

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
 Data File : BG061675.D
 Acq On : 24 Jun 2024 14:50
 Operator : MA/JU
 Sample : P2775-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4BQ3

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: BG061675.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	27	31	35	rVB5	17020	26370	1.03%	0.092%
2	3.737	72	76	83	rBV	75774	113088	4.42%	0.396%
3	3.901	99	104	107	rVV4	8934	11588	0.45%	0.041%
4	4.007	117	122	128	rVV	35550	51849	2.02%	0.181%
5	4.078	131	134	139	rVV4	7581	9056	0.35%	0.032%
6	4.131	139	143	147	rVB5	3734	5273	0.21%	0.018%
7	4.195	151	154	155	rBV2	3809	4329	0.17%	0.015%
8	4.236	157	161	167	rVB2	30411	49240	1.92%	0.172%
9	4.383	182	186	187	rBV	5776	5166	0.20%	0.018%
10	4.742	243	247	251	rVV5	5673	9443	0.37%	0.033%
11	4.789	251	255	261	rVB3	17179	29191	1.14%	0.102%
12	4.930	277	279	285	rVB4	4680	5630	0.22%	0.020%
13	5.029	294	296	299	rBV4	5076	4288	0.17%	0.015%
14	5.153	310	317	323	rBV	46218	78400	3.06%	0.274%
15	5.235	327	331	339	rVB5	10298	18382	0.72%	0.064%
16	6.134	479	484	490	rVB2	14607	26053	1.02%	0.091%
17	7.021	632	635	639	rBV4	7191	11142	0.44%	0.039%
18	7.215	661	668	679	rVB	299293	515984	20.15%	1.806%
19	7.397	692	699	706	rBV	234538	394030	15.38%	1.379%
20	7.603	728	734	740	rBV2	287545	494630	19.31%	1.731%
21	7.662	740	744	754	rVV	59478	97986	3.83%	0.343%
22	7.744	756	758	763	rVB3	3024	4643	0.18%	0.016%
23	8.079	806	815	821	rBV	256631	445986	17.41%	1.561%
24	8.132	821	824	829	rVB3	10732	15595	0.61%	0.055%
25	8.543	892	894	897	rBV2	4520	5789	0.23%	0.020%
26	8.766	924	932	940	rBV	330215	562215	21.95%	1.968%
27	8.896	948	954	962	rVB3	25065	45348	1.77%	0.159%
28	9.236	1005	1012	1019	rBV	326214	572935	22.37%	2.005%
29	9.324	1025	1027	1034	rVV5	10017	12688	0.50%	0.044%
30	9.395	1035	1039	1044	rVB4	5186	8038	0.31%	0.028%
31	9.483	1053	1054	1057	rBV2	4489	4584	0.18%	0.016%
32	9.636	1076	1080	1084	rBV3	3819	6248	0.24%	0.022%
33	9.794	1103	1107	1113	rVB3	35836	54543	2.13%	0.191%
34	9.959	1128	1135	1145	rVV3	275184	499202	19.49%	1.747%
35	10.071	1148	1154	1156	rBV4	10811	20449	0.80%	0.072%
36	10.159	1166	1169	1175	rVB4	7132	12791	0.50%	0.045%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
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37	10.500	1219	1227	1235	rBV	426220	769749	30.05%	2.694%
38	10.652	1250	1253	1255	rBV2	5065	5503	0.21%	0.019%
39	10.887	1285	1293	1300	rBV	449297	787660	30.75%	2.757%
40	11.017	1308	1315	1325	rBV2	170981	316439	12.36%	1.108%
41	11.410	1379	1382	1388	rVB3	22335	32062	1.25%	0.112%
42	11.622	1410	1418	1430	rBV3	130153	246567	9.63%	0.863%
43	11.745	1437	1439	1443	rVB2	4895	5695	0.22%	0.020%
44	12.180	1512	1513	1516	rBV	4038	4484	0.18%	0.016%
45	12.233	1521	1522	1527	rVB4	4527	5285	0.21%	0.018%
46	12.321	1535	1537	1539	rBV3	4580	5249	0.20%	0.018%
47	12.462	1559	1561	1568	rVB4	13320	17905	0.70%	0.063%
48	12.532	1571	1573	1574	rBV2	4951	4371	0.17%	0.015%
49	12.762	1607	1612	1613	rBV3	3654	4312	0.17%	0.015%
50	12.944	1638	1643	1644	rBV3	4798	6131	0.24%	0.021%
51	13.249	1692	1695	1698	rVB4	4880	7112	0.28%	0.025%
52	13.320	1703	1707	1708	rBV3	4883	6226	0.24%	0.022%
53	13.537	1742	1744	1748	rVB4	8257	7218	0.28%	0.025%
54	13.602	1753	1755	1759	rBV4	5313	5948	0.23%	0.021%
55	13.737	1775	1778	1781	rVB4	3581	4440	0.17%	0.016%
56	13.825	1788	1793	1797	rBV	12885	23629	0.92%	0.083%
57	14.013	1822	1825	1827	rBV2	4415	5387	0.21%	0.019%
58	14.107	1835	1841	1848	rVB	877792	1251550	48.87%	4.380%
59	14.160	1848	1850	1854	rVB3	6056	5324	0.21%	0.019%
60	14.236	1858	1863	1864	rBV3	5088	7148	0.28%	0.025%
61	14.342	1875	1881	1885	rBV5	29821	49203	1.92%	0.172%
62	14.395	1885	1890	1899	rVB2	899911	1480476	57.80%	5.182%
63	14.595	1922	1924	1927	rVB2	7539	7201	0.28%	0.025%
64	14.701	1933	1942	1949	rBV	769748	1205534	47.07%	4.219%
65	14.765	1950	1953	1959	rVB2	22148	36182	1.41%	0.127%
66	14.871	1965	1971	1977	rBV	367846	643795	25.14%	2.253%
67	15.035	1995	1999	2006	rVB4	32043	48901	1.91%	0.171%
68	15.100	2006	2010	2015	rBV4	31002	52009	2.03%	0.182%
69	15.406	2060	2062	2065	rBV4	6653	7025	0.27%	0.025%
70	15.458	2069	2071	2074	rBV4	4232	4512	0.18%	0.016%
71	15.482	2074	2075	2076	rBV	7266	4670	0.18%	0.016%
72	15.693	2104	2111	2117	rBV	1293360	1930846	75.39%	6.758%
73	15.746	2117	2120	2123	rVV3	29542	39478	1.54%	0.138%
74	15.793	2123	2128	2135	rVB	406189	591055	23.08%	2.069%
75	15.911	2146	2148	2150	rBV3	6241	4758	0.19%	0.017%
76	16.046	2167	2171	2176	rVB6	10773	19374	0.76%	0.068%

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77	16.105	2179	2181	2182	rBV2	8395	5562	0.22%	0.019%
78	16.228	2200	2202	2206	rVB4	18237	18812	0.73%	0.066%
79	16.410	2229	2233	2240	rBV9	16743	37791	1.48%	0.132%
80	16.745	2288	2290	2295	rVB6	13388	15675	0.61%	0.055%
81	16.804	2298	2300	2302	rBV3	8435	5266	0.21%	0.018%
82	16.986	2326	2331	2335	rBV7	21193	34374	1.34%	0.120%
83	17.145	2354	2358	2361	rBV5	20118	32890	1.28%	0.115%
84	17.444	2403	2409	2413	rBV	1014707	1556686	60.78%	5.448%
85	17.485	2413	2416	2421	rVV	364236	535850	20.92%	1.876%
86	17.544	2421	2426	2438	rVB	1383326	2123929	82.93%	7.434%
87	17.844	2475	2477	2484	rVB3	39580	42008	1.64%	0.147%
88	18.114	2519	2523	2527	rBV2	49820	66256	2.59%	0.232%
89	18.332	2554	2560	2564	rBV2	396324	575416	22.47%	2.014%
90	18.514	2586	2591	2594	rBV3	79480	132637	5.18%	0.464%
91	18.813	2640	2642	2645	rBV	39258	43676	1.71%	0.153%
92	18.854	2646	2649	2653	rVB2	61433	81245	3.17%	0.284%
93	19.489	2753	2757	2763	rVB	820649	1153853	45.05%	4.039%
94	19.601	2772	2776	2782	rBV4	137113	196576	7.68%	0.688%
95	19.824	2809	2814	2818	rBV	1581808	2561174	100.00%	8.964%
96	21.363	3073	3076	3083	rVB3	313969	488566	19.08%	1.710%
97	21.610	3115	3118	3122	rVB	208085	267688	10.45%	0.937%
98	21.710	3127	3135	3140	rBV2	995718	2075775	81.05%	7.265%
99	24.718	3642	3647	3655	rVB2	505068	1155358	45.11%	4.044%
100	24.941	3677	3685	3695	rVB2	537027	1471357	57.45%	5.150%

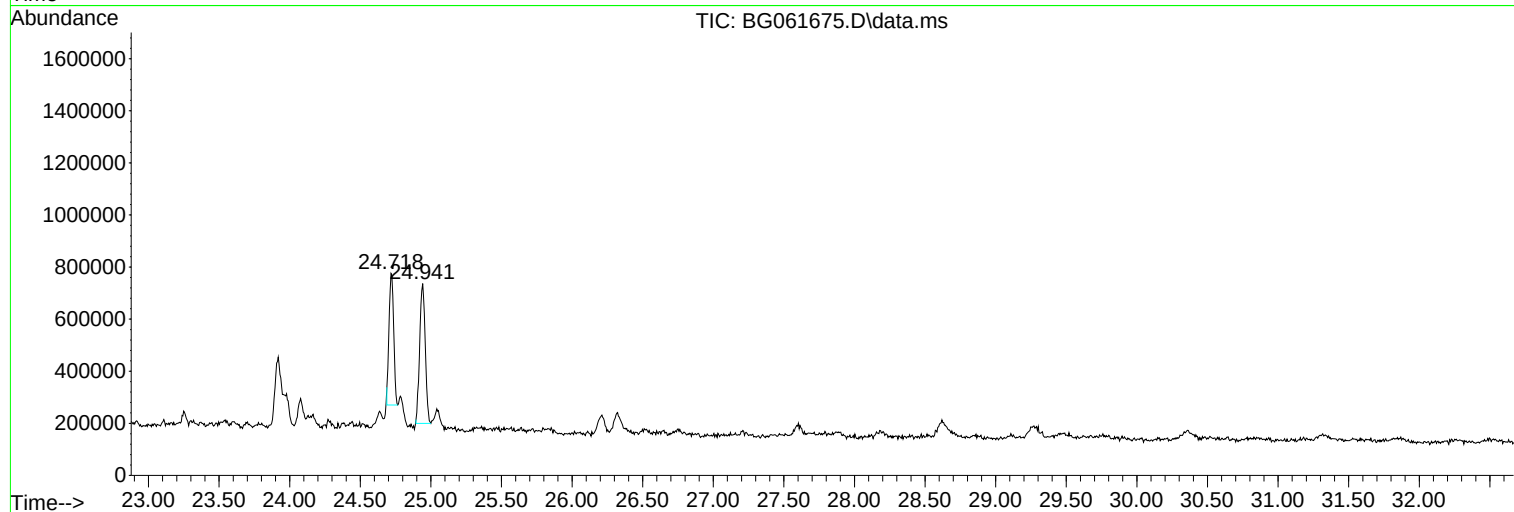
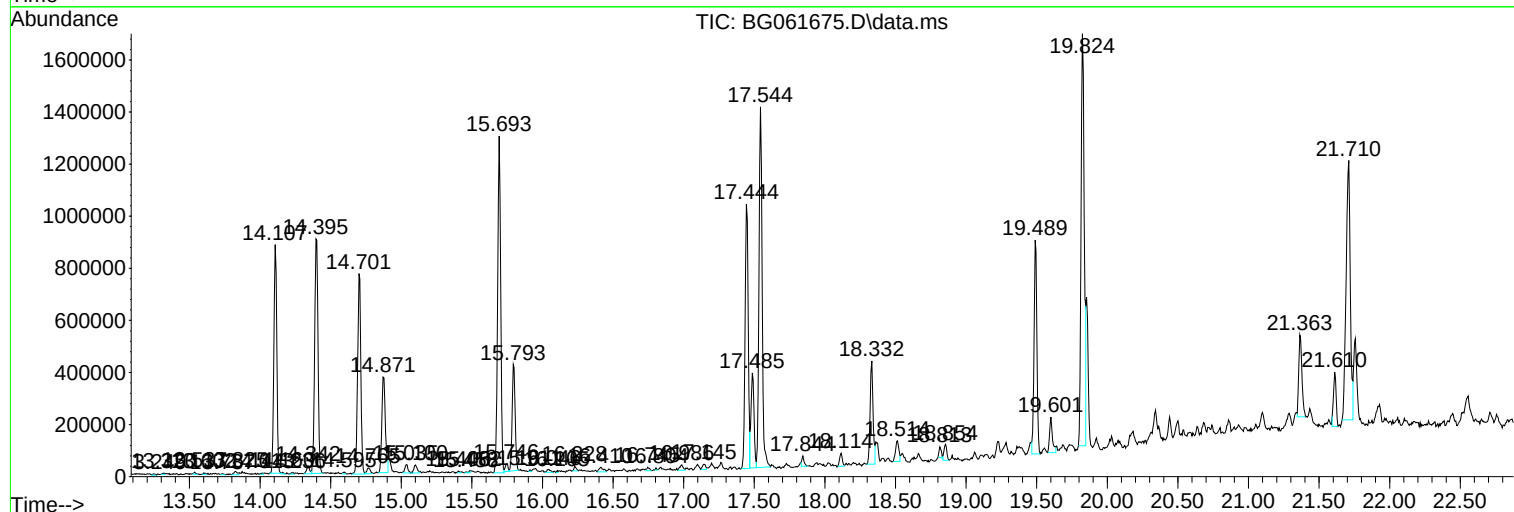
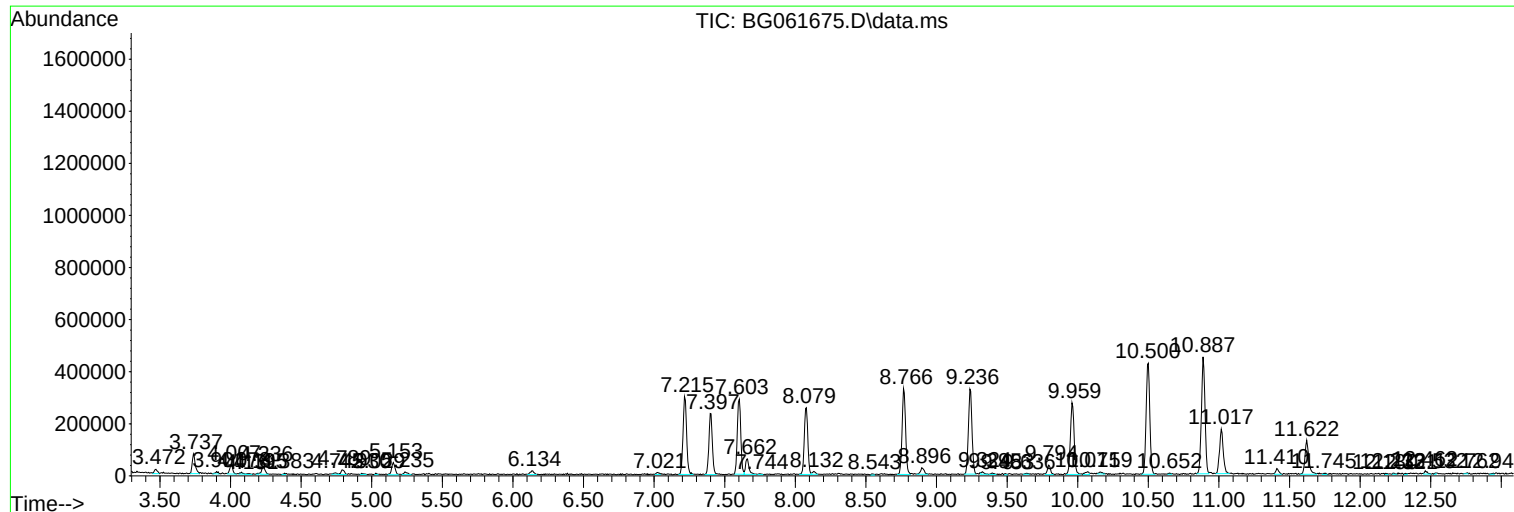
Sum of corrected areas: 28571005

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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BQ3

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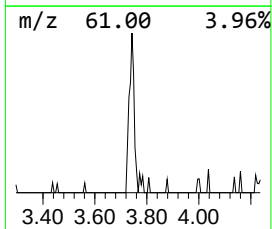
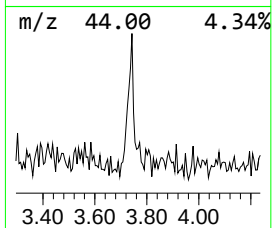
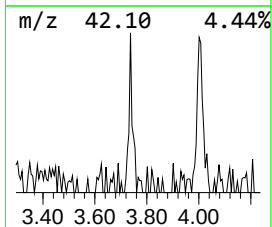
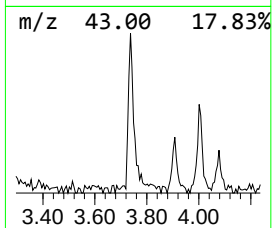
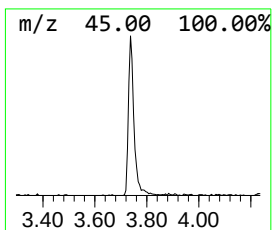
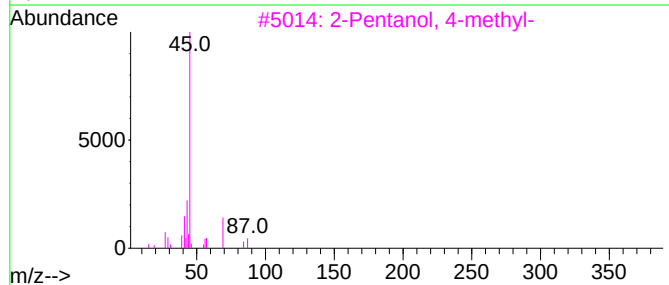
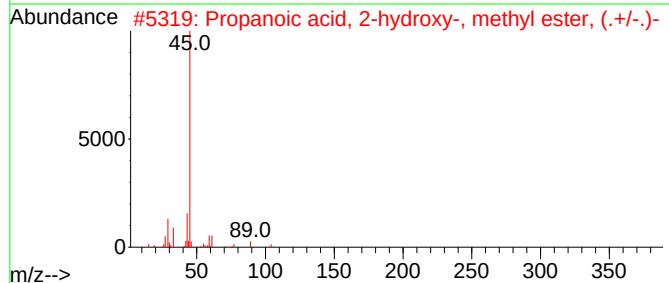
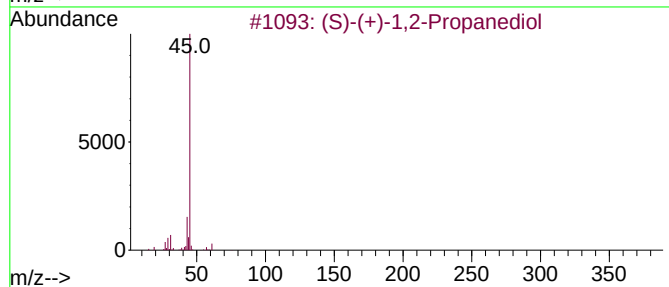
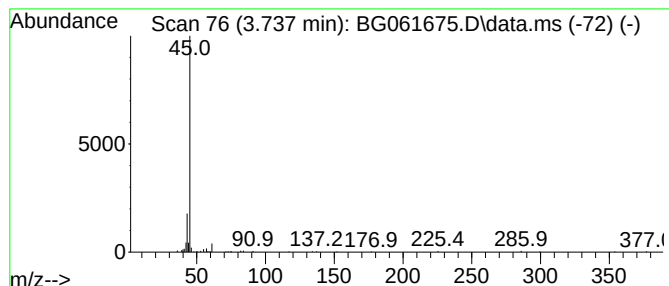
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.737	5.07 ng/ul	113088	1,4-Dichlorobenzene-d4	8.079

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	40
2	Propanoic acid, 2-hydroxy-, meth...			104	C4H8O3	000547-64-8	9
3	2-Pentanol, 4-methyl-			102	C6H14O	000108-11-2	9
4	Methoxyacetyl chloride			108	C3H5ClO2	038870-89-2	9
5	Isopropyl Alcohol			60	C3H8O	000067-63-0	9



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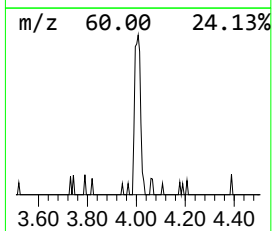
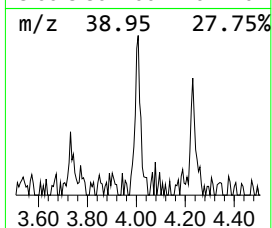
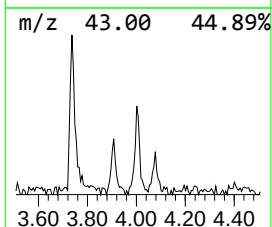
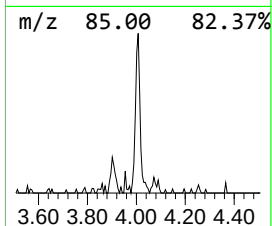
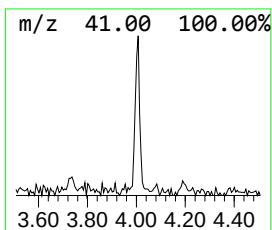
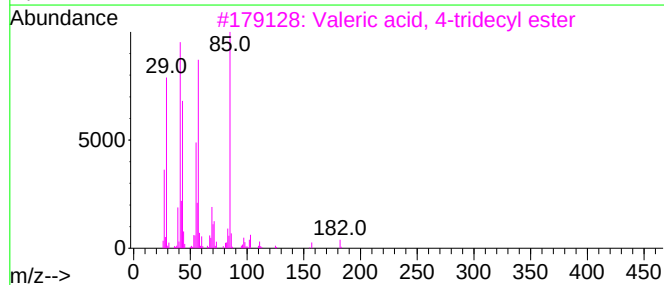
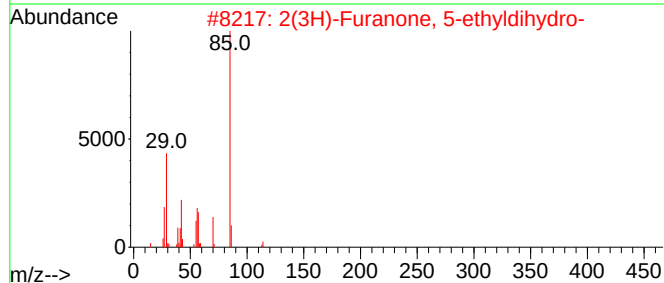
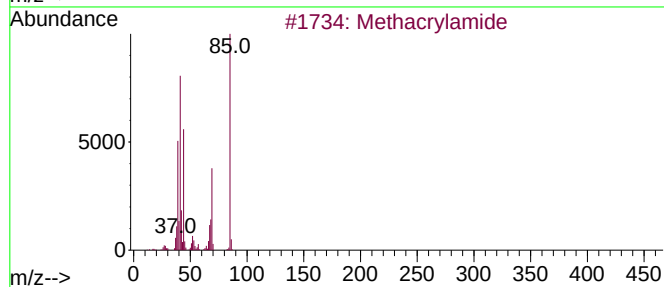
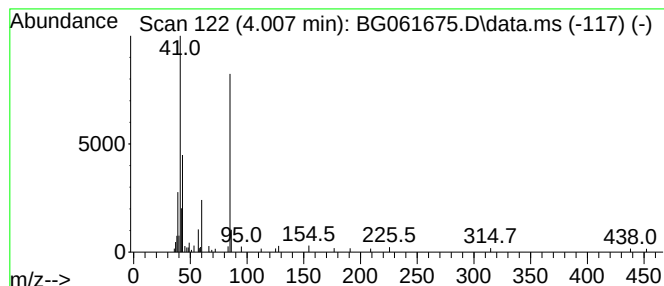
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.007	2.33 ng/ul	51849	1,4-Dichlorobenzene-d4	8.079

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	47
2			2(3H)-Furanone, 5-ethyldihydro-	114	C6H10O2	000695-06-7	45
3			Valeric acid, 4-tridecyl ester	284	C18H36O2	1000281-99-0	38
4			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
5			Methacrylamide	85	C4H7NO	000079-39-0	33



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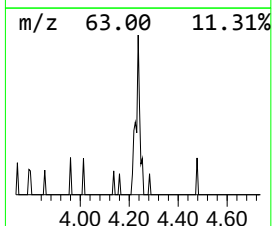
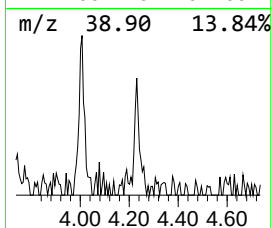
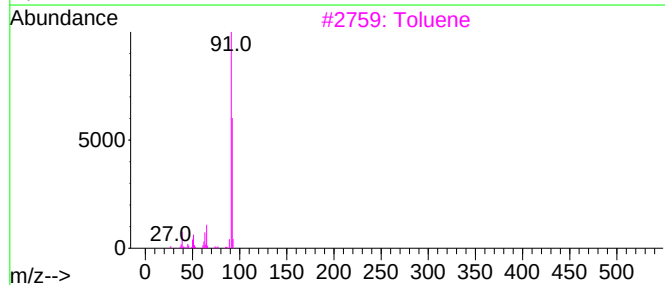
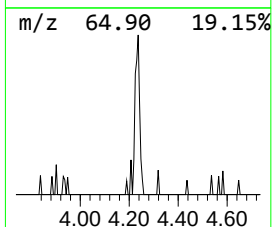
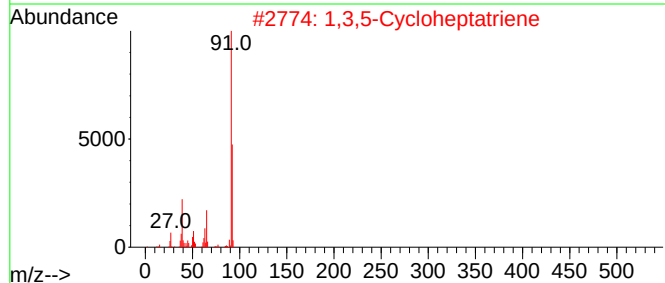
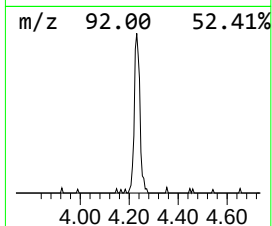
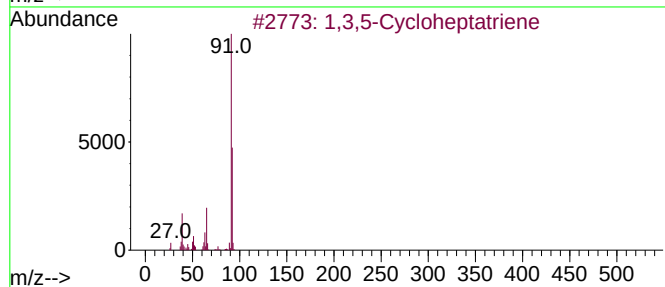
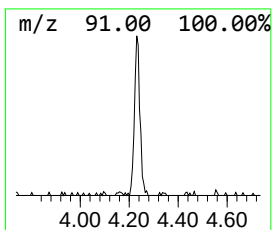
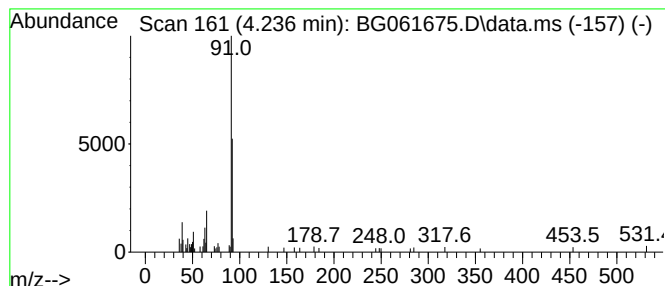
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,3,5-Cycloheptatriene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.236	2.21 ng/ul	49240	1,4-Dichlorobenzene-d4	8.079

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	68
2	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	68
3	Toluene			92	C7H8	000108-88-3	68
4	Toluene			92	C7H8	000108-88-3	64
5	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061675.D
Acq On : 24 Jun 2024 14:50
Operator : MA/JU
Sample : P2775-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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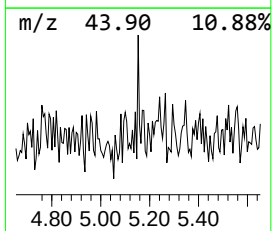
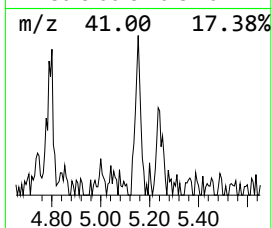
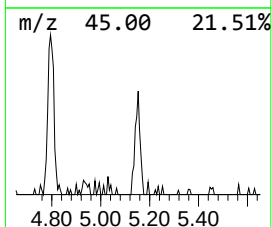
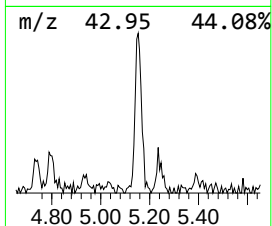
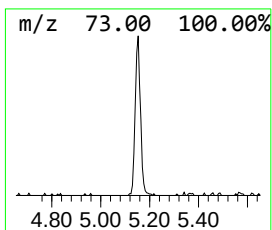
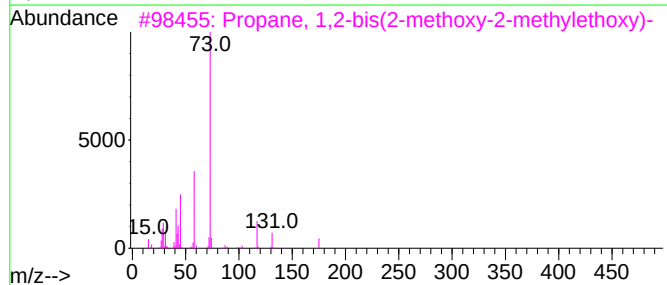
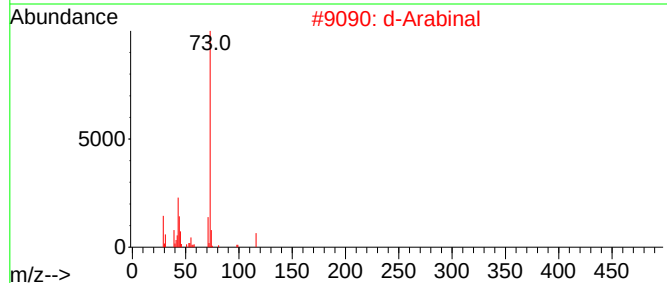
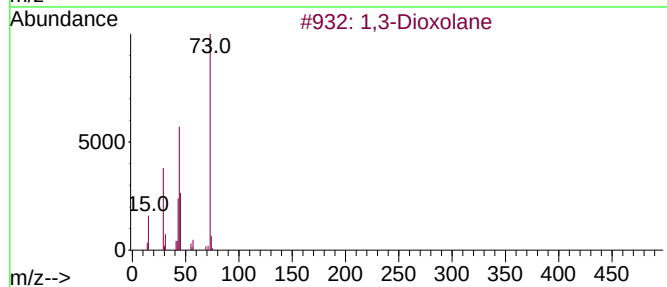
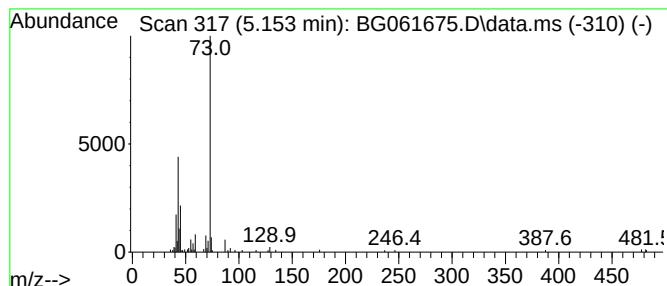
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.153	3.52 ng/ul	78400	1,4-Dichlorobenzene-d4	8.079

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane	74	C3H6O2	000646-06-0	43
2			d-Arabinol	116	C5H8O3	000496-61-7	32
3			Propane, 1,2-bis(2-methoxy-2-met...	220	C11H24O4	042769-21-1	28
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	25
5			1,3-Dioxolane, 2-(1-methylpropyl)-	130	C7H14O2	014447-25-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061675.D
Acq On : 24 Jun 2024 14:50
Operator : MA/JU
Sample : P2775-01
Misc :
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Instrument :
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ClientSampleId :
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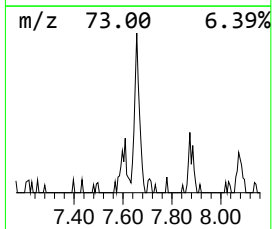
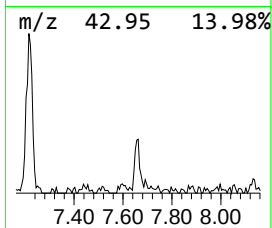
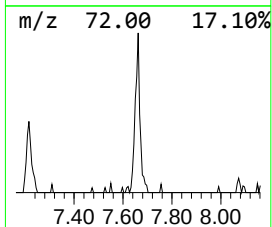
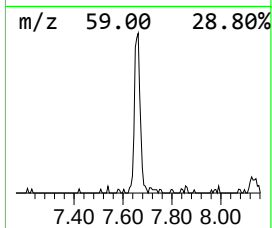
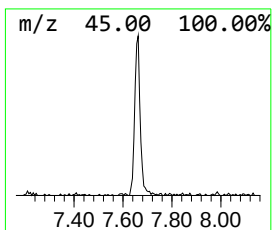
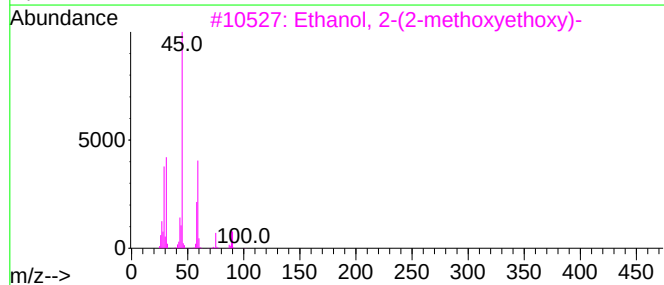
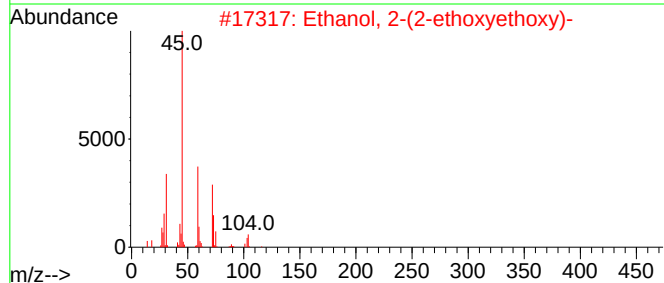
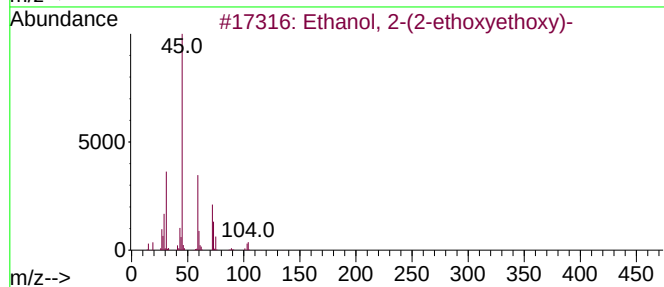
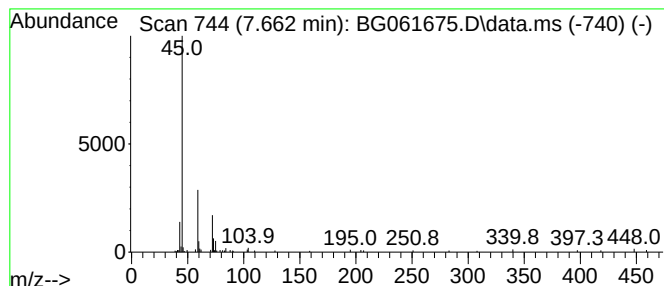
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.662	4.39 ng/ul	97986	1,4-Dichlorobenzene-d4	8.079

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	56
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	40
3			Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	39
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	38
5			Trichloroacetic acid, 2-methoxye...	220	C5H7Cl3O3	1000330-88-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061675.D
Acq On : 24 Jun 2024 14:50
Operator : MA/JU
Sample : P2775-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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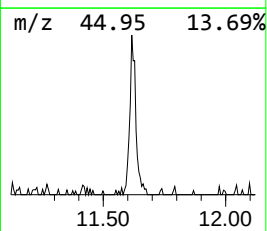
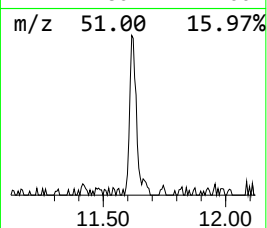
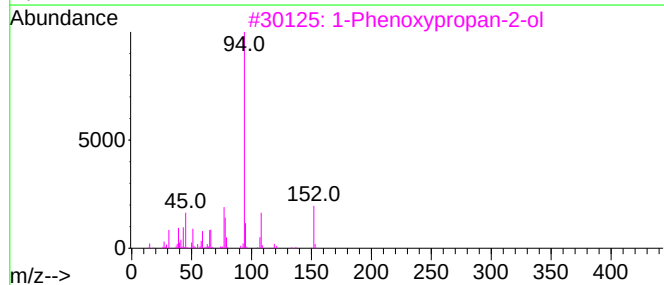
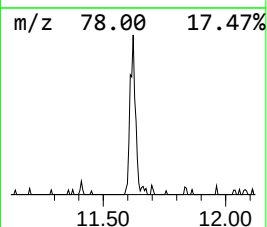
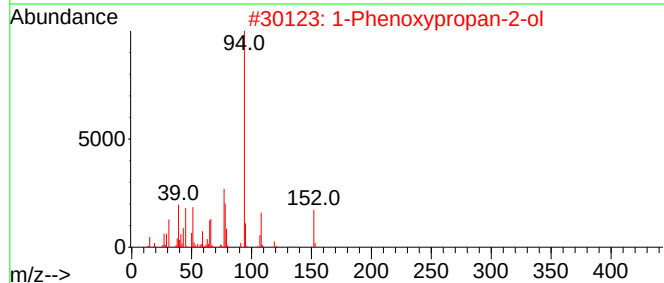
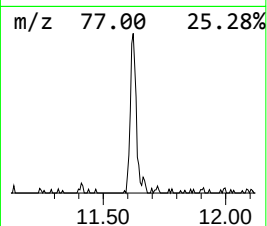
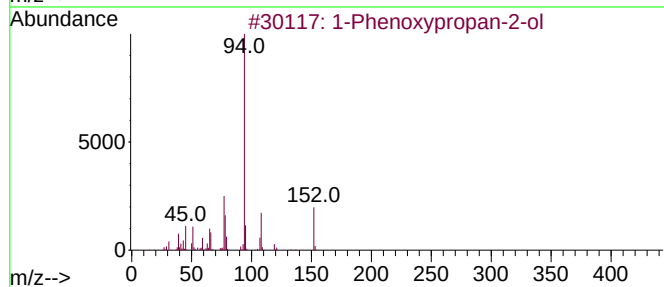
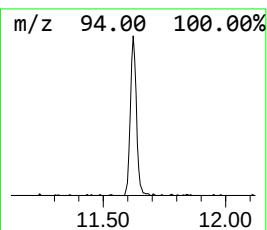
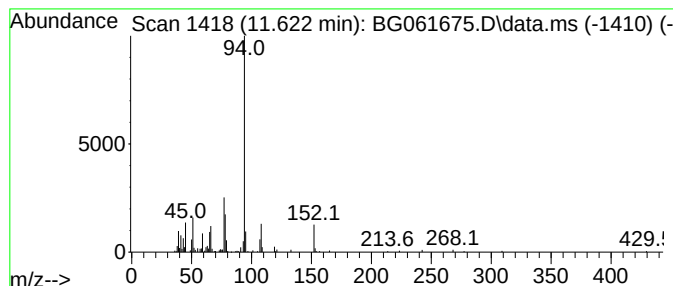
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.622	6.26 ng/ul	246567	Naphthalene-d8	10.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061675.D
Acq On : 24 Jun 2024 14:50
Operator : MA/JU
Sample : P2775-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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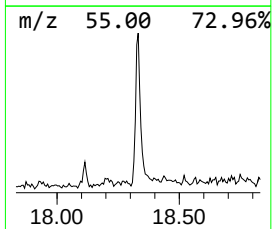
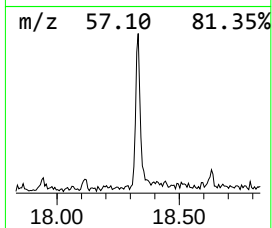
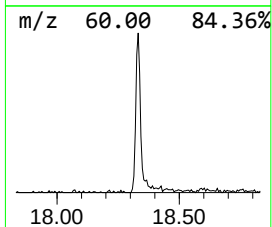
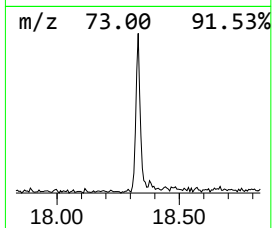
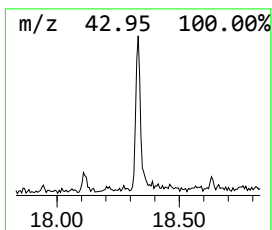
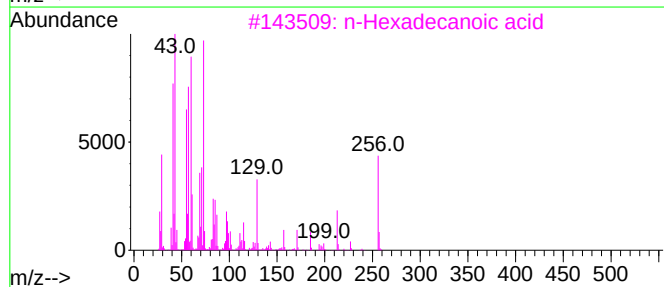
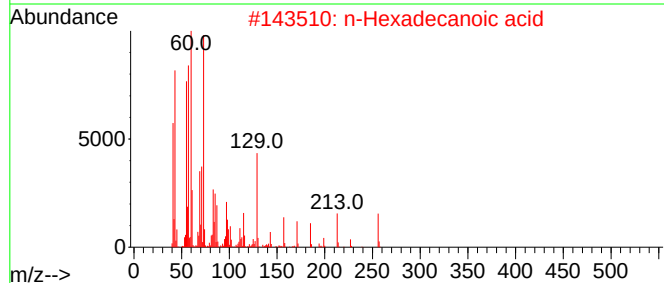
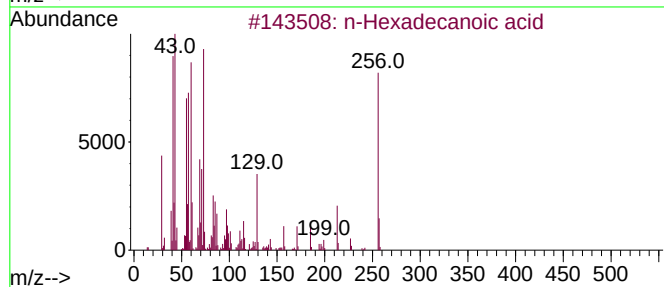
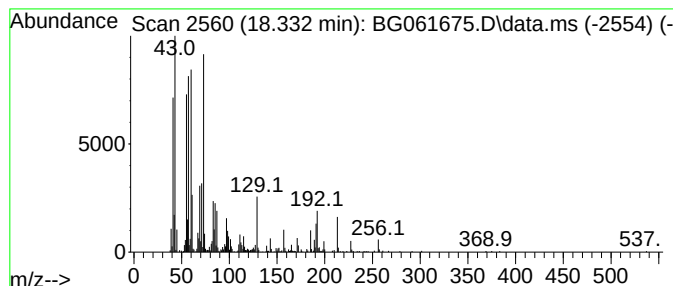
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.331	7.39 ng/ul	575416	Phenanthrene-d10	17.444

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
5			Dodecanoic acid	200	C12H24O2	000143-07-7	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061675.D
Acq On : 24 Jun 2024 14:50
Operator : MA/JU
Sample : P2775-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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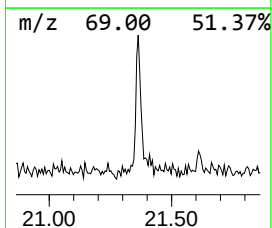
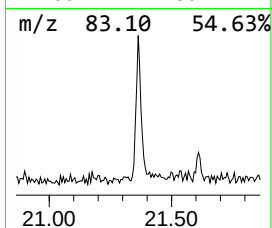
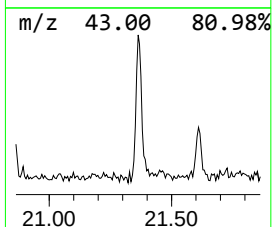
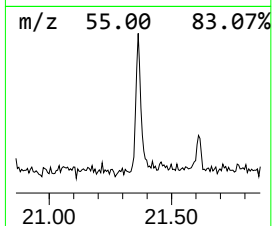
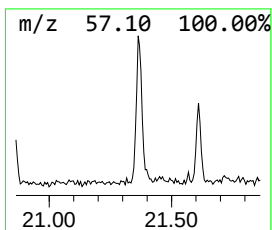
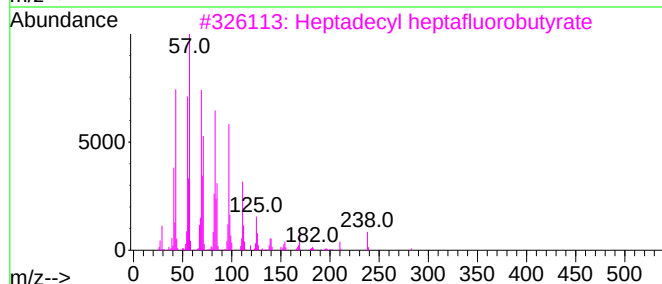
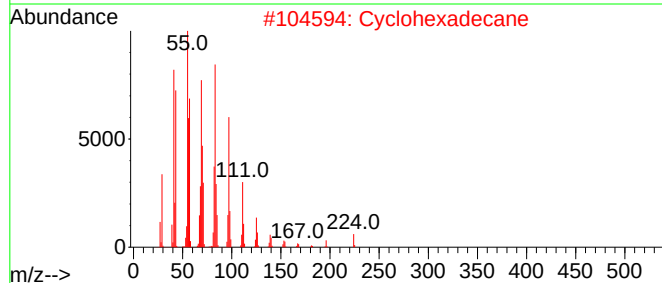
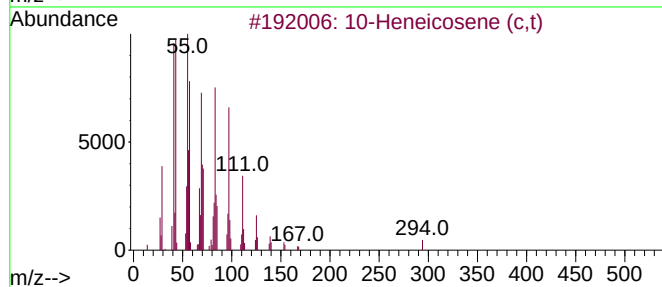
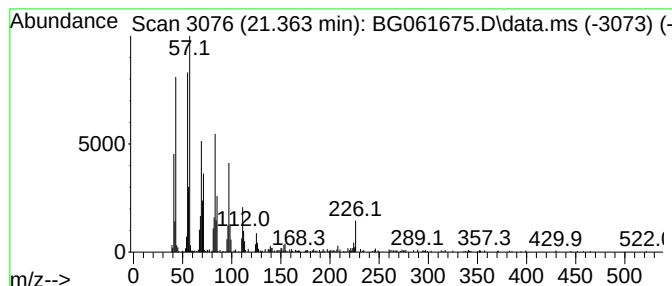
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.363	4.71 ng/ul	488566	Chrysene-d12	21.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Heneicosene (c,t)			294	C21H42	095008-11-0	93
2	Cyclohexadecane			224	C16H32	000295-65-8	91
3	Heptadecyl heptafluorobutyrate			452	C21H35F7O2	959085-66-6	90
4	Eicosyl trifluoroacetate			394	C22H41F3O2	1000351-75-2	90
5	Cyclooctacosane			392	C28H56	000297-24-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061675.D
Acq On : 24 Jun 2024 14:50
Operator : MA/JU
Sample : P2775-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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unknown-02	4.007	2.3	ng/ul	51849	1	8.079	445986	20.0
1,3,5-Cyclohept...	4.236	2.2	ng/ul	49240	1	8.079	445986	20.0
unknown-03	5.153	3.5	ng/ul	78400	1	8.079	445986	20.0
Ethanol, 2-(2-e...	7.662	4.4	ng/ul	97986	1	8.079	445986	20.0
1-Phenoxypropan...	11.622	6.3	ng/ul	246567	2	10.887	787660	20.0
n-Hexadecanoic ...	18.331	7.4	ng/ul	575416	4	17.444	1556690	20.0
(DEL) Alkane: S...	21.363	4.7	ng/ul	488566	5	21.710	2075780	20.0