

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
 Data File : BG061645.D
 Acq On : 22 Jun 2024 12:20
 Operator : MA/JU
 Sample : P2775-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4C95

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-BG062024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061645.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.472	26	31	35	rBV3	15729	21841	0.92%	0.076%
2	3.736	74	76	80	rBV3	3824	6747	0.28%	0.024%
3	3.895	100	103	108	rBV	27039	37226	1.57%	0.130%
4	4.000	118	121	125	rBV4	10812	16037	0.68%	0.056%
5	4.188	149	153	156	rVV4	6534	8913	0.38%	0.031%
6	4.229	156	160	166	rVB6	6764	13800	0.58%	0.048%
7	4.447	193	197	201	rVB5	8118	13184	0.56%	0.046%
8	4.741	244	247	250	rBV4	7082	9554	0.40%	0.033%
9	4.794	252	256	262	rVB2	17348	25133	1.06%	0.088%
10	5.152	310	317	323	rVB	50344	75264	3.17%	0.263%
11	5.457	364	369	371	rBV3	4706	6588	0.28%	0.023%
12	5.722	409	414	415	rBV4	4627	6558	0.28%	0.023%
13	6.045	464	469	471	rBV5	3783	5831	0.25%	0.020%
14	6.133	481	484	489	rBV4	14744	22675	0.96%	0.079%
15	6.750	584	589	594	rVB6	8141	12879	0.54%	0.045%
16	7.026	629	636	641	rVV4	9204	14030	0.59%	0.049%
17	7.214	660	668	677	rBV	340297	587451	24.75%	2.056%
18	7.402	691	700	710	rVB	224437	371545	15.65%	1.301%
19	7.514	717	719	724	rVB4	5107	6318	0.27%	0.022%
20	7.602	727	734	741	rBV2	306834	521847	21.98%	1.827%
21	7.667	741	745	749	rBV2	9398	16140	0.68%	0.056%
22	7.761	758	761	764	rBV3	5254	6897	0.29%	0.024%
23	8.072	808	814	821	rBV	210309	353939	14.91%	1.239%
24	8.137	821	825	829	rVB4	8717	14902	0.63%	0.052%
25	8.313	850	855	858	rBV4	5908	10000	0.42%	0.035%
26	8.613	901	906	907	rBV3	5167	6420	0.27%	0.022%
27	8.765	924	932	940	rBV	382166	662653	27.92%	2.320%
28	8.900	947	955	964	rVB2	24801	49940	2.10%	0.175%
29	9.235	1004	1012	1023	rBV	299493	531060	22.37%	1.859%
30	9.318	1023	1026	1030	rBV4	6219	7844	0.33%	0.027%
31	9.394	1035	1039	1046	rVB6	5966	10443	0.44%	0.037%
32	9.653	1078	1083	1086	rVB5	5251	7191	0.30%	0.025%
33	9.799	1101	1108	1117	rVB2	36881	62718	2.64%	0.220%
34	9.964	1127	1136	1142	rBV2	246968	460140	19.38%	1.611%
35	10.070	1150	1154	1158	rBV4	15514	20553	0.87%	0.072%
36	10.158	1165	1169	1179	rVB6	10330	19733	0.83%	0.069%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
 Title : SVOA CALIBRATION

37	10.493	1218	1226	1234	rBV	410771	731361	30.81%	2.560%
38	10.646	1249	1252	1254	rBV2	4994	5466	0.23%	0.019%
39	10.728	1264	1266	1270	rVB5	4628	5246	0.22%	0.018%
40	10.886	1284	1293	1299	rBV	323899	577677	24.34%	2.022%

41	11.016	1307	1315	1324	rVB2	233593	440789	18.57%	1.543%
42	11.409	1379	1382	1389	rVB5	16760	29702	1.25%	0.104%
43	11.468	1389	1392	1395	rBV3	5350	5955	0.25%	0.021%
44	11.627	1412	1419	1429	rVB2	56082	124116	5.23%	0.434%
45	12.291	1528	1532	1535	rBV5	5684	7485	0.32%	0.026%

46	12.461	1557	1561	1568	rVB5	9160	14935	0.63%	0.052%
47	12.537	1568	1574	1578	rBV4	10699	16855	0.71%	0.059%
48	12.755	1607	1611	1617	rVB3	8261	12371	0.52%	0.043%
49	13.066	1661	1664	1668	rVV5	6170	8264	0.35%	0.029%
50	13.607	1755	1756	1762	rVB5	7411	9028	0.38%	0.032%

51	13.665	1762	1766	1773	rVB5	14288	23568	0.99%	0.082%
52	13.736	1773	1778	1781	rBV6	4851	8951	0.38%	0.031%
53	13.859	1793	1799	1800	rBV4	7348	8534	0.36%	0.030%
54	13.883	1801	1803	1807	rVB3	9188	8196	0.35%	0.029%
55	14.024	1823	1827	1833	rVB4	16986	28666	1.21%	0.100%

56	14.106	1833	1841	1847	rBV	713318	1053260	44.37%	3.687%
57	14.259	1864	1867	1872	rVB6	6660	10584	0.45%	0.037%
58	14.341	1877	1881	1883	rBV4	13765	19230	0.81%	0.067%
59	14.394	1883	1890	1900	rVB2	777098	1285317	54.15%	4.499%
60	14.600	1921	1925	1928	rBV6	5129	6848	0.29%	0.024%

61	14.653	1931	1934	1935	rBV2	9439	7831	0.33%	0.027%
62	14.700	1935	1942	1948	rVB	534543	820443	34.56%	2.872%
63	14.764	1948	1953	1959	rVB	74153	126554	5.33%	0.443%
64	14.870	1966	1971	1978	rBV	374960	589037	24.81%	2.062%
65	14.976	1986	1989	1993	rVB4	45055	57386	2.42%	0.201%

66	15.099	2005	2010	2016	rVB2	48522	80061	3.37%	0.280%
67	15.375	2053	2057	2061	rVB6	8921	13502	0.57%	0.047%
68	15.693	2104	2111	2116	rBV	1045136	1576758	66.42%	5.519%
69	15.745	2116	2120	2123	rVV3	63992	92487	3.90%	0.324%
70	15.792	2123	2128	2134	rVB	288569	423893	17.86%	1.484%

71	16.045	2168	2171	2177	rVB5	23757	39041	1.64%	0.137%
72	16.116	2180	2183	2184	rBV2	8694	7156	0.30%	0.025%
73	16.133	2184	2186	2189	rBV3	12528	10588	0.45%	0.037%
74	16.186	2192	2195	2200	rVB6	16265	26640	1.12%	0.093%
75	16.315	2213	2217	2225	rVB4	40470	62690	2.64%	0.219%

76	16.380	2225	2228	2230	rBV3	13556	18668	0.79%	0.065%
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 Title : SVOA CALIBRATION

77	16.539	2253	2255	2257	rBV3	9241	8335	0.35%	0.029%
78	17.091	2346	2349	2355	rVB	41049	60515	2.55%	0.212%
79	17.167	2358	2362	2363	rBV4	15697	19456	0.82%	0.068%
80	17.261	2374	2378	2384	rVB	48114	74662	3.15%	0.261%
81	17.320	2386	2388	2391	rBV4	15008	19423	0.82%	0.068%
82	17.443	2403	2409	2413	rBV	652582	1043761	43.97%	3.654%
83	17.485	2413	2416	2421	rVB	814022	1154351	48.63%	4.041%
84	17.543	2421	2426	2431	rBV	1073447	1608312	67.75%	5.630%
85	17.843	2474	2477	2481	rVB	75089	96565	4.07%	0.338%
86	18.025	2505	2508	2518	rVB	27935	66151	2.79%	0.232%
87	18.272	2547	2550	2554	rBV4	43843	66749	2.81%	0.234%
88	18.331	2554	2560	2564	rBV3	538388	807573	34.02%	2.827%
89	18.513	2586	2591	2595	rBV3	157650	278875	11.75%	0.976%
90	18.754	2629	2632	2636	rVB	124701	138363	5.83%	0.484%
91	18.848	2645	2648	2654	rVB	131973	196074	8.26%	0.686%
92	18.924	2658	2661	2666	rVB4	115658	144665	6.09%	0.506%
93	19.494	2752	2758	2764	rVB	1621311	2339364	98.55%	8.189%
94	19.600	2773	2776	2786	rVB6	230869	494916	20.85%	1.732%
95	19.823	2806	2814	2817	rBV2	1293658	2373748	100.00%	8.309%
96	19.852	2817	2819	2827	rVB	1063779	1278148	53.85%	4.474%
97	20.440	2916	2919	2922	rVB	132966	141377	5.96%	0.495%
98	21.362	3073	3076	3083	rVB4	403429	647960	27.30%	2.268%
99	21.697	3127	3133	3139	rBV4	830315	2146368	90.42%	7.513%
100	21.756	3140	3143	3152	rVB	616727	942994	39.73%	3.301%

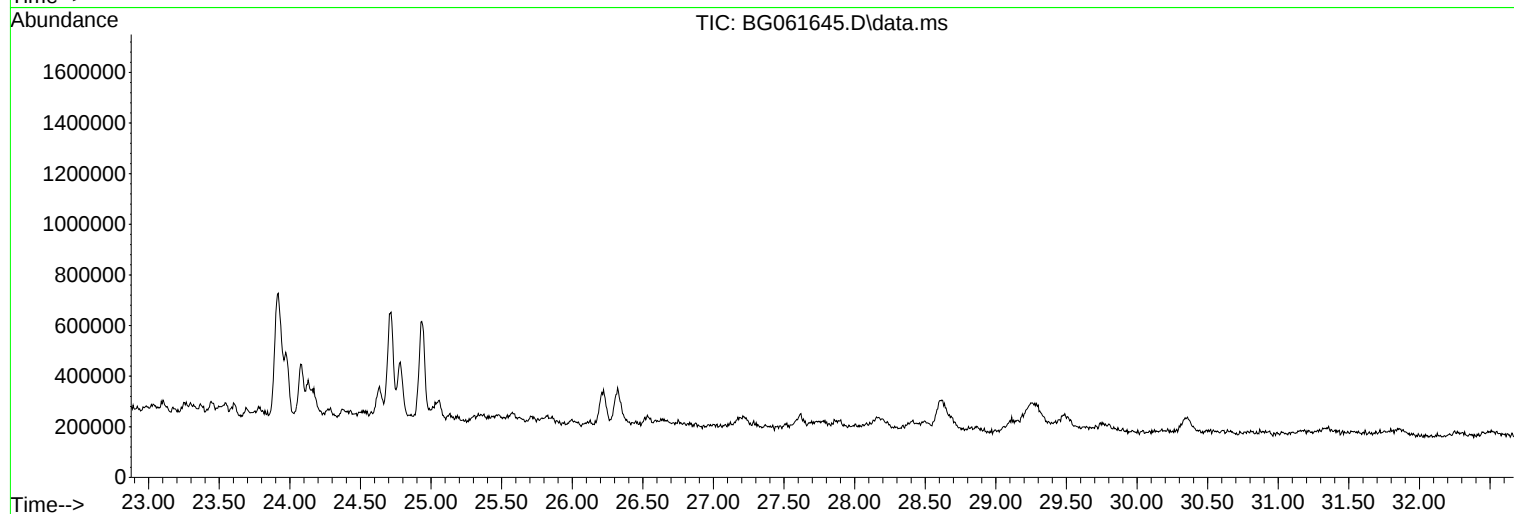
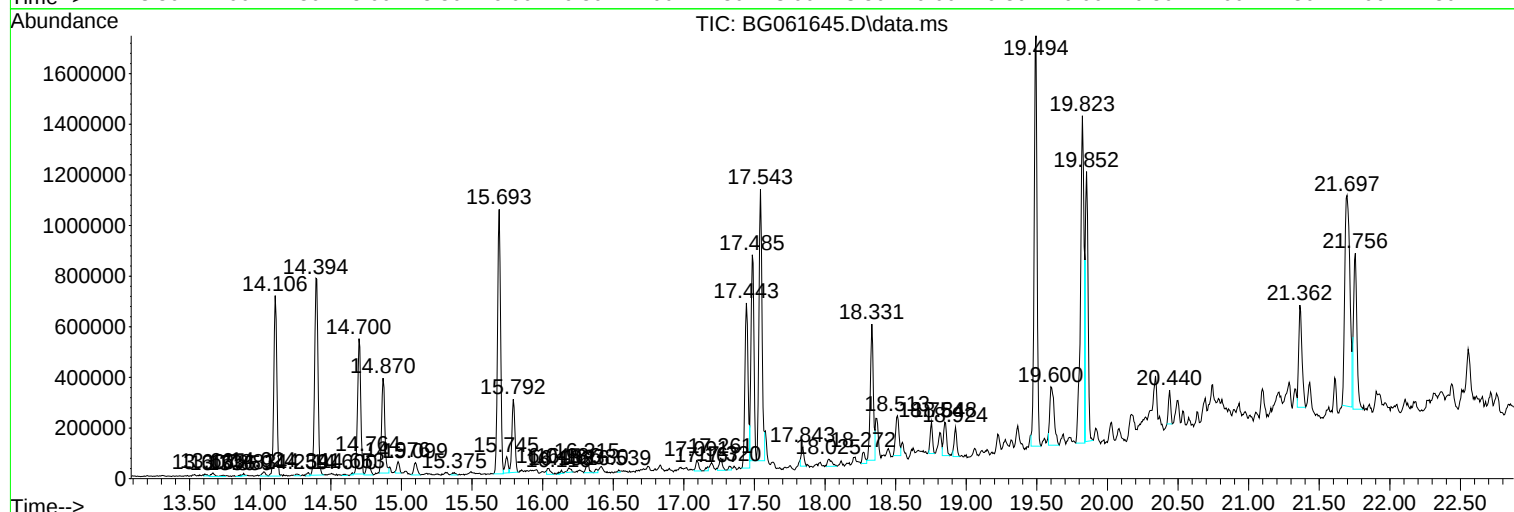
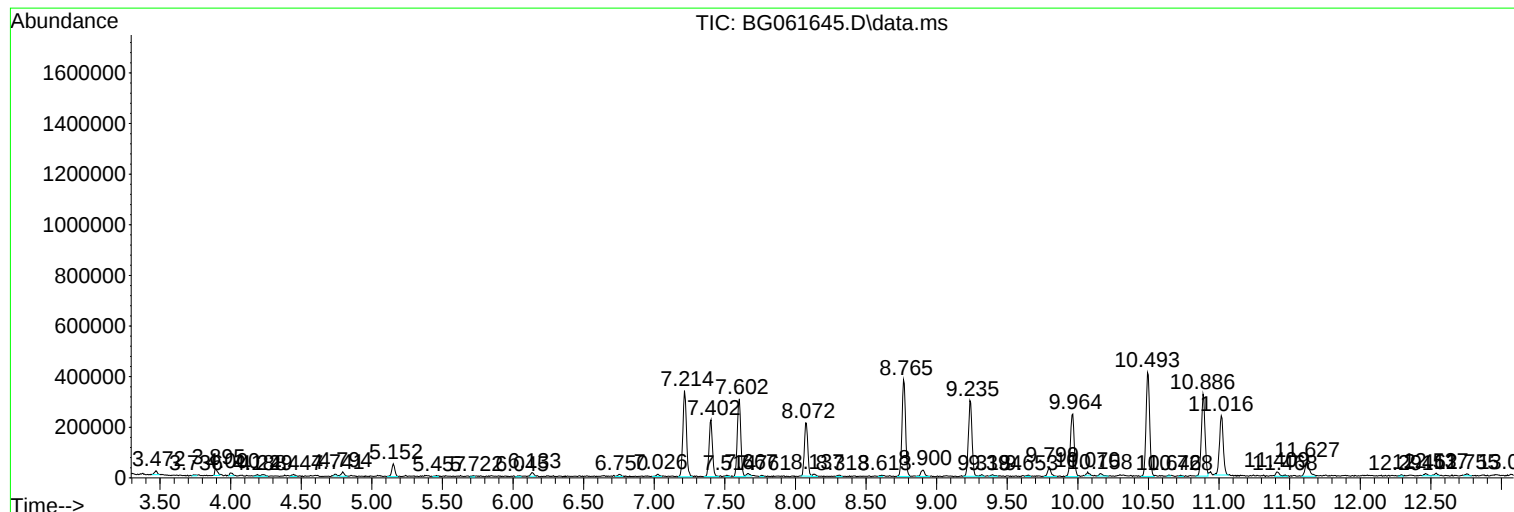
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C95

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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



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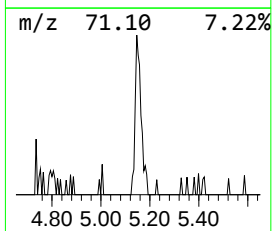
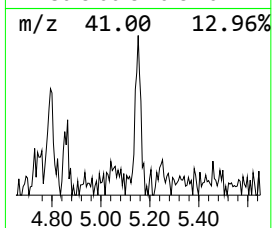
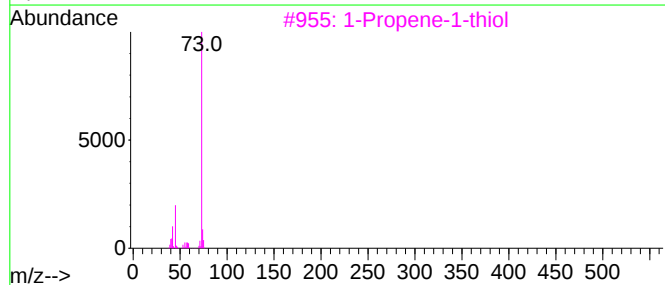
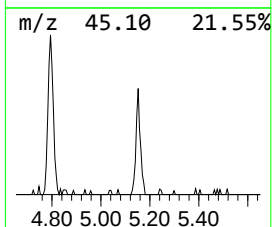
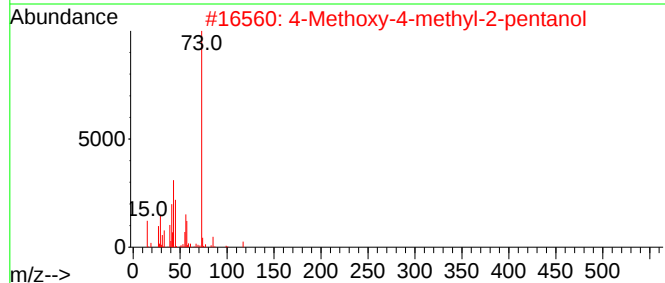
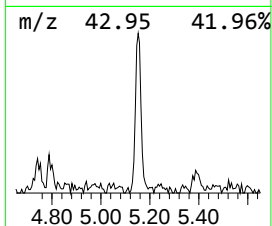
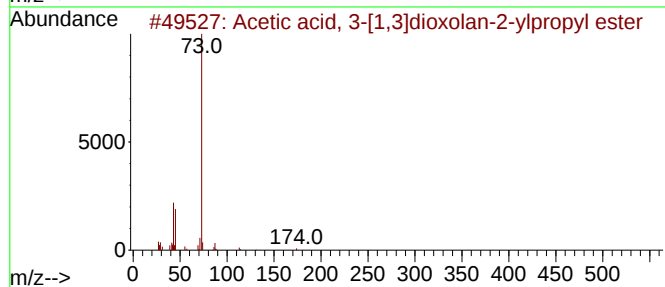
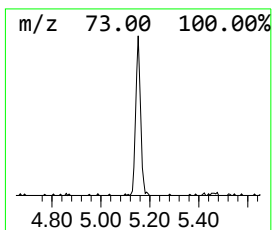
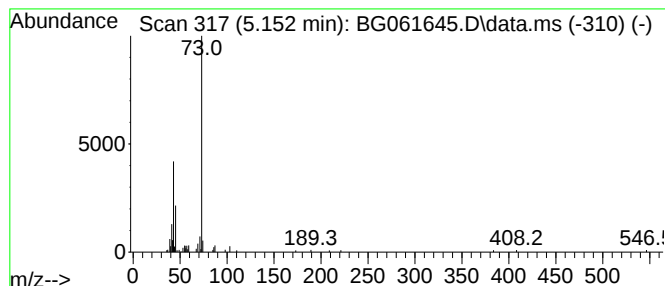
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.152	4.25 ng/ul	75264	1,4-Dichlorobenzene-d4	8.078

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	42
2			4-Methoxy-4-methyl-2-pentanol	132	C7H16O2	000141-73-1	33
3			1-Propene-1-thiol	74	C3H6S	000925-89-3	33
4			1,3-Dioxolane, 2-(3-bromo-5,5,5-...	352	C10H16BrCl3O2	1000115-31-4	33
5			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	33



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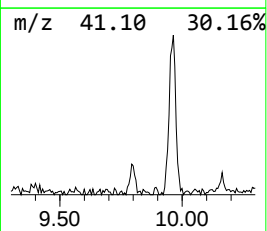
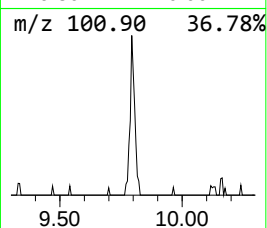
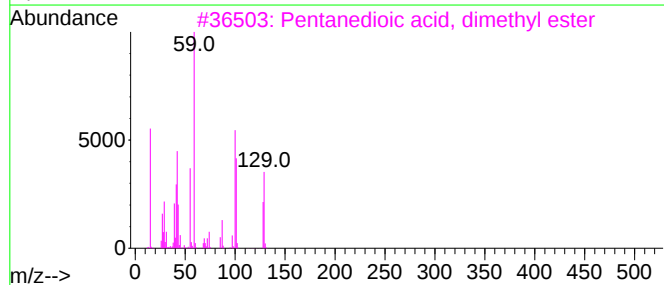
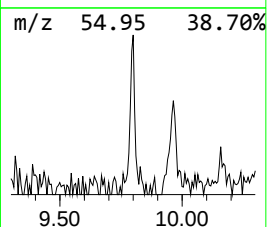
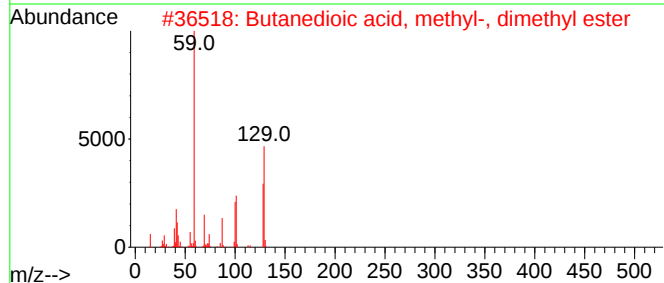
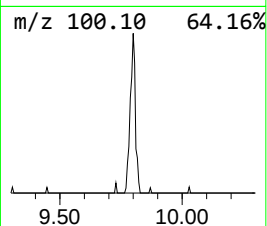
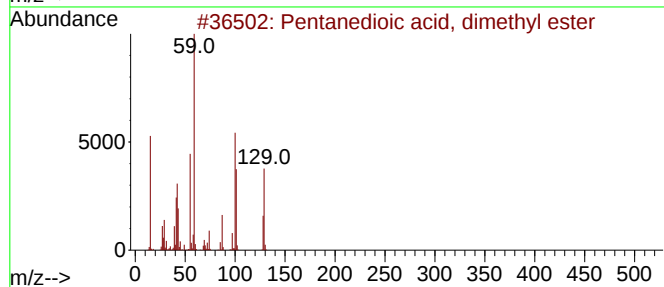
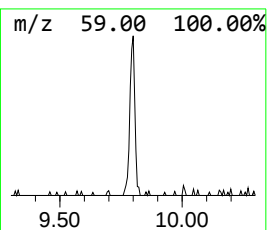
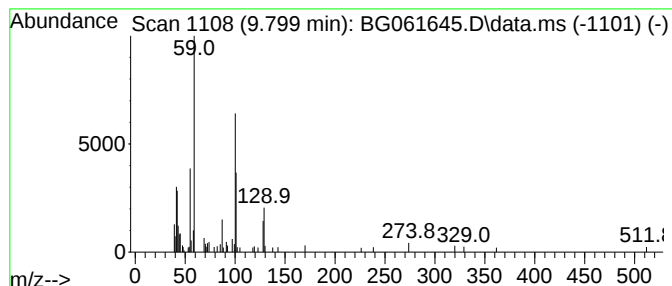
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Pentanedioic acid, dimethyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.799	2.17 ng/ul	62718	Naphthalene-d8	10.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	56
2			Butanedioic acid, methyl-, dimet...	160	C7H12O4	001604-11-1	53
3			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	47
4			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	47
5			Butanedioic acid, methyl-, dimet...	160	C7H12O4	001604-11-1	36



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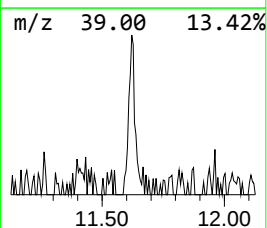
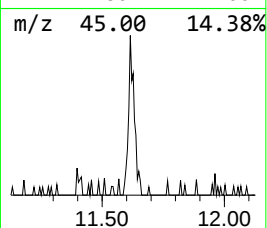
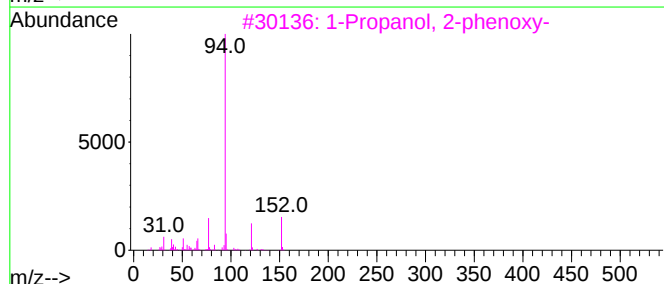
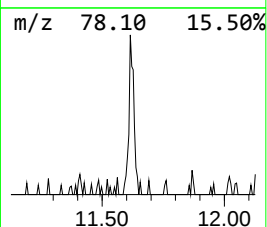
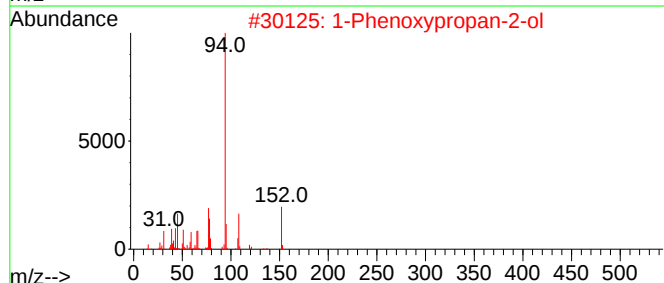
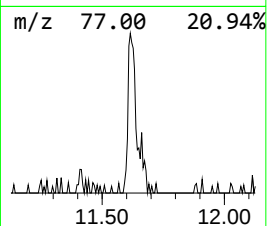
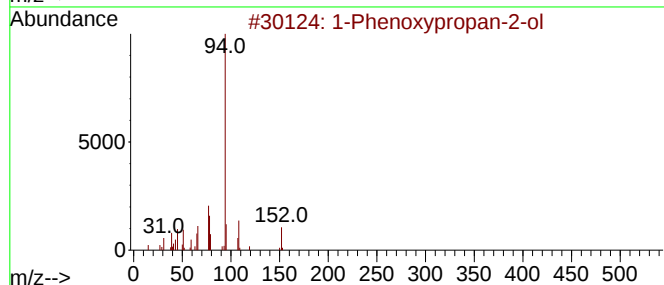
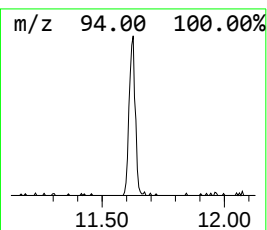
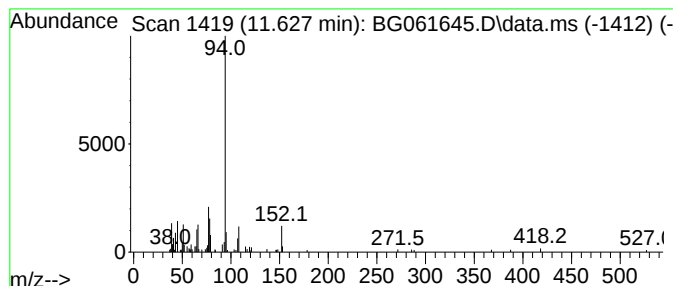
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.627	4.30 ng/ul	124116	Naphthalene-d8	10.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
3			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	72
4			1,2-Propanediol, 3-phenoxy-	168	C9H12O3	000538-43-2	64
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
Operator : MA/JU
Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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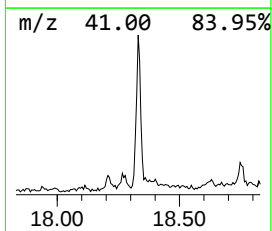
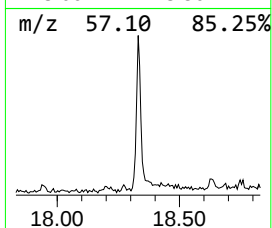
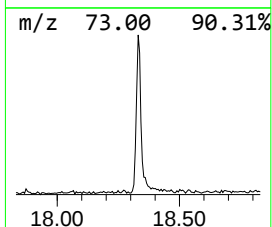
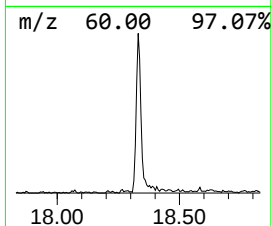
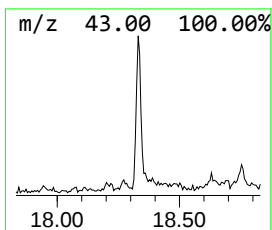
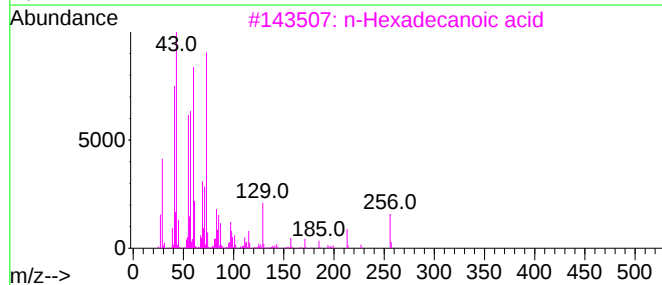
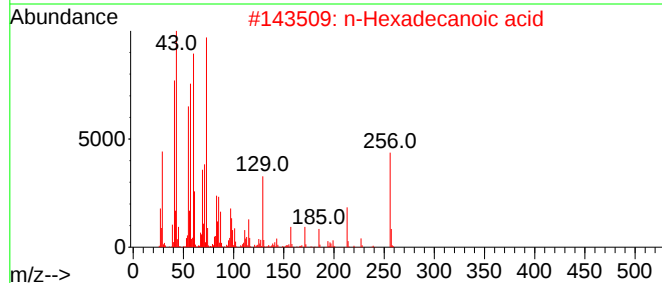
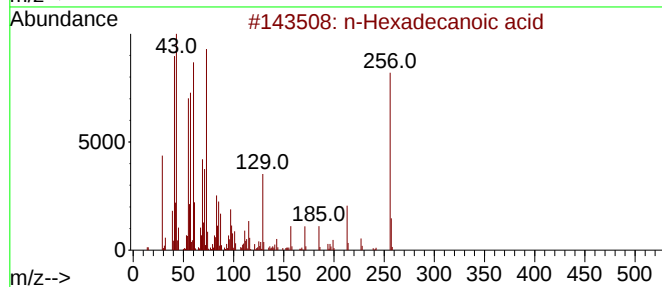
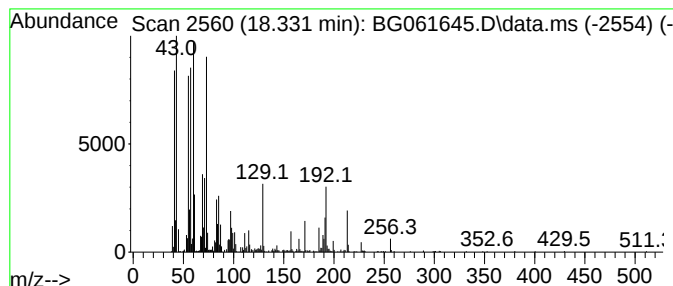
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.331	15.47 ng/ul	807573	Phenanthrene-d10	17.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	62
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
Operator : MA/JU
Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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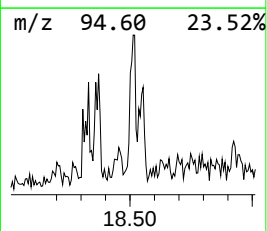
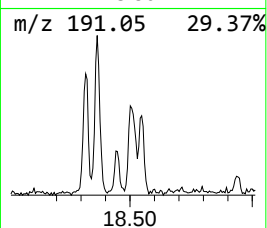
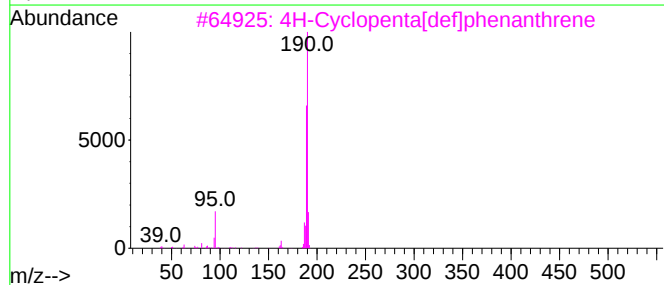
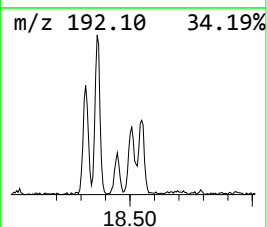
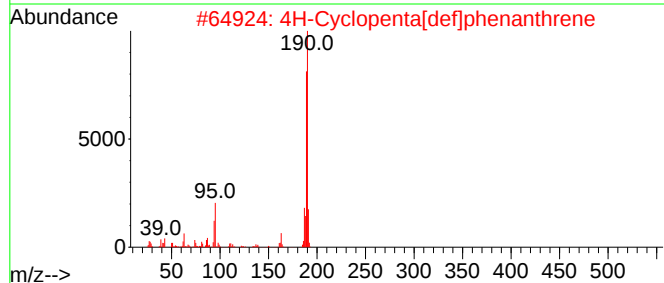
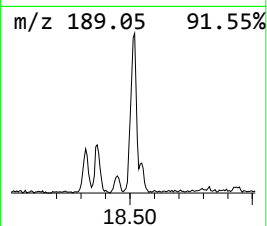
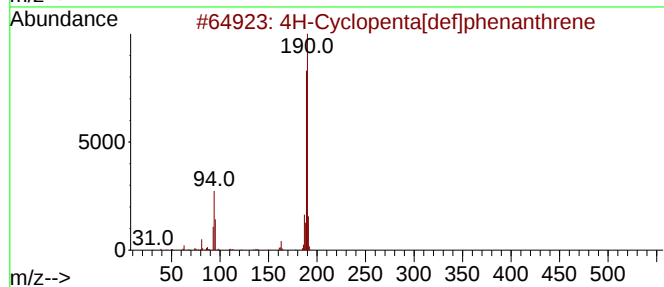
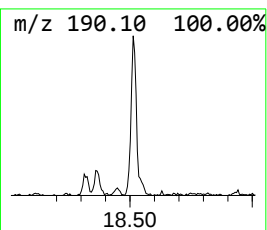
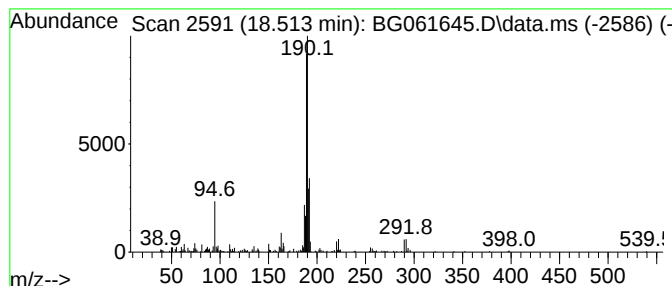
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.513	5.34 ng/ul	278875	Phenanthrene-d10	17.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	72
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	59
4			Methyl diselenide	190	C2H6Se2	007101-31-7	55
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
Operator : MA/JU
Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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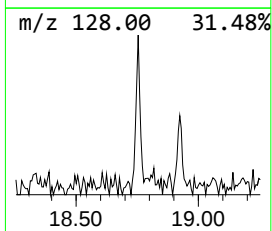
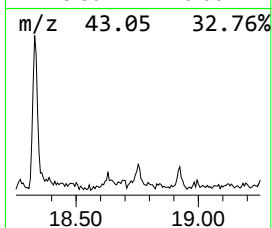
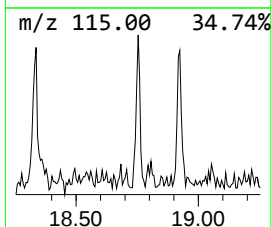
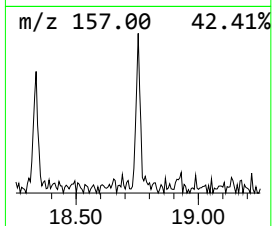
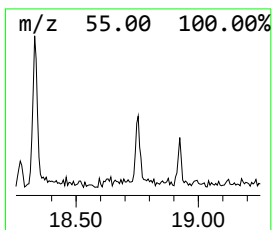
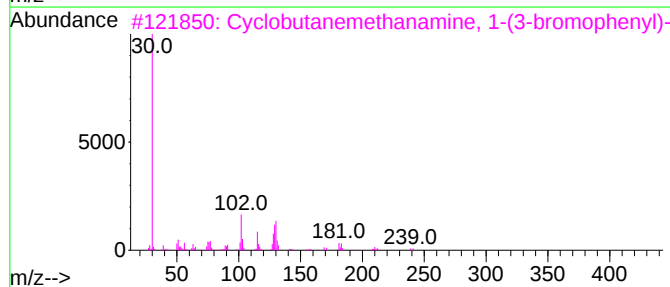
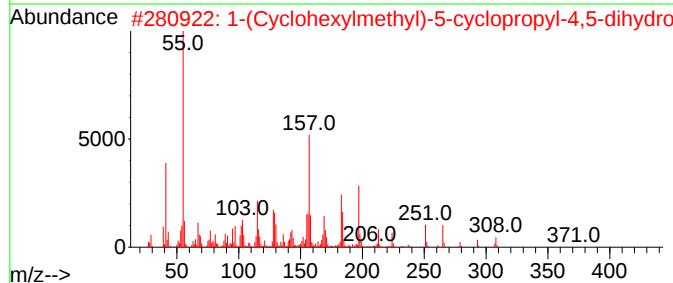
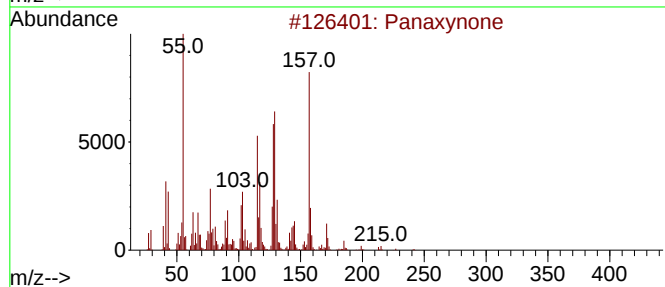
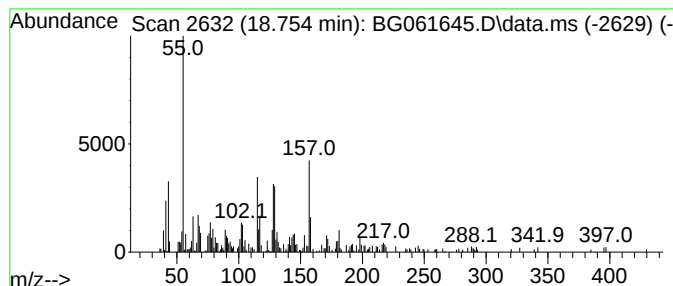
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Panaxynone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.754	2.65 ng/ul	138363	Phenanthrene-d10	17.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Panaxynone	242	C17H22O	004117-11-7	58
2			1-(Cyclohexylmethyl)-5-cycloprop...	372	C19H24N4O2S	1000481-83-1	25
3			Cyclobutanemethanamine, 1-(3-bro...	239	C11H14BrN	915690-61-8	11
4			Diethyl adipate	202	C10H18O4	000141-28-6	9
5			Propanedioic acid, methyl-, diet...	174	C8H14O4	000609-08-5	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
Operator : MA/JU
Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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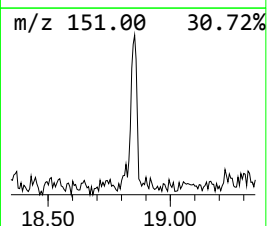
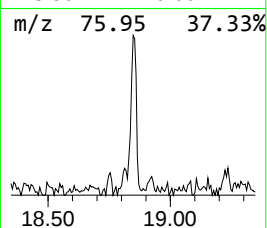
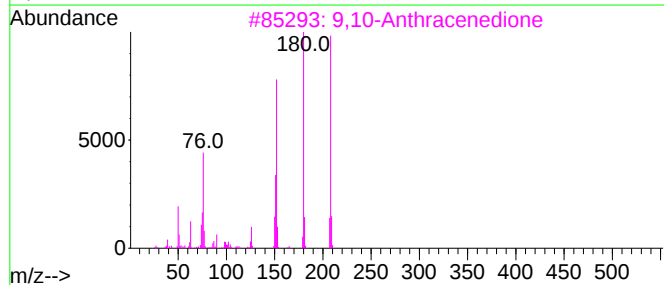
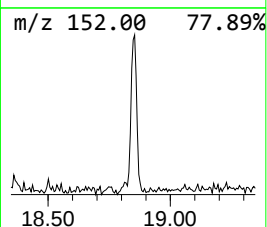
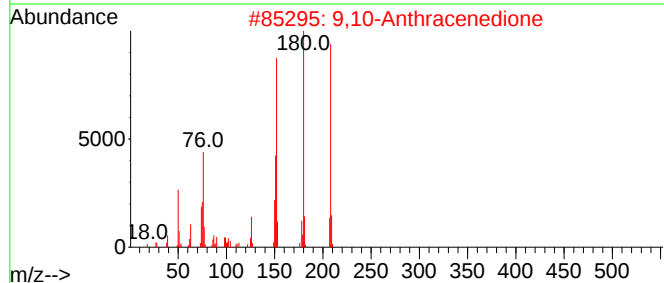
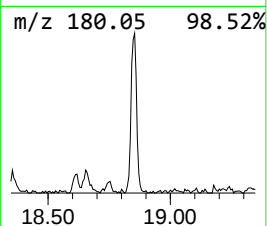
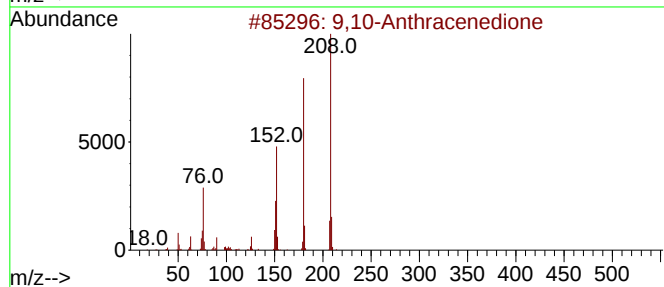
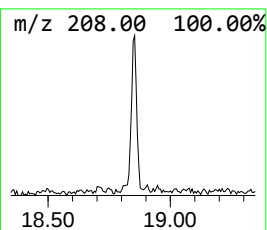
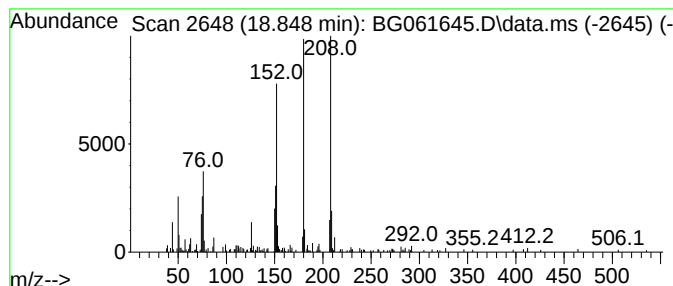
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.848	3.76 ng/ul	196074	Phenanthrene-d10	17.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	91
5			9H-Fluoren-9-one	180	C13H8O	000486-25-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
Operator : MA/JU
Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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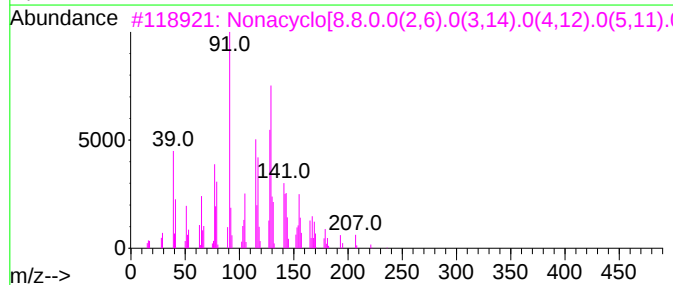
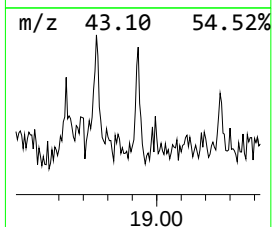
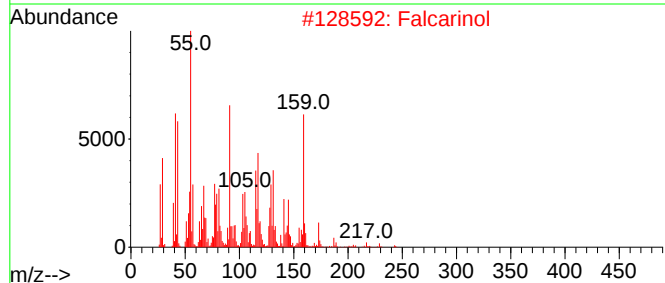
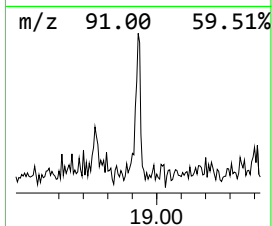
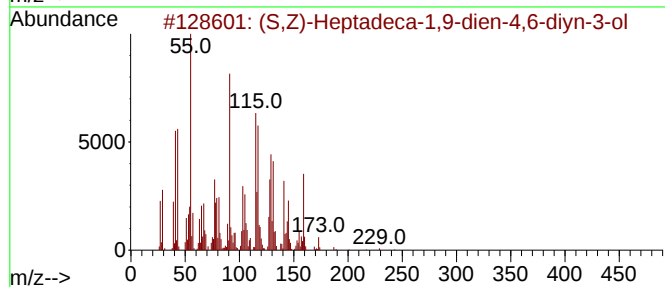
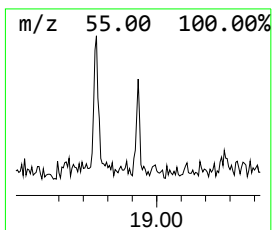
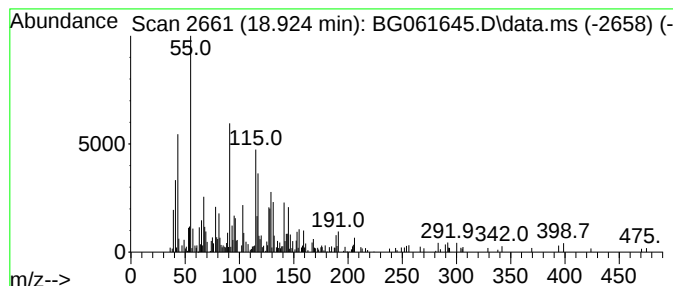
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (S,Z)-Heptadeca-1,9-dien-4,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.924	2.77 ng/ul	144665	Phenanthrene-d10	17.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S,Z)-Heptadeca-1,9-dien-4,6-diy...	244	C17H24O	081203-57-8	72		
2	Falcarinol	244	C17H24O	021852-80-2	38		
3	Nonacyclo[8.8.0.0(2,6).0(3,14).0...	236	C18H20	1000436-96-5	32		
4	Octanal, 2-(phenylmethylene)-	216	C15H20O	000101-86-0	32		
5	trans-2-Methyl-.beta.-methyl-.be...	177	C10H11NO2	147102-54-3	32		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
Operator : MA/JU
Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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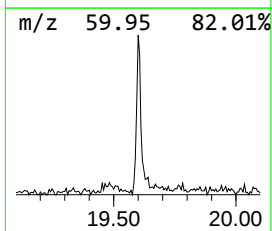
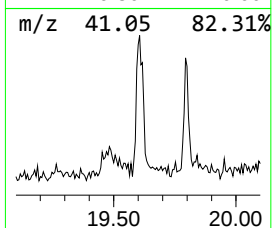
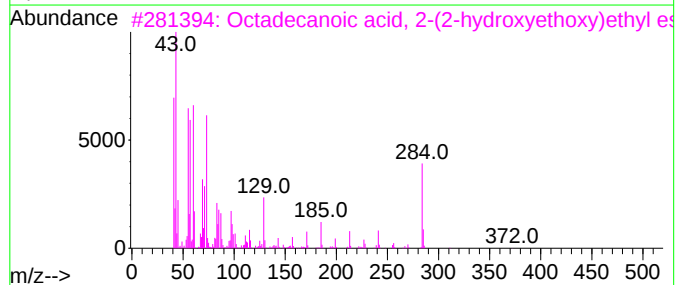
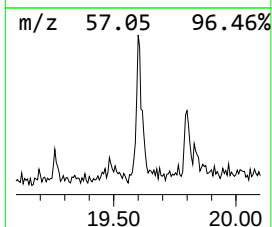
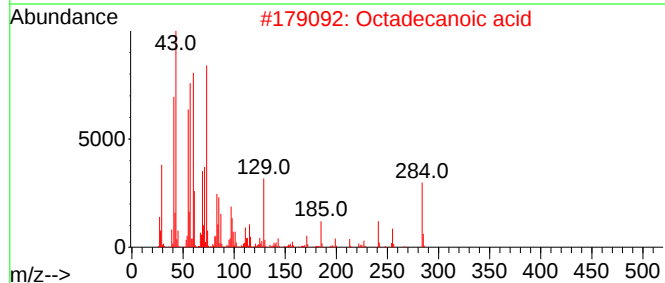
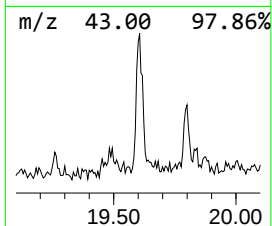
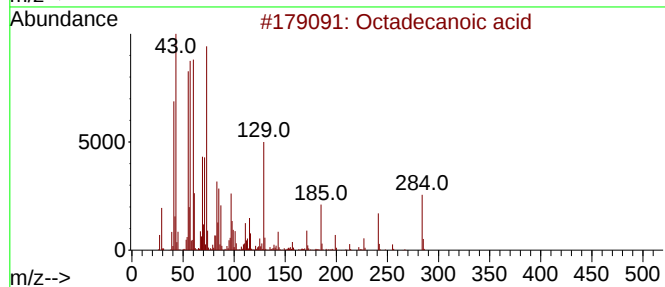
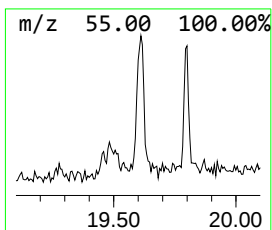
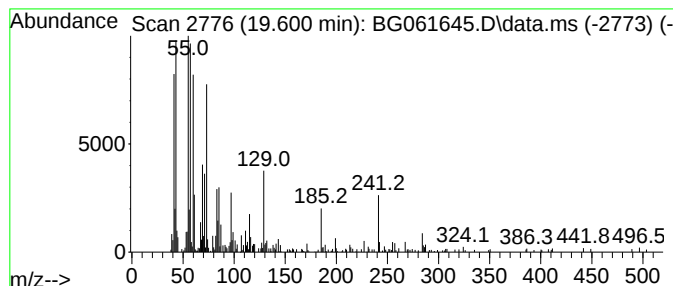
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.600	4.61 ng/ul	494916	Chrysene-d12	21.709

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	93
2			Octadecanoic acid	284	C18H36O2	000057-11-4	89
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	78
4			Octadecanoic acid	284	C18H36O2	000057-11-4	64
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
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Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

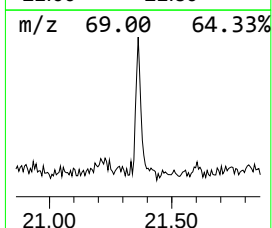
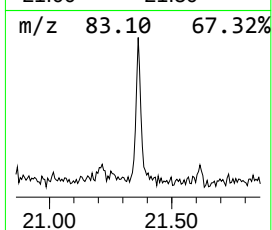
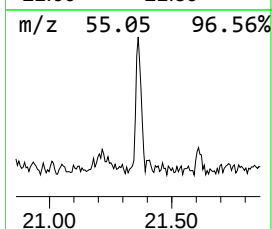
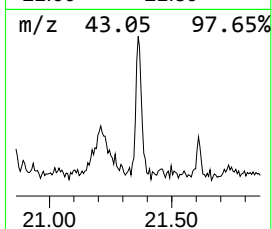
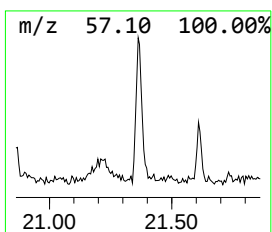
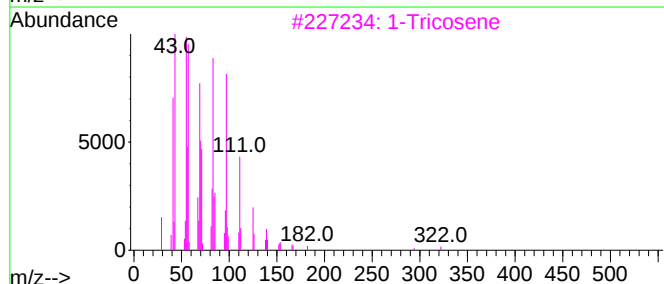
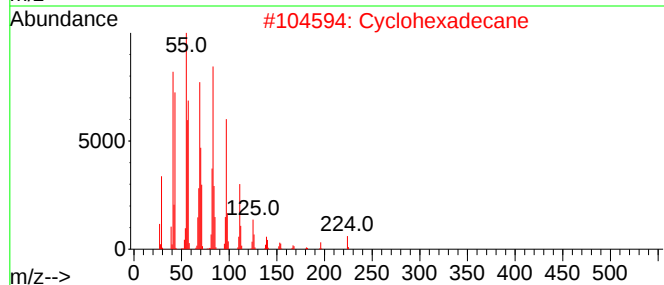
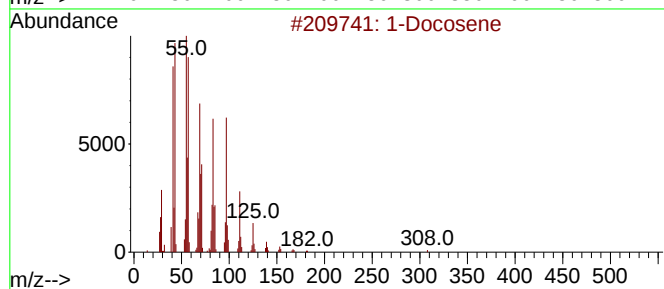
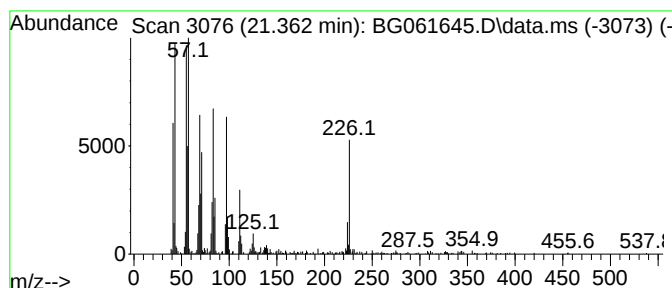
Instrument :
BNA_G
ClientSampleId :
A4C95

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.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 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.362	6.04 ng/ul	647960	Chrysene-d12	21.709	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	95
2	Cyclohexadecane	224	C16H32	000295-65-8	89
3	1-Tricosene	322	C23H46	018835-32-0	83
4	Cyclotriacontane	420	C30H60	000297-35-8	78
5	Cetene	224	C16H32	000629-73-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061645.D
Acq On : 22 Jun 2024 12:20
Operator : MA/JU
Sample : P2775-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C95

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	5.152	4.3	ng/ul	75264	1	8.078	353939	20.0
Pentanedioic ac...	9.799	2.2	ng/ul	62718	2	10.886	577677	20.0
1-Phenoxypropan...	11.627	4.3	ng/ul	124116	2	10.886	577677	20.0
n-Hexadecanoic ...	18.331	15.5	ng/ul	807573	4	17.443	1043760	20.0
4H-Cyclopenta[d...	18.513	5.3	ng/ul	278875	4	17.443	1043760	20.0
Panaxynone	18.754	2.6	ng/ul	138363	4	17.443	1043760	20.0
9,10-Anthracene...	18.848	3.8	ng/ul	196074	4	17.443	1043760	20.0
(S,Z)-Heptadeca...	18.924	2.8	ng/ul	144665	4	17.443	1043760	20.0
Octadecanoic acid	19.600	4.6	ng/ul	494916	5	21.709	2146370	20.0
(DEL) Alkane: S...	21.362	6.0	ng/ul	647960	5	21.709	2146370	20.0