A4DK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.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\BG072224\
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Instrument :
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A4DK8

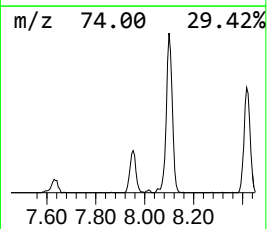
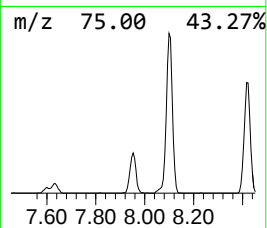
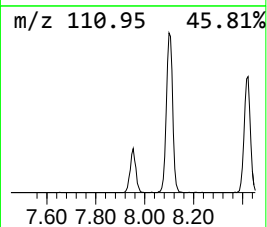
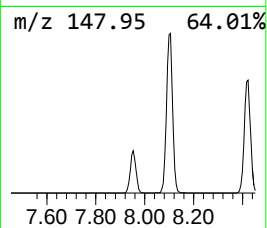
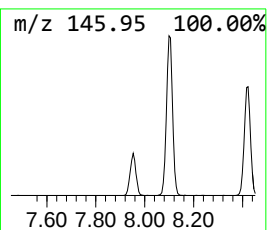
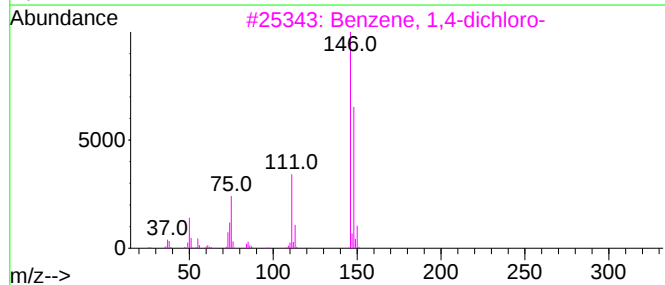
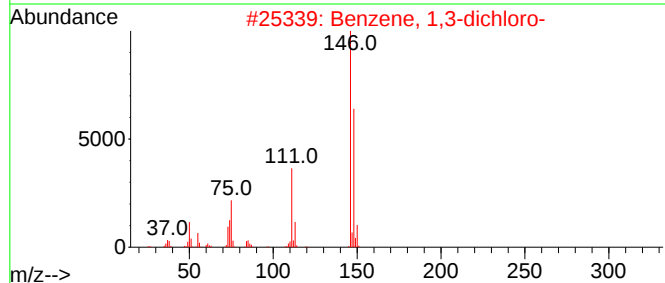
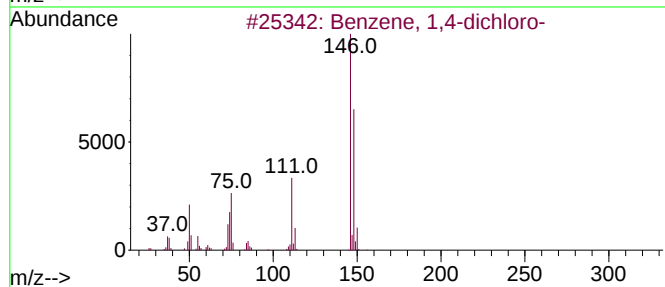
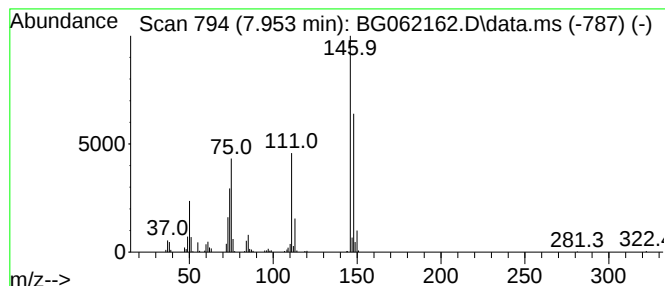
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,4-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.953	4.92 ng/ul	354592	1,4-Dichlorobenzene-d4	8.064

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93



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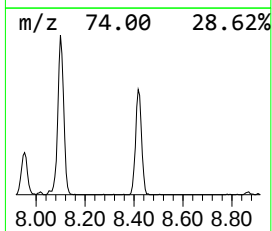
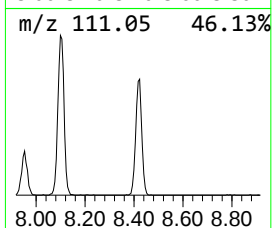
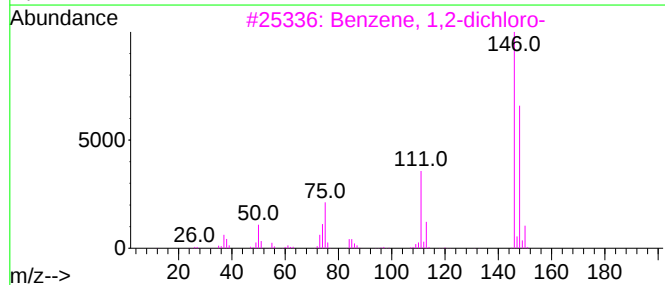
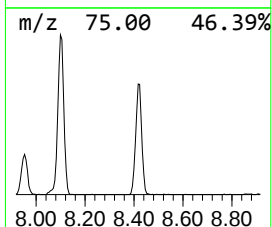
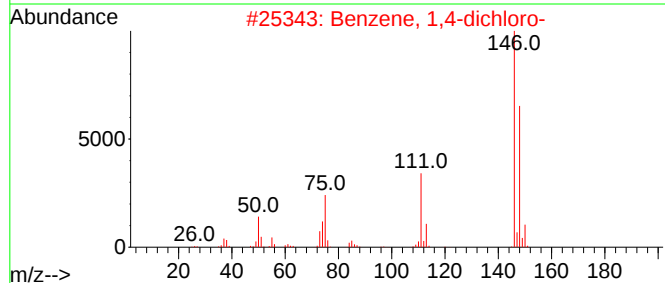
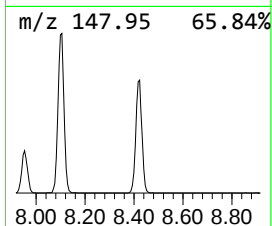
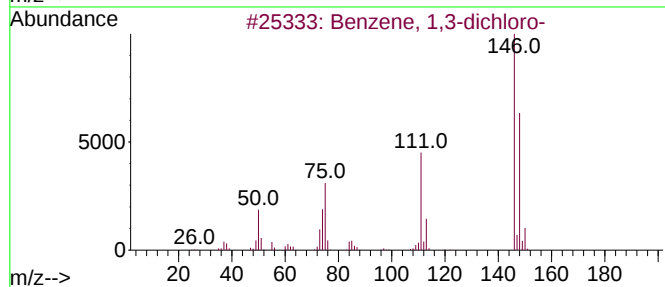
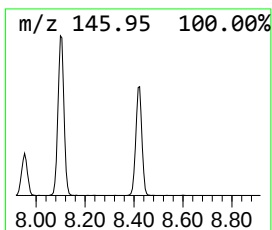
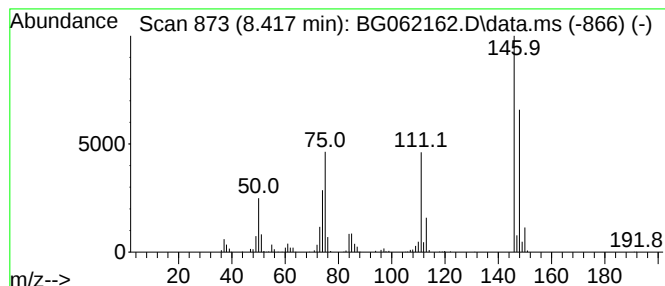
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.417	13.86 ng/ul	998328	1,4-Dichlorobenzene-d4	8.064

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96



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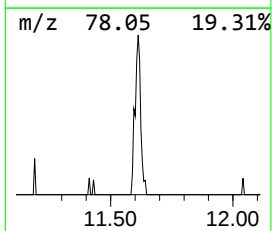
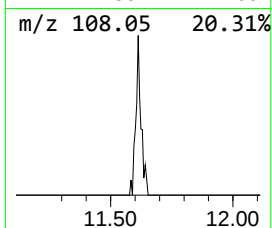
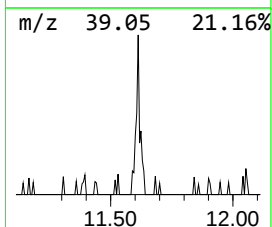
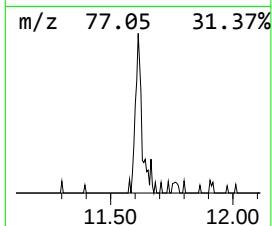
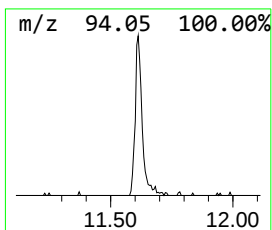
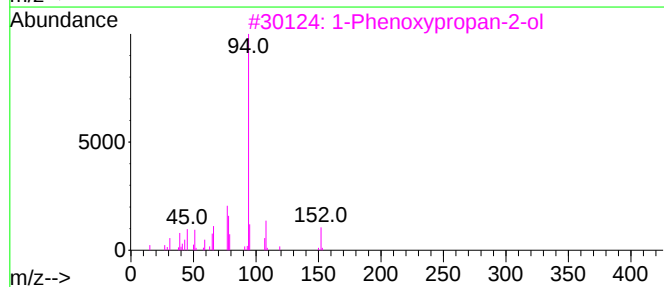
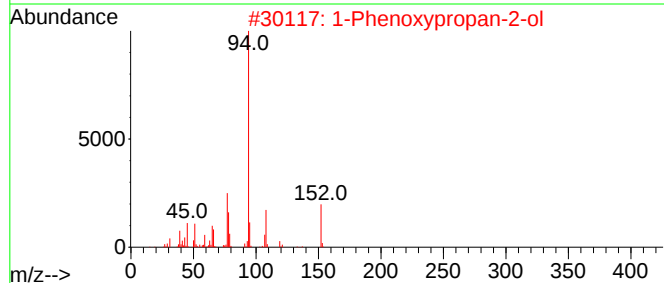
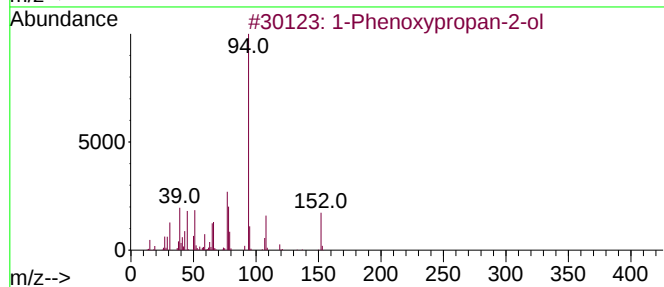
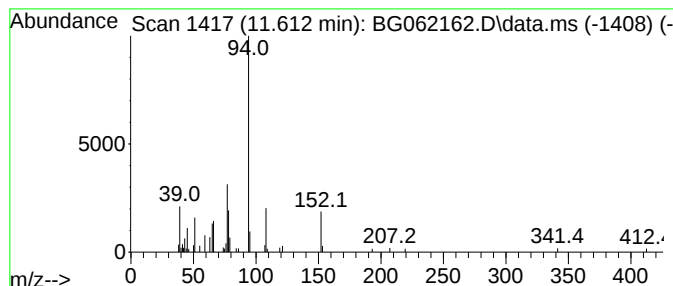
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.612	3.99 ng/ul	72871	Naphthalene-d8	10.878

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	83
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	80
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	72
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062162.D
Acq On : 22 Jul 2024 15:09
Operator : MA/JU
Sample : P3263-24
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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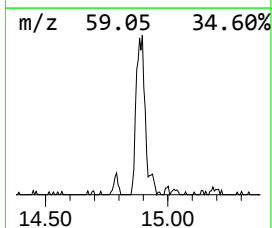
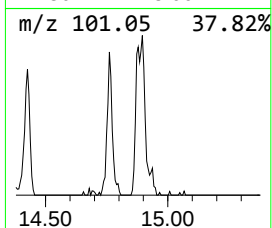
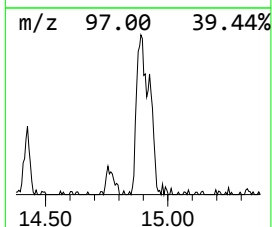
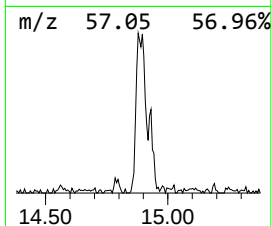
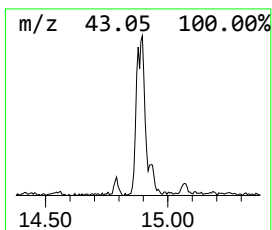
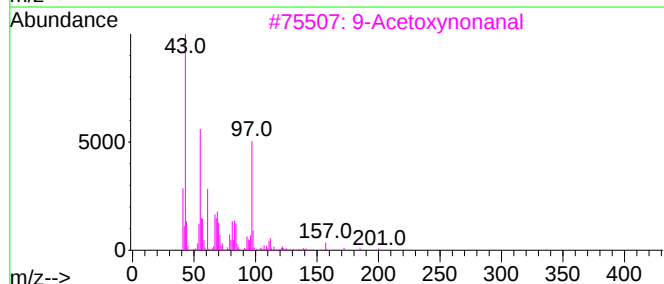
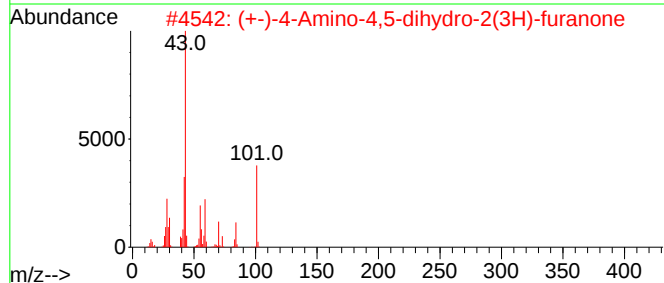
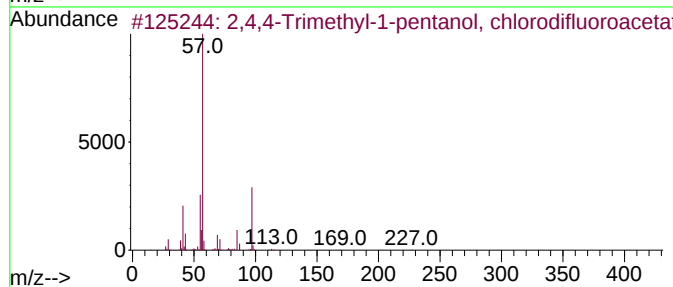
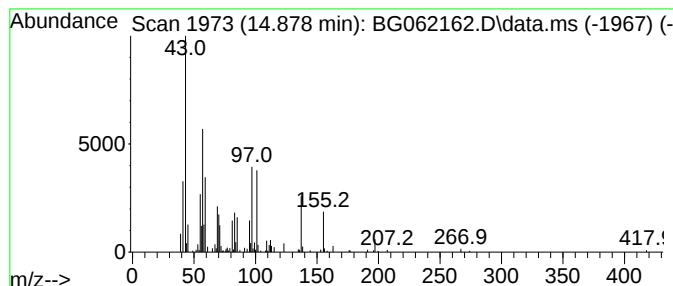
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.878	6.98 ng/ul	196398	Acenaphthene-d10	14.696

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4-Trimethyl-1-pentanol, chlo...	242	C10H17ClF2O2	1000447-27-2	22		
2	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	18		
3	9-Acetoxy-nonanal	200	C11H20O3	029541-97-7	14		
4	N-Hexadecylpyridinium bromide	383	C21H38BrN	000140-72-7	10		
5	Acetic acid, 3-hydroxy-2-nitroph...	197	C8H7NO5	1000269-41-1	10		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062162.D
Acq On : 22 Jul 2024 15:09
Operator : MA/JU
Sample : P3263-24
Misc :
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Instrument :
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ClientSampleId :
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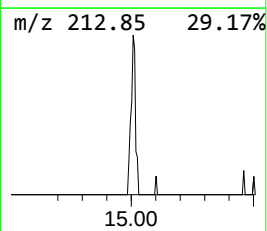
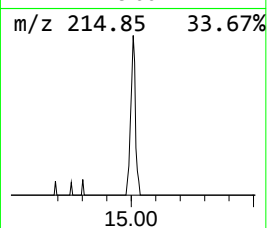
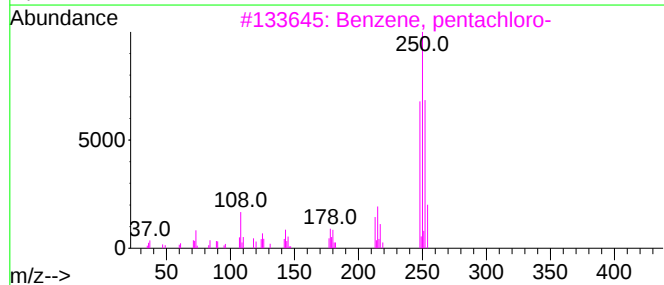
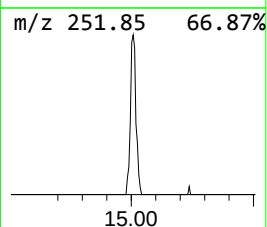
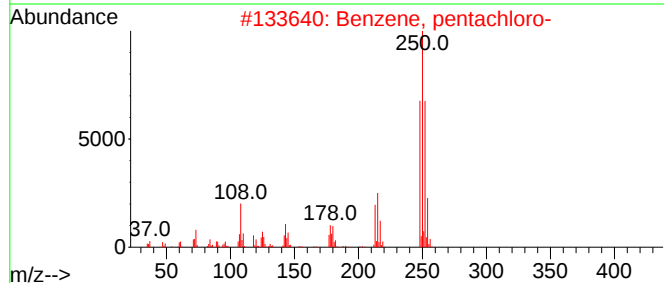
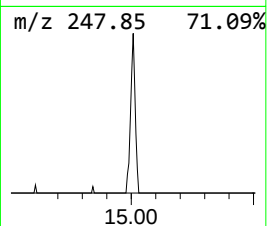
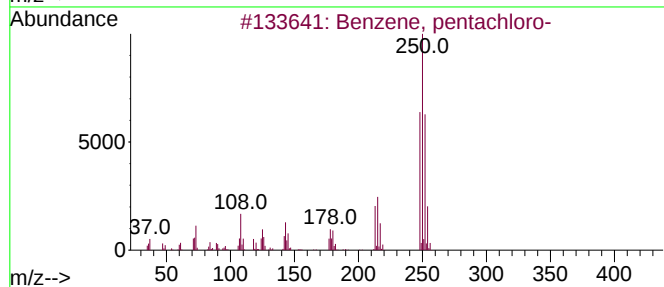
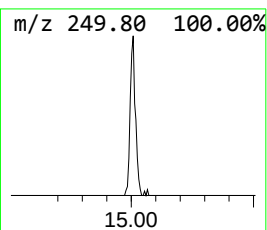
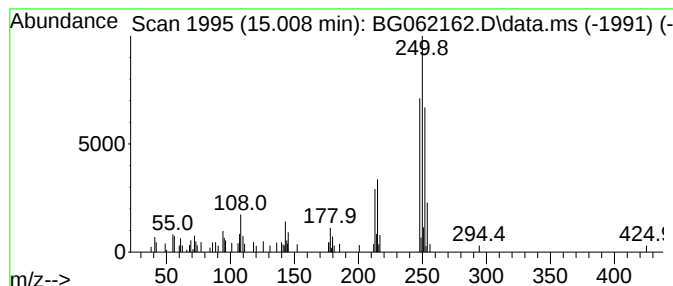
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, pentachloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.008	2.79 ng/ul	78453	Acenaphthene-d10	14.696

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentachloro-	248	C6HCl5	000608-93-5	95
2			Benzene, pentachloro-	248	C6HCl5	000608-93-5	94
3			Benzene, pentachloro-	248	C6HCl5	000608-93-5	94
4			Benzene, pentachloro-	248	C6HCl5	000608-93-5	93
5			Benzene, pentachloro-	248	C6HCl5	000608-93-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062162.D
Acq On : 22 Jul 2024 15:09
Operator : MA/JU
Sample : P3263-24
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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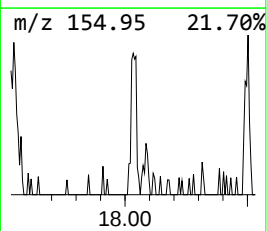
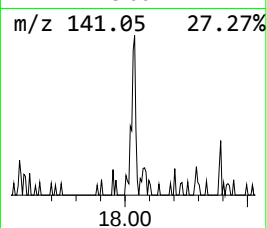
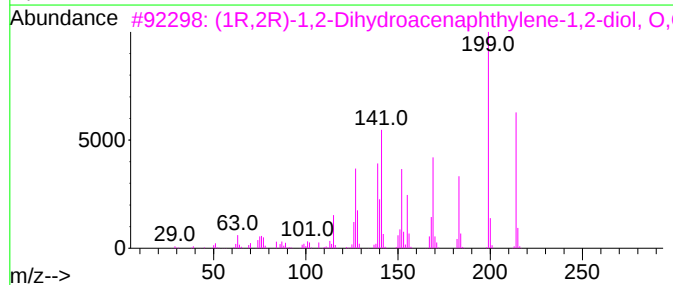
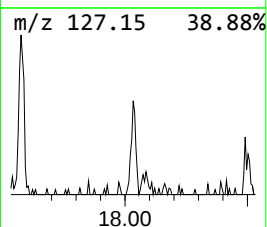
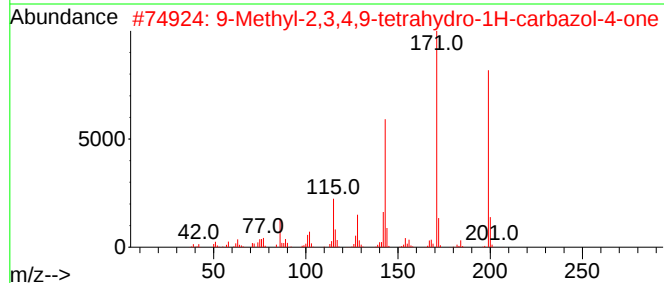
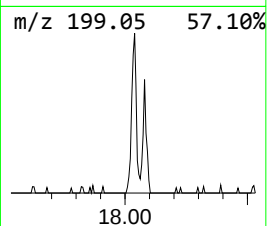
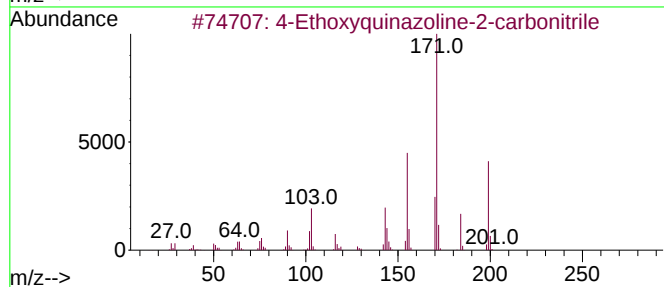
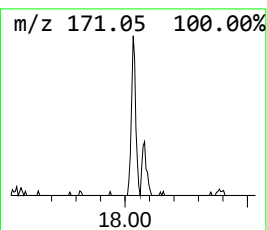
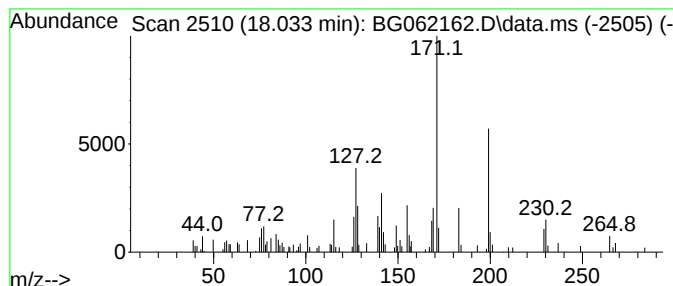
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.033	2.30 ng/ul	88018	Phenanthrene-d10	17.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Ethoxyquinazoline-2-carbonitrile	199	C11H9N3O	099973-07-6	43
2			9-Methyl-2,3,4,9-tetrahydro-1H-c...	199	C13H13NO	027387-31-1	38
3			(1R,2R)-1,2-Dihydroacenaphthylene...	214	C14H14O2	1000496-44-2	35
4			1H-Pyrido[3,4-b]indole, 1-ethyl-...	200	C13H16N2	006678-86-0	30
5			2,5-Difluorophenyl isothiocyanate	171	C7H3F2NS	206559-57-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062162.D
Acq On : 22 Jul 2024 15:09
Operator : MA/JU
Sample : P3263-24
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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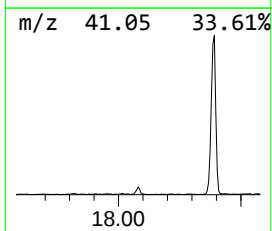
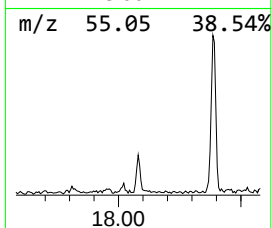
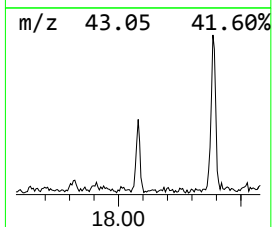
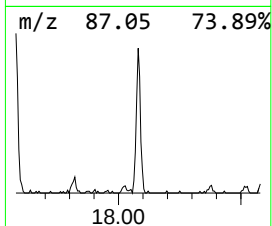
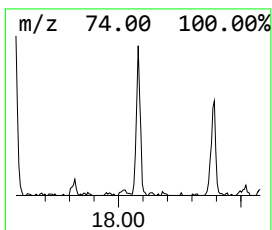
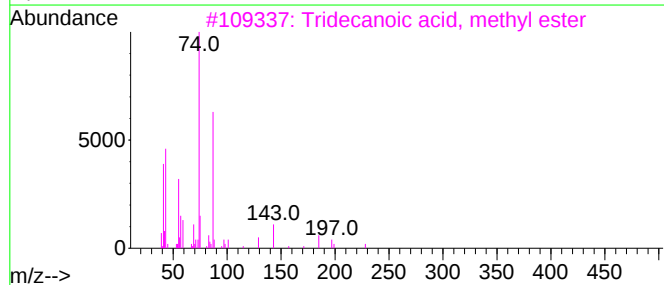
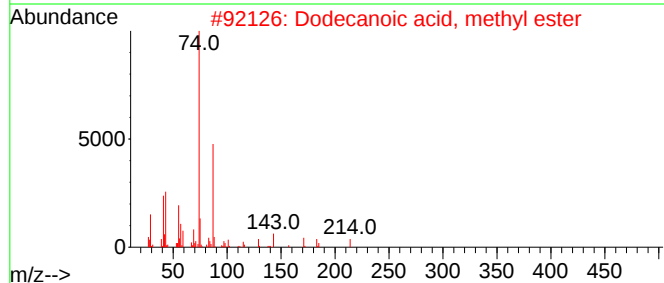
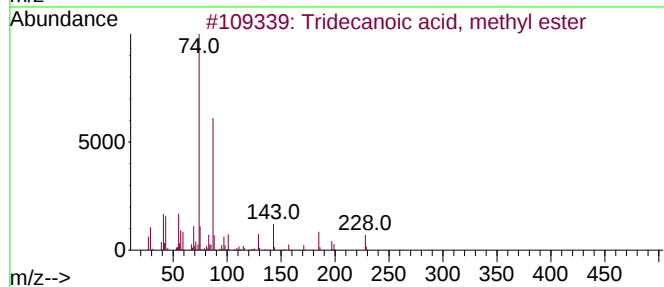
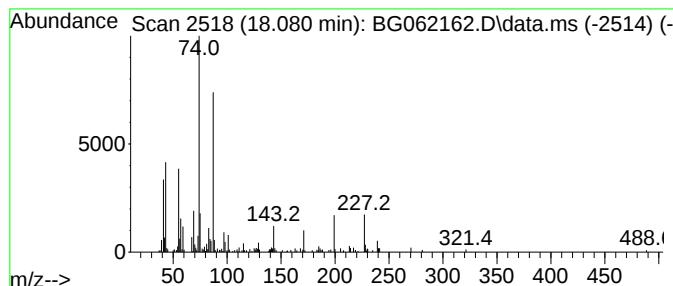
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tridecanoic acid, methyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.080	3.90 ng/ul	149248	Phenanthrene-d10	17.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	83
2			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	83
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	81
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	80
5			Heneicosanoic acid, methyl ester	340	C22H44O2	006064-90-0	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062162.D
Acq On : 22 Jul 2024 15:09
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Sample : P3263-24
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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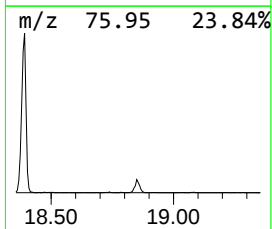
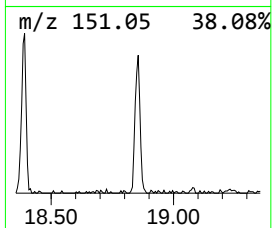
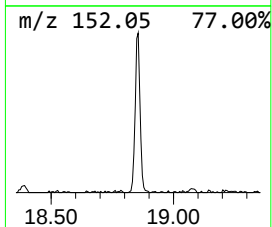
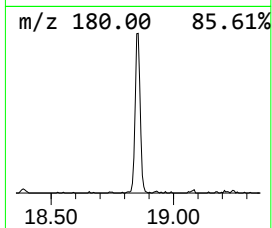
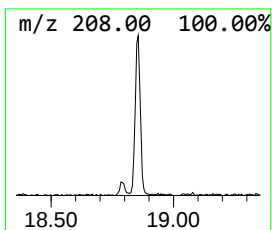
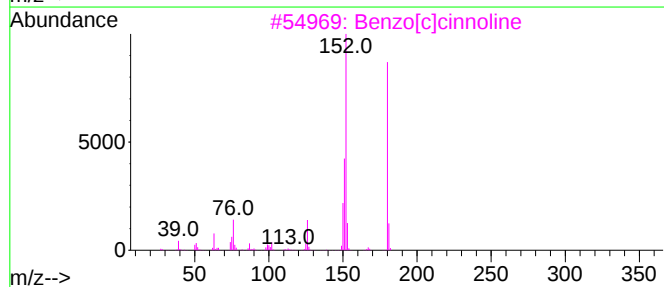
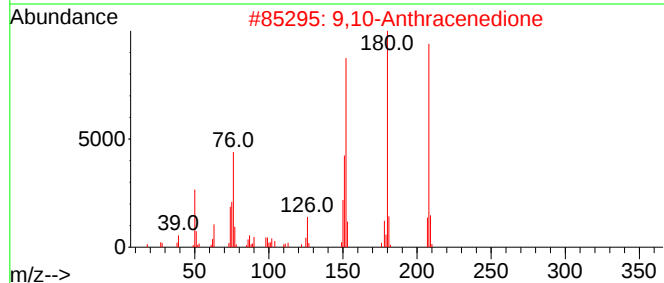
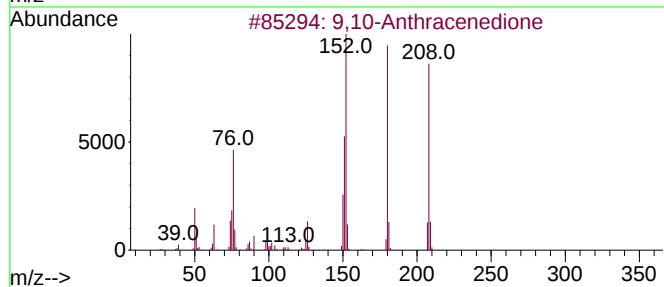
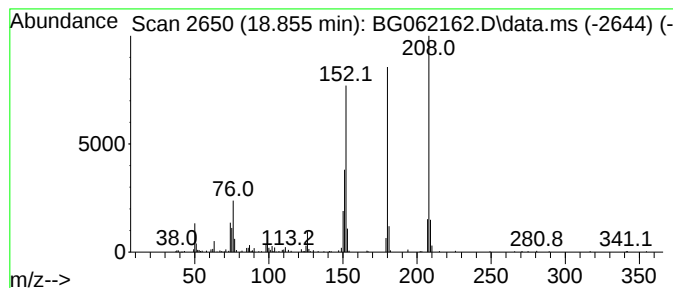
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.855	8.78 ng/ul	335693	Phenanthrene-d10	17.440

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	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	96
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
5	9,10-Anthracenedione			208	C14H8O2	000084-65-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062162.D
Acq On : 22 Jul 2024 15:09
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Sample : P3263-24
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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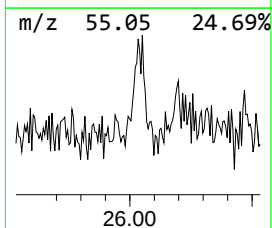
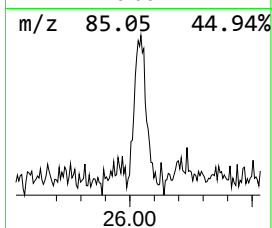
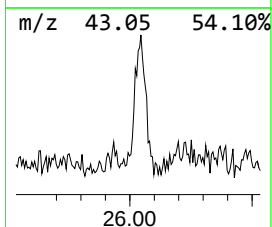
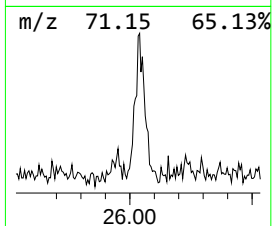
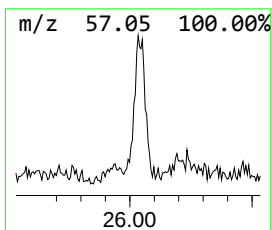
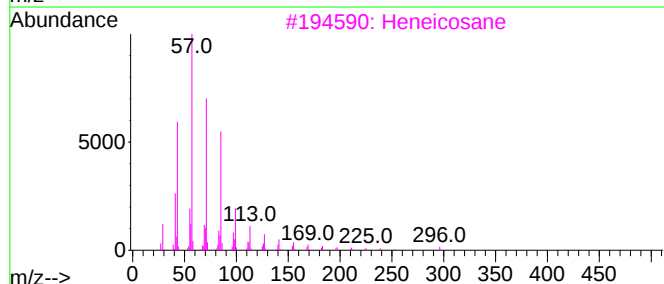
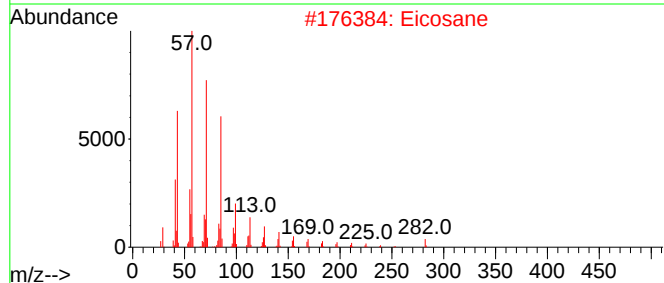
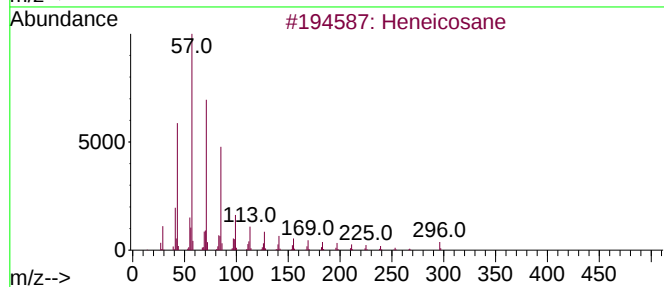
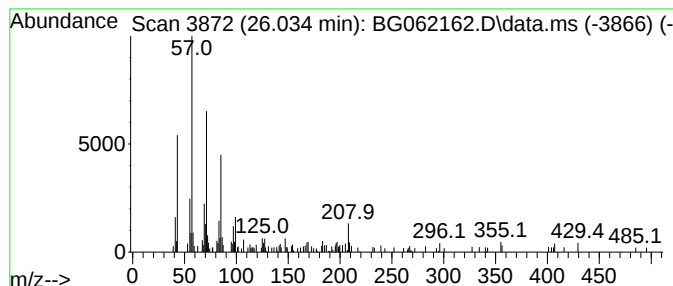
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.034	3.13 ng/ul	144710	Perylene-d12	24.935

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	86
2			Eicosane	282	C20H42	000112-95-8	70
3			Heneicosane	296	C21H44	000629-94-7	62
4			Heneicosane	296	C21H44	000629-94-7	58
5			Hexatriacontane	507	C36H74	000630-06-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062162.D
Acq On : 22 Jul 2024 15:09
Operator : MA/JU
Sample : P3263-24
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4DK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.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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Benzene, 1,3-di...	8.417	13.9	ng/ul	998328	1	8.064	1440640	20.0
1-Phenoxypropan...	11.612	4.0	ng/ul	72871	2	10.878	364956	20.0
unknown-01	14.878	7.0	ng/ul	196398	3	14.696	562659	20.0
Benzene, pentac...	15.008	2.8	ng/ul	78453	3	14.696	562659	20.0
unknown-02	18.033	2.3	ng/ul	88018	4	17.440	764752	20.0
Tridecanoic aci...	18.080	3.9	ng/ul	149248	4	17.440	764752	20.0
9,10-Anthracene...	18.855	8.8	ng/ul	335693	4	17.440	764752	20.0
(DEL) Alkane: S...	26.034	3.1	ng/ul	144710	6	24.935	923509	20.0