

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
 Data File : BG061862.D
 Acq On : 3 Jul 2024 13:14
 Operator : MA/JU
 Sample : P3065-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4DN0

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-BG070224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061862.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.752	76	79	87	rVB4	12072	20674	0.24%	0.016%
2	4.252	158	164	173	rVB2	25973	45383	0.53%	0.035%
3	5.721	406	414	417	rBV2	15630	27358	0.32%	0.021%
4	7.278	666	679	689	rBV2	861199	1688128	19.89%	1.304%
5	7.419	694	703	711	rVV	99857	167833	1.98%	0.130%
6	7.513	713	719	727	rVB	193725	305053	3.59%	0.236%
7	7.660	739	744	754	rVB	486447	850059	10.02%	0.656%
8	7.983	791	799	808	rBV	262655	431757	5.09%	0.333%
9	8.094	811	818	820	rVV	174224	268626	3.17%	0.207%
10	8.135	820	825	833	rVB	989508	1672417	19.71%	1.292%
11	8.453	865	879	885	rBV	690406	1192814	14.05%	0.921%
12	8.799	932	938	947	rBV	183172	318365	3.75%	0.246%
13	8.899	947	955	963	rBV	818022	1375929	16.21%	1.063%
14	9.181	995	1003	1010	rBV	537189	949850	11.19%	0.734%
15	9.264	1010	1017	1020	rVV	148442	272282	3.21%	0.210%
16	9.305	1020	1024	1031	rVB	251174	436214	5.14%	0.337%
17	10.016	1133	1145	1153	rVB3	307945	768168	9.05%	0.593%
18	10.309	1188	1195	1202	rVB	432678	699100	8.24%	0.540%
19	10.550	1225	1236	1243	rBV2	399116	1031681	12.16%	0.797%
20	10.915	1290	1298	1301	rBV	261275	479179	5.65%	0.370%
21	10.962	1301	1306	1314	rVV	1494406	2747428	32.37%	2.122%
22	11.044	1314	1320	1330	rVB	168883	309261	3.64%	0.239%
23	11.643	1416	1422	1428	rBV5	22313	46739	0.55%	0.036%
24	12.178	1504	1513	1526	rBV	1059600	1757078	20.70%	1.357%
25	12.953	1639	1645	1654	rBV4	9756	20482	0.24%	0.016%
26	13.159	1673	1680	1686	rBV	707622	1118049	13.17%	0.863%
27	13.230	1686	1692	1703	rVV	1602762	2583373	30.44%	1.995%
28	13.606	1748	1756	1767	rVB	850076	1410592	16.62%	1.089%
29	13.776	1780	1785	1790	rBV5	17207	18918	0.22%	0.015%
30	13.999	1817	1823	1829	rBV5	20532	34633	0.41%	0.027%
31	14.128	1838	1845	1851	rBV	529280	804678	9.48%	0.621%
32	14.305	1866	1875	1881	rBV	2533980	3981424	46.91%	3.075%
33	14.358	1881	1884	1889	rVV5	14599	31983	0.38%	0.025%
34	14.452	1889	1900	1910	rVB2	2339706	4438449	52.30%	3.428%
35	14.722	1937	1946	1951	rBV	454004	695380	8.19%	0.537%
36	14.787	1951	1957	1967	rVB	2061727	3333762	39.28%	2.575%

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Title : SVOA CALIBRATION

37	14.939	1972	1983	1994	rBV2	1741910	3313051	39.04%	2.559%
38	15.033	1994	1999	2001	rVV4	43980	68017	0.80%	0.053%
39	15.086	2001	2008	2019	rVB	1542495	2356065	27.76%	1.820%
40	15.339	2046	2051	2056	rBV6	34074	52523	0.62%	0.041%
41	15.539	2077	2085	2090	rBV	3613555	5358317	63.13%	4.138%
42	15.709	2107	2114	2118	rBV	717551	1100656	12.97%	0.850%
43	15.762	2118	2123	2129	rVV2	3656057	5451842	64.24%	4.210%
44	15.821	2129	2133	2144	rVB	308036	471679	5.56%	0.364%
45	15.944	2150	2154	2161	rVB10	8886	18262	0.22%	0.014%
46	16.014	2161	2166	2172	rBV3	19266	34307	0.40%	0.026%
47	16.249	2201	2206	2211	rBV5	19037	32460	0.38%	0.025%
48	16.302	2211	2215	2220	rVB4	25705	40034	0.47%	0.031%
49	16.408	2229	2233	2234	rBV2	16831	19159	0.23%	0.015%
50	16.432	2234	2237	2242	rVV6	33763	60364	0.71%	0.047%
51	16.602	2261	2266	2273	rVB2	45557	63871	0.75%	0.049%
52	16.767	2286	2294	2301	rBV	4001429	6439237	75.87%	4.973%
53	16.978	2325	2330	2337	rVB9	13329	23085	0.27%	0.018%
54	17.113	2343	2353	2360	rBV	2666468	4073260	47.99%	3.146%
55	17.242	2371	2375	2377	rBV5	34432	51098	0.60%	0.039%
56	17.466	2402	2413	2416	rVV	543101	894071	10.53%	0.690%
57	17.507	2416	2420	2425	rVV	1572759	2297291	27.07%	1.774%
58	17.566	2425	2430	2432	rVV2	831306	1342604	15.82%	1.037%
59	17.601	2432	2436	2442	rVB	2091468	3029817	35.70%	2.340%
60	17.730	2453	2458	2463	rBV5	17748	29986	0.35%	0.023%
61	17.859	2473	2480	2485	rVB5	25875	53035	0.62%	0.041%
62	18.059	2509	2514	2518	rVV4	57557	93815	1.11%	0.072%
63	18.112	2518	2523	2530	rVB	146151	204971	2.42%	0.158%
64	18.335	2558	2561	2568	rVB5	33524	43966	0.52%	0.034%
65	18.418	2568	2575	2581	rBV	4205317	5605445	66.05%	4.329%
66	18.517	2589	2592	2596	rBV6	14147	18718	0.22%	0.014%
67	18.623	2609	2610	2614	rVV4	16401	20334	0.24%	0.016%
68	18.705	2617	2624	2629	rVV8	39443	99335	1.17%	0.077%
69	18.758	2629	2633	2636	rVV5	30317	47810	0.56%	0.037%
70	18.817	2636	2643	2647	rVV3	30059	64576	0.76%	0.050%
71	18.870	2647	2652	2660	rVB	245103	369863	4.36%	0.286%
72	19.199	2705	2708	2711	rBV5	12344	18801	0.22%	0.015%
73	19.281	2719	2722	2725	rVV	87944	105954	1.25%	0.082%
74	19.305	2725	2726	2730	rVB4	20306	20551	0.24%	0.016%
75	19.410	2739	2744	2747	rBV2	61759	75716	0.89%	0.058%
76	19.516	2753	2762	2769	rBV	4475390	6746436	79.49%	5.210%

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 Title : SVOA CALIBRATION

77	19.751	2798	2802	2806	rVB	51343	65914	0.78%	0.051%
78	19.839	2811	2817	2819	rVV	1062352	1549027	18.25%	1.196%
79	19.869	2819	2822	2828	rVV	2307624	3351131	39.48%	2.588%
80	19.928	2829	2832	2837	rVV6	30469	48486	0.57%	0.037%
81	19.969	2837	2839	2843	rVB5	15215	21211	0.25%	0.016%
82	20.116	2862	2864	2868	rVB5	22501	22352	0.26%	0.017%
83	20.497	2926	2929	2933	rBV2	76657	82142	0.97%	0.063%
84	20.662	2955	2957	2960	rBV4	16508	23024	0.27%	0.018%
85	20.844	2985	2988	2993	rVB7	29858	52169	0.61%	0.040%
86	21.367	3072	3077	3088	rBV3	148230	303807	3.58%	0.235%
87	21.520	3100	3103	3108	rVB	133824	173719	2.05%	0.134%
88	21.614	3112	3119	3124	rVV	5934099	8487152	100.00%	6.554%
89	21.714	3128	3136	3142	rVV2	3007321	6032215	71.07%	4.659%
90	21.784	3142	3148	3157	rVB	4082274	6925109	81.60%	5.348%
91	22.530	3272	3275	3279	rBV3	40542	65394	0.77%	0.051%
92	22.736	3306	3310	3318	rVB3	103444	161319	1.90%	0.125%
93	24.023	3517	3529	3539	rBV	2803935	7538930	88.83%	5.822%
94	24.334	3580	3582	3588	rVB5	40099	56268	0.66%	0.043%
95	24.751	3642	3653	3662	rBV	533485	1497059	17.64%	1.156%
96	24.969	3682	3690	3705	rVB	324957	979390	11.54%	0.756%
97	26.161	3886	3893	3903	rVB5	98957	286255	3.37%	0.221%
98	26.285	3912	3914	3915	rBV2	23412	20591	0.24%	0.016%
99	28.770	4318	4337	4349	rVB	1188141	5181472	61.05%	4.002%
100	29.170	4402	4405	4406	rBV3	17578	19835	0.23%	0.015%

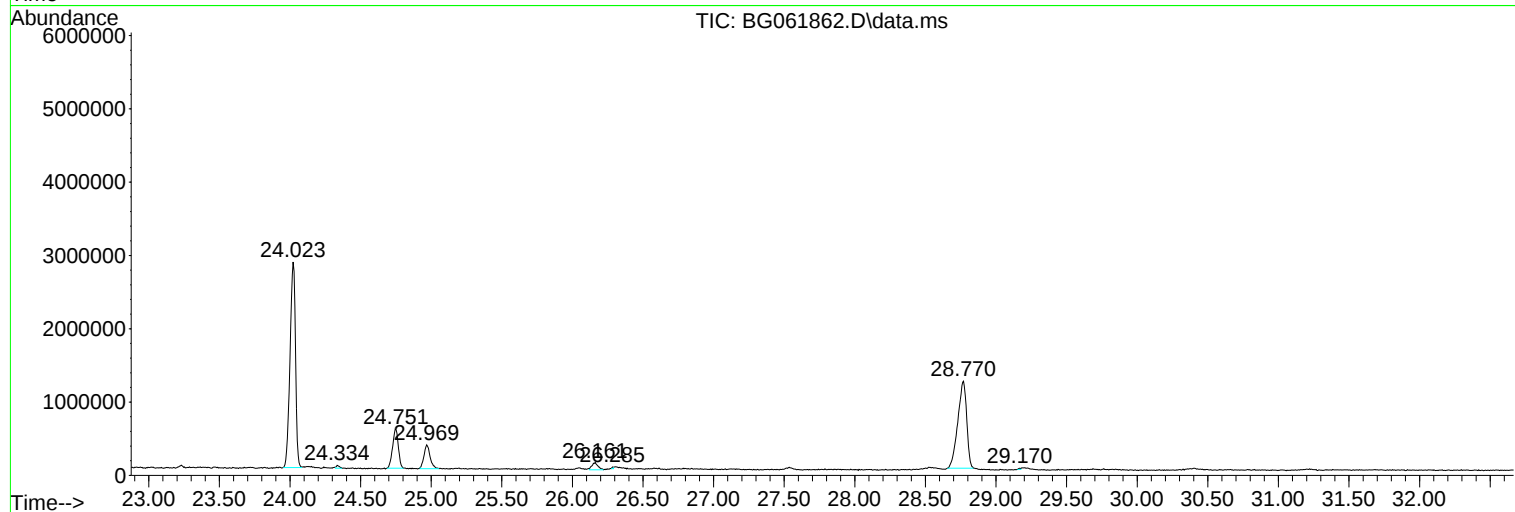
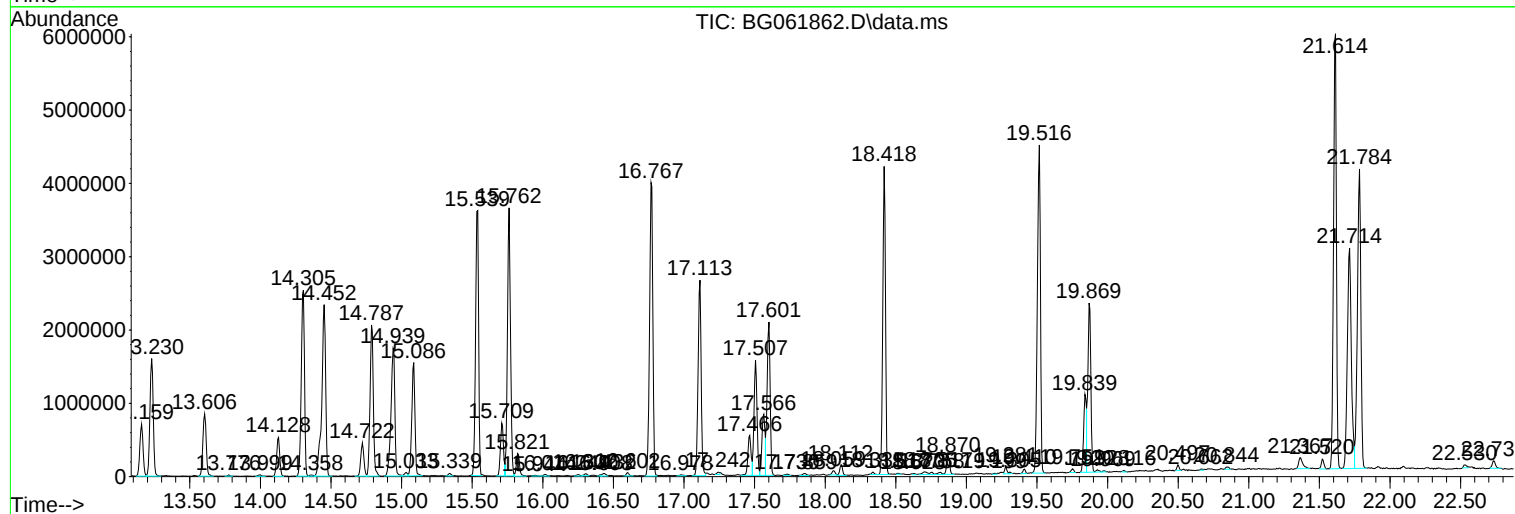
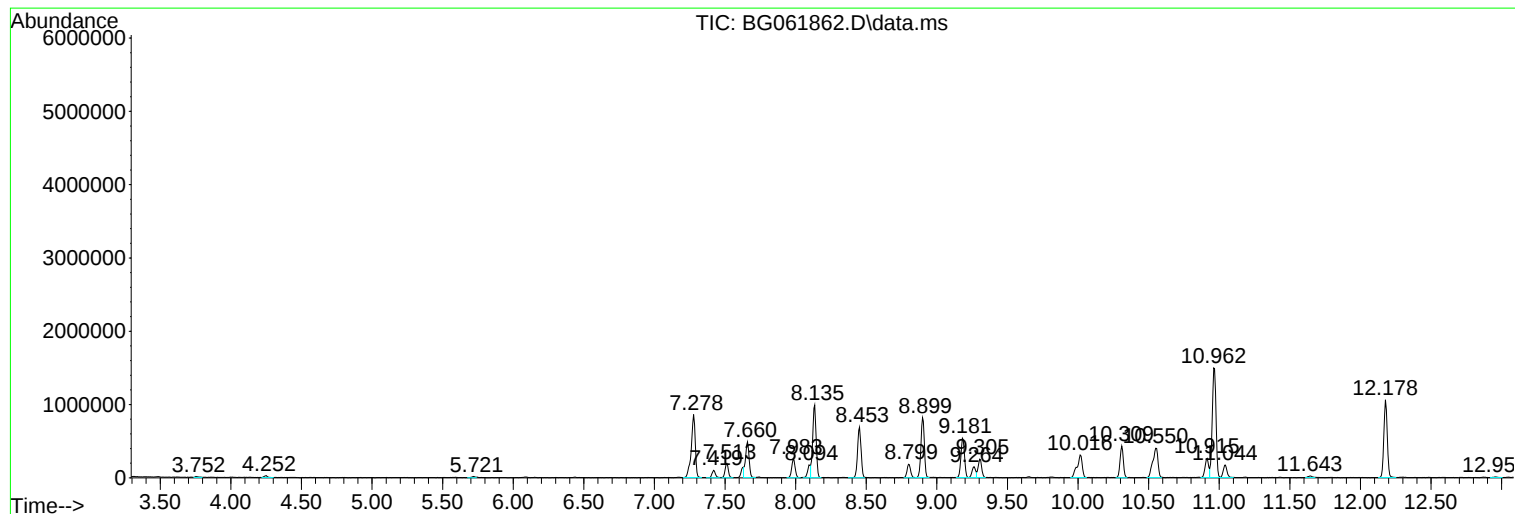
Sum of corrected areas: 129487452

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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



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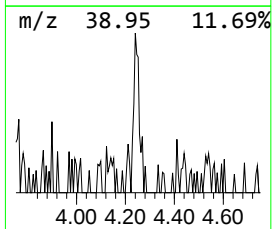
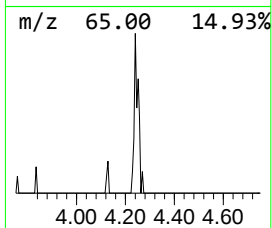
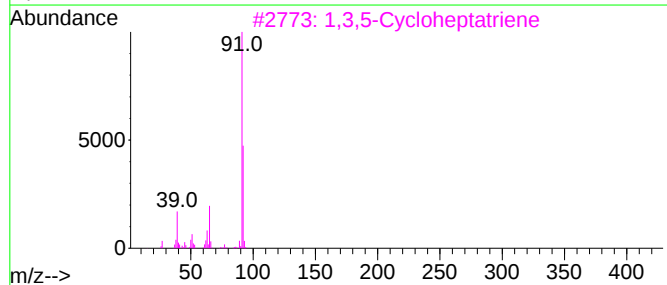
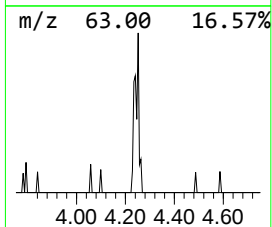
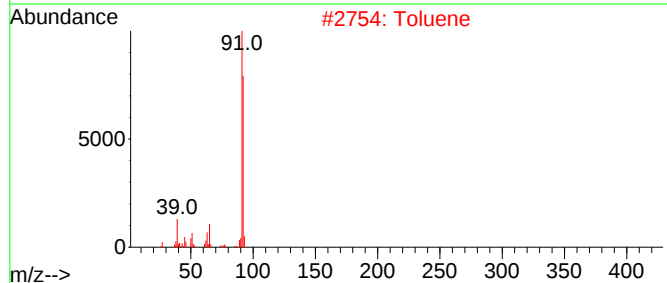
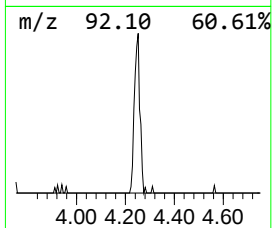
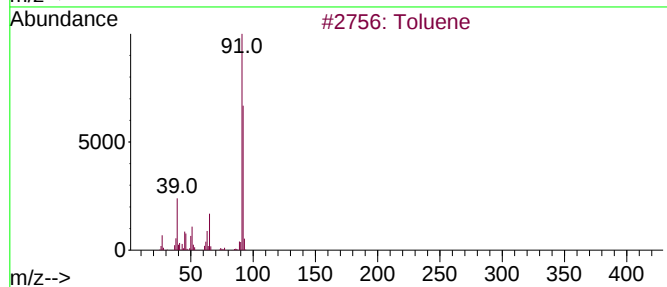
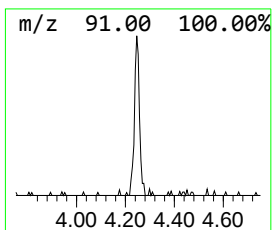
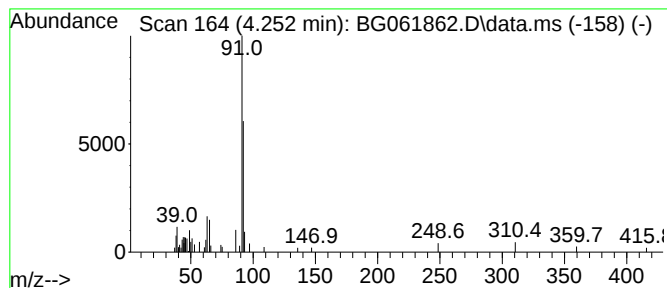
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Toluene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.252	3.38 ng/ul	45383	1,4-Dichlorobenzene-d4	8.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	53
2			Toluene	92	C7H8	000108-88-3	49
3			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	49
4			Toluene	92	C7H8	000108-88-3	47
5			Toluene	92	C7H8	000108-88-3	47



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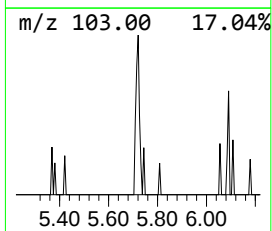
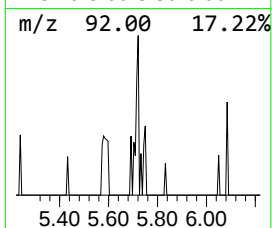
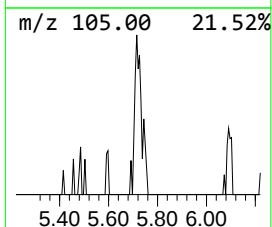
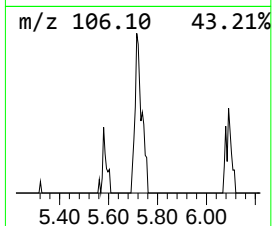
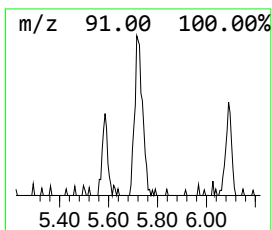
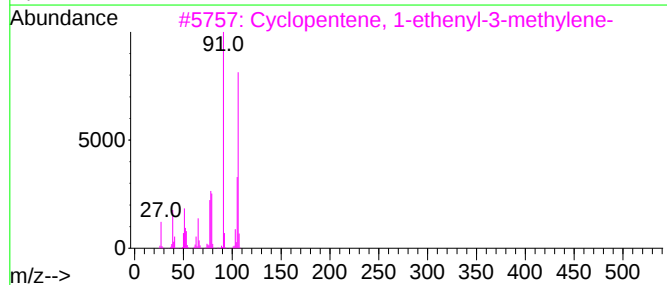
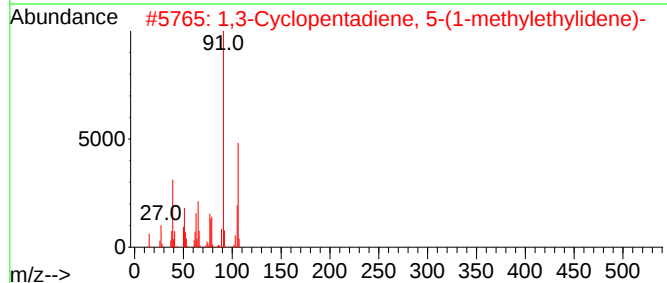
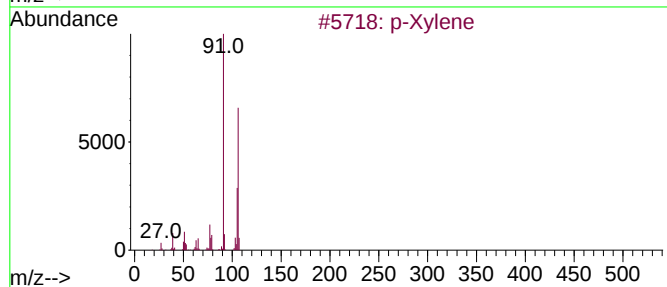
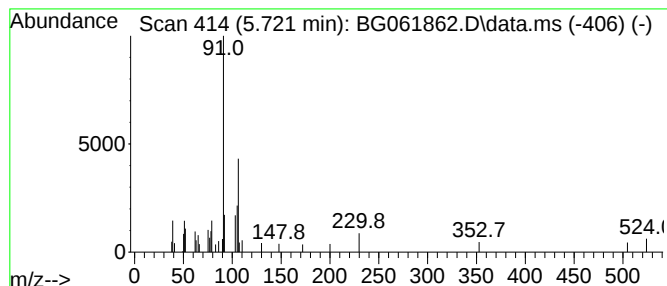
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.721	2.04 ng/ul	27358	1,4-Dichlorobenzene-d4	8.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Xylene	106	C8H10	000106-42-3	43
2			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	43
3			Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	38
4			p-Xylene	106	C8H10	000106-42-3	38
5			Benzaldehyde 3-benzylisothiosemi...	269	C15H15N3S	1000295-63-3	38



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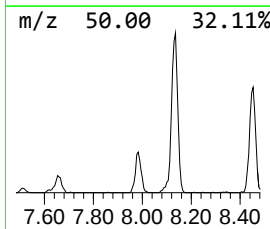
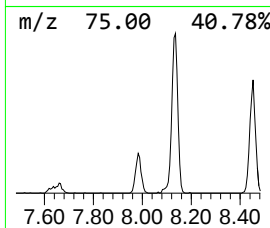
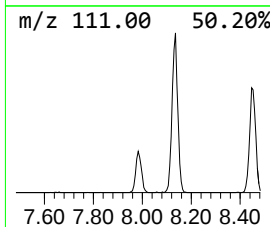
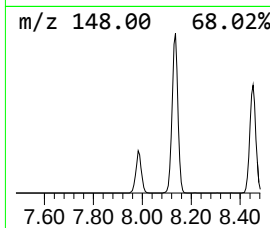
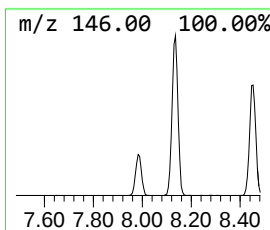
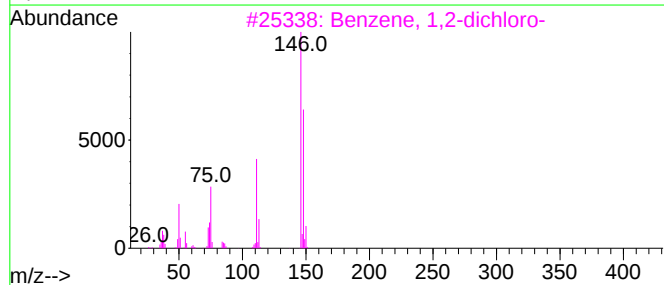
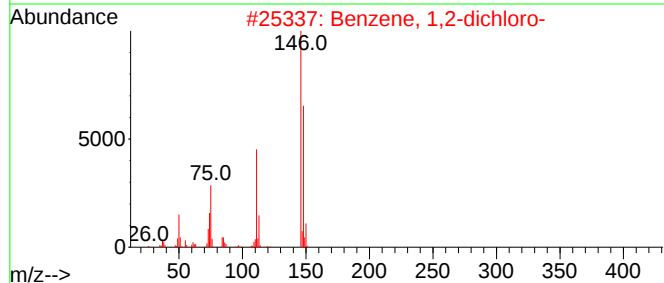
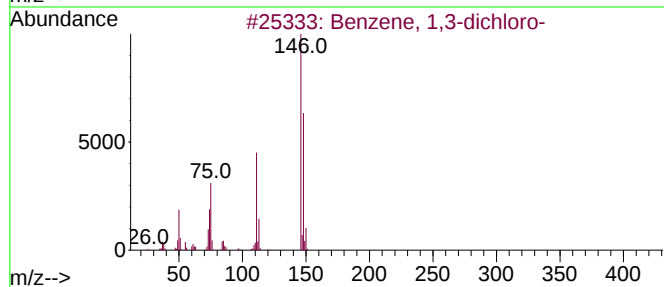
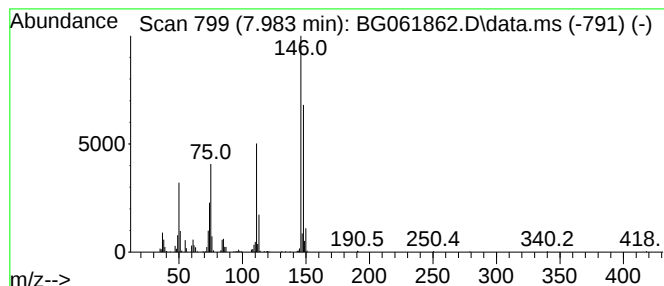
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.983	32.15 ng/ul	431757	1,4-Dichlorobenzene-d4	8.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061862.D
Acq On : 3 Jul 2024 13:14
Operator : MA/JU
Sample : P3065-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4DN0

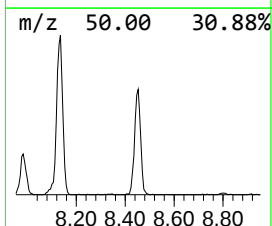
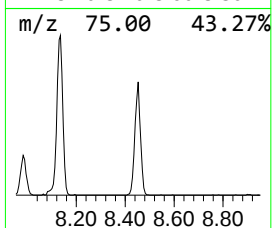
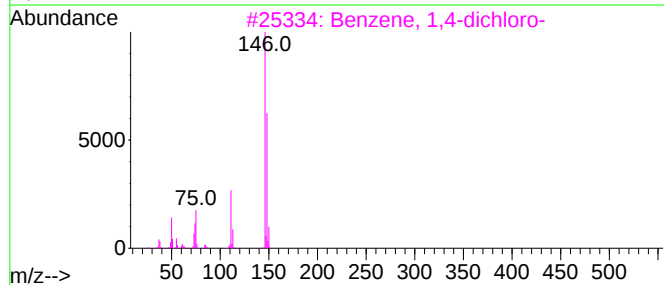
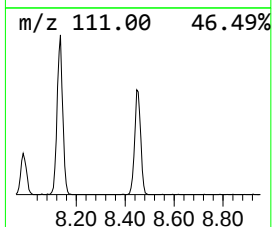
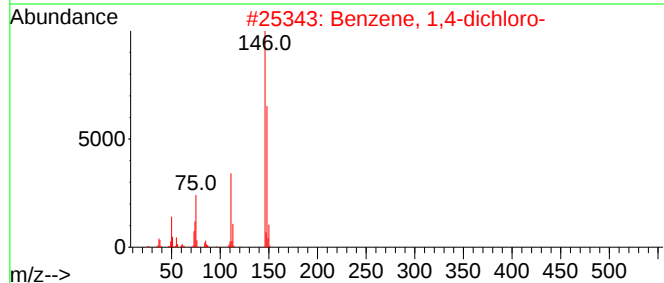
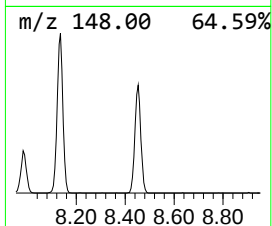
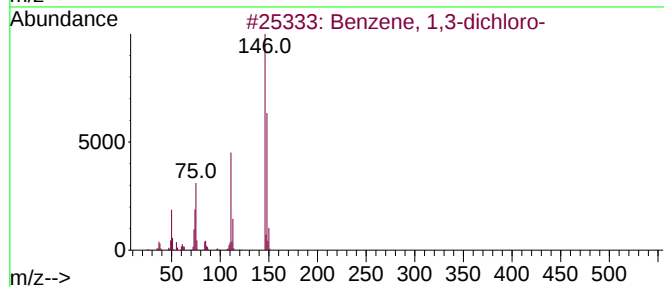
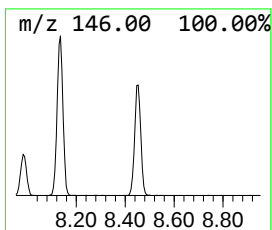
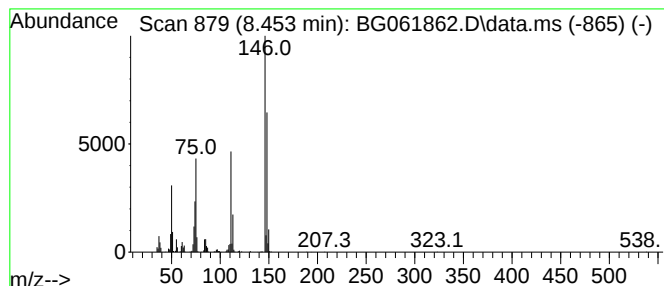
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.453	88.81 ng/ul	1192810	1,4-Dichlorobenzene-d4	8.094

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, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061862.D
Acq On : 3 Jul 2024 13:14
Operator : MA/JU
Sample : P3065-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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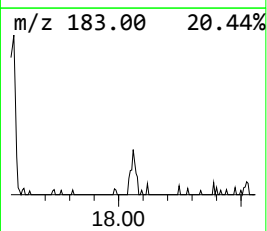
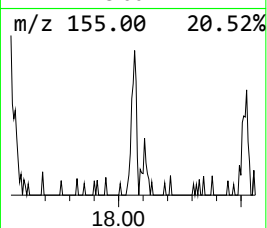
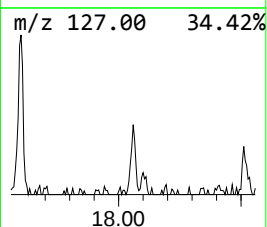
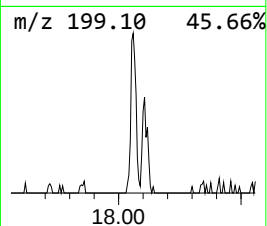
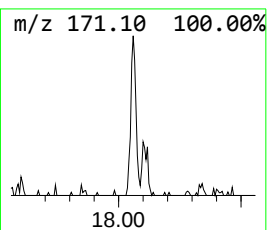
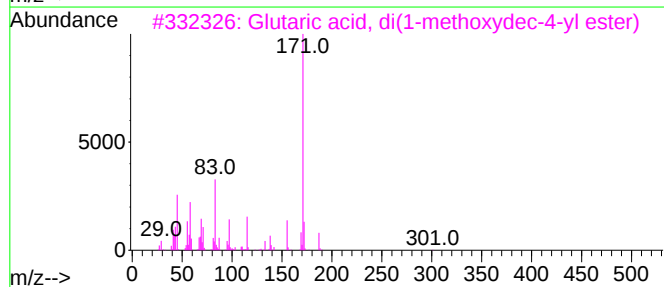
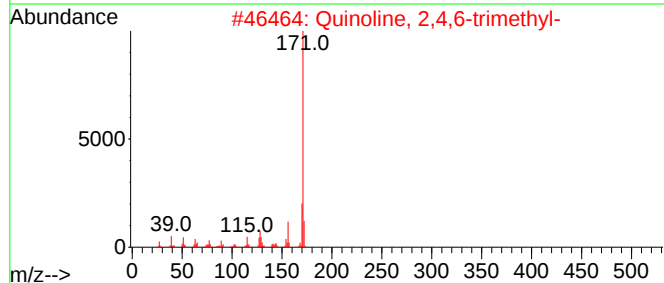
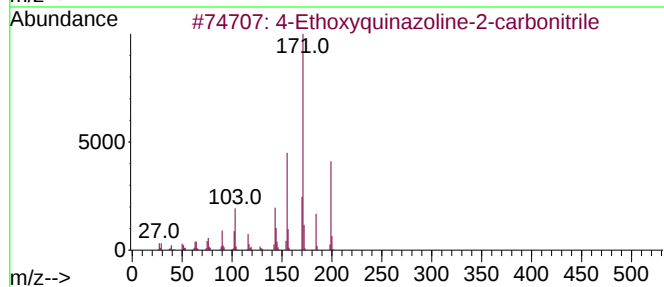
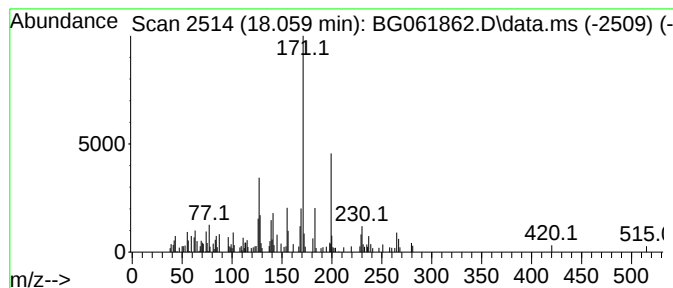
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.059	2.10 ng/ul	93815	Phenanthrene-d10	17.466

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Ethoxyquinazoline-2-carbonitrile	199	C11H9N3O	099973-07-6	35
2			Quinoline, 2,4,6-trimethyl-	171	C12H13N	002243-89-2	27
3			Glutaric acid, di(1-methoxydec-4...	472	C27H52O6	1000358-48-0	25
4			2,4(1H,3H)-Pyrimidinedione, 6-me...	171	C5H5N3O4	016632-21-6	22
5			3,5-Bis(methylthio)pyridine	171	C7H9NS2	070999-08-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061862.D
Acq On : 3 Jul 2024 13:14
Operator : MA/JU
Sample : P3065-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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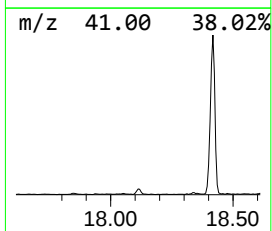
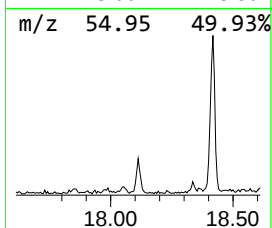
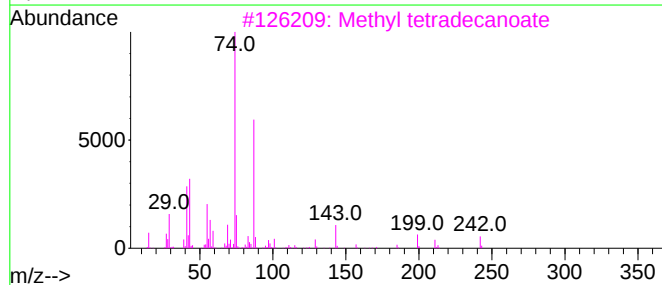
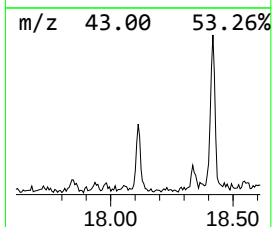
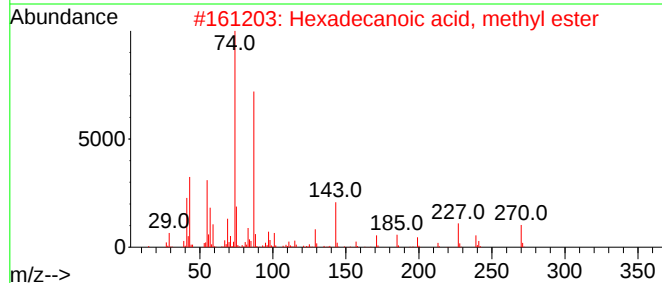
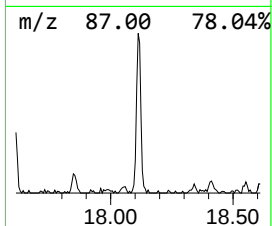
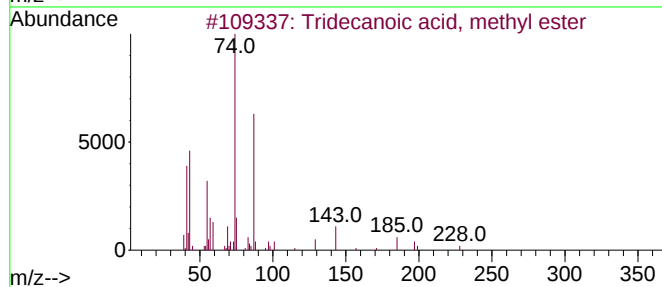
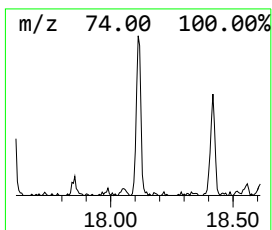
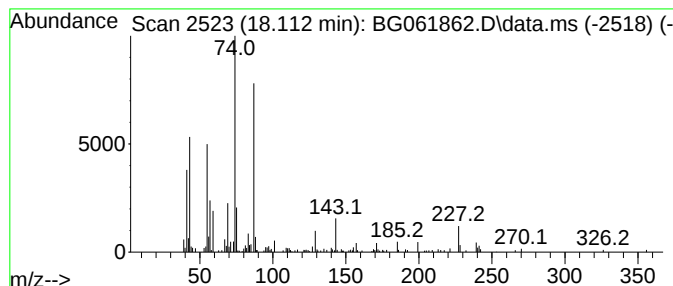
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Tridecanoic acid, methyl ester Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.112	4.59 ng/ul	204971	Phenanthrene-d10	17.466

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	86
3			Methyl tetradecanoate	242	C15H30O2	000124-10-7	80
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	72
5			Methyl 8-methyl-nonanoate	186	C11H22O2	1000452-00-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061862.D
Acq On : 3 Jul 2024 13:14
Operator : MA/JU
Sample : P3065-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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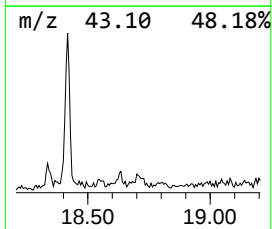
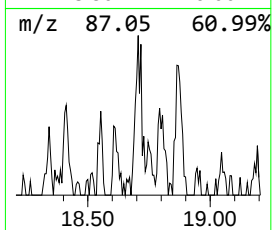
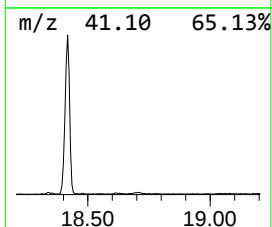
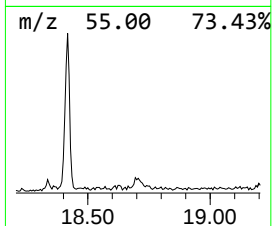
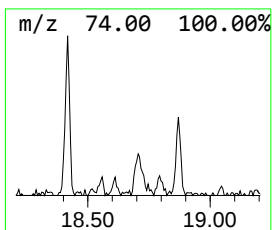
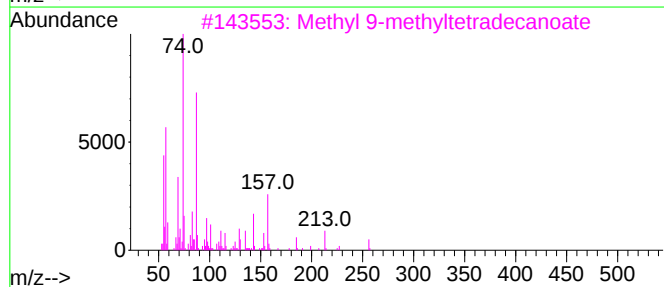
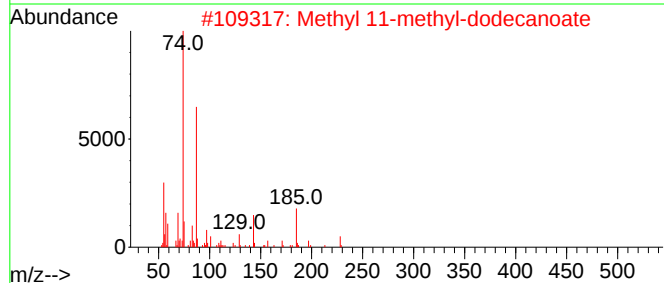
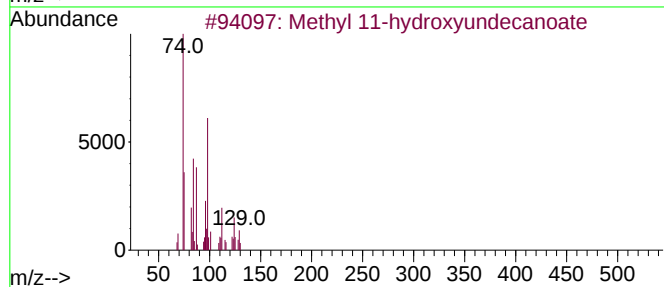
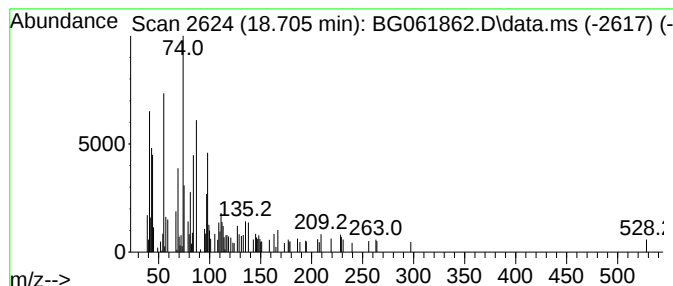
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Methyl 11-hydroxyundecanoate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.705	2.22 ng/ul	99335	Phenanthrene-d10	17.466

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 11-hydroxyundecanoate	216	C12H24O3	1000427-32-2	59
2			Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	43
3			Methyl 9-methyltetradecanoate	256	C16H32O2	213617-69-7	43
4			Tetradecanoic acid, 12-methyl-, ...	256	C16H32O2	062691-05-8	43
5			Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061862.D
Acq On : 3 Jul 2024 13:14
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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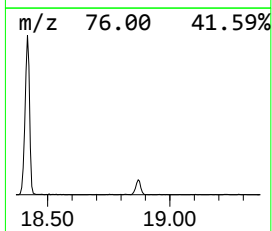
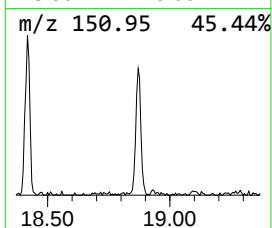
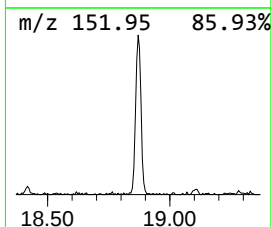
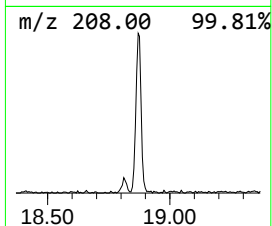
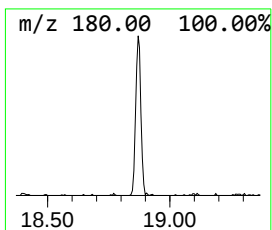
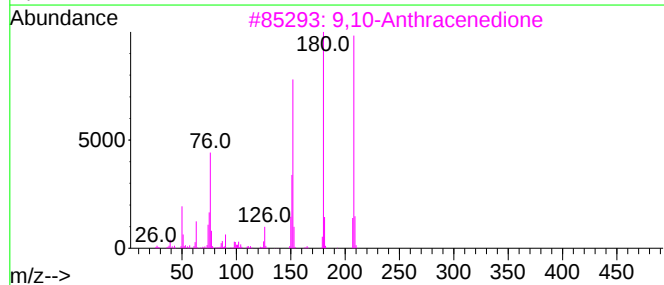
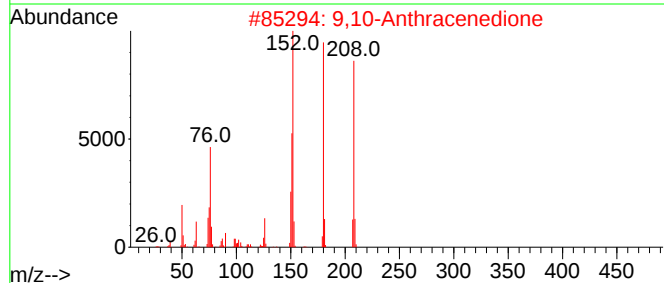
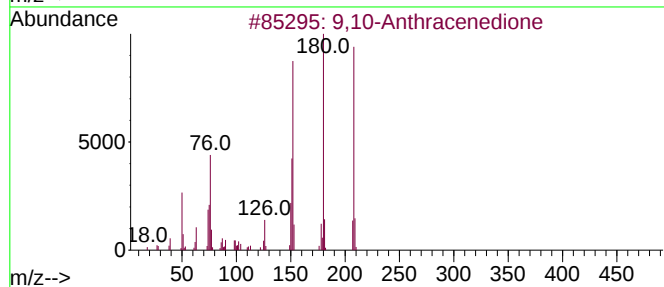
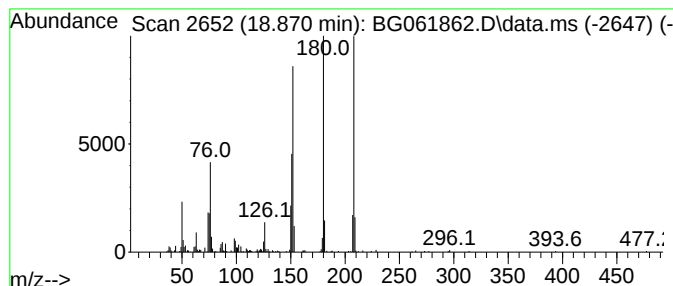
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.870	8.27 ng/ul	369863	Phenanthrene-d10	17.466

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
5			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
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Misc :
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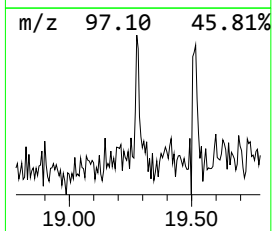
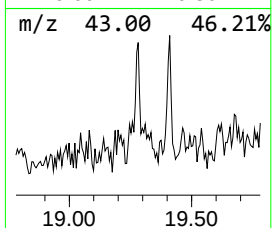
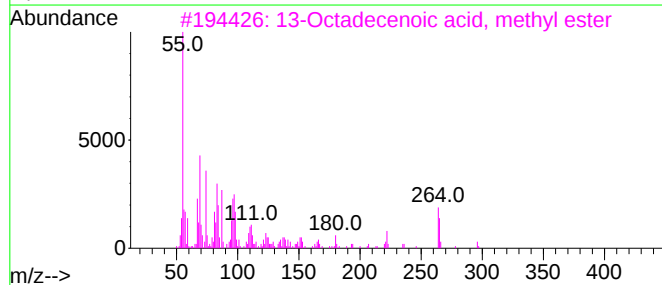
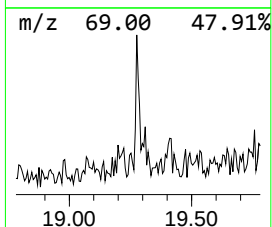
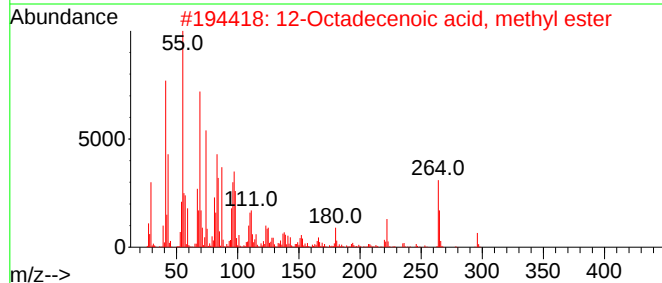
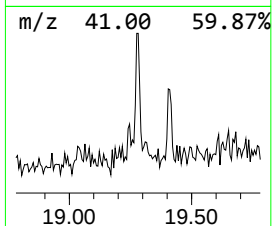
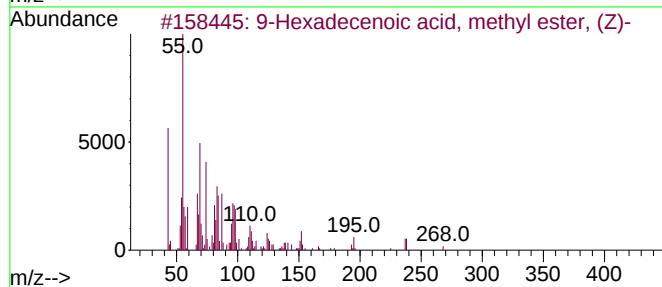
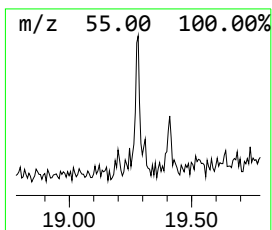
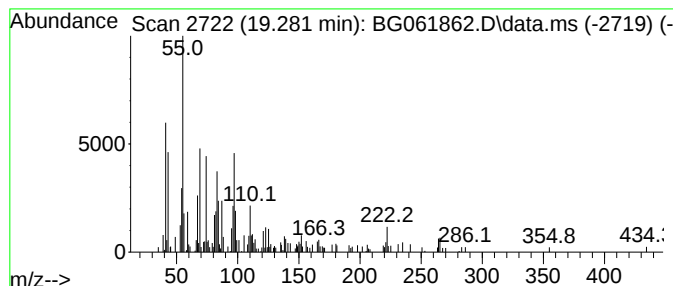
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 9-Hexadecenoic acid, methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.281	2.37 ng/ul	105954	Phenanthrene-d10	17.466	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	70
2	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	64
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	59
4	Methyl myristoleate	240	C15H28O2	056219-06-8	58
5	9-Octadecenoic acid, 12-hydroxy-...	312	C19H36O3	000141-24-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061862.D
Acq On : 3 Jul 2024 13:14
Operator : MA/JU
Sample : P3065-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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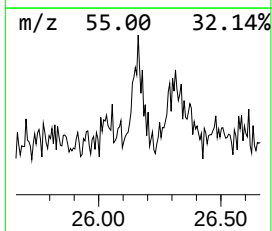
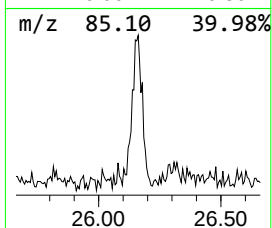
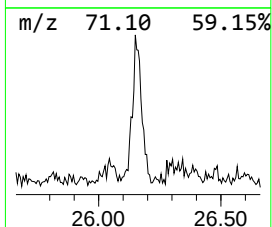
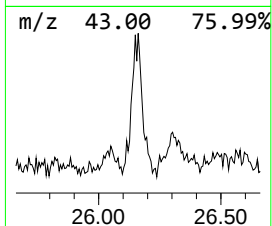
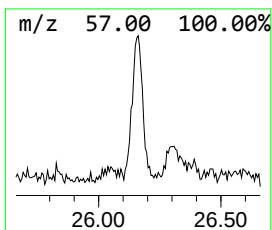
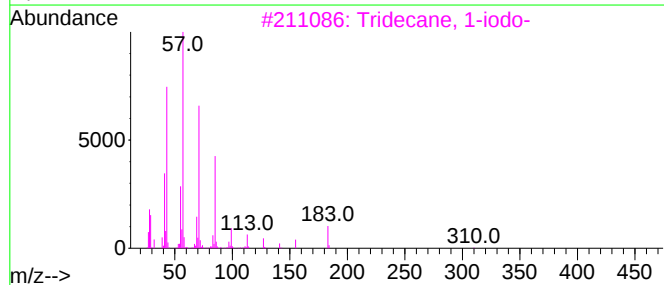
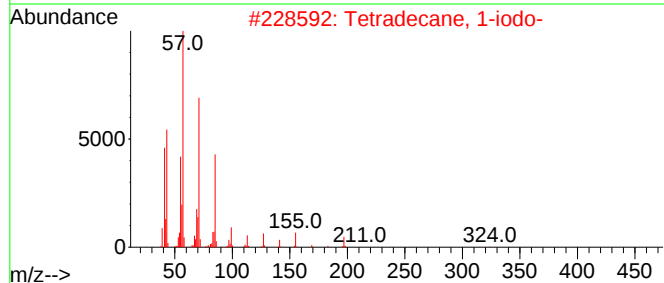
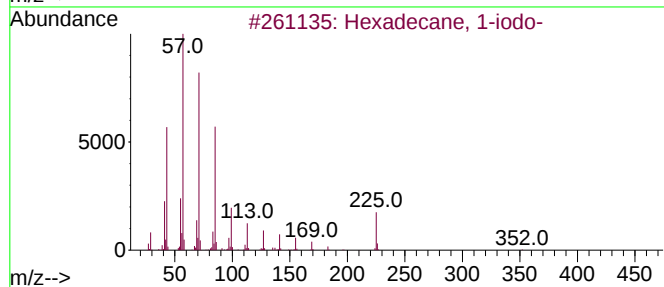
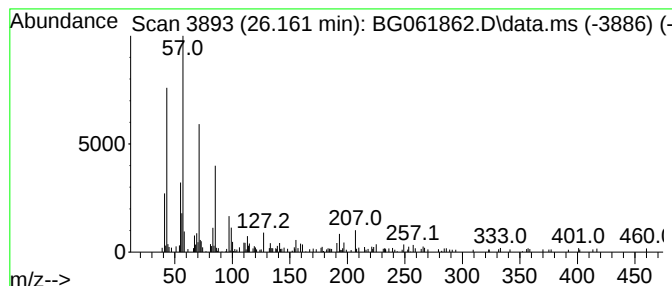
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Hexadecane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.161	5.85 ng/ul	286255	Perylene-d12	24.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	76
2			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	64
3			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64
4			Hexacosane	366	C26H54	000630-01-3	64
5			Undecane, 5-methyl-	170	C12H26	001632-70-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061862.D
Acq On : 3 Jul 2024 13:14
Operator : MA/JU
Sample : P3065-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4DN0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.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
Toluene	4.252	3.4	ng/ul	45383	1	8.094	268626	20.0
unknown-01	5.721	2.0	ng/ul	27358	1	8.094	268626	20.0
Benzene, 1,3-di...	7.983	32.1	ng/ul	431757	1	8.094	268626	20.0
Benzene, 1,4-di...	8.453	88.8	ng/ul	1192810	1	8.094	268626	20.0
unknown-02	18.059	2.1	ng/ul	93815	4	17.466	894071	20.0
Tridecanoic aci...	18.112	4.6	ng/ul	204971	4	17.466	894071	20.0
Methyl 11-hydro...	18.705	2.2	ng/ul	99335	4	17.466	894071	20.0
9,10-Anthracene...	18.870	8.3	ng/ul	369863	4	17.466	894071	20.0
9-Hexadecenoic ...	19.281	2.4	ng/ul	105954	4	17.466	894071	20.0
Hexadecane, 1-i...	26.161	5.8	ng/ul	286255	6	24.969	979390	20.0