

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
 Data File : BG062160.D
 Acq On : 22 Jul 2024 13:26
 Operator : MA/JU
 Sample : P3238-23
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4DN3

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-BG071924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062160.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.401	526	530	535	rBV4	6784	12485	0.17%	0.014%
2	7.265	664	677	684	rBV2	313868	761873	10.54%	0.824%
3	7.388	692	698	705	rBV	86412	146045	2.02%	0.158%
4	7.482	707	714	722	rVB	400846	637751	8.83%	0.690%
5	7.600	726	734	736	rBV	118389	199240	2.76%	0.216%
6	7.635	736	740	748	rVB	221022	389295	5.39%	0.421%
7	7.952	786	794	802	rBV	379971	619184	8.57%	0.670%
8	8.064	807	813	815	rVV	130194	207190	2.87%	0.224%
9	8.099	815	819	831	rVB	321587	555916	7.69%	0.601%
10	8.422	865	874	883	rVB	1260377	2166404	29.98%	2.344%
11	8.786	929	936	943	rBV	172995	296209	4.10%	0.320%
12	8.875	943	951	958	rBV	990596	1702274	23.56%	1.841%
13	9.239	1006	1013	1015	rBV2	127282	224248	3.10%	0.243%
14	9.280	1015	1020	1028	rVB	612567	1094943	15.15%	1.184%
15	9.615	1071	1077	1083	rVB6	10018	17716	0.25%	0.019%
16	9.991	1128	1141	1150	rBV3	667921	1396496	19.32%	1.511%
17	10.279	1182	1190	1199	rVB	544500	885342	12.25%	0.958%
18	10.531	1220	1233	1244	rBV2	1028631	2142989	29.65%	2.318%
19	10.878	1286	1292	1299	rVB	192071	334685	4.63%	0.362%
20	11.019	1309	1316	1322	rVB2	91369	166231	2.30%	0.180%
21	11.154	1334	1339	1343	rVB4	8554	12598	0.17%	0.014%
22	11.618	1409	1418	1423	rBV2	39844	78042	1.08%	0.084%
23	12.164	1502	1511	1520	rBV	1198694	2050486	28.37%	2.218%
24	12.928	1638	1641	1647	rVB3	12392	16211	0.22%	0.018%
25	13.134	1669	1676	1683	rBV	644819	1035355	14.33%	1.120%
26	13.210	1683	1689	1700	rVB	875434	1456732	20.16%	1.576%
27	13.609	1753	1757	1764	rBV2	10766	18684	0.26%	0.020%
28	13.962	1810	1817	1822	rBV5	6970	12245	0.17%	0.013%
29	14.097	1833	1840	1846	rVB	397555	569562	7.88%	0.616%
30	14.273	1863	1870	1876	rBV	906871	1387126	19.20%	1.501%
31	14.326	1876	1879	1884	rVV6	8065	15722	0.22%	0.017%
32	14.391	1884	1890	1899	rVB	403259	652164	9.02%	0.706%
33	14.696	1935	1942	1947	rBV2	355928	539265	7.46%	0.583%
34	14.761	1947	1953	1963	rVB	2404792	3846690	53.23%	4.161%
35	14.878	1968	1973	1974	rBV	237636	278007	3.85%	0.301%
36	14.902	1974	1977	1978	rVV2	281108	364173	5.04%	0.394%

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Smoothing : OFF

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

37	14.931	1978	1982	1990	rVB3	478113	896271	12.40%	0.970%
38	15.072	1998	2006	2018	rVB	1565669	2523989	34.93%	2.730%
39	15.325	2044	2049	2056	rVB7	17777	28250	0.39%	0.031%
40	15.501	2073	2079	2084	rVV	1022411	1378075	19.07%	1.491%
41	15.554	2084	2088	2096	rVB3	14262	24982	0.35%	0.027%
42	15.689	2104	2111	2115	rBV	546929	869699	12.03%	0.941%
43	15.748	2115	2121	2127	rVV	2009700	2996501	41.47%	3.242%
44	15.806	2127	2131	2138	rVB	170016	254962	3.53%	0.276%
45	15.912	2146	2149	2155	rVB6	6778	12748	0.18%	0.014%
46	15.988	2157	2162	2170	rBV3	31219	55408	0.77%	0.060%
47	16.223	2197	2202	2208	rBV6	16959	30196	0.42%	0.033%
48	16.411	2226	2234	2244	rVB9	19409	46202	0.64%	0.050%
49	16.564	2257	2260	2264	rVV2	38217	51670	0.72%	0.056%
50	16.623	2264	2270	2276	rVB	1230794	1715722	23.74%	1.856%
51	16.952	2320	2326	2331	rBV7	9363	18591	0.26%	0.020%
52	17.093	2342	2350	2361	rBV	1422545	2222920	30.76%	2.405%
53	17.210	2366	2370	2383	rVB	30623	82332	1.14%	0.089%
54	17.439	2403	2409	2412	rBV	470509	740809	10.25%	0.801%
55	17.486	2412	2417	2422	rVV	2313247	3442592	47.64%	3.724%
56	17.539	2422	2426	2429	rVV2	683176	1057024	14.63%	1.143%
57	17.575	2429	2432	2440	rVB	774271	1069325	14.80%	1.157%
58	17.815	2468	2473	2480	rBV8	15936	34851	0.48%	0.038%
59	17.951	2491	2496	2500	rBV8	9307	12709	0.18%	0.014%
60	18.021	2503	2508	2509	rBV4	20065	25657	0.36%	0.028%
61	18.039	2509	2511	2514	rVV4	23194	24698	0.34%	0.027%
62	18.080	2514	2518	2522	rVB	120502	146521	2.03%	0.159%
63	18.385	2564	2570	2577	rBV	3194519	4293422	59.41%	4.645%
64	18.509	2584	2591	2595	rBV8	16516	36390	0.50%	0.039%
65	18.673	2615	2619	2625	rBV8	15157	27269	0.38%	0.029%
66	18.855	2645	2650	2657	rBV	119651	188135	2.60%	0.204%
67	18.996	2671	2674	2676	rBV4	10948	15007	0.21%	0.016%
68	19.037	2679	2681	2684	rVB4	10896	10888	0.15%	0.012%
69	19.219	2708	2712	2713	rBV3	19028	20871	0.29%	0.023%
70	19.243	2713	2716	2720	rVB3	68323	81189	1.12%	0.088%
71	19.378	2732	2739	2744	rBV2	58062	89124	1.23%	0.096%
72	19.425	2744	2747	2750	rVV3	24837	34667	0.48%	0.038%
73	19.490	2751	2758	2766	rVB	2270765	3173296	43.91%	3.433%
74	19.725	2793	2798	2804	rBV	104265	161755	2.24%	0.175%
75	19.860	2816	2821	2826	rVB	4808003	7226463	100.00%	7.817%
76	20.083	2856	2859	2861	rBV4	11288	12108	0.17%	0.013%

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 Title : SVOA CALIBRATION

77	20.171	2872	2874	2877	rBV4	13244	14338	0.20%	0.016%
78	20.465	2921	2924	2928	rBV2	68160	70385	0.97%	0.076%
79	20.723	2962	2968	2972	rBV	2128696	2553886	35.34%	2.763%
80	20.817	2982	2984	2989	rVB5	19542	17051	0.24%	0.018%
81	21.328	3066	3071	3074	rBV3	88866	148111	2.05%	0.160%
82	21.481	3093	3097	3101	rVB2	101912	129244	1.79%	0.140%
83	21.569	3106	3112	3118	rBV	2934135	3884484	53.75%	4.202%
84	21.687	3125	3132	3138	rBV2	2116720	4065978	56.27%	4.399%
85	21.751	3138	3143	3151	rVB	1269788	2028974	28.08%	2.195%
86	22.039	3190	3192	3198	rVB7	27839	38333	0.53%	0.041%
87	22.474	3263	3266	3269	rBV4	29495	33574	0.46%	0.036%
88	22.679	3297	3301	3307	rVB2	77044	137726	1.91%	0.149%
89	22.785	3311	3319	3327	rVB	2672175	4527157	62.65%	4.897%
90	23.907	3501	3510	3515	rBV2	631508	1569895	21.72%	1.698%
91	23.978	3515	3522	3532	rVB	1177154	2783640	38.52%	3.011%
92	24.712	3639	3647	3653	rBV	396576	1104705	15.29%	1.195%
93	24.788	3653	3660	3669	rVB	433948	1159686	16.05%	1.255%
94	24.935	3674	3685	3695	rBV	294643	861176	11.92%	0.932%
95	25.940	3852	3856	3857	rBV4	15389	18262	0.25%	0.020%
96	26.040	3866	3873	3882	rVB4	62791	190017	2.63%	0.206%
97	26.192	3893	3899	3900	rBV6	20432	33620	0.47%	0.036%
98	26.363	3926	3928	3937	rVB10	11886	17255	0.24%	0.019%
99	26.827	4005	4007	4012	rBV6	8631	10707	0.15%	0.012%
100	28.718	4300	4329	4348	rVV2	1053437	5618805	77.75%	6.078%

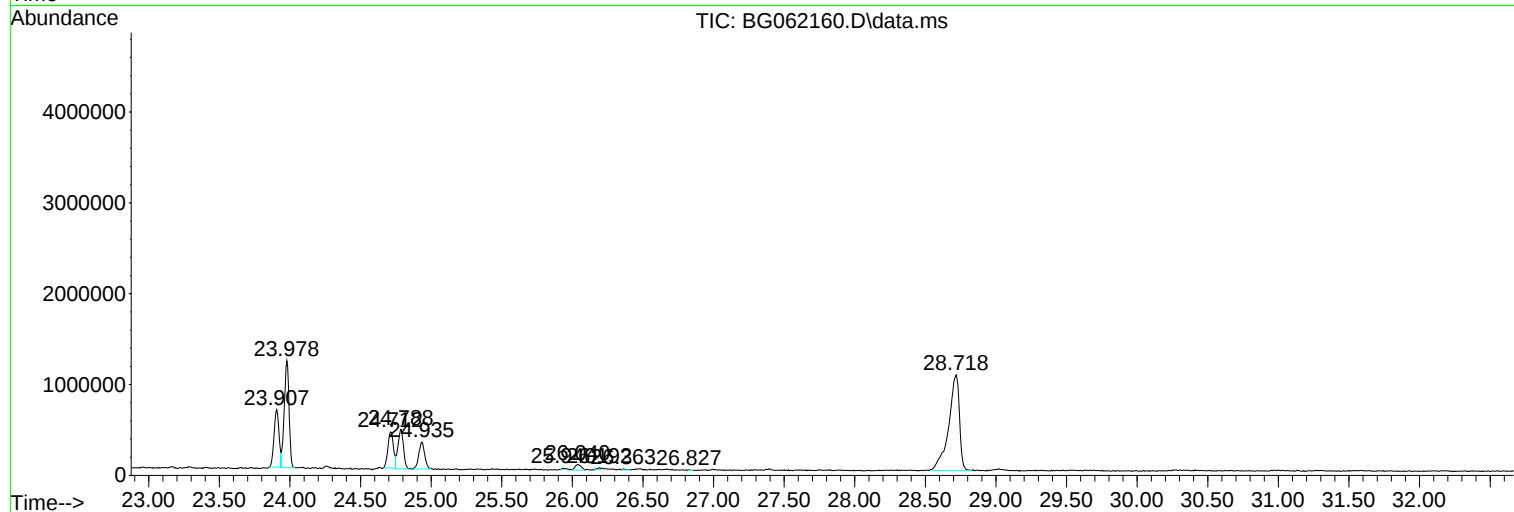
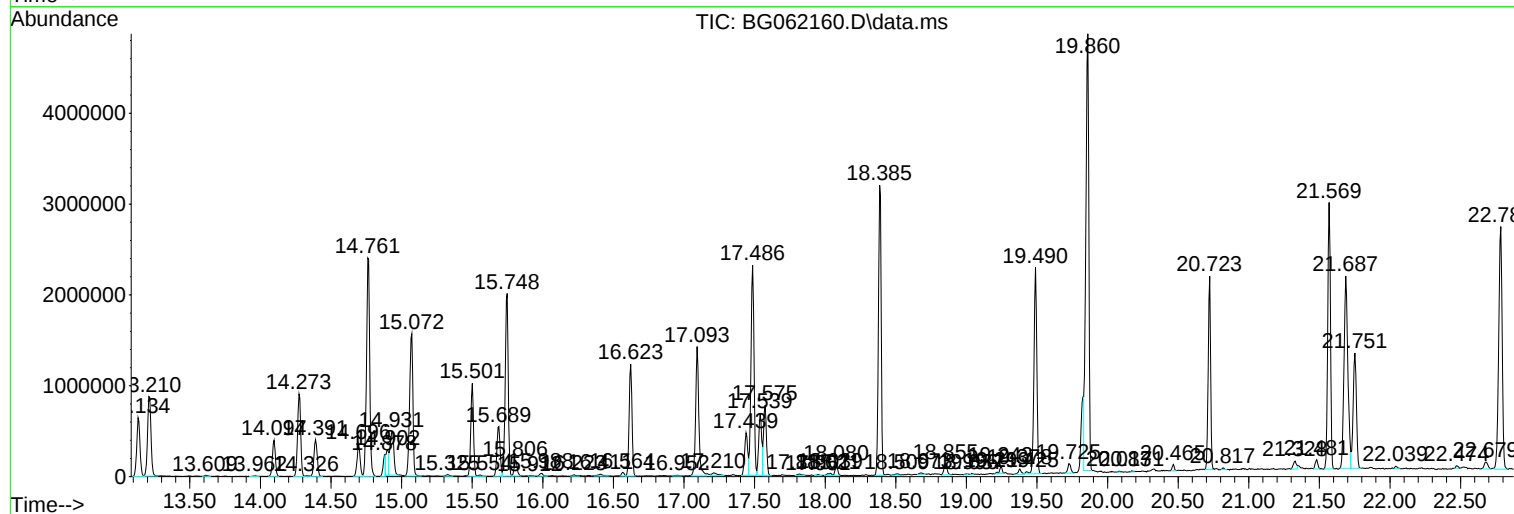
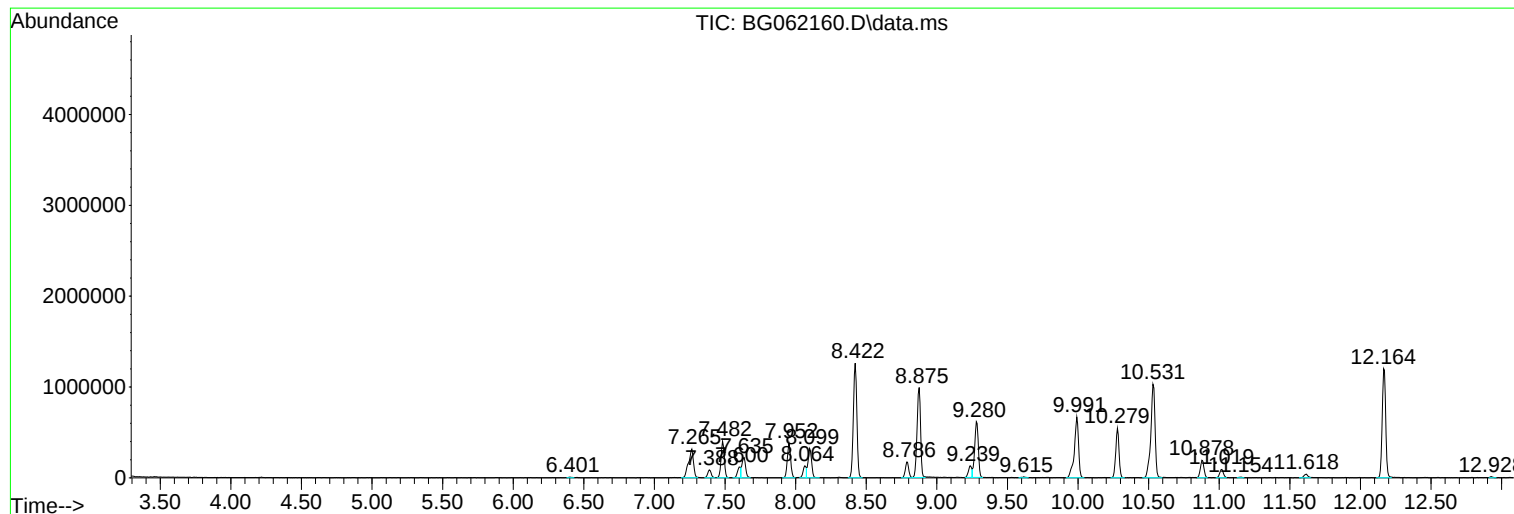
Sum of corrected areas: 92439875

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Instrument :
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ClientSampleId :
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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



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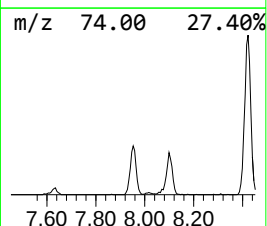
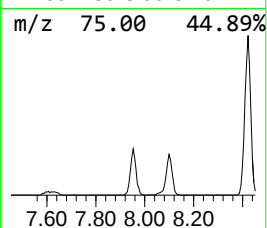
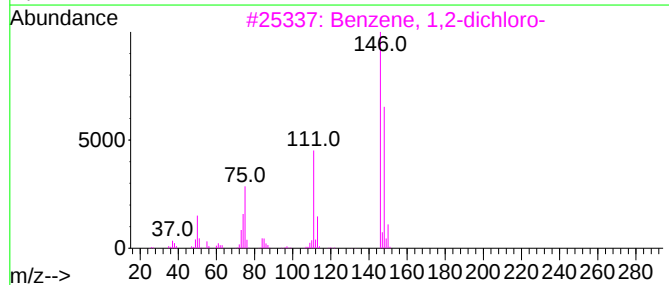
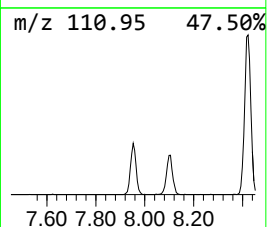
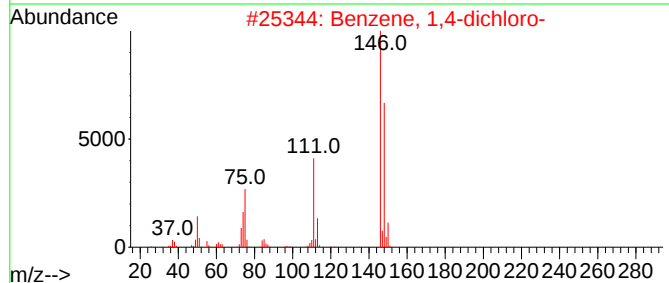
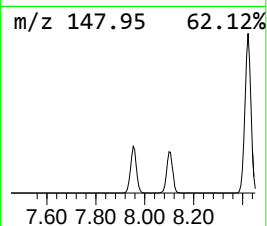
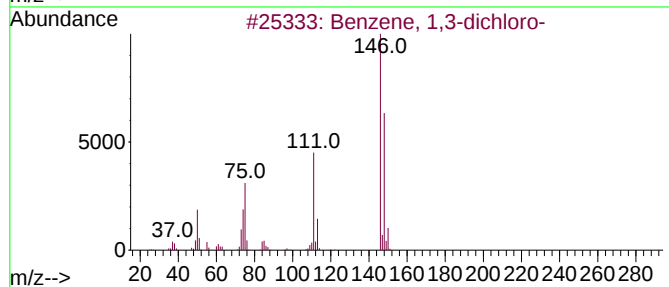
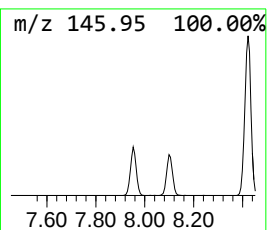
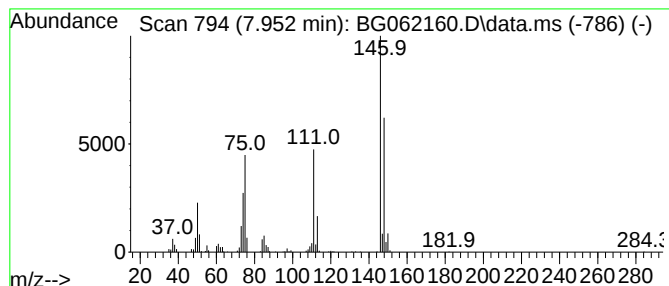
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.952	59.77 ng/ul	619184	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	96
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	93
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93



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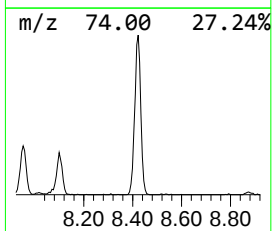
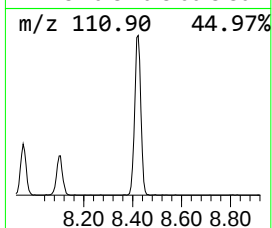
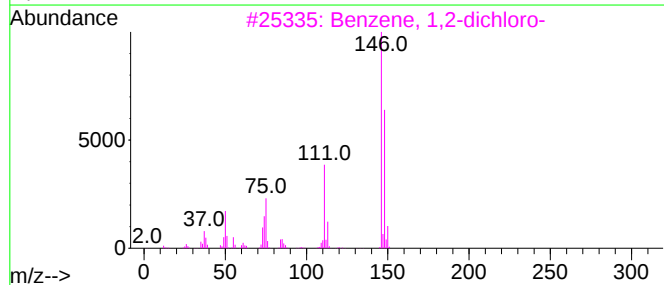
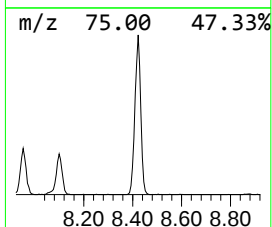
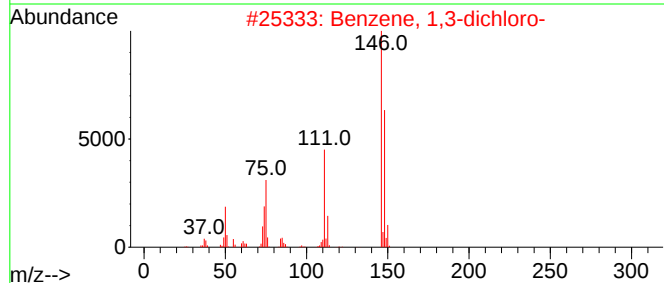
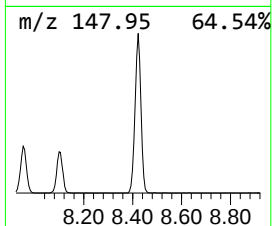
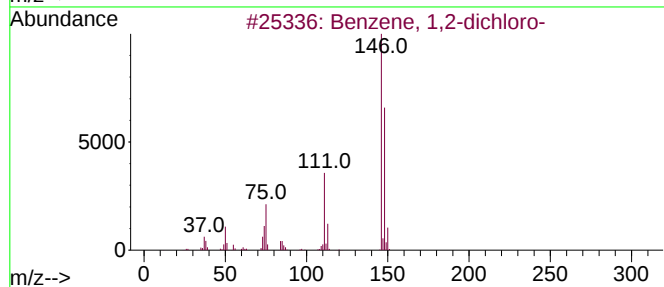
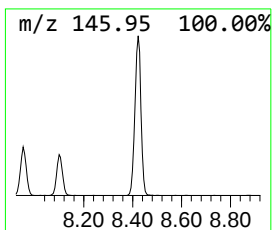
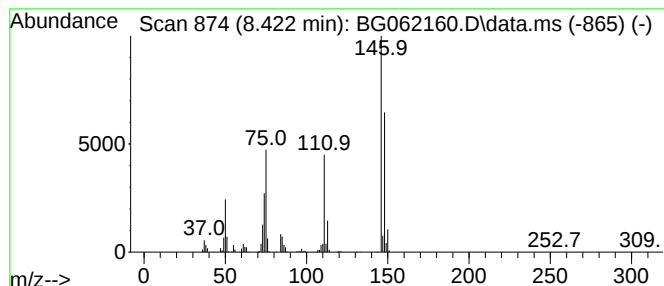
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.422	209.12 ng/ul	2166400	1,4-Dichlorobenzene-d4	8.064

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



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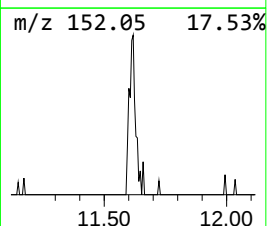
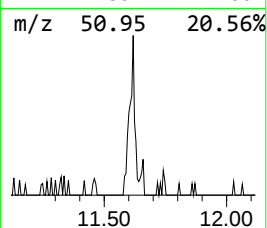
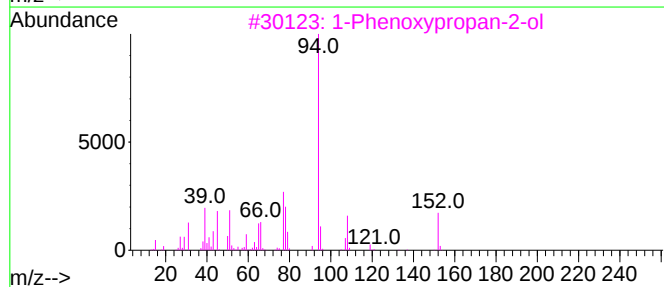
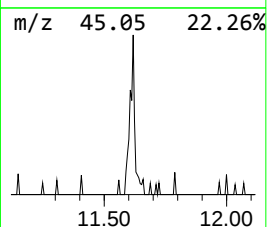
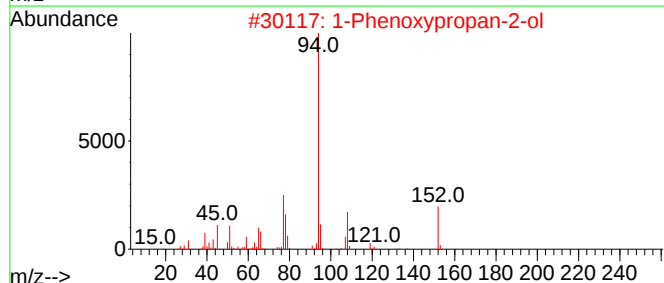
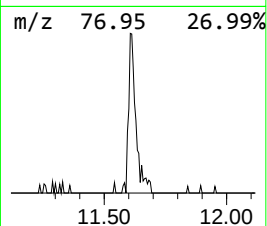
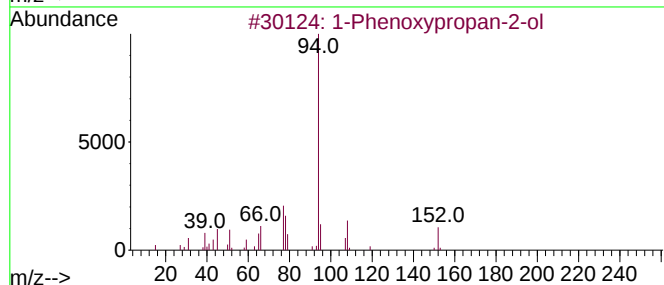
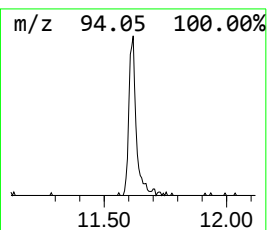
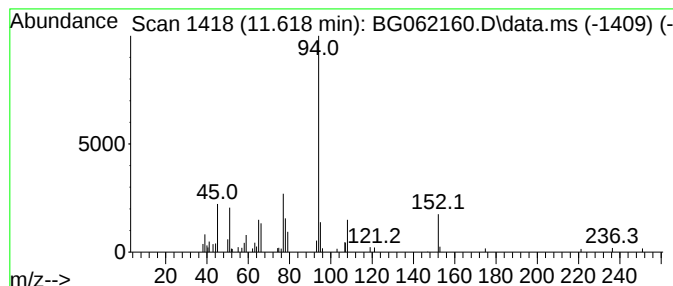
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.618	4.66 ng/ul	78042	Naphthalene-d8	10.878

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
Acq On : 22 Jul 2024 13:26
Operator : MA/JU
Sample : P3238-23
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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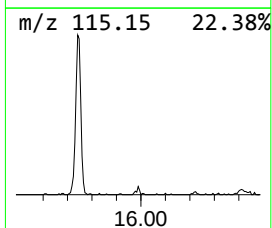
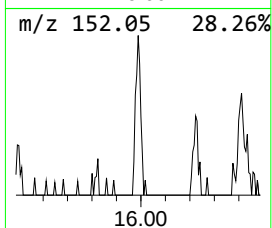
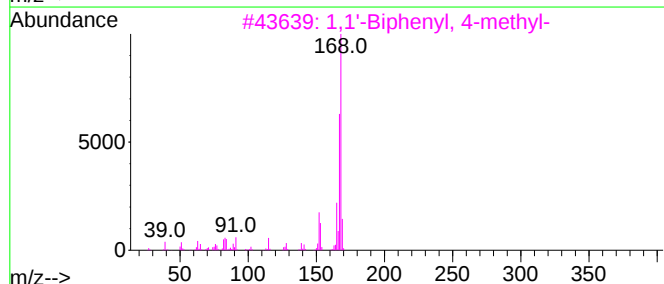
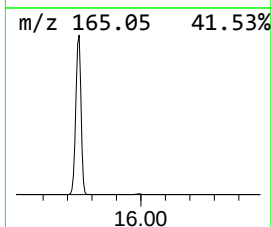
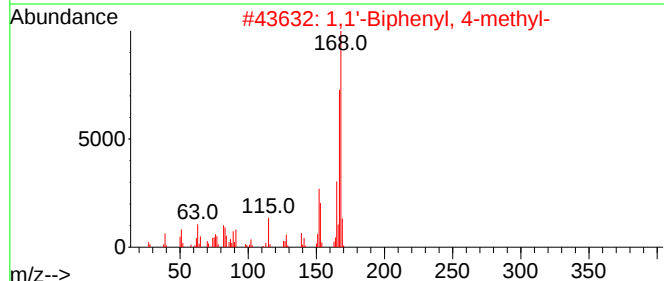
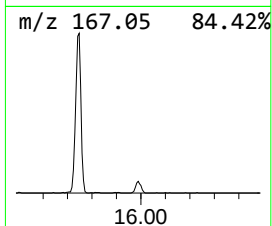
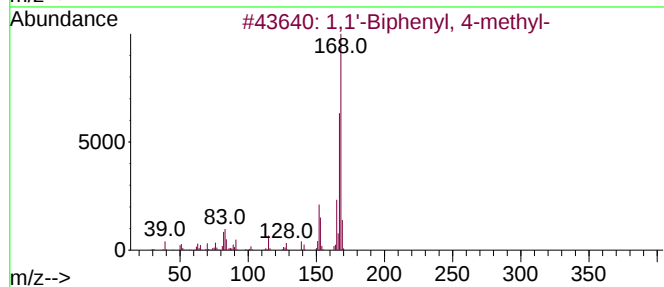
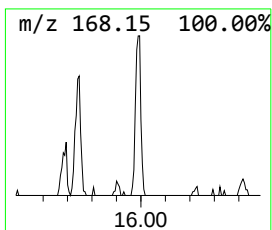
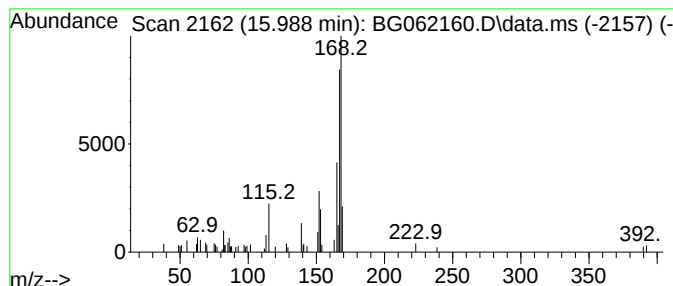
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.989	2.05 ng/ul	55408	Acenaphthene-d10	14.696

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	74
2			1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	74
3			1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	74
4			Nordiphenamid	225	C15H15NO	000954-21-2	72
5			1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
Acq On : 22 Jul 2024 13:26
Operator : MA/JU
Sample : P3238-23
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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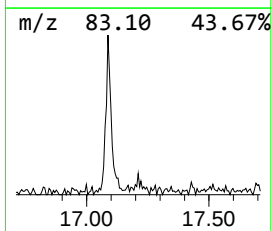
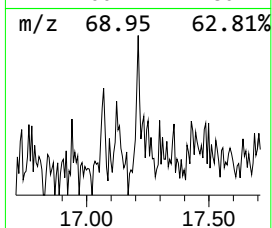
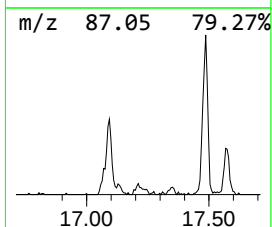
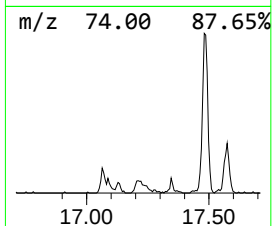
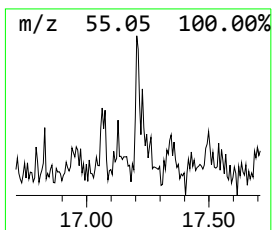
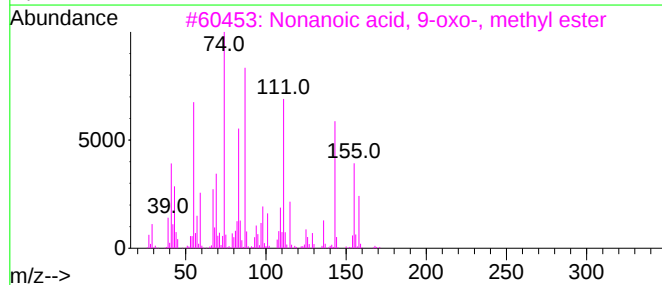
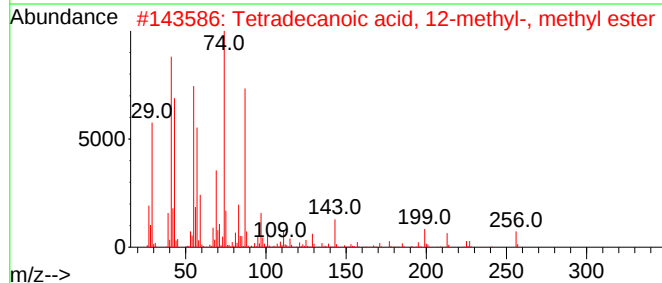
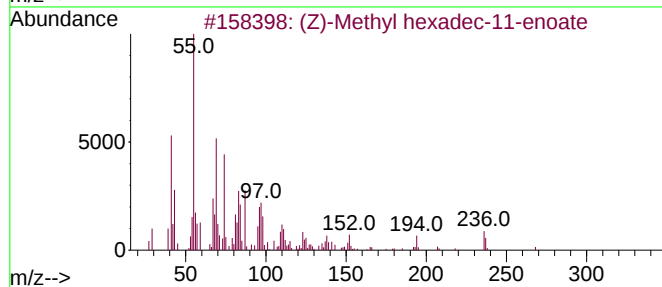
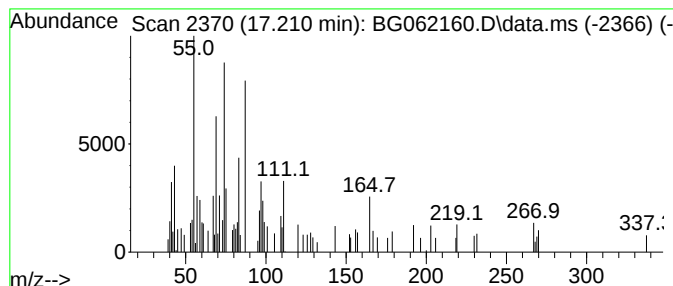
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (Z)-Methyl hexadec-11-enoate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.210	2.22 ng/ul	82332	Phenanthrene-d10	17.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(Z)-Methyl hexadec-11-enoate	268	C17H32O2	000822-05-9	50
2			Tetradecanoic acid, 12-methyl-, ...	256	C16H32O2	005129-66-8	47
3			Nonanoic acid, 9-oxo-, methyl ester	186	C10H18O3	001931-63-1	47
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	43
5			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
Acq On : 22 Jul 2024 13:26
Operator : MA/JU
Sample : P3238-23
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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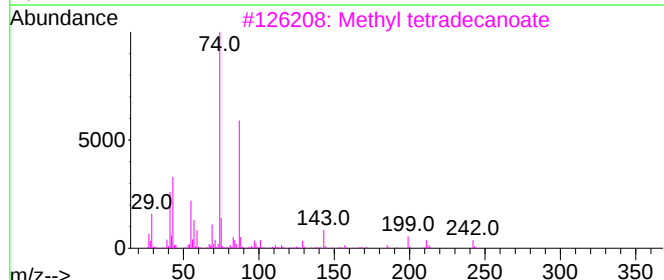
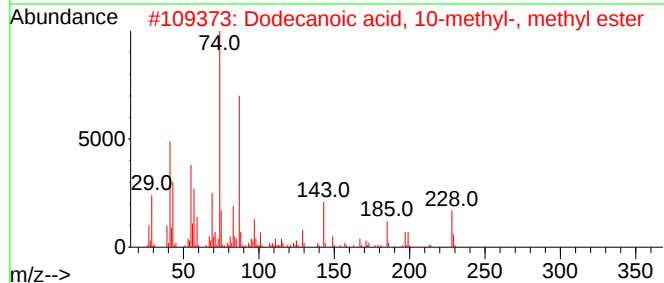
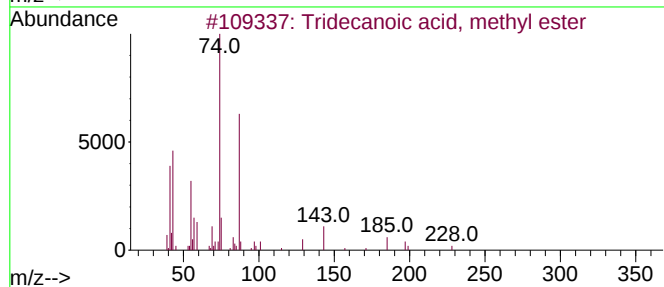
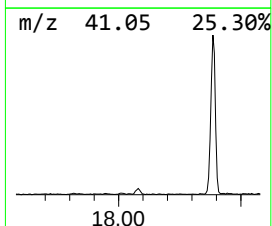
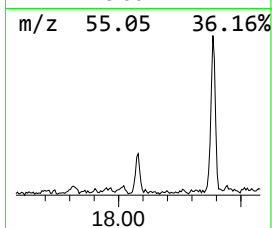
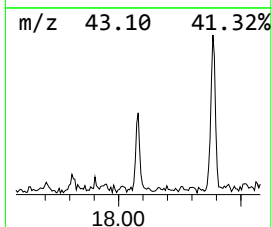
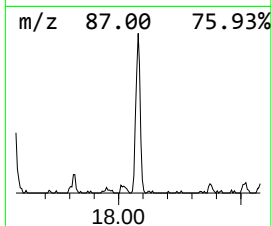
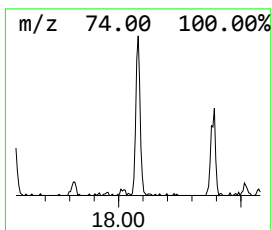
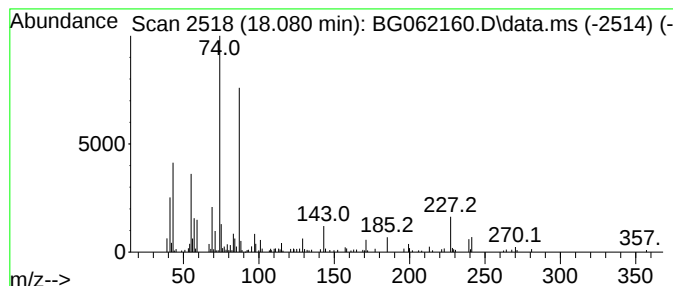
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Tridecanoic acid, methyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.080	3.96 ng/ul	146521	Phenanthrene-d10	17.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
2			Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	83
3			Methyl tetradecanoate	242	C15H30O2	000124-10-7	80
4			Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	80
5			Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
Acq On : 22 Jul 2024 13:26
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Misc :
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Instrument :
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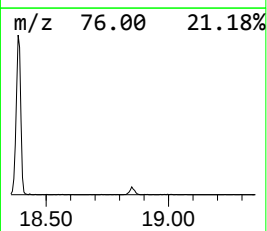
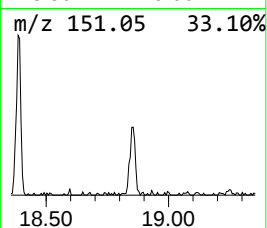
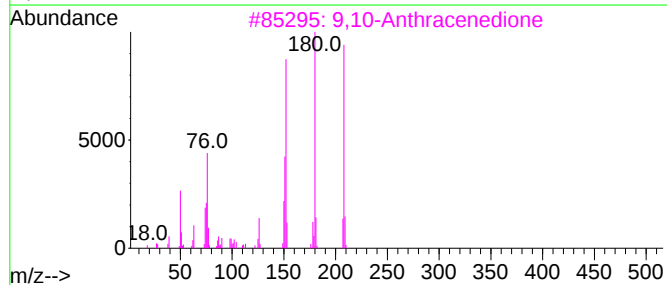
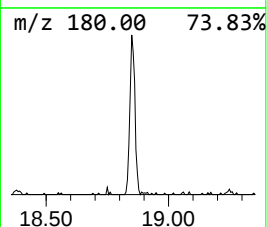
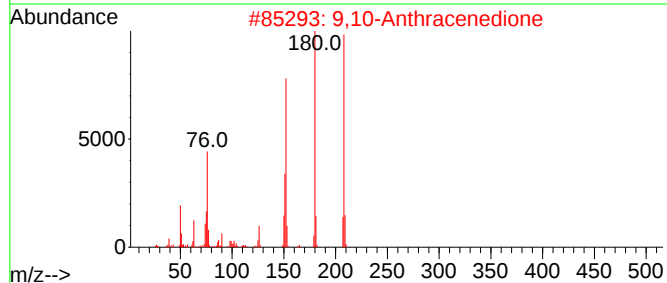
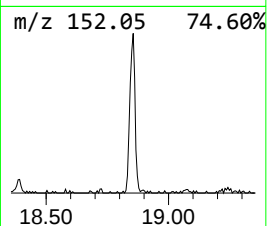
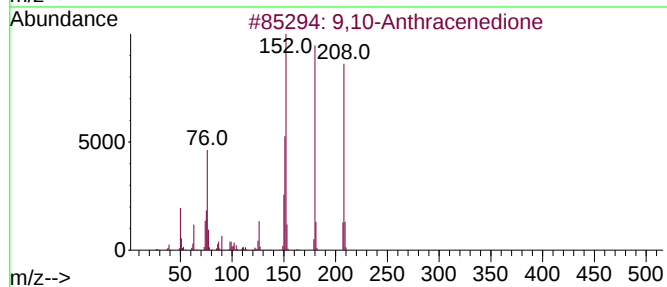
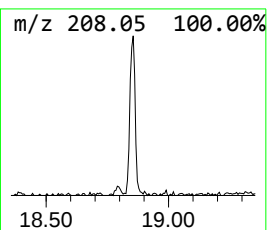
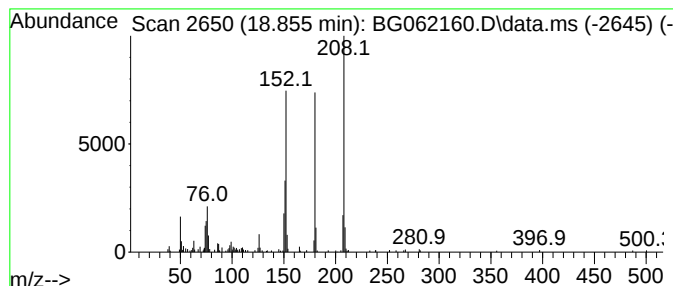
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.855	5.08 ng/ul	188135	Phenanthrene-d10	17.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	97
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	90
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	89
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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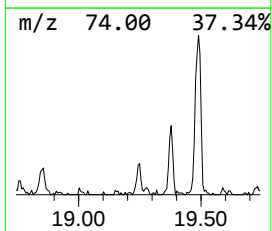
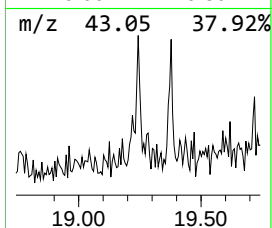
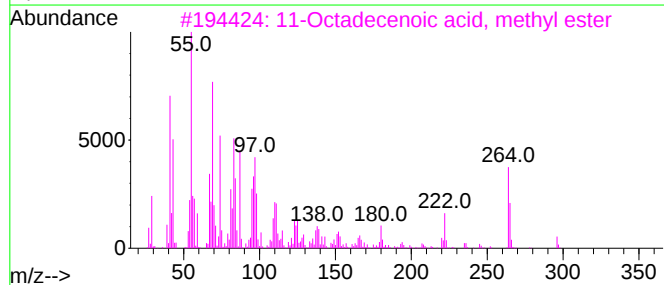
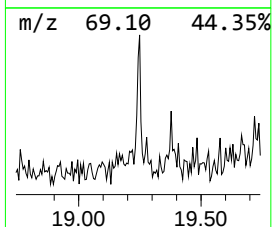
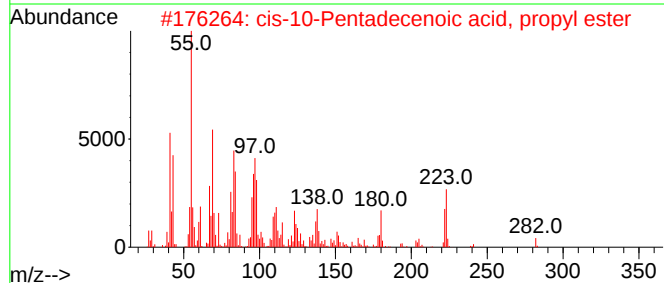
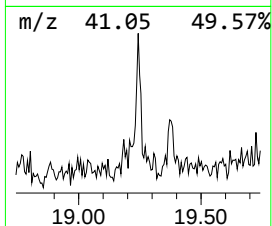
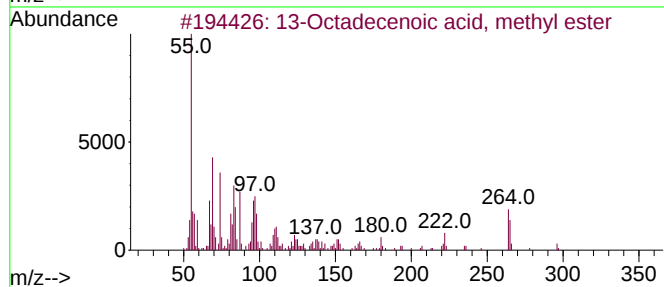
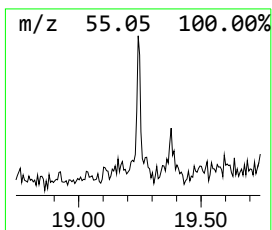
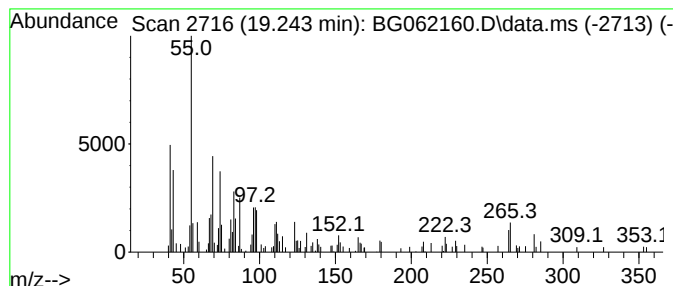
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 13-Octadecenoic acid, methy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.243	2.19 ng/ul	81189	Phenanthrene-d10	17.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	72
2			cis-10-Pentadecenoic acid, propy...	282	C18H34O2	1000405-14-4	70
3			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	58
4			15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	58
5			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
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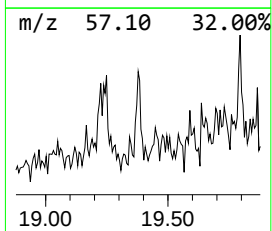
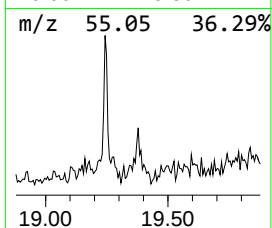
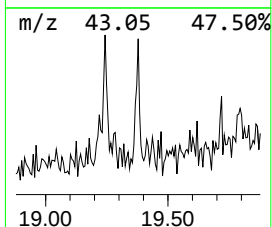
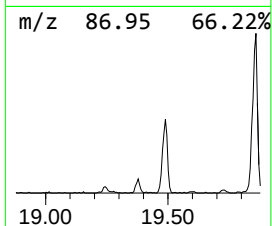
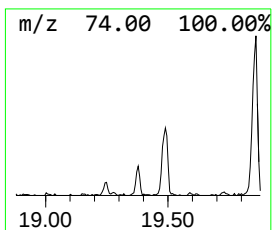
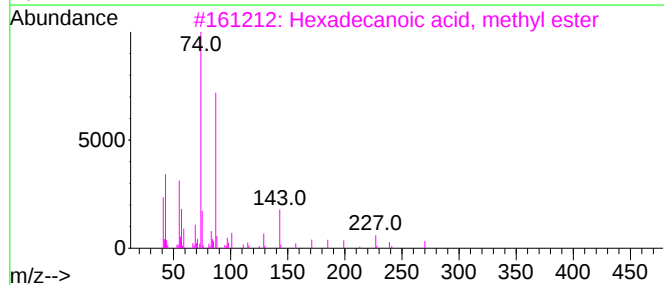
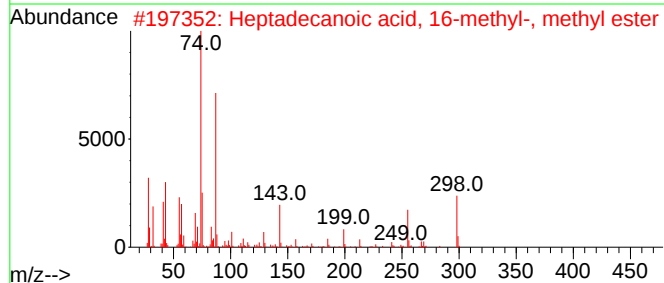
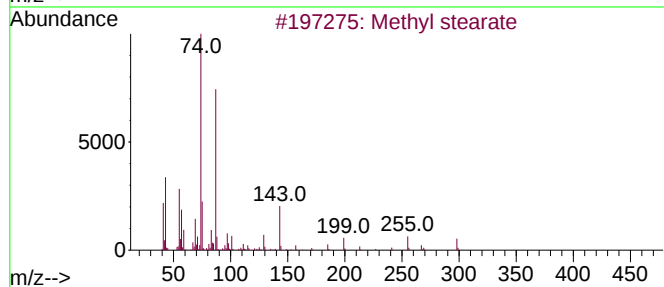
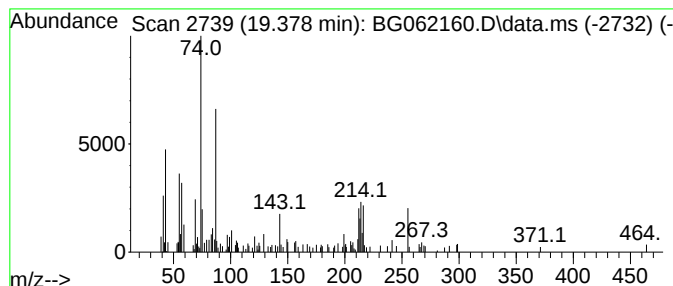
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Methyl stearate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.378	2.41 ng/ul	89124	Phenanthrene-d10	17.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	90
2			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	70
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	50
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	50
5			Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
Acq On : 22 Jul 2024 13:26
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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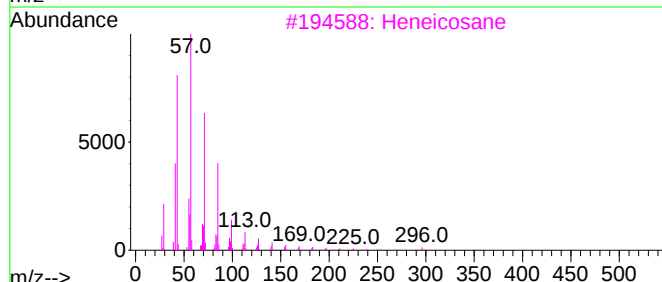
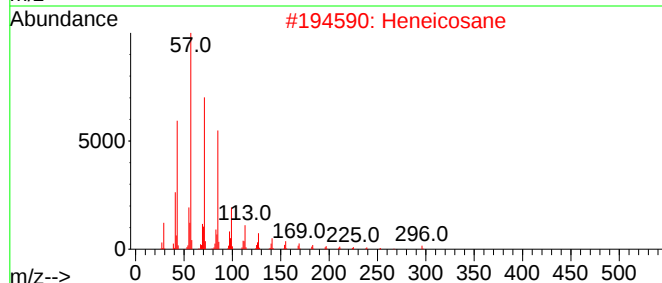
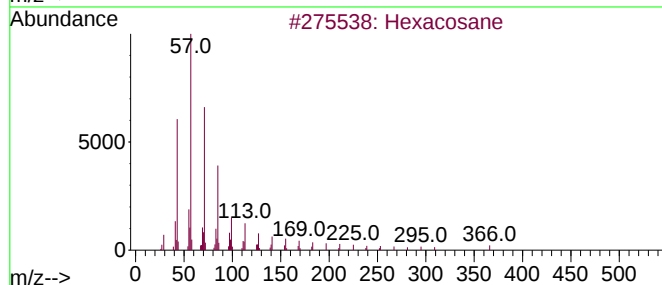
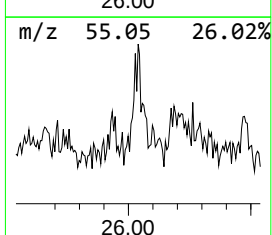
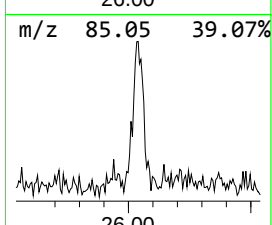
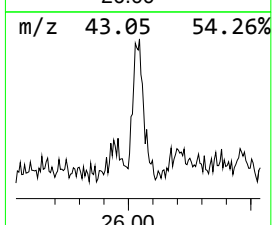
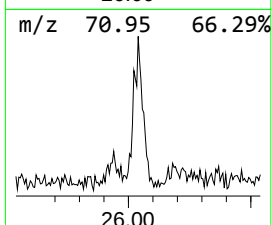
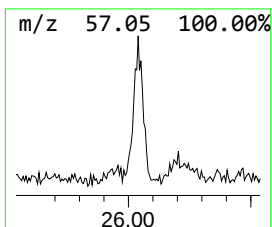
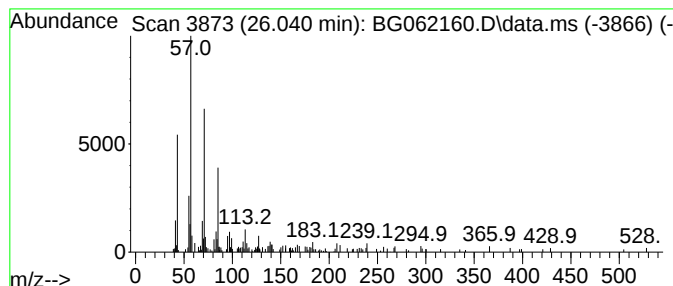
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.040	4.41 ng/ul	190017	Perylene-d12	24.929

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Heneicosane	296	C21H44	000629-94-7	90
3			Heneicosane	296	C21H44	000629-94-7	87
4			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	80
5			Heptadecane, 2,3-dimethyl-	268	C19H40	061868-03-9	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072224\
Data File : BG062160.D
Acq On : 22 Jul 2024 13:26
Operator : MA/JU
Sample : P3238-23
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4DN3

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
Benzene, 1,3-di...	7.952	59.8	ng/ul	619184	1	8.064	207190	20.0
Benzene, 1,2-di...	8.422	209.1	ng/ul	2166400	1	8.064	207190	20.0
1-Phenoxypropan...	11.618	4.7	ng/ul	78042	2	10.878	334685	20.0
1,1'-Biphenyl, ...	15.989	2.0	ng/ul	55408	3	14.696	539265	20.0
(Z)-Methyl hexa...	17.210	2.2	ng/ul	82332	4	17.439	740809	20.0
Tridecanoic aci...	18.080	4.0	ng/ul	146521	4	17.439	740809	20.0
9,10-Anthracene...	18.855	5.1	ng/ul	188135	4	17.439	740809	20.0
13-Octadecenoic...	19.243	2.2	ng/ul	81189	4	17.439	740809	20.0
Methyl stearate	19.378	2.4	ng/ul	89124	4	17.439	740809	20.0
(DEL) Alkane: S...	26.040	4.4	ng/ul	190017	6	24.929	861176	20.0