

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
 Data File : BG061387.D
 Acq On : 31 May 2024 10:11
 Operator : MA/JU
 Sample : P2570-22
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH654

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-BG052124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061387.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.083	9	16	34	rVB	270662	581224	55.39%	4.770%
2	3.341	55	60	67	rVB5	7472	13559	1.29%	0.111%
3	3.617	100	107	115	rBV	14121	23595	2.25%	0.194%
4	3.746	124	129	141	rBV2	71649	127323	12.13%	1.045%
5	4.075	181	185	191	rBV3	2415	4308	0.41%	0.035%
6	5.956	501	505	511	rVB2	4797	7888	0.75%	0.065%
7	7.060	685	693	700	rBV	146330	247871	23.62%	2.034%
8	7.207	710	718	726	rBV	91260	145461	13.86%	1.194%
9	7.413	747	753	760	rBV	143946	230861	22.00%	1.894%
10	7.466	760	762	767	rVV2	3510	5350	0.51%	0.044%
11	7.877	825	832	838	rBV	127128	209255	19.94%	1.717%
12	7.936	838	842	845	rVB3	4782	6855	0.65%	0.056%
13	8.594	947	954	967	rVV	161665	276237	26.32%	2.267%
14	9.029	1021	1028	1037	rVV	137603	239236	22.80%	1.963%
15	9.452	1095	1100	1106	rVV2	2933	4671	0.45%	0.038%
16	9.587	1118	1123	1128	rBV2	16928	24567	2.34%	0.202%
17	9.751	1143	1151	1162	rBV2	123520	217750	20.75%	1.787%
18	10.303	1238	1245	1254	rBV	176192	307880	29.34%	2.526%
19	10.439	1263	1268	1278	rBV2	19073	34469	3.28%	0.283%
20	10.668	1299	1307	1315	rBV	159977	280577	26.74%	2.302%
21	10.803	1323	1330	1339	rBV	140086	241771	23.04%	1.984%
22	11.197	1390	1397	1402	rBV3	4727	6227	0.59%	0.051%
23	11.408	1425	1433	1442	rBV2	15481	31050	2.96%	0.255%
24	12.254	1573	1577	1582	rBV2	2228	4027	0.38%	0.033%
25	13.329	1757	1760	1767	rVB	4818	6698	0.64%	0.055%
26	13.400	1767	1772	1779	rBV2	3325	6040	0.58%	0.050%
27	13.917	1854	1860	1866	rBV	305537	422022	40.22%	3.463%
28	14.205	1902	1909	1916	rVB	319227	497126	47.37%	4.079%
29	14.410	1940	1944	1949	rVB2	3256	4663	0.44%	0.038%
30	14.510	1953	1961	1968	rBV	253154	384443	36.63%	3.155%
31	14.657	1982	1986	1990	rBV2	2349	3278	0.31%	0.027%
32	14.734	1990	1999	2002	rBV4	163331	413560	39.41%	3.394%
33	14.763	2002	2004	2020	rVV	146124	264906	25.24%	2.174%
34	14.869	2020	2022	2027	rVV5	5081	9186	0.88%	0.075%
35	14.904	2027	2028	2031	rVV2	4394	4179	0.40%	0.034%
36	15.356	2100	2105	2109	rVB5	4159	6247	0.60%	0.051%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
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37	15.503	2122	2130	2137	rVV	417391	626622	59.71%	5.142%
38	15.638	2147	2153	2162	rBV	97603	142729	13.60%	1.171%
39	16.150	2235	2240	2245	rVB3	3016	6085	0.58%	0.050%
40	16.226	2249	2253	2261	rVV3	4444	9831	0.94%	0.081%

41	17.260	2421	2429	2433	rBV	314718	461850	44.01%	3.790%
42	17.301	2433	2436	2440	rVV	43272	60218	5.74%	0.494%
43	17.360	2440	2446	2456	rVB	454407	669619	63.81%	5.495%
44	17.660	2495	2497	2501	rVB3	3673	5214	0.50%	0.043%
45	17.765	2509	2515	2517	rBV4	2741	5033	0.48%	0.041%

46	17.924	2538	2542	2549	rVB2	6816	9500	0.91%	0.078%
47	18.159	2570	2582	2585	rBV5	29192	59893	5.71%	0.491%
48	18.224	2591	2593	2597	rVB4	4274	4308	0.41%	0.035%
49	18.329	2606	2611	2615	rBV3	13899	24423	2.33%	0.200%
50	18.365	2615	2617	2621	rVB2	7417	8319	0.79%	0.068%

51	18.453	2626	2632	2635	rBV4	4479	5374	0.51%	0.044%
52	18.494	2635	2639	2642	rBV5	3655	5811	0.55%	0.048%
53	18.635	2659	2663	2668	rVV4	6291	10400	0.99%	0.085%
54	18.682	2668	2671	2676	rVB4	11671	16237	1.55%	0.133%
55	18.882	2702	2705	2707	rVV3	5204	5947	0.57%	0.049%

56	18.935	2711	2714	2717	rVV4	3898	4501	0.43%	0.037%
57	18.970	2717	2720	2723	rVV4	4178	4903	0.47%	0.040%
58	19.099	2739	2742	2747	rVB5	24590	32795	3.13%	0.269%
59	19.146	2747	2750	2755	rVB6	3582	5516	0.53%	0.045%
60	19.205	2755	2760	2764	rBV5	10990	23631	2.25%	0.194%

61	19.316	2770	2779	2785	rBV	145431	217398	20.72%	1.784%
62	19.375	2787	2789	2792	rVB4	5461	4094	0.39%	0.034%
63	19.434	2794	2799	2804	rBV6	6427	12364	1.18%	0.101%
64	19.516	2809	2813	2817	rBV6	3551	6792	0.65%	0.056%
65	19.557	2819	2820	2823	rBV3	3991	4217	0.40%	0.035%

66	19.651	2830	2836	2849	rBV2	597143	1049413	100.00%	8.611%
67	19.757	2849	2854	2857	rBV4	8730	14233	1.36%	0.117%
68	19.857	2863	2871	2878	rBV3	9512	24151	2.30%	0.198%
69	19.916	2878	2881	2887	rBV5	7225	12397	1.18%	0.102%
70	19.998	2889	2895	2903	rBV7	13165	37486	3.57%	0.308%

71	20.139	2916	2919	2920	rBV3	4821	5183	0.49%	0.043%
72	20.274	2938	2942	2946	rVV2	12461	16566	1.58%	0.136%
73	20.327	2948	2951	2956	rVV3	13030	22458	2.14%	0.184%
74	20.468	2973	2975	2979	rBV4	10191	9003	0.86%	0.074%
75	20.515	2981	2983	2986	rVV4	8737	10502	1.00%	0.086%

76	20.568	2989	2992	2996	rVB3	18884	23665	2.26%	0.194%
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 Title : SVOA CALIBRATION

77	20.680	3006	3011	3016	rBV2	23191	33954	3.24%	0.279%
78	20.791	3028	3030	3032	rBV2	5622	4156	0.40%	0.034%
79	20.926	3050	3053	3058	rVB	19476	33862	3.23%	0.278%
80	20.967	3058	3060	3063	rBV4	7581	8986	0.86%	0.074%
81	21.020	3067	3069	3071	rVV3	6127	3484	0.33%	0.029%
82	21.103	3081	3083	3086	rVB3	10342	10869	1.04%	0.089%
83	21.173	3092	3095	3104	rVB3	48035	90317	8.61%	0.741%
84	21.402	3130	3134	3139	rVB	50062	67939	6.47%	0.558%
85	21.514	3144	3153	3157	rVV2	364846	704494	67.13%	5.781%
86	21.555	3157	3160	3166	rVB	71243	109470	10.43%	0.898%
87	21.702	3181	3185	3190	rBV3	31623	43201	4.12%	0.355%
88	21.902	3217	3219	3220	rBV2	10430	7764	0.74%	0.064%
89	22.284	3280	3284	3290	rBV6	26659	42103	4.01%	0.345%
90	22.924	3387	3393	3401	rBV3	102076	224726	21.41%	1.844%
91	23.094	3420	3422	3424	rBV3	6660	5722	0.55%	0.047%
92	23.623	3505	3512	3519	rBV	59821	185959	17.72%	1.526%
93	24.311	3623	3629	3633	rBV2	30728	74258	7.08%	0.609%
94	24.387	3634	3642	3650	rVV2	330010	792863	75.55%	6.506%
95	24.598	3668	3678	3687	rBV2	231636	666867	63.55%	5.472%
96	25.674	3857	3861	3868	rVB7	18980	44264	4.22%	0.363%
97	25.973	3910	3912	3913	rBV2	6298	4257	0.41%	0.035%
98	26.919	4071	4073	4074	rBV2	5987	3687	0.35%	0.030%
99	27.965	4250	4251	4253	rBV2	5649	3200	0.30%	0.026%
100	28.089	4264	4272	4275	rBV3	17896	46745	4.45%	0.384%

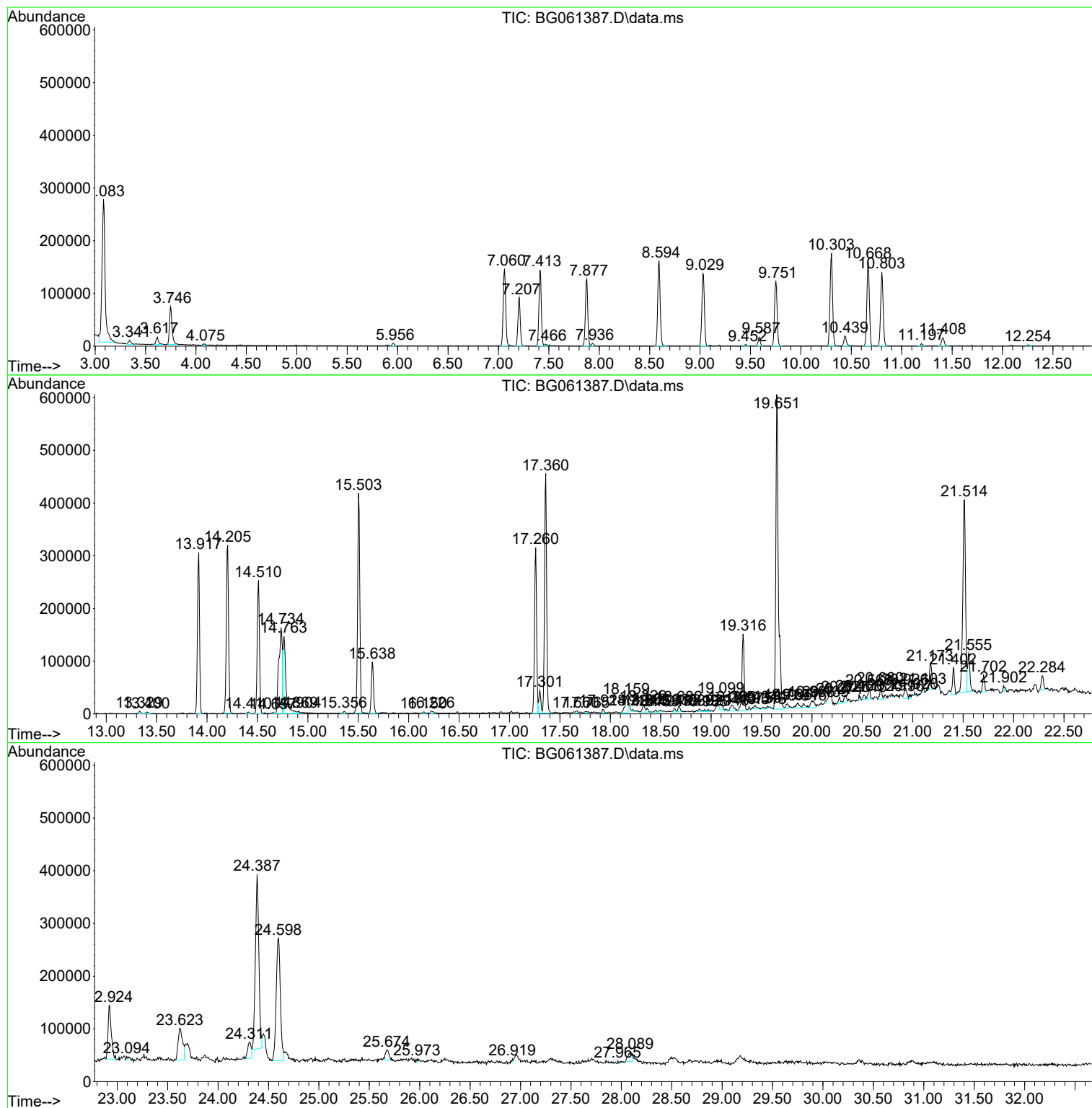
Sum of corrected areas: 12186188

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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



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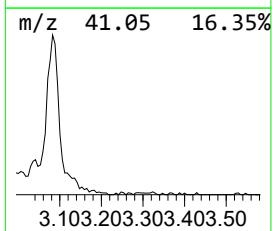
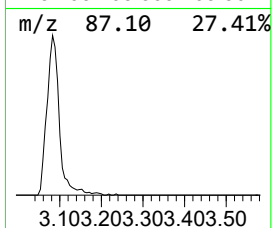
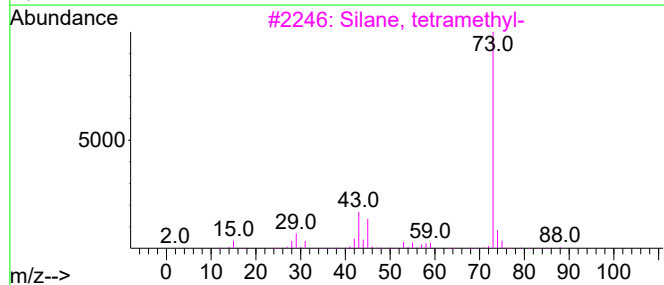
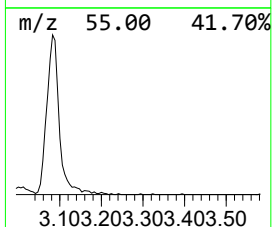
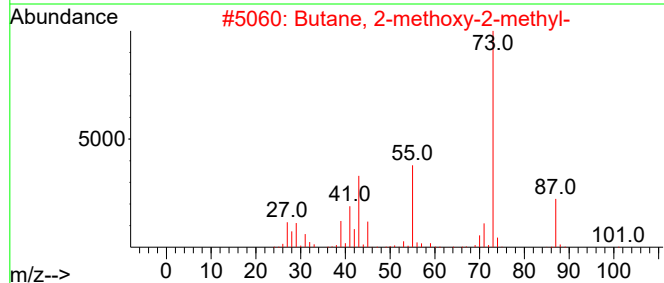
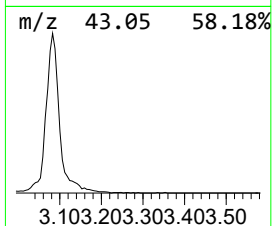
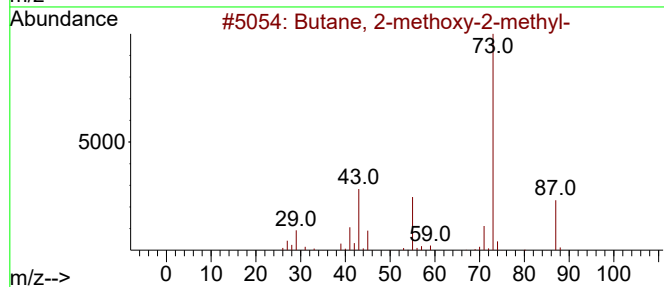
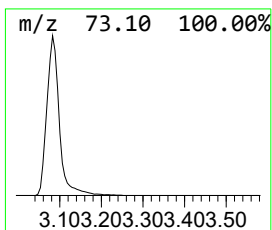
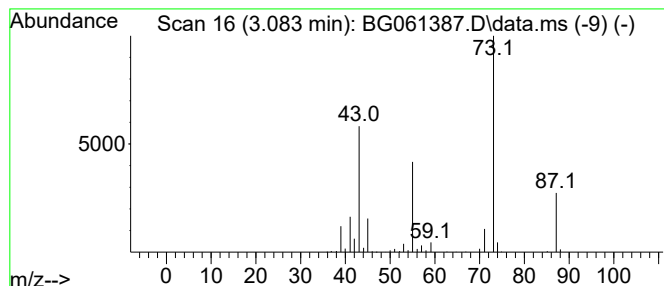
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.083	55.55 ng/ul	581224	1,4-Dichlorobenzene-d4	7.877

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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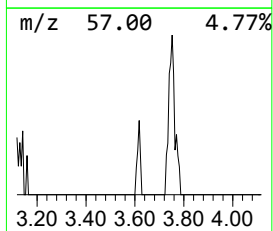
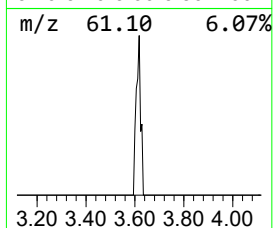
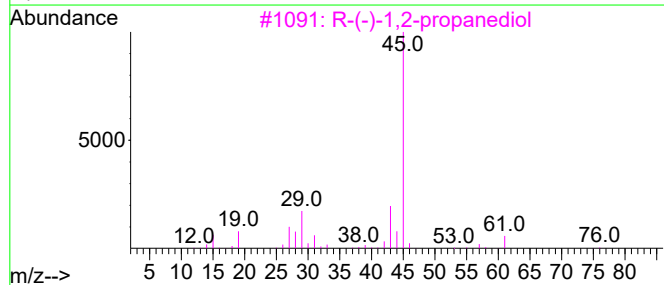
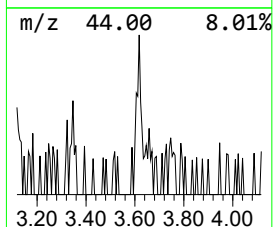
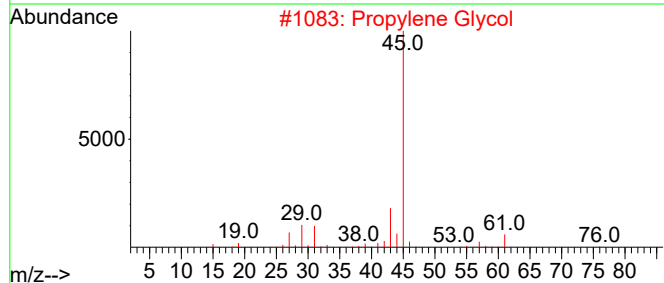
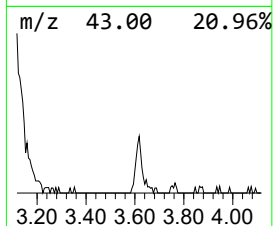
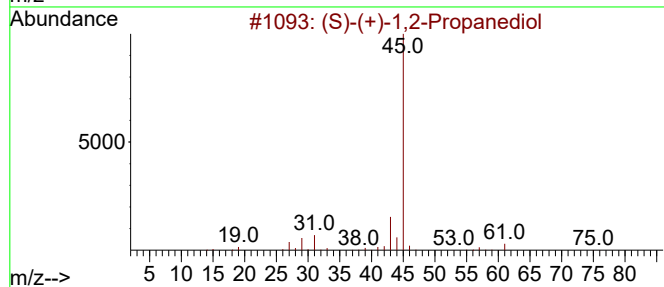
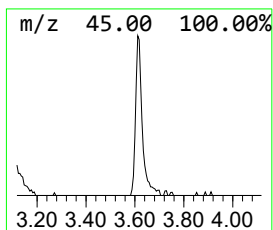
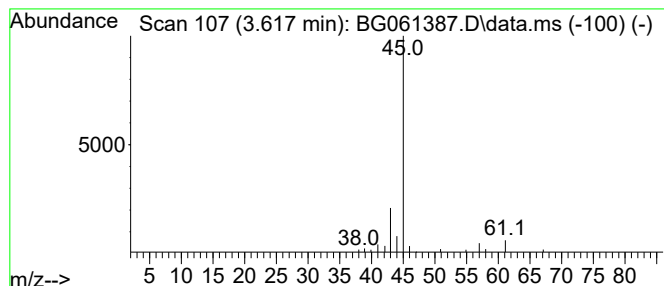
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.617	2.26 ng/ul	23595	1,4-Dichlorobenzene-d4	7.877

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	43
2	Propylene Glycol			76	C3H8O2	000057-55-6	43
3	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	43
4	Propanoic acid, 2-hydroxy-, meth...			104	C4H8O3	000547-64-8	40
5	Propylene Glycol			76	C3H8O2	000057-55-6	9



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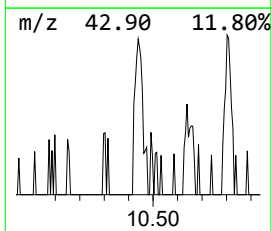
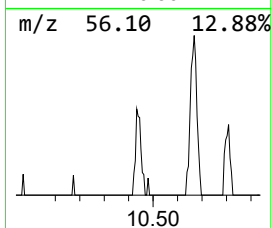
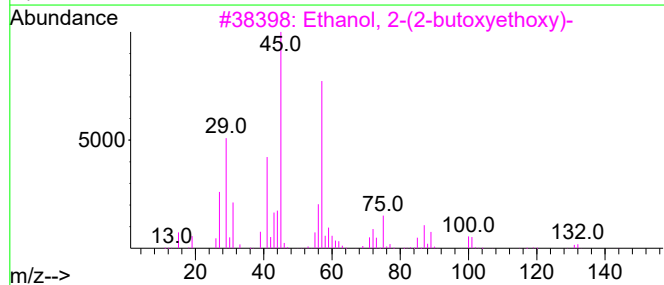
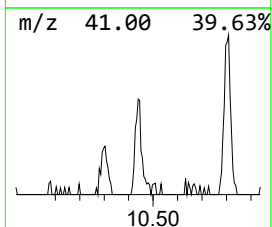
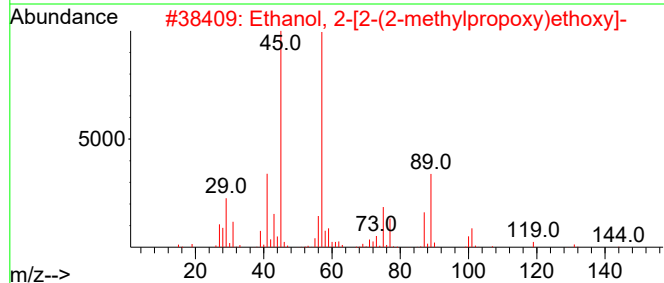
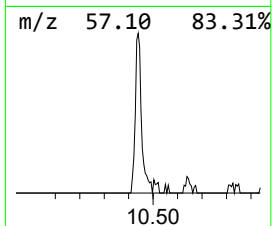
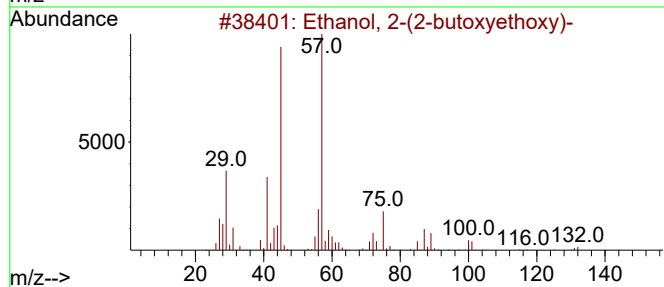
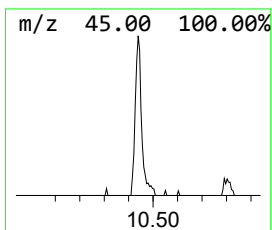
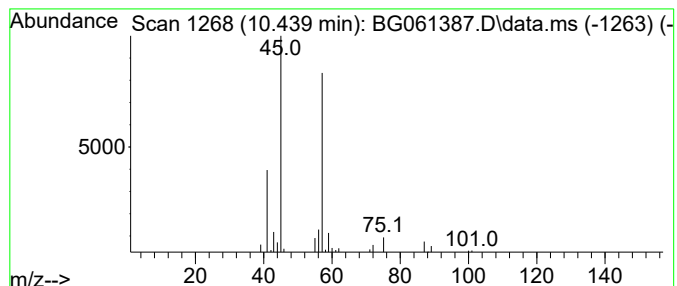
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.439	2.46 ng/ul	34469	Naphthalene-d8	10.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	74
2			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	47
4			1-Butene, 4-methoxy	86	C5H10O	004696-30-4	47
5			Di-sec-Butyl ether	130	C8H18O	006863-58-7	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061387.D
Acq On : 31 May 2024 10:11
Operator : MA/JU
Sample : P2570-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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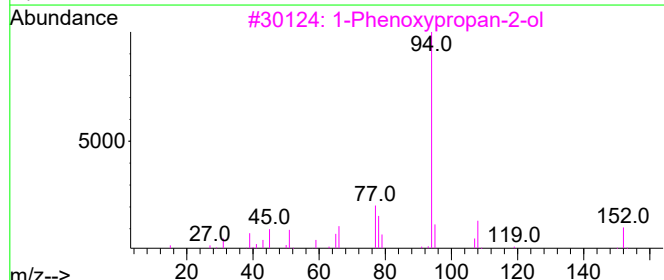
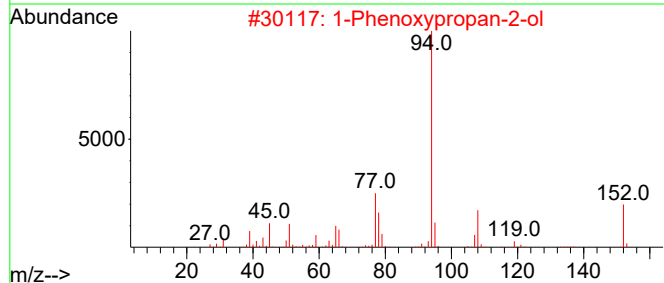
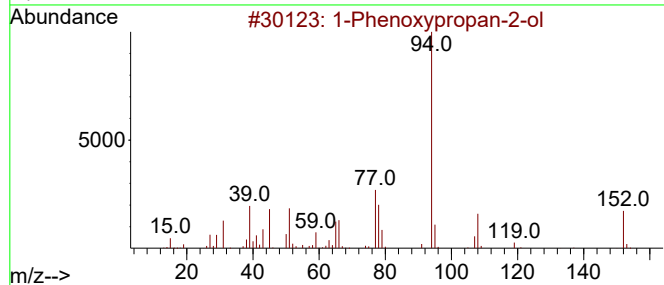
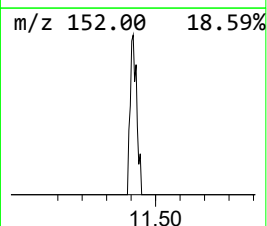
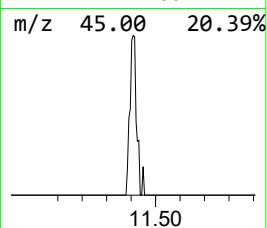
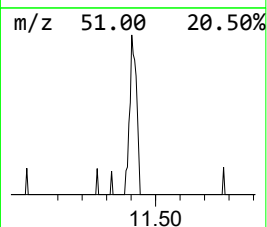
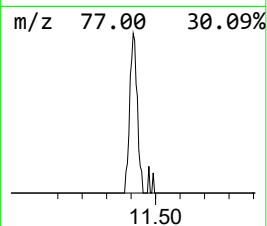
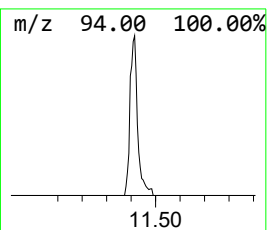
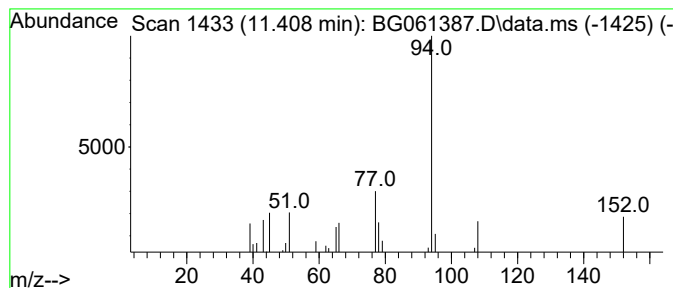
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.408	2.21 ng/ul	31050	Naphthalene-d8	10.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	81
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061387.D
Acq On : 31 May 2024 10:11
Operator : MA/JU
Sample : P2570-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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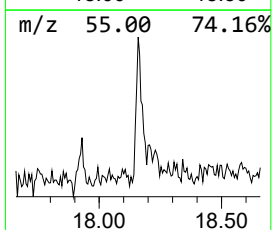
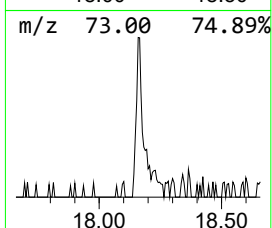
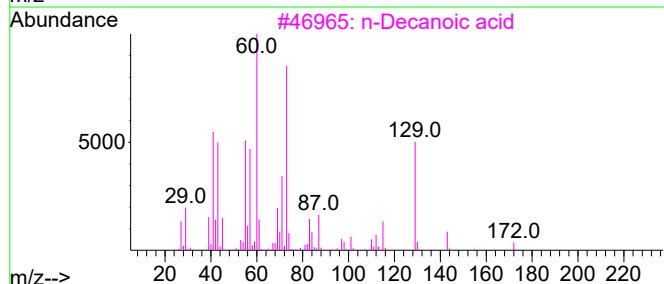
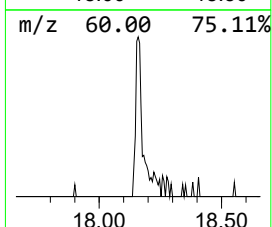
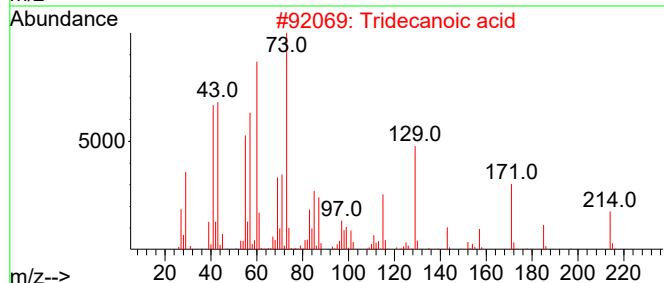
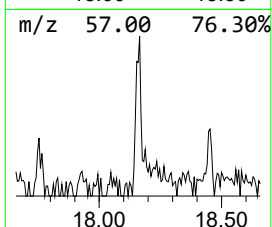
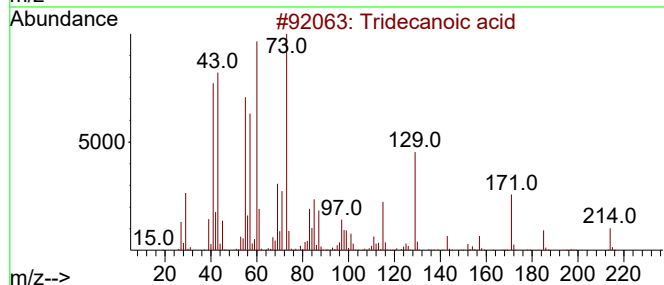
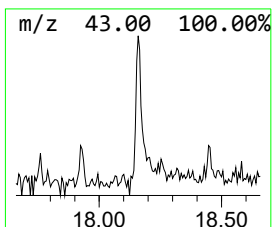
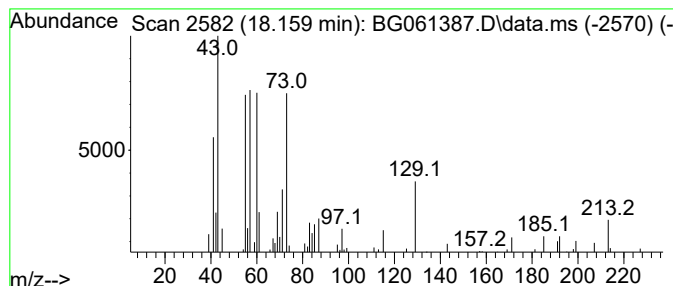
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Tridecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.159	2.59 ng/ul	59893	Phenanthrene-d10	17.260

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid	214	C13H26O2	000638-53-9	83
2			Tridecanoic acid	214	C13H26O2	000638-53-9	64
3			n-Decanoic acid	172	C10H20O2	000334-48-5	53
4			n-Decanoic acid	172	C10H20O2	000334-48-5	53
5			Tridecanoic acid	214	C13H26O2	000638-53-9	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061387.D
Acq On : 31 May 2024 10:11
Operator : MA/JU
Sample : P2570-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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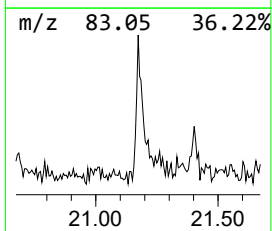
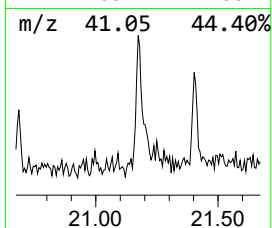
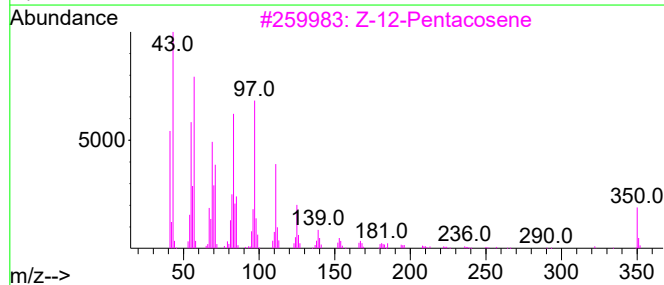
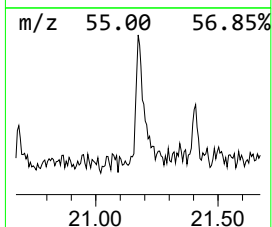
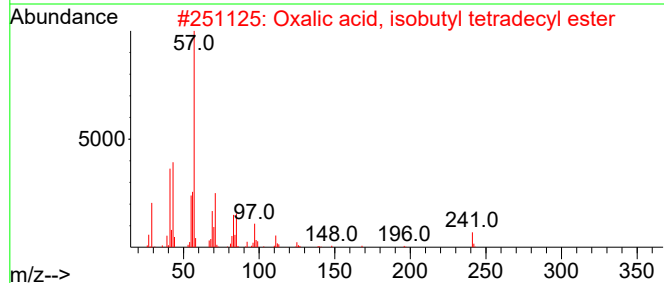
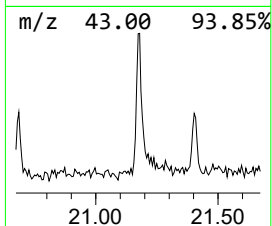
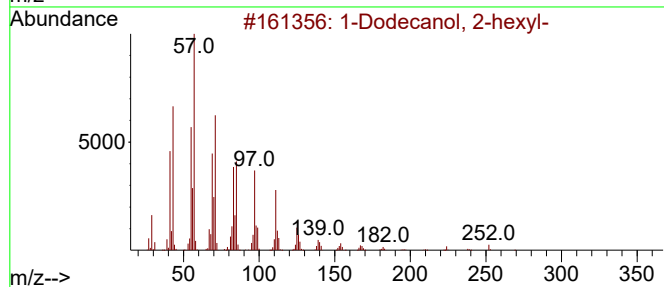
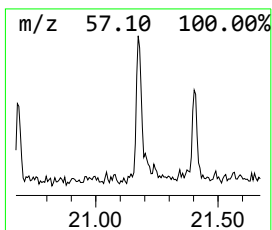
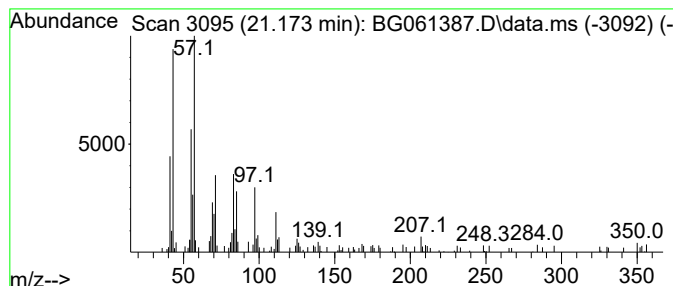
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Dodecanol, 2-hexyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.173	2.56 ng/ul	90317	Chrysene-d12	21.514

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Dodecanol, 2-hexyl-			270	C18H38O	110225-00-8	81
2	Oxalic acid, isobutyl tetradecyl...			342	C20H38O4	1000309-37-9	80
3	Z-12-Pentacosene			350	C25H50	1000131-09-4	76
4	Tetratriacontyl heptafluorobutyrate			691	C38H69F7O2	1000351-84-1	74
5	1-Dodecanol, 2-octyl-			298	C20H42O	005333-42-6	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061387.D
Acq On : 31 May 2024 10:11
Operator : MA/JU
Sample : P2570-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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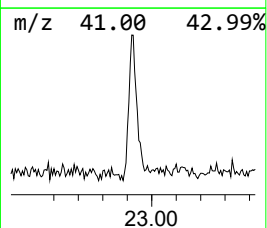
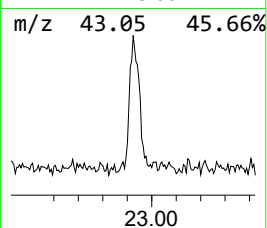
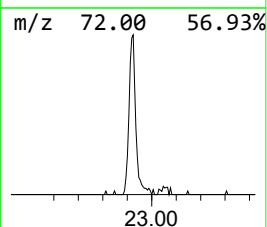
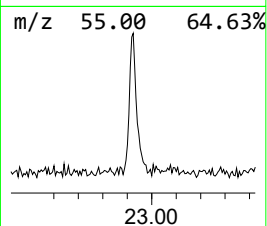
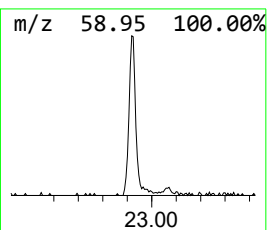
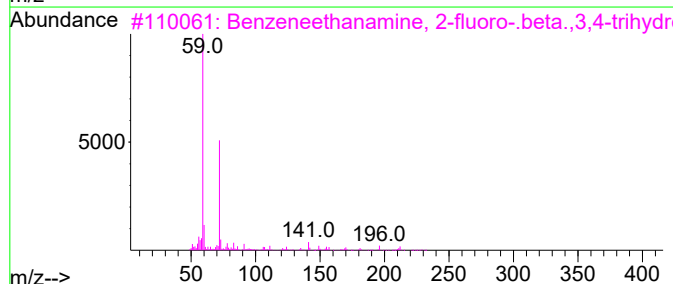
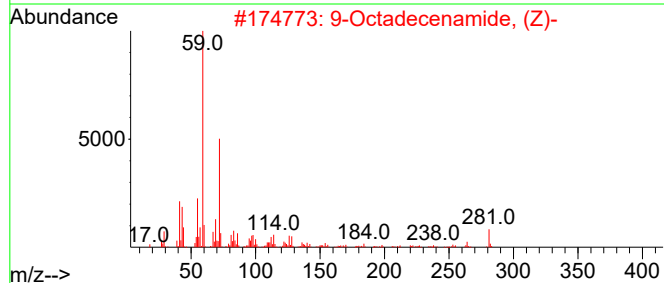
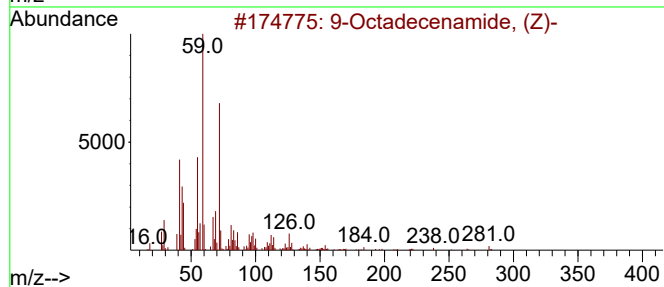
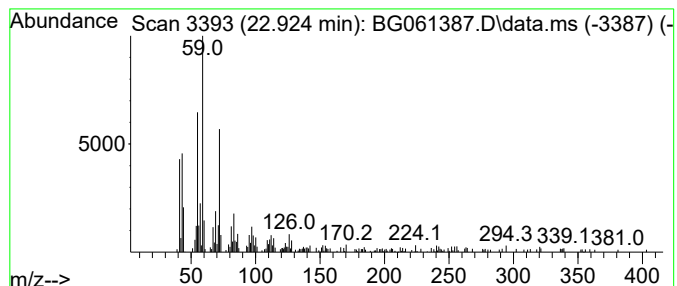
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9-Octadecenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.924	6.38 ng/ul	224726	Chrysene-d12	21.514

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	74
3			Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	72
4			1-Phenyl-9-thioureido-3,6-diazah...	302	C16H22N4S	153464-71-2	64
5			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061387.D
Acq On : 31 May 2024 10:11
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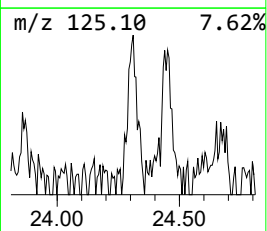
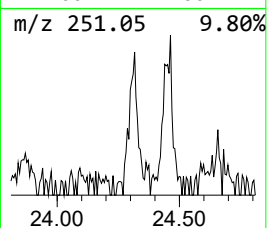
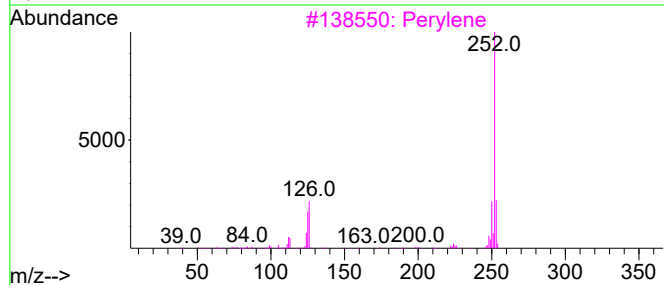
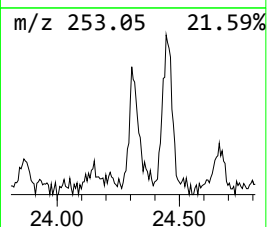
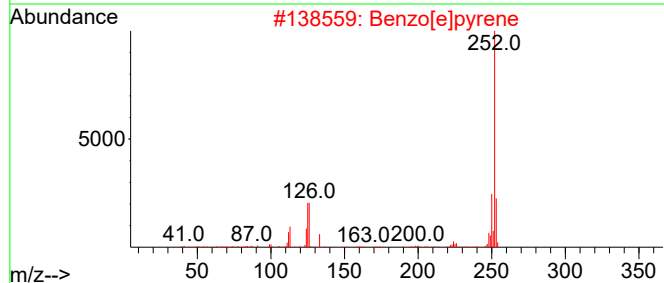
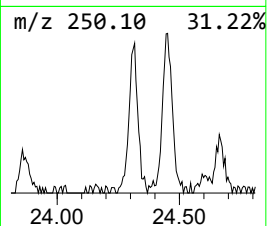
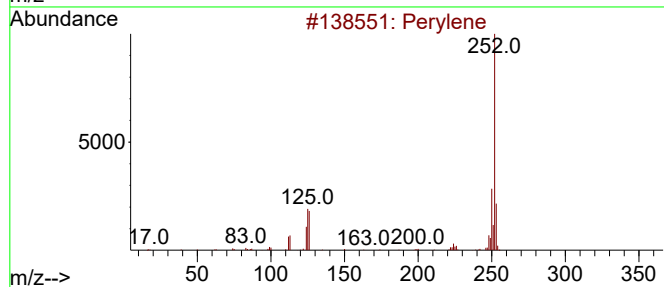
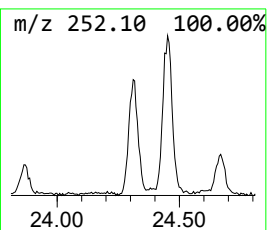
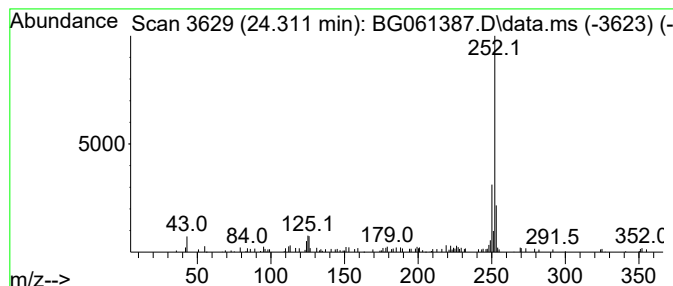
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Perylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.311	2.23 ng/ul	74258	Perylene-d12	24.599

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	94
2			Benzo[e]pyrene	252	C20H12	000192-97-2	90
3			Perylene	252	C20H12	000198-55-0	90
4			Perylene	252	C20H12	000198-55-0	90
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061387.D
Acq On : 31 May 2024 10:11
Operator : MA/JU
Sample : P2570-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown-01	3.617	2.3	ng/ul	23595	1	7.877	209255	20.0
Ethanol, 2-(2-b...	10.439	2.5	ng/ul	34469	2	10.668	280577	20.0
1-Phenoxypropan...	11.408	2.2	ng/ul	31050	2	10.668	280577	20.0
Tridecanoic acid	18.159	2.6	ng/ul	59893	4	17.260	461850	20.0
1-Dodecanol, 2-...	21.173	2.6	ng/ul	90317	5	21.514	704494	20.0
9-Octadecenamid...	22.924	6.4	ng/ul	224726	5	21.514	704494	20.0
Perylene	24.311	2.2	ng/ul	74258	6	24.599	666867	20.0