

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
 Data File : BG063103.D
 Acq On : 28 Sep 2024 11:49
 Operator : RC/JU
 Sample : P3910-05DL 10X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDC07DL

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-BG092524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063103.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.964	312	319	325	rBV	75389	110411	4.62%	0.686%
2	5.875	471	474	480	rVB4	15758	22127	0.93%	0.137%
3	6.163	517	523	531	rVV	150412	251491	10.53%	1.562%
4	6.410	561	565	570	rVB4	9043	12928	0.54%	0.080%
5	6.733	616	620	627	rVB5	5144	11371	0.48%	0.071%
6	6.839	635	638	645	rBV7	7867	15472	0.65%	0.096%
7	6.897	645	648	654	rBV7	6236	11610	0.49%	0.072%
8	7.015	661	668	675	rBV5	29514	61553	2.58%	0.382%
9	7.174	691	695	699	rBV3	15720	23689	0.99%	0.147%
10	7.385	724	731	738	rVB3	23892	43008	1.80%	0.267%
11	7.479	741	747	756	rBV3	42367	86938	3.64%	0.540%
12	7.849	801	810	815	rBV	243333	430673	18.04%	2.675%
13	7.914	817	821	827	rVB3	33353	56936	2.38%	0.354%
14	8.078	845	849	853	rVV4	14358	21173	0.89%	0.132%
15	8.272	876	882	884	rBV4	10544	17701	0.74%	0.110%
16	8.390	896	902	907	rVB7	9886	19848	0.83%	0.123%
17	8.495	913	920	921	rBV4	14098	25480	1.07%	0.158%
18	8.554	925	930	937	rVB3	29296	55403	2.32%	0.344%
19	8.619	937	941	944	rBV4	7232	11829	0.50%	0.073%
20	8.725	955	959	962	rBV4	7833	10944	0.46%	0.068%
21	8.848	977	980	986	rVB4	8274	16289	0.68%	0.101%
22	8.966	992	1000	1002	rBV5	12637	24796	1.04%	0.154%
23	9.007	1002	1007	1013	rVB4	16595	34361	1.44%	0.213%
24	9.077	1013	1019	1024	rVB3	39842	70908	2.97%	0.440%
25	9.148	1028	1031	1032	rBV3	12835	12678	0.53%	0.079%
26	9.530	1094	1096	1098	rVV2	9904	10564	0.44%	0.066%
27	9.565	1098	1102	1106	rVV3	11919	20898	0.88%	0.130%
28	9.723	1121	1129	1136	rBV7	33182	66397	2.78%	0.412%
29	9.782	1136	1139	1143	rBV4	8788	12397	0.52%	0.077%
30	9.935	1161	1165	1169	rVB5	8361	13448	0.56%	0.084%
31	9.994	1172	1175	1178	rVV5	9150	10617	0.44%	0.066%
32	10.141	1192	1200	1206	rVV3	34653	72706	3.05%	0.452%
33	10.264	1216	1221	1228	rVB2	48867	77484	3.25%	0.481%
34	10.405	1243	1245	1251	rVB5	8873	14301	0.60%	0.089%
35	10.511	1255	1263	1267	rBV4	22286	46745	1.96%	0.290%
36	10.640	1276	1285	1291	rBV	360158	670957	28.10%	4.168%

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37	10.775	1300	1308	1313	rBV6	35194	88021	3.69%	0.547%
38	10.934	1330	1335	1339	rVB7	13232	24102	1.01%	0.150%
39	11.369	1407	1409	1412	rVV4	8640	11232	0.47%	0.070%
40	11.668	1456	1460	1464	rVV5	12160	22068	0.92%	0.137%
41	11.739	1468	1472	1478	rVB3	37755	58024	2.43%	0.360%
42	12.068	1520	1528	1533	rBV6	27606	58077	2.43%	0.361%
43	12.144	1538	1541	1545	rVV5	12846	16938	0.71%	0.105%
44	12.297	1564	1567	1575	rVB5	17463	31857	1.33%	0.198%
45	12.520	1601	1605	1608	rBV5	9797	13425	0.56%	0.083%
46	12.585	1613	1616	1623	rVB6	16686	27570	1.15%	0.171%
47	13.308	1735	1739	1747	rVB3	37928	65412	2.74%	0.406%
48	13.396	1747	1754	1755	rBV3	5830	12384	0.52%	0.077%
49	13.519	1770	1775	1776	rBV5	10059	12261	0.51%	0.076%
50	13.642	1794	1796	1797	rBV2	13878	10894	0.46%	0.068%
51	13.813	1819	1825	1828	rBV4	16940	35688	1.49%	0.222%
52	13.889	1833	1838	1843	rVB	95292	150659	6.31%	0.936%
53	13.971	1850	1852	1856	rVB3	13986	15411	0.65%	0.096%
54	14.112	1874	1876	1879	rVB4	9198	11125	0.47%	0.069%
55	14.177	1882	1887	1892	rBV2	95667	164295	6.88%	1.021%
56	14.389	1918	1923	1928	rVB2	152966	195040	8.17%	1.212%
57	14.483	1931	1939	1946	rBV	755124	1192652	49.96%	7.409%
58	14.571	1949	1954	1960	rBV4	44964	74824	3.13%	0.465%
59	14.629	1960	1964	1969	rVB3	55655	82794	3.47%	0.514%
60	14.694	1971	1975	1981	rBV7	26988	55059	2.31%	0.342%
61	14.788	1987	1991	1996	rBV7	19353	32277	1.35%	0.201%
62	14.888	2005	2008	2014	rVB7	19498	24008	1.01%	0.149%
63	15.011	2026	2029	2032	rVB5	10404	13652	0.57%	0.085%
64	15.047	2032	2035	2039	rBV6	11920	17386	0.73%	0.108%
65	15.111	2039	2046	2055	rBV4	106227	214554	8.99%	1.333%
66	15.246	2064	2069	2077	rBV2	126762	209524	8.78%	1.302%
67	15.340	2082	2085	2090	rVB2	60266	79310	3.32%	0.493%
68	15.481	2104	2109	2116	rVB	150346	213018	8.92%	1.323%
69	15.587	2121	2127	2135	rBV5	41854	94014	3.94%	0.584%
70	15.752	2150	2155	2159	rVB7	29257	52201	2.19%	0.324%
71	15.840	2168	2170	2177	rVB7	24144	42195	1.77%	0.262%
72	15.899	2177	2180	2181	rBV3	10928	13233	0.55%	0.082%
73	16.045	2201	2205	2206	rBV4	18365	23100	0.97%	0.144%
74	16.116	2213	2217	2222	rVB4	66427	94162	3.94%	0.585%
75	16.204	2228	2232	2236	rBV4	55102	98293	4.12%	0.611%
76	16.416	2266	2268	2271	rVB2	17175	18874	0.79%	0.117%

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77	16.551	2288	2291	2295	rVB6	33543	43112	1.81%	0.268%
78	16.886	2343	2348	2358	rBV	1208410	1990835	83.39%	12.367%
79	16.997	2364	2367	2371	rVV5	53178	60666	2.54%	0.377%
80	17.050	2373	2376	2382	rVB8	40416	62197	2.61%	0.386%
81	17.238	2401	2408	2414	rBV	1066635	1641080	68.74%	10.195%
82	17.332	2419	2424	2428	rVV	180154	272537	11.42%	1.693%
83	17.379	2428	2432	2435	rVB6	35791	48478	2.03%	0.301%
84	17.526	2453	2457	2464	rVB7	40749	87820	3.68%	0.546%
85	17.738	2489	2493	2500	rVB3	54331	79699	3.34%	0.495%
86	17.814	2502	2506	2509	rBV2	72500	89866	3.76%	0.558%
87	18.355	2593	2598	2599	rBV5	19093	26307	1.10%	0.163%
88	18.431	2607	2611	2618	rVB5	38048	54764	2.29%	0.340%
89	19.036	2710	2714	2715	rBV4	13072	17270	0.72%	0.107%
90	19.065	2716	2719	2726	rVB7	32798	54931	2.30%	0.341%
91	19.171	2734	2737	2742	rBV7	21977	37934	1.59%	0.236%
92	19.265	2751	2753	2756	rBV4	12539	13185	0.55%	0.082%
93	19.630	2810	2815	2822	rBV	247057	357895	14.99%	2.223%
94	19.941	2866	2868	2870	rBV3	11820	11114	0.47%	0.069%
95	20.176	2905	2908	2910	rBV4	21216	27397	1.15%	0.170%
96	20.587	2975	2978	2980	rBV4	14225	18127	0.76%	0.113%
97	21.163	3073	3076	3079	rVB4	46248	55254	2.31%	0.343%
98	21.486	3125	3131	3139	rBV	1329117	2163406	90.62%	13.439%
99	24.318	3608	3613	3624	rVB	137259	350502	14.68%	2.177%
100	24.530	3640	3649	3662	rVB	881608	2387324	100.00%	14.830%

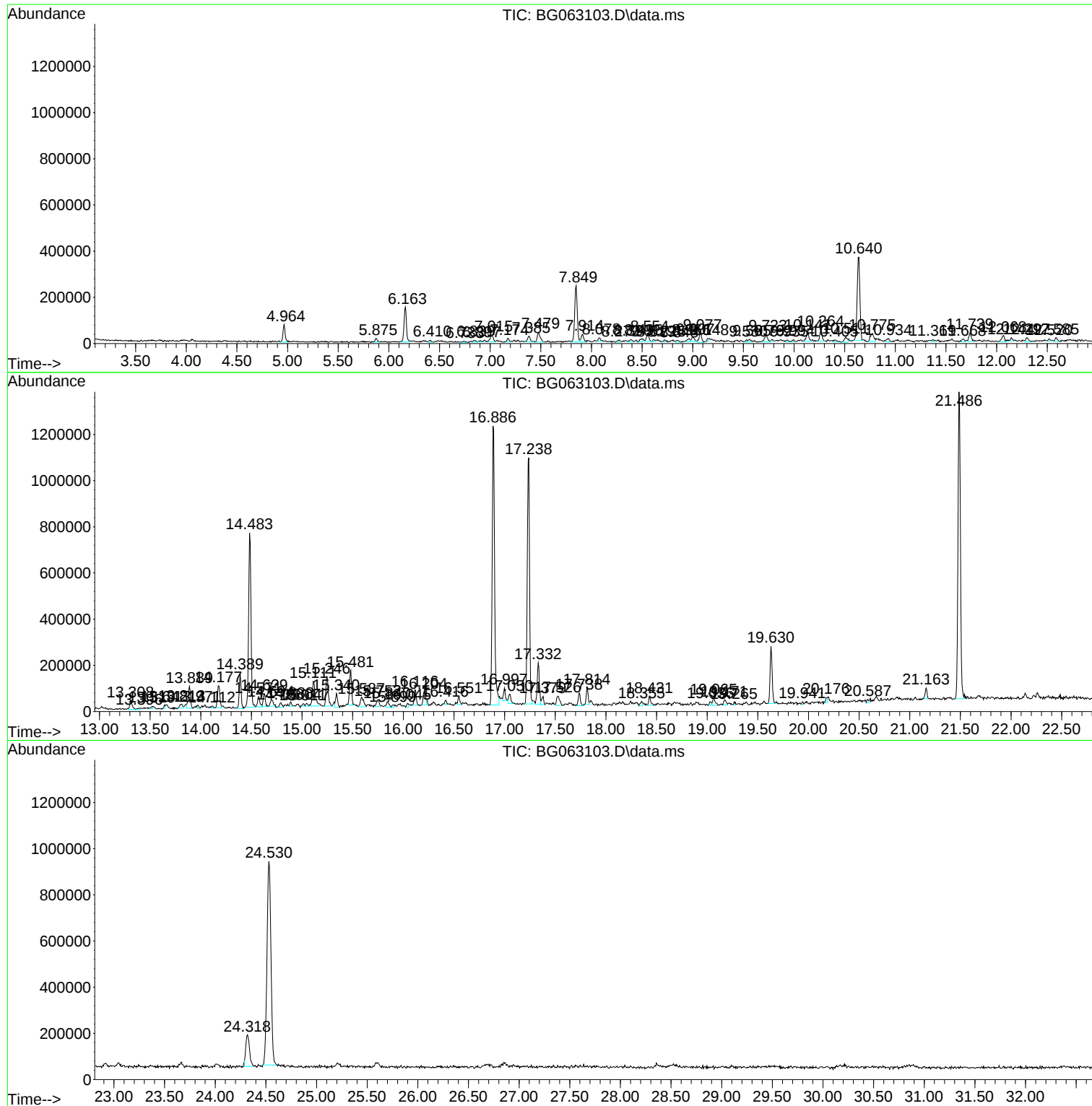
Sum of corrected areas: 16097444

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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



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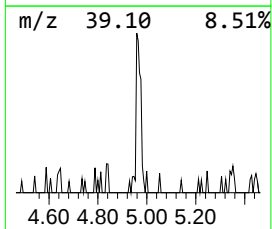
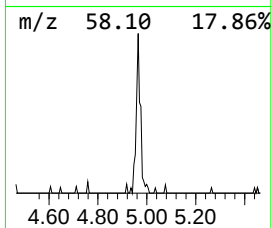
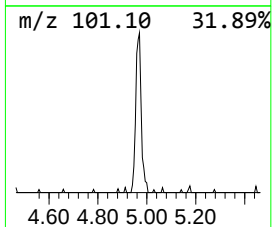
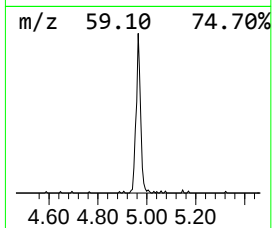
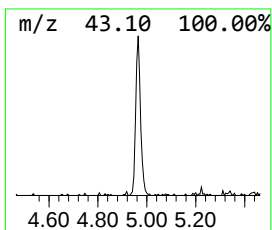
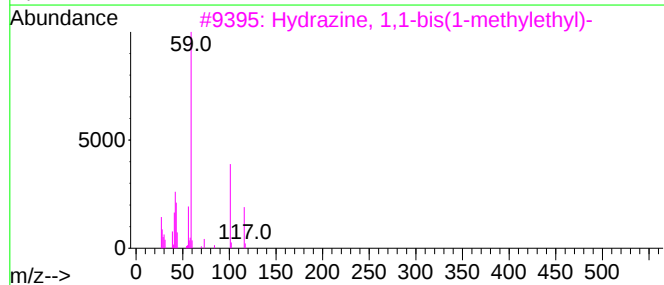
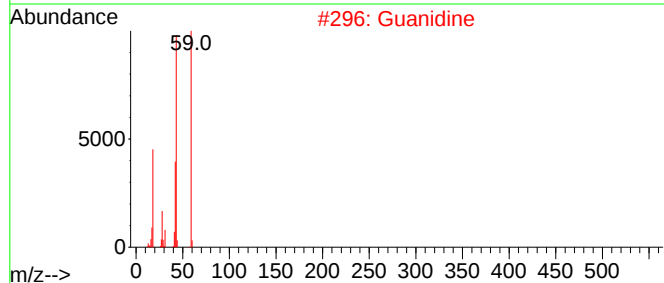
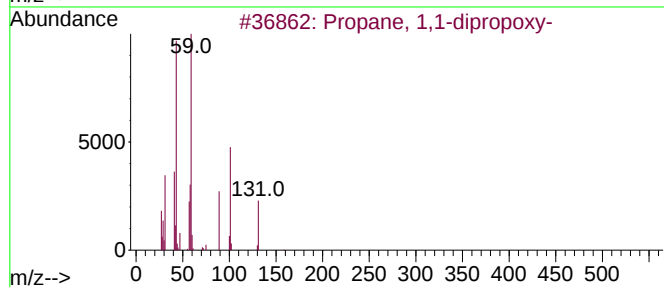
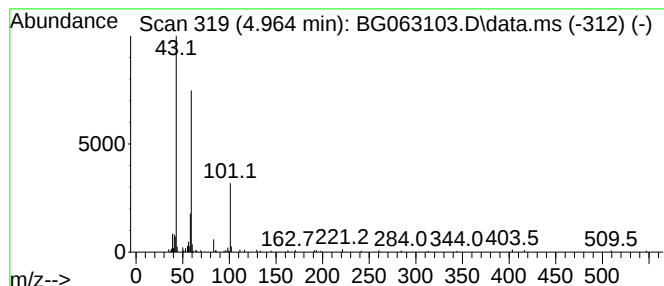
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propane, 1,1-dipropoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.964	5.13 ng/ul	110411	1,4-Dichlorobenzene-d4	7.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	72
2			Guanidine	59	CH5N3	000113-00-8	53
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	53
4			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
5			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40



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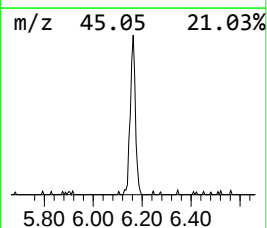
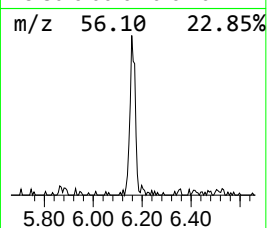
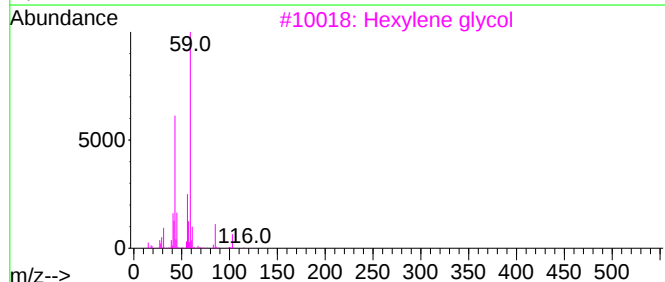
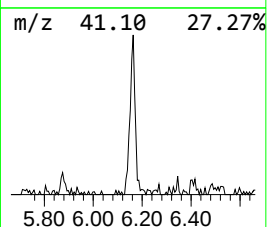
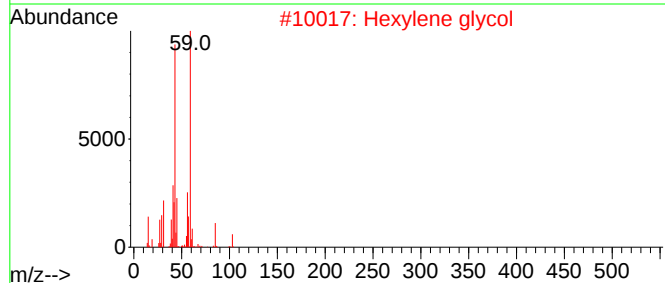
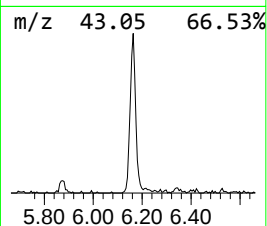
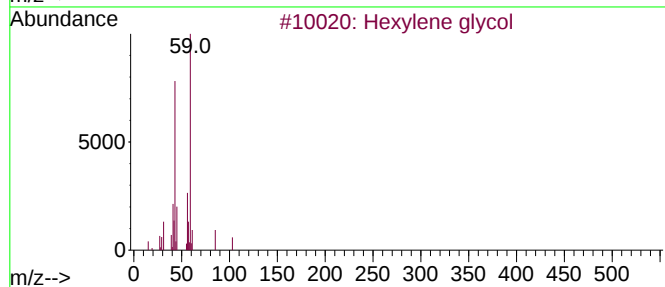
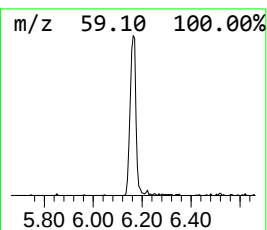
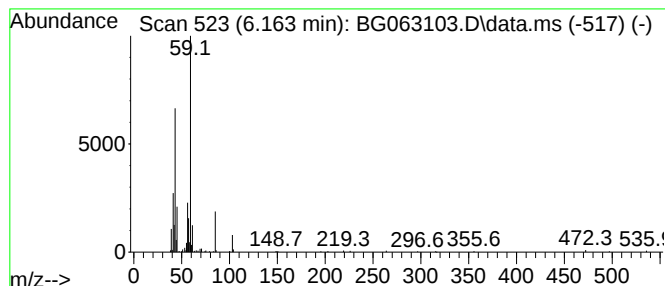
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexylene glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.163	11.68 ng/ul	251491	1,4-Dichlorobenzene-d4	7.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	86
2			Hexylene glycol	118	C6H14O2	000107-41-5	83
3			Hexylene glycol	118	C6H14O2	000107-41-5	78
4			Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	64
5			2-Butanol, 3-methoxy-	104	C5H12O2	053778-72-6	50



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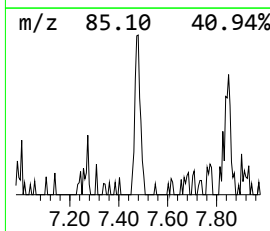
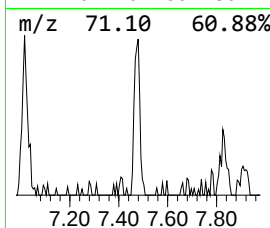
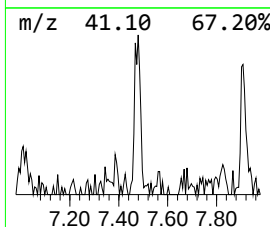
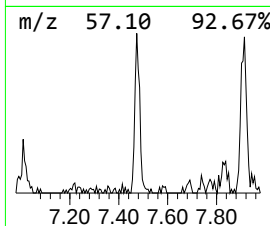
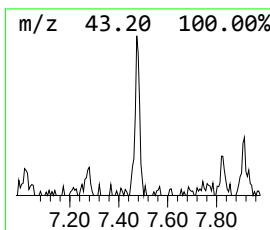
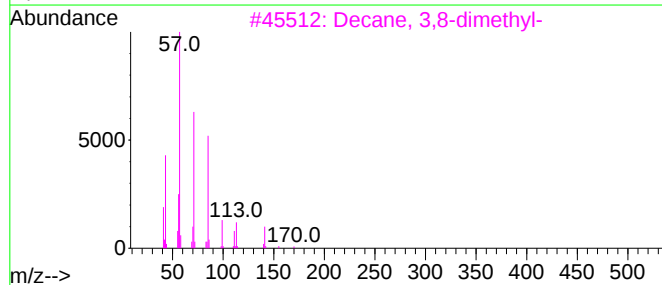
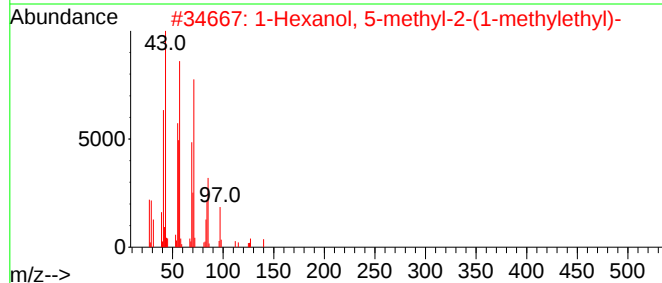
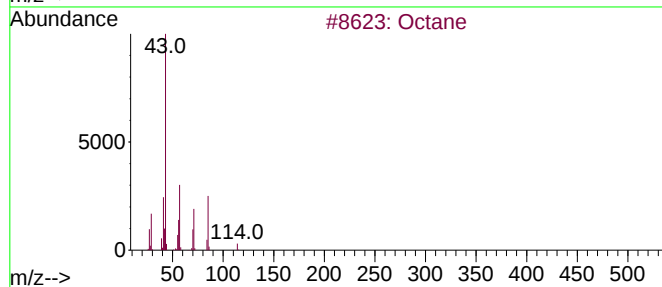
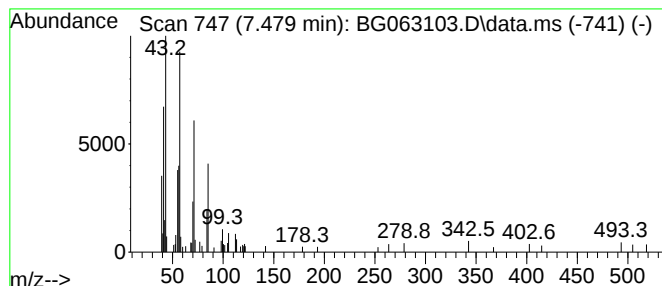
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.479	4.04 ng/ul	86938	1,4-Dichlorobenzene-d4	7.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane	114	C8H18	000111-65-9	59
2			1-Hexanol, 5-methyl-2-(1-methyle...	158	C10H22O	002051-33-4	53
3			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	47
4			L-Homoserine lactone, N,N-dimethyl-	129	C6H11NO2	1010374-45-7	47
5			Octadecane, 2-methyl-	268	C19H40	001560-88-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063103.D
Acq On : 28 Sep 2024 11:49
Operator : RC/JU
Sample : P3910-05DL 10X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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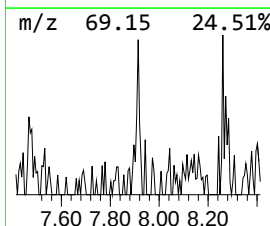
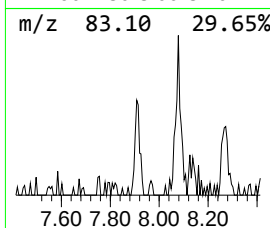
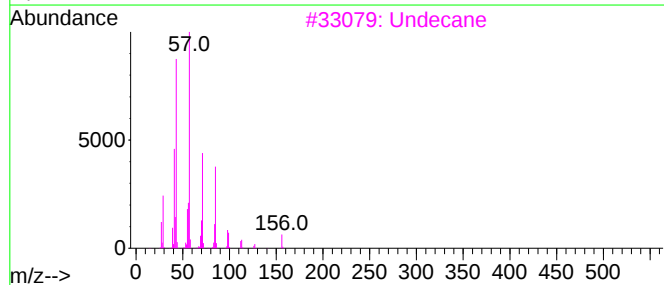
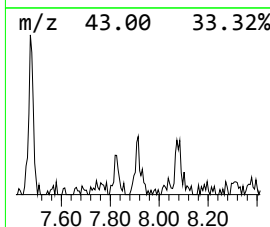
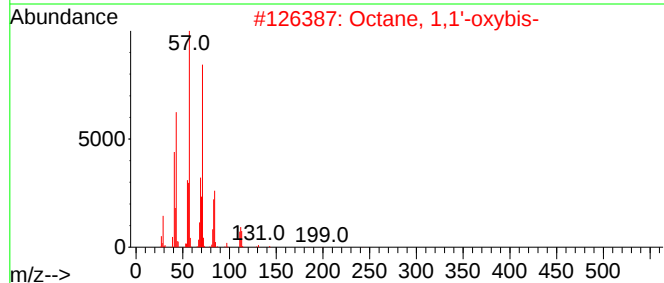
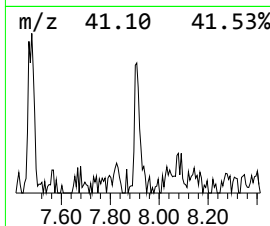
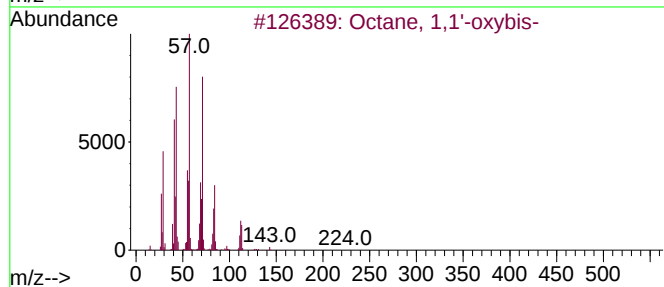
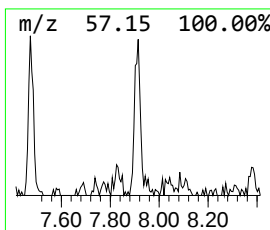
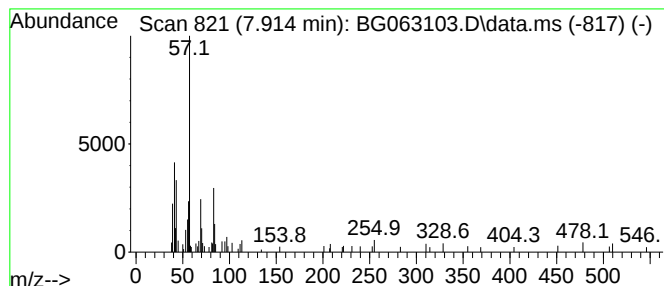
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.914	2.64 ng/ul	56936	1,4-Dichlorobenzene-d4	7.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 1,1'-oxybis-		242	C16H34O	000629-82-3	10	
2	Octane, 1,1'-oxybis-		242	C16H34O	000629-82-3	10	
3	Undecane		156	C11H24	001120-21-4	9	
4	Octane, 1,1'-oxybis-		242	C16H34O	000629-82-3	9	
5	Pentadecane		212	C15H32	000629-62-9	9	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063103.D
Acq On : 28 Sep 2024 11:49
Operator : RC/JU
Sample : P3910-05DL 10X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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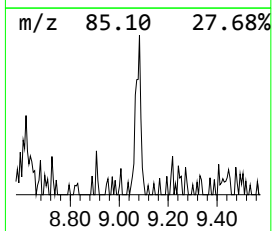
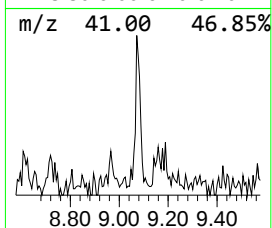
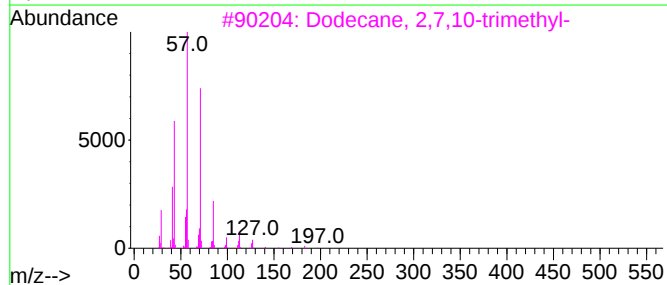
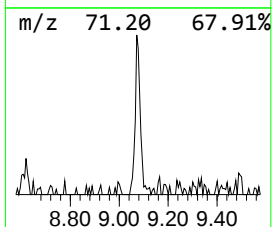
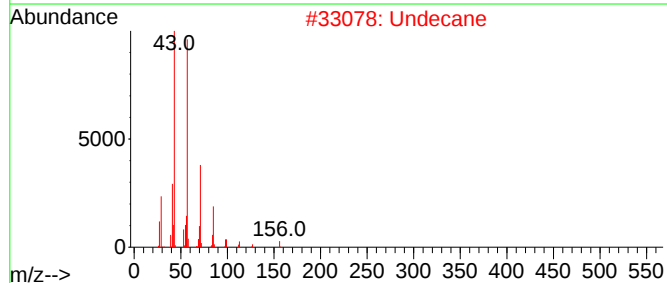
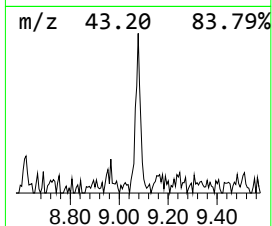
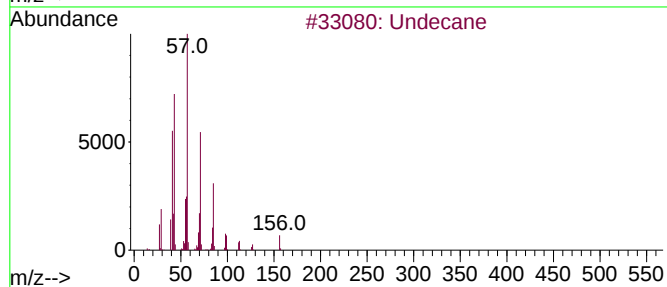
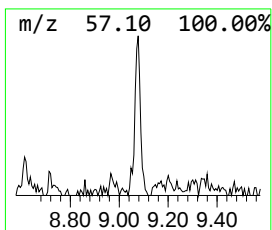
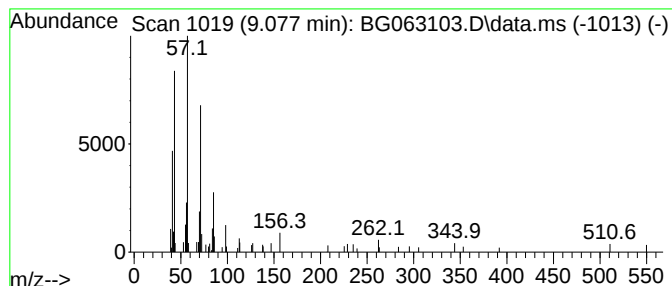
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.077	3.29 ng/ul	70908	1,4-Dichlorobenzene-d4	7.843

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	62
2	Undecane			156	C11H24	001120-21-4	59
3	Dodecane, 2,7,10-trimethyl-			212	C15H32	074645-98-0	53
4	Dodecane, 4,6-dimethyl-			198	C14H30	061141-72-8	53
5	Decane, 3,7-dimethyl-			170	C12H26	017312-54-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063103.D
Acq On : 28 Sep 2024 11:49
Operator : RC/JU
Sample : P3910-05DL 10X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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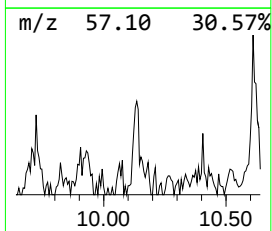
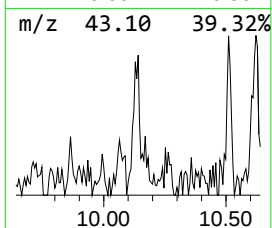
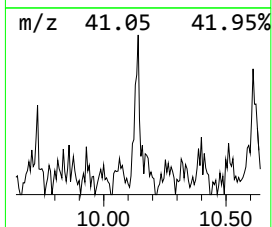
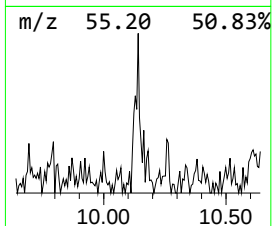
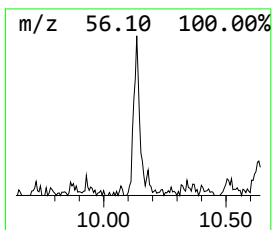
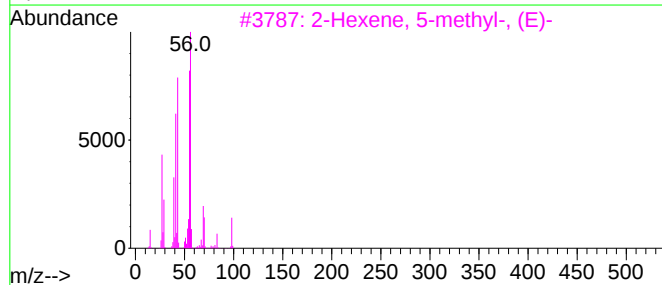
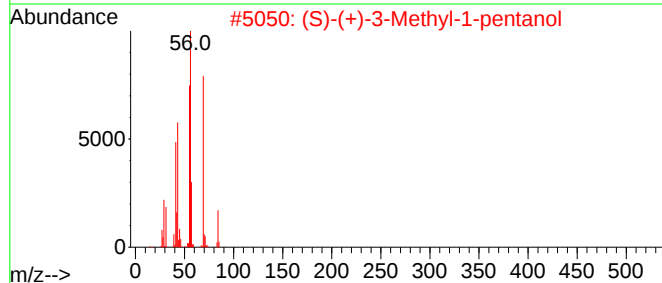
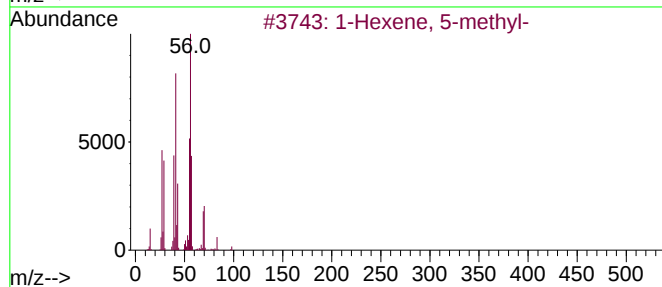
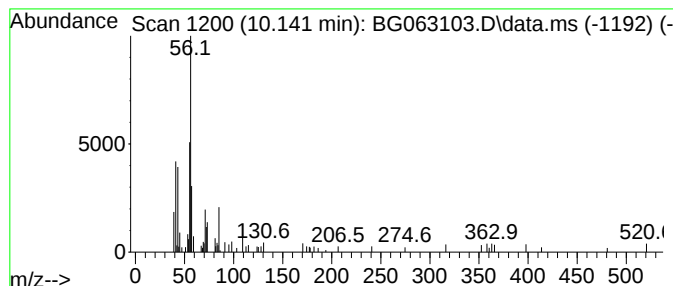
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.141	2.17 ng/ul	72706	Naphthalene-d8	10.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexene, 5-methyl-	98	C7H14	003524-73-0	47
2			(S)-(+)-3-Methyl-1-pentanol	102	C6H14O	042072-39-9	38
3			2-Hexene, 5-methyl-, (E)-	98	C7H14	007385-82-2	38
4			1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	38
5			trans-2-Aminocyclohexanol	115	C6H13NO	006982-39-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063103.D
Acq On : 28 Sep 2024 11:49
Operator : RC/JU
Sample : P3910-05DL 10X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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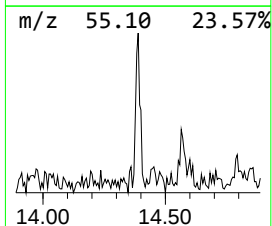
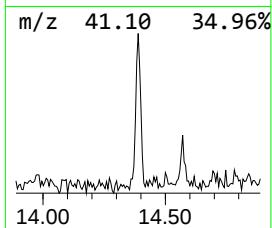
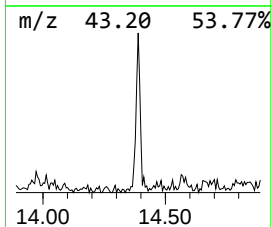
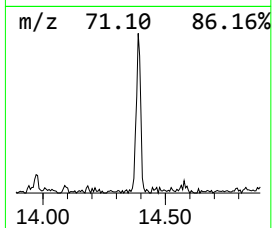
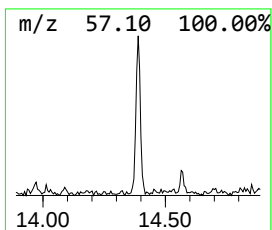
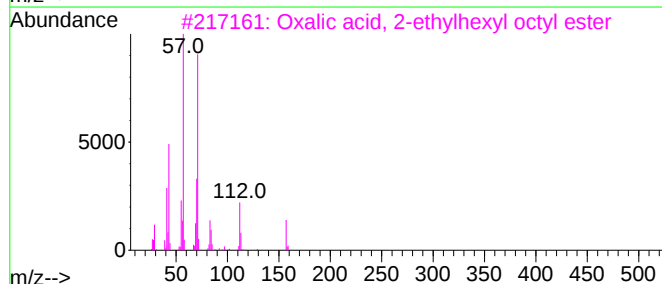
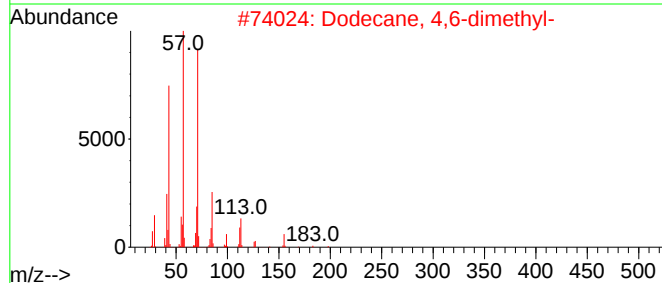
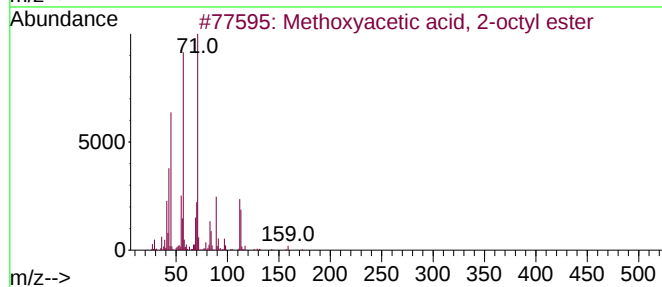
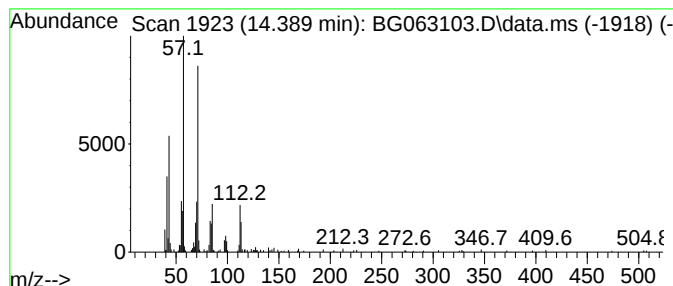
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Methoxyacetic acid, 2-octyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.389	3.27 ng/ul	195040	Acenaphthene-d10	14.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methoxyacetic acid, 2-octyl ester	202	C11H22O3	1000282-69-6	80
2			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	64
3			Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	64
4			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	64
5			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063103.D
Acq On : 28 Sep 2024 11:49
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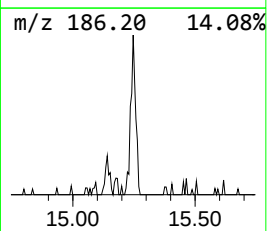
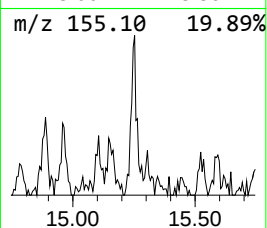
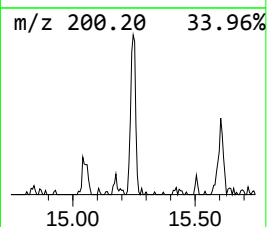
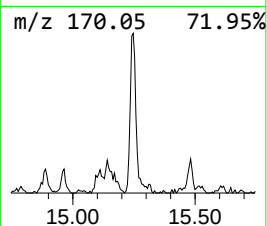
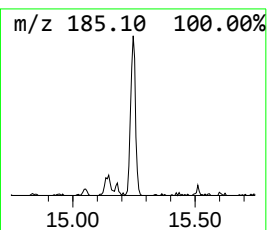
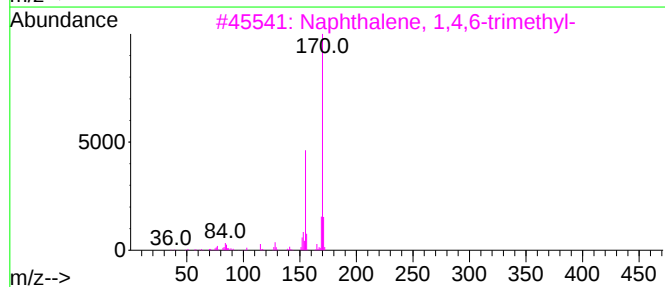
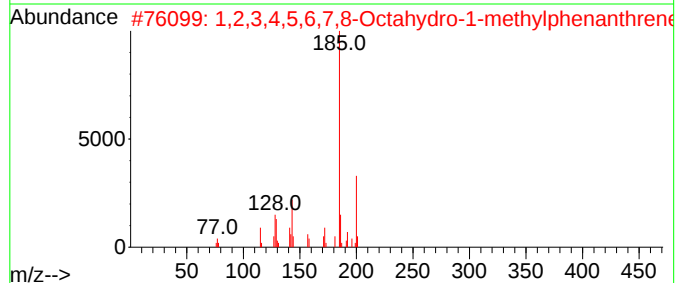
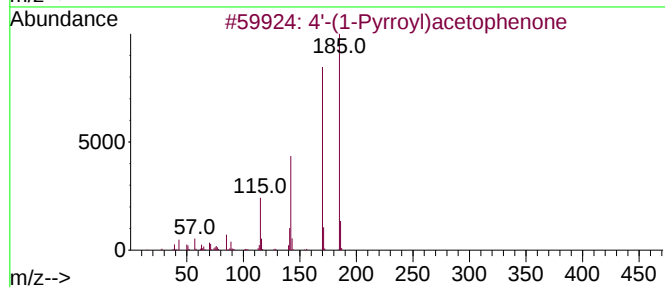
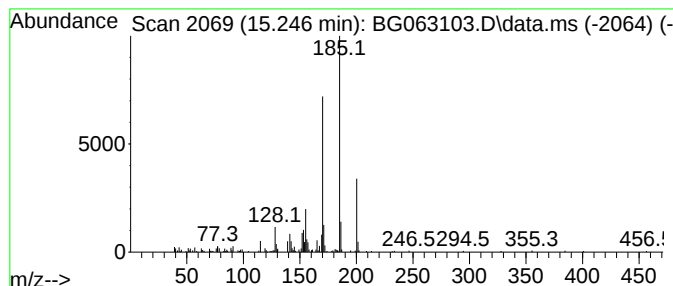
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 4'-(1-Pyrrolyl)acetophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.246	3.51 ng/ul	209524	Acenaphthene-d10	14.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	52
2			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	47
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	38
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	38
5			Indole, 3-(1,3,4-oxadiazol-2-yl)-	185	C10H7N3O	082380-80-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063103.D
Acq On : 28 Sep 2024 11:49
Operator : RC/JU
Sample : P3910-05DL 10X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.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
Propane, 1,1-di...	4.964	5.1	ng/ul	110411	1	7.843	430673	20.0
Hexylene glycol	6.163	11.7	ng/ul	251491	1	7.843	430673	20.0
(DEL) Alkane: S...	7.479	4.0	ng/ul	86938	1	7.843	430673	20.0
unknown-01	7.914	2.6	ng/ul	56936	1	7.843	430673	20.0
(DEL) Alkane: S...	9.077	3.3	ng/ul	70908	1	7.843	430673	20.0
(DEL) Alkane: S...	10.141	2.2	ng/ul	72706	2	10.640	670957	20.0
Methoxyacetic a...	14.389	3.3	ng/ul	195040	3	14.483	1192650	20.0
4'-(1-Pyrroyl)a...	15.246	3.5	ng/ul	209524	3	14.483	1192650	20.0