

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
 Data File : BG062442.D
 Acq On : 16 Aug 2024 13:29
 Operator : MA/JU
 Sample : P3555-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCXV2

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-BG081424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062442.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.422	19	23	30	rVB5	9749	14498	0.62%	0.080%
2	3.681	63	67	73	rBV	35164	59107	2.54%	0.328%
3	3.939	106	111	116	rVB2	15154	22055	0.95%	0.122%
4	4.163	146	149	155	rVB3	4735	7135	0.31%	0.040%
5	4.380	181	186	197	rVB	37506	61115	2.63%	0.339%
6	4.715	238	243	246	rBV4	2717	4689	0.20%	0.026%
7	4.785	250	255	262	rVB	19395	29438	1.27%	0.163%
8	4.956	281	284	291	rVB5	3897	6739	0.29%	0.037%
9	5.067	298	303	309	rBV	9394	14573	0.63%	0.081%
10	5.155	314	318	322	rVB2	5124	5779	0.25%	0.032%
11	5.954	449	454	457	rBV6	2037	3231	0.14%	0.018%
12	6.036	464	468	474	rVB3	3410	6846	0.29%	0.038%
13	6.224	494	500	504	rBV5	6167	9676	0.42%	0.054%
14	7.112	643	651	664	rBV	134428	262003	11.27%	1.454%
15	7.282	674	680	689	rBV	95043	158855	6.83%	0.882%
16	7.487	709	715	721	rVV	120522	207598	8.93%	1.152%
17	7.546	721	725	732	rVV	19751	31648	1.36%	0.176%
18	7.734	753	757	759	rBV4	1452	2350	0.10%	0.013%
19	7.957	788	795	802	rVV	235566	405805	17.45%	2.252%
20	8.022	802	806	810	rVV4	3905	7075	0.30%	0.039%
21	8.328	851	858	862	rVB3	1763	3677	0.16%	0.020%
22	8.651	907	913	920	rBV	145929	251513	10.82%	1.396%
23	8.774	928	934	945	rVV2	30955	61285	2.64%	0.340%
24	8.886	949	953	962	rVB2	3895	7726	0.33%	0.043%
25	9.027	970	977	979	rBV3	3949	6589	0.28%	0.037%
26	9.109	983	991	999	rVV	122436	210715	9.06%	1.169%
27	9.215	1004	1009	1014	rVV3	3059	4173	0.18%	0.023%
28	9.561	1063	1068	1070	rVB3	1788	2640	0.11%	0.015%
29	9.667	1080	1086	1092	rBV2	14715	22655	0.97%	0.126%
30	9.767	1098	1103	1106	rBV4	1992	3060	0.13%	0.017%
31	9.831	1106	1114	1122	rVB	95633	173156	7.45%	0.961%
32	9.966	1130	1137	1148	rBV	14824	32110	1.38%	0.178%
33	10.043	1148	1150	1157	rVB6	2508	5027	0.22%	0.028%
34	10.184	1168	1174	1180	rBV2	9426	14231	0.61%	0.079%
35	10.366	1199	1205	1214	rBV	163512	285461	12.28%	1.584%
36	10.525	1227	1232	1239	rVB3	1314	2997	0.13%	0.017%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BG081424.MA.M
Title : SVOA CALIBRATION

37	10.754	1263	1271	1280	rVB	363778	645803	27.77%	3.584%
38	10.883	1285	1293	1303	rVB3	25041	47850	2.06%	0.266%
39	11.276	1356	1360	1367	rVB5	13478	19539	0.84%	0.108%
40	11.482	1388	1395	1401	rBV	64237	117287	5.04%	0.651%

41	12.334	1536	1540	1545	rVB3	3531	4233	0.18%	0.023%
42	12.845	1624	1627	1632	rBV2	1596	2478	0.11%	0.014%
43	13.192	1682	1686	1690	rVV4	2541	3253	0.14%	0.018%
44	13.415	1720	1724	1729	rVB4	3053	4553	0.20%	0.025%
45	13.473	1732	1734	1739	rBV3	3693	5634	0.24%	0.031%

46	13.979	1814	1820	1826	rBV	306704	439442	18.90%	2.439%
47	14.225	1854	1862	1863	rBV4	5225	7446	0.32%	0.041%
48	14.267	1863	1869	1879	rVV	338663	549615	23.64%	3.050%
49	14.484	1902	1906	1910	rVB4	2404	3118	0.13%	0.017%
50	14.578	1915	1922	1930	rVB	633846	980580	42.17%	5.441%

51	14.801	1946	1960	1974	rBV	950733	2325415	100.00%	12.904%
52	14.913	1974	1979	1984	rVB5	3538	5727	0.25%	0.032%
53	14.983	1987	1991	1992	rBV2	5643	5140	0.22%	0.029%
54	15.083	2002	2008	2012	rBV4	2071	3857	0.17%	0.021%
55	15.377	2054	2058	2062	rVV5	3800	6516	0.28%	0.036%

56	15.430	2064	2067	2071	rVV3	2722	3583	0.15%	0.020%
57	15.565	2084	2090	2098	rVB	480842	715206	30.76%	3.969%
58	15.671	2102	2108	2115	rBV	112467	171267	7.37%	0.950%
59	15.735	2115	2119	2121	rVB4	2486	3676	0.16%	0.020%
60	16.117	2180	2184	2187	rBV3	5192	7803	0.34%	0.043%

61	16.246	2201	2206	2209	rBV5	1843	2782	0.12%	0.015%
62	16.293	2209	2214	2219	rBV4	4457	8549	0.37%	0.047%
63	16.422	2234	2236	2241	rBV5	4671	7137	0.31%	0.040%
64	16.716	2282	2286	2290	rBV3	13441	17034	0.73%	0.095%
65	16.881	2309	2314	2320	rBV8	4476	8450	0.36%	0.047%

66	16.969	2323	2329	2334	rVB4	14323	25205	1.08%	0.140%
67	17.016	2334	2337	2339	rBV4	5017	5209	0.22%	0.029%
68	17.074	2343	2347	2348	rBV3	5374	5813	0.25%	0.032%
69	17.215	2364	2371	2378	rBV	61580	99712	4.29%	0.553%
70	17.315	2378	2388	2394	rVV	788850	1270398	54.63%	7.050%

71	17.356	2394	2395	2398	rVV2	12920	12870	0.55%	0.071%
72	17.415	2398	2405	2414	rVV2	482463	754149	32.43%	4.185%
73	17.486	2414	2417	2424	rVB6	15530	22658	0.97%	0.126%
74	17.821	2471	2474	2477	rVB5	3524	4584	0.20%	0.025%
75	17.944	2490	2495	2499	rVV6	8787	16549	0.71%	0.092%

76	17.991	2499	2503	2508	rVB3	15999	20984	0.90%	0.116%
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 Title : SVOA CALIBRATION

77	18.091	2514	2520	2526	rBV4	29795	60285	2.59%	0.335%
78	18.150	2526	2530	2536	rBV3	30120	40885	1.76%	0.227%
79	18.214	2536	2541	2550	rBV	298246	430751	18.52%	2.390%
80	18.508	2587	2591	2596	rBV4	31515	56191	2.42%	0.312%
81	18.637	2611	2613	2617	rBV5	8474	7749	0.33%	0.043%
82	18.772	2632	2636	2641	rBV4	14144	28008	1.20%	0.155%
83	18.925	2659	2662	2665	rBV4	7676	8663	0.37%	0.048%
84	19.001	2672	2675	2680	rBV2	11039	15294	0.66%	0.085%
85	19.072	2685	2687	2688	rBV2	7008	4636	0.20%	0.026%
86	19.160	2699	2702	2706	rVB3	38593	42497	1.83%	0.236%
87	19.336	2729	2732	2734	rBV4	17831	23407	1.01%	0.130%
88	19.366	2734	2737	2739	rBV2	55255	62835	2.70%	0.349%
89	19.483	2753	2757	2763	rBV2	94544	135847	5.84%	0.754%
90	19.694	2788	2793	2797	rBV	611726	965754	41.53%	5.359%
91	19.730	2797	2799	2810	rVB	188375	304407	13.09%	1.689%
92	21.228	3050	3054	3067	rBV	237132	381481	16.40%	2.117%
93	21.557	3104	3110	3117	rBV2	831075	1319591	56.75%	7.323%
94	22.373	3243	3249	3261	rBV	262967	554641	23.85%	3.078%
95	23.807	3488	3493	3497	rBV2	67430	129955	5.59%	0.721%
96	23.859	3498	3502	3517	rVB4	80382	215375	9.26%	1.195%
97	24.435	3591	3600	3611	rVB2	272764	724444	31.15%	4.020%
98	24.652	3627	3637	3645	rBV	494220	1359133	58.45%	7.542%
99	25.792	3826	3831	3840	rVB2	72954	187916	8.08%	1.043%
100	27.090	4044	4052	4062	rBV2	55963	192620	8.28%	1.069%

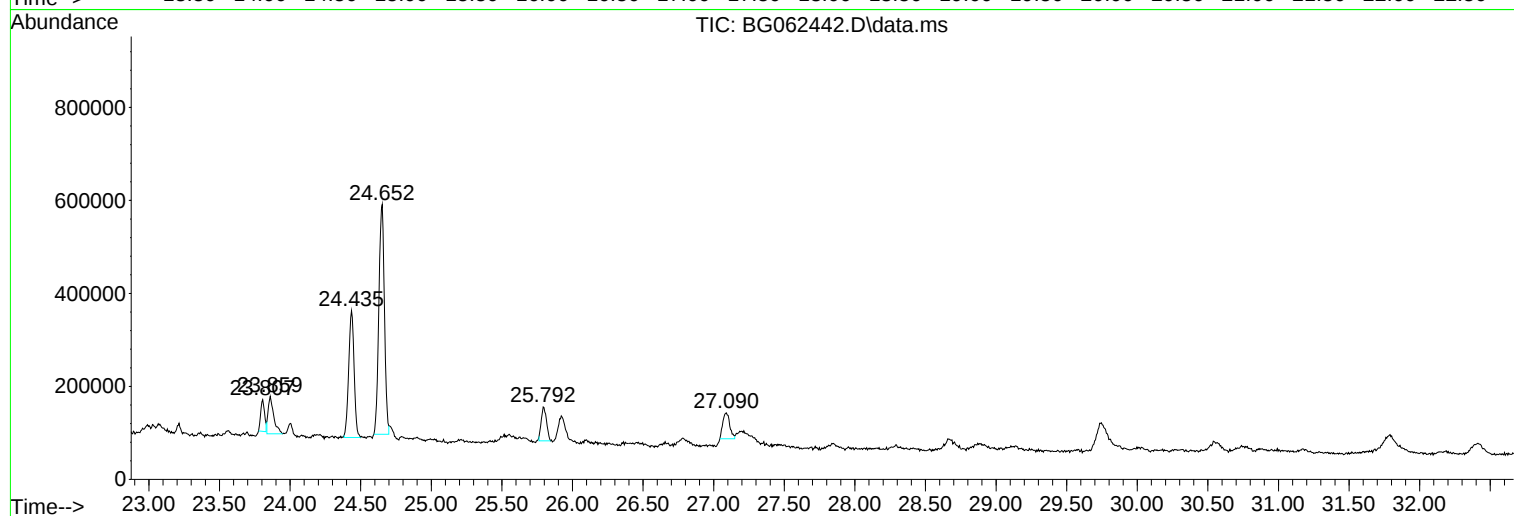
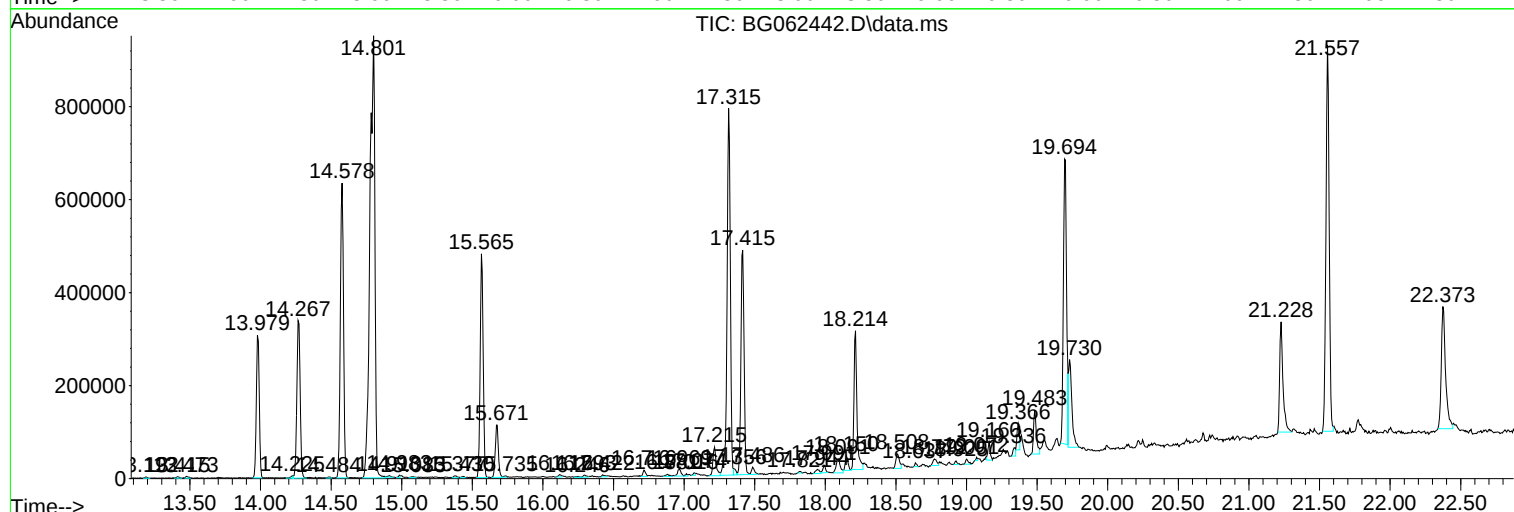
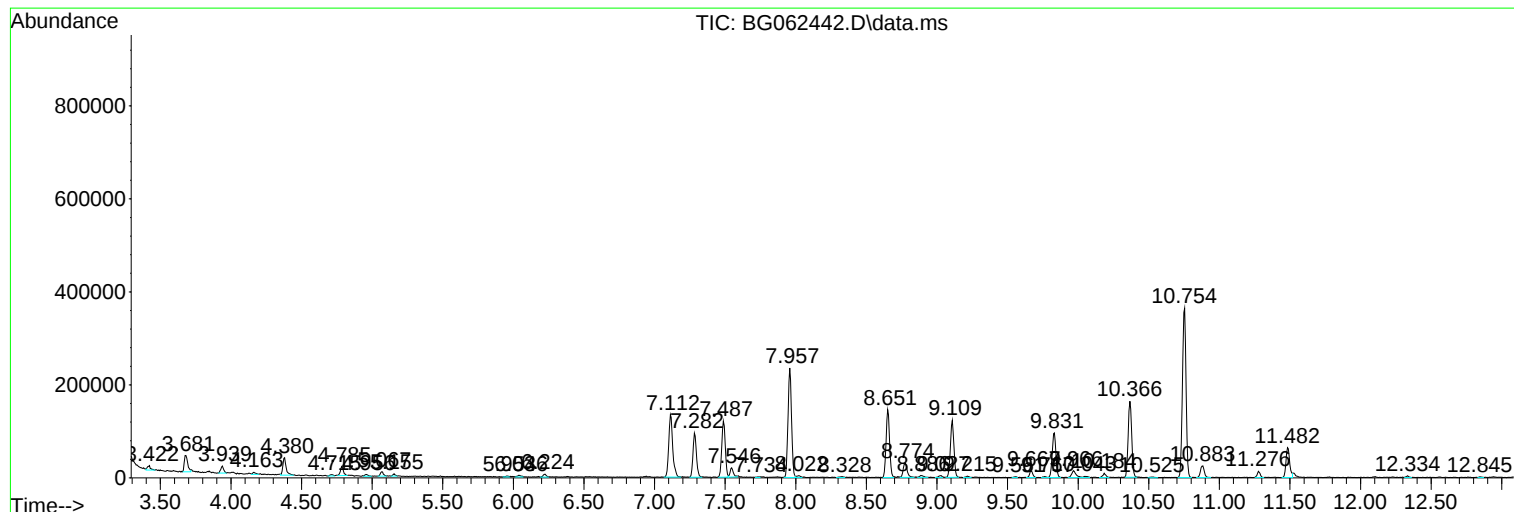
Sum of corrected areas: 18020699

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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCXV2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
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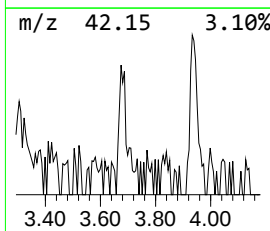
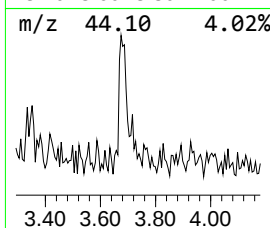
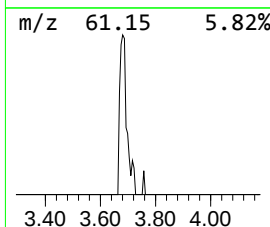
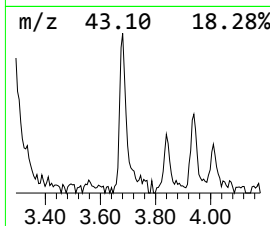
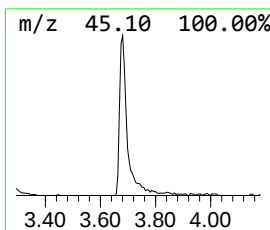
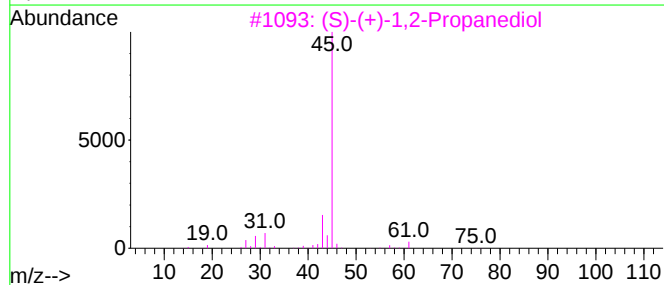
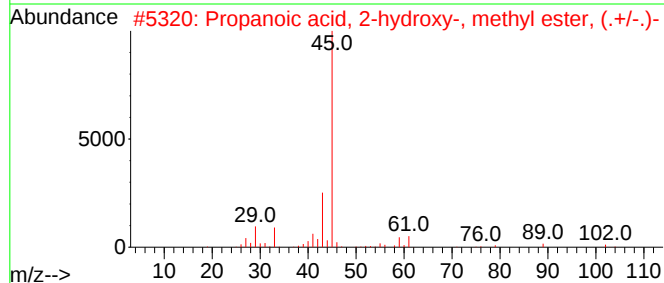
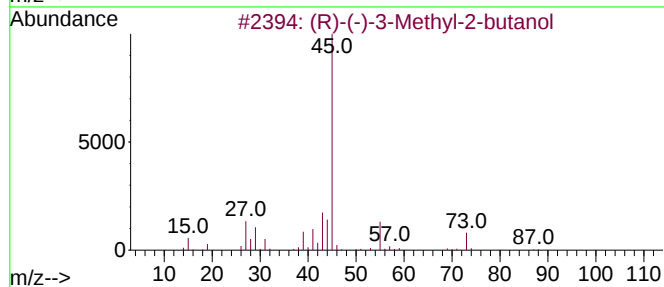
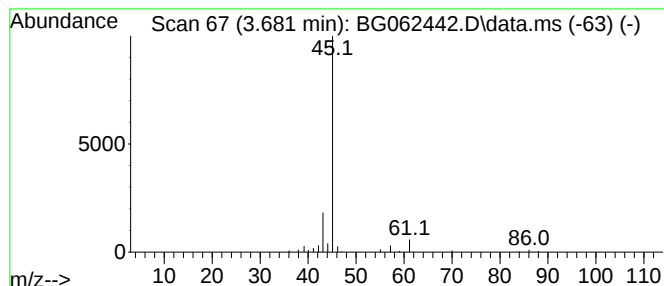
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (R)-(-)-3-Methyl-2-butanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.681	2.91 ng/ul	59107	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(R)-(-)-3-Methyl-2-butanol			88	C5H12O	001572-93-6	64
2	Propanoic acid, 2-hydroxy-, meth...			104	C4H8O3	000547-64-8	39
3	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	39
4	Butane, 1-methoxy-			88	C5H12O	000628-28-4	9
5	1-Butene, 4-methoxy			86	C5H10O	004696-30-4	9



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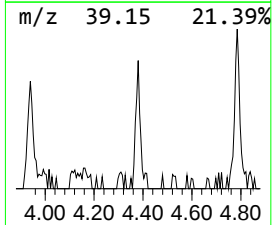
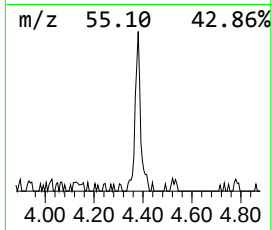
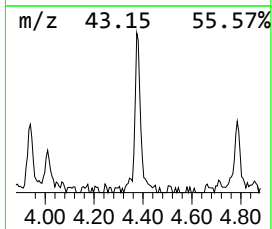
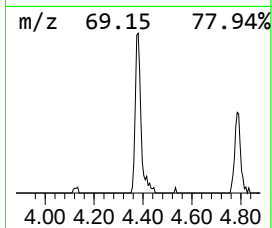
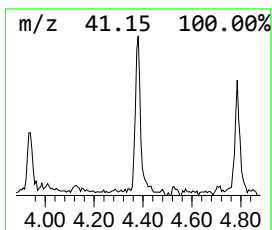
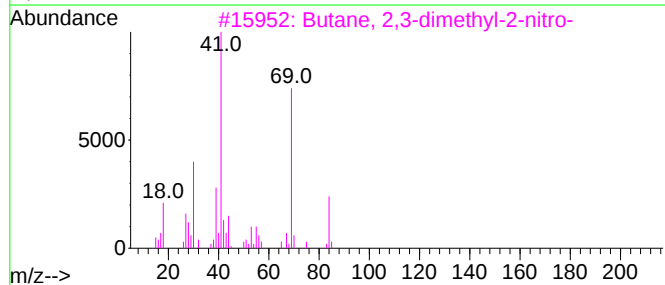
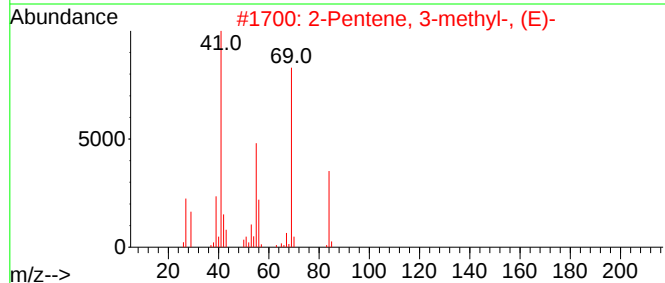
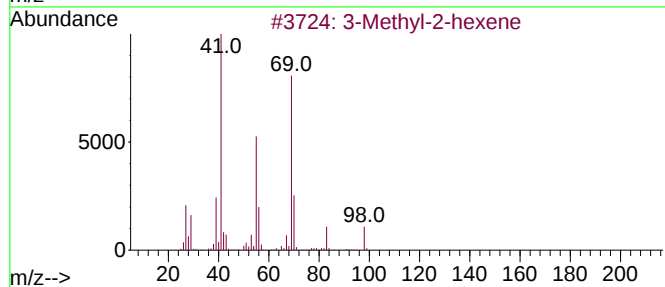
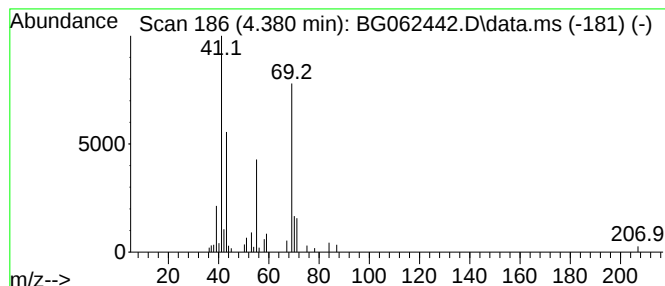
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.380	3.01 ng/ul	61115	1,4-Dichlorobenzene-d4	7.957

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Methyl-2-hexene	98	C7H14	017618-77-8	72
2		2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	64
3		Butane, 2,3-dimethyl-2-nitro-	131	C6H13NO2	034075-28-0	59
4		2-Pentene, 4-methyl-	84	C6H12	004461-48-7	59
5		2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	59



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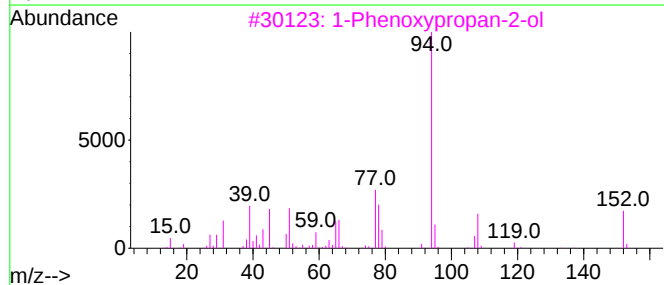
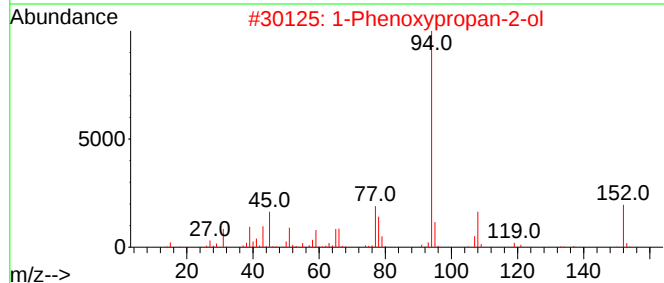
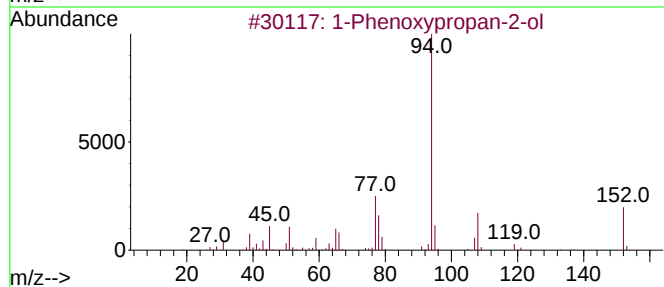
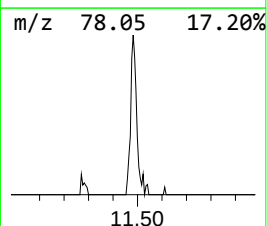
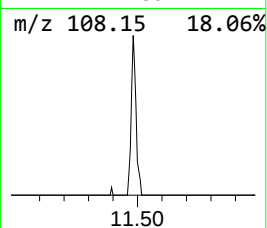
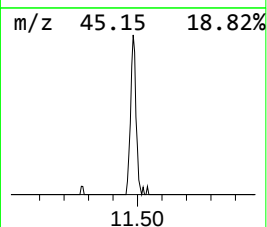
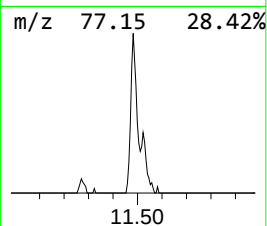
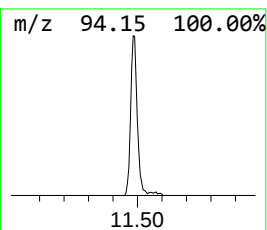
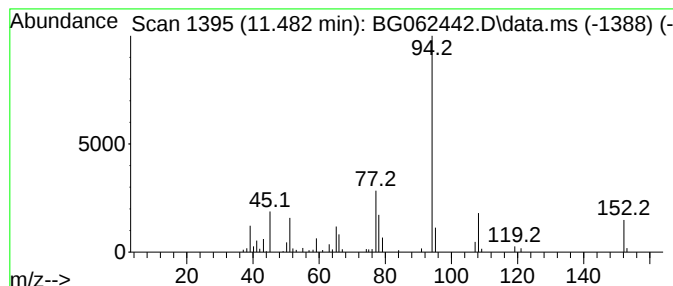
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.482	3.63 ng/ul	117287	Naphthalene-d8	10.754

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	90
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
Data File : BG062442.D
Acq On : 16 Aug 2024 13:29
Operator : MA/JU
Sample : P3555-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

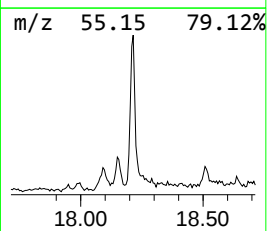
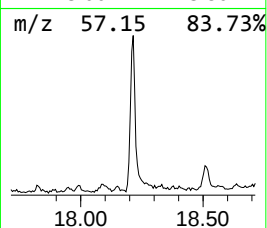
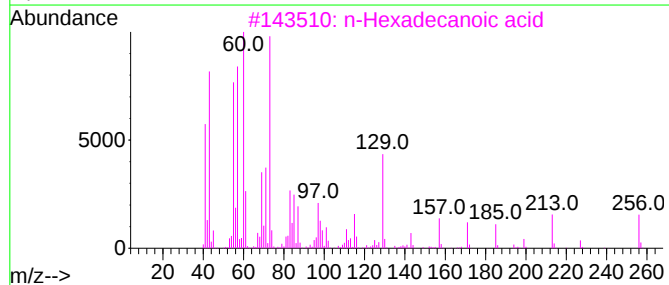
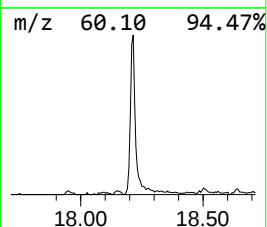
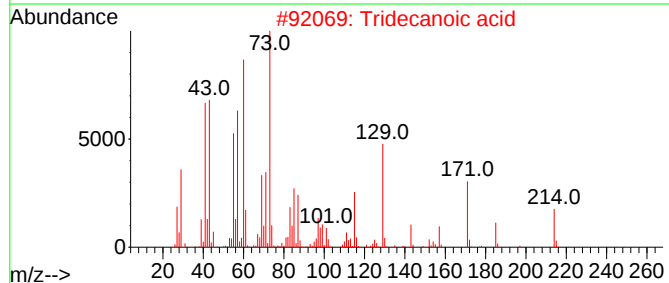
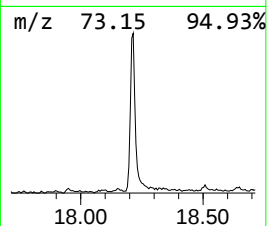
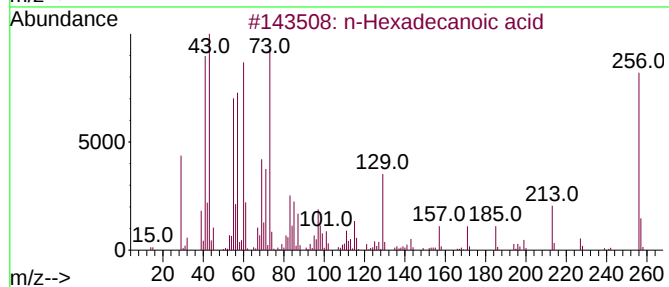
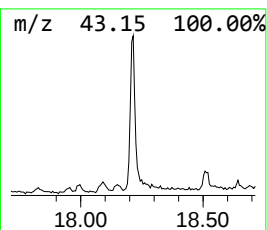
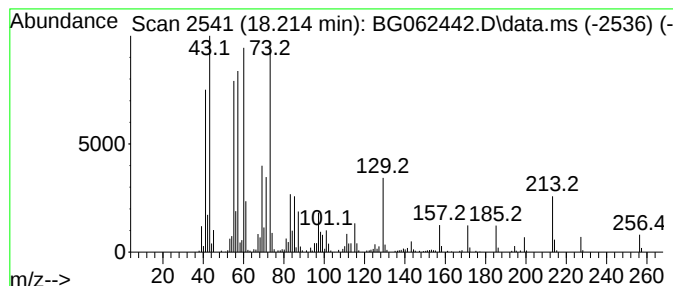
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.214	6.78 ng/ul	430751	Phenanthrene-d10	17.315	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
4	Tridecanoic acid	214	C13H26O2	000638-53-9	90
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
Data File : BG062442.D
Acq On : 16 Aug 2024 13:29
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Sample : P3555-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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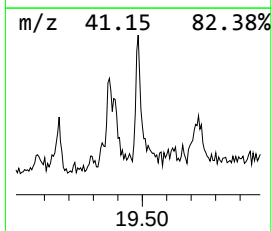
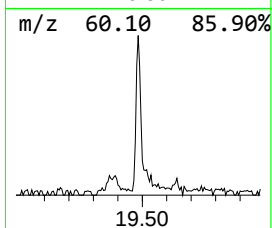
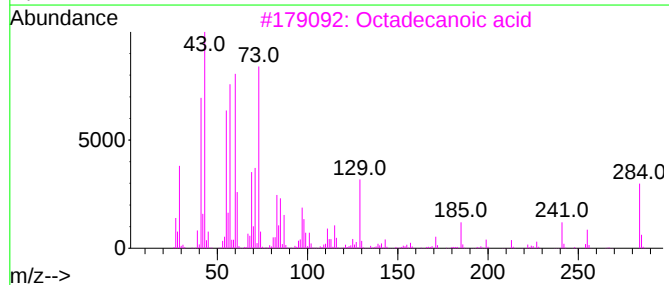
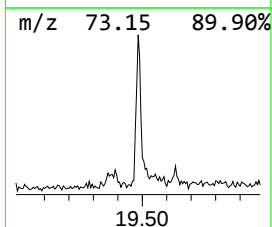
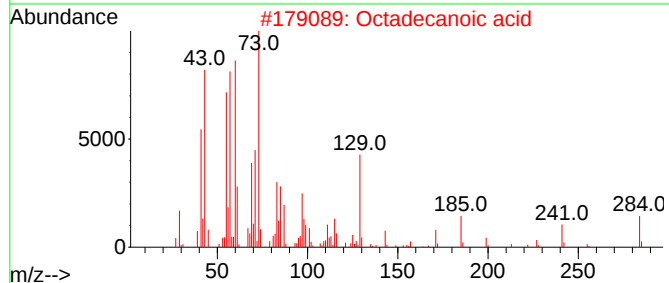
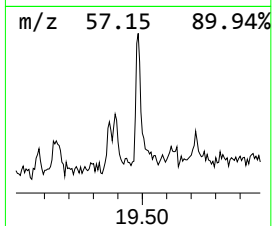
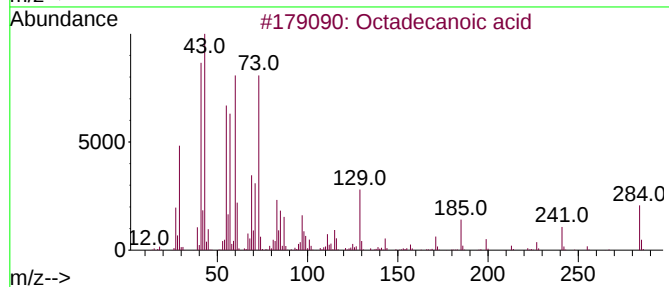
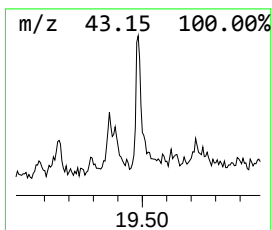
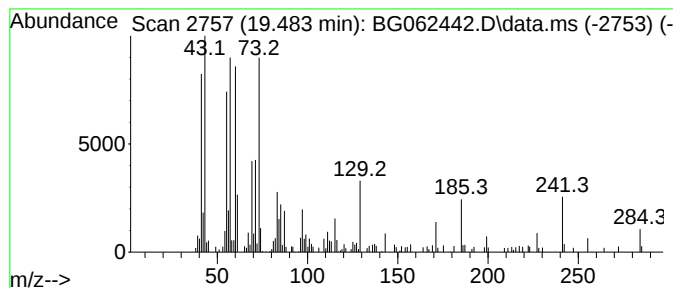
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.483	2.06 ng/ul	135847	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	97
2			Octadecanoic acid	284	C18H36O2	000057-11-4	94
3			Octadecanoic acid	284	C18H36O2	000057-11-4	93
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	80
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
Data File : BG062442.D
Acq On : 16 Aug 2024 13:29
Operator : MA/JU
Sample : P3555-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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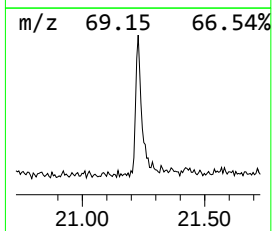
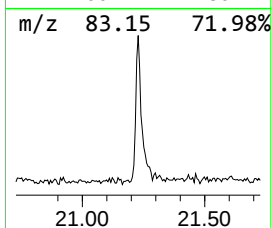
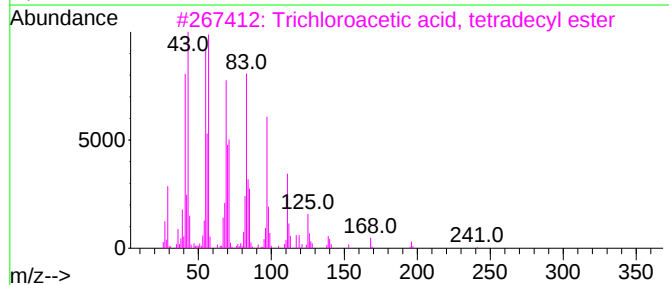
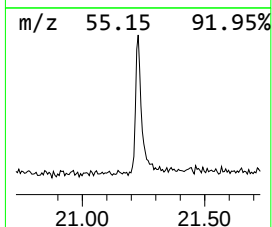
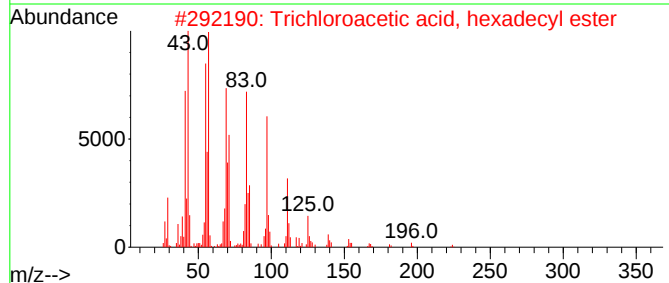
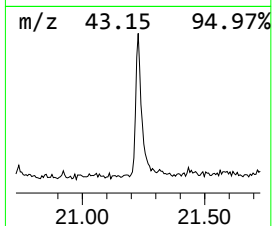
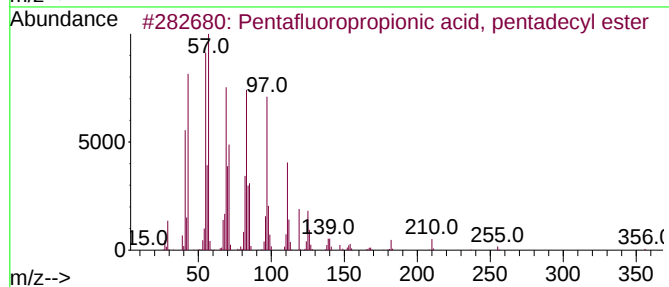
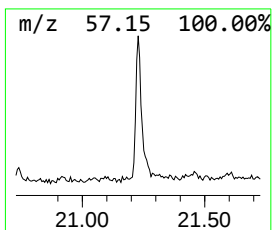
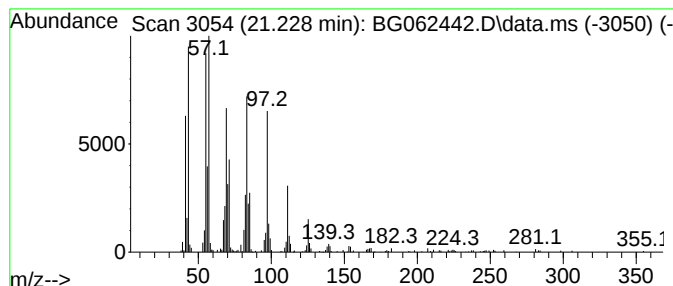
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Pentafluoropropionic acid, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.228	5.78 ng/ul	381481	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
3			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	90
4			1-Tetracosene	336	C24H48	010192-32-2	90
5			1-Heptadecene	238	C17H34	006765-39-5	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
Data File : BG062442.D
Acq On : 16 Aug 2024 13:29
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Sample : P3555-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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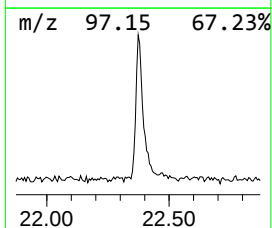
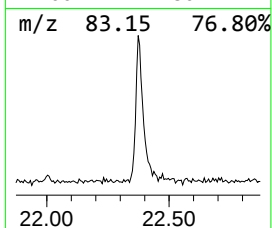
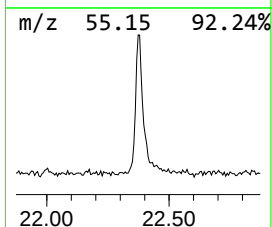
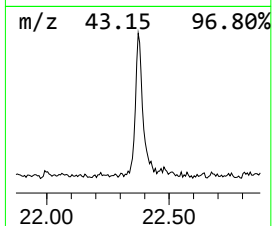
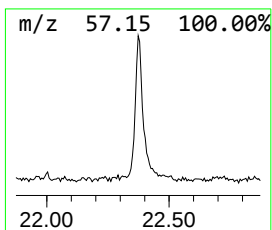
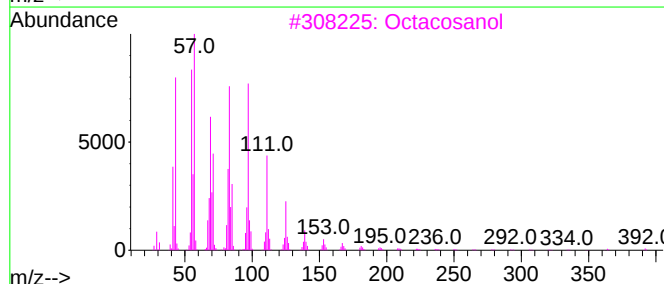
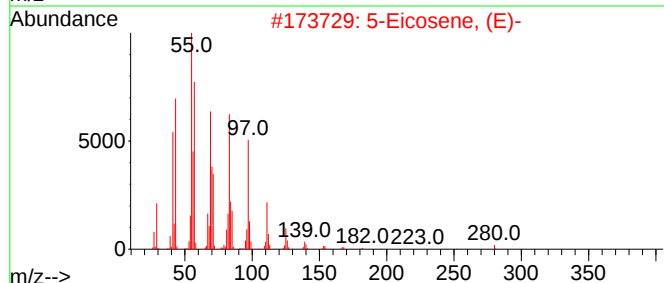
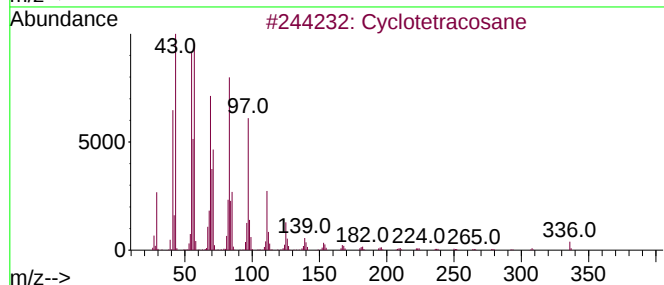
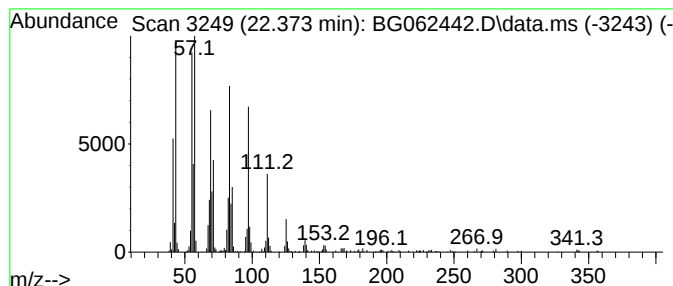
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic22.373 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.373	8.41 ng/ul	554641	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	94
2			5-Eicosene, (E)-	280	C20H40	074685-30-6	91
3			Octacosanol	410	C28H58O	000557-61-9	91
4			1-Hexacosene	364	C26H52	018835-33-1	91
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
Data File : BG062442.D
Acq On : 16 Aug 2024 13:29
Operator : MA/JU
Sample : P3555-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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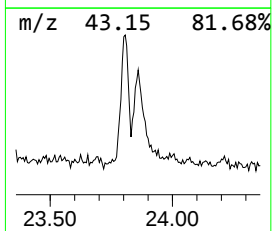
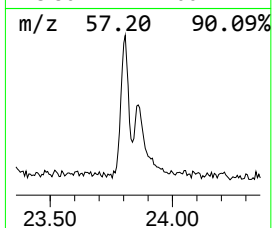
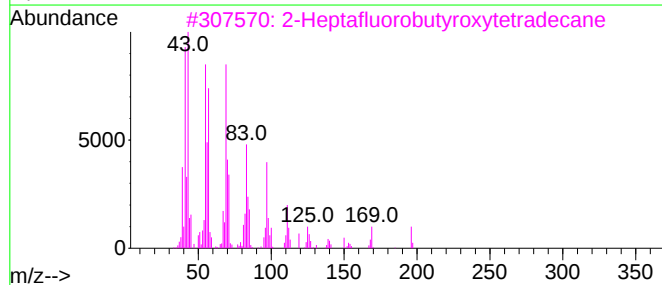
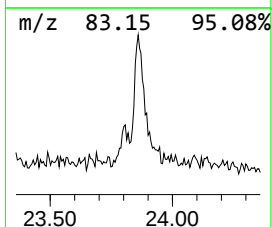
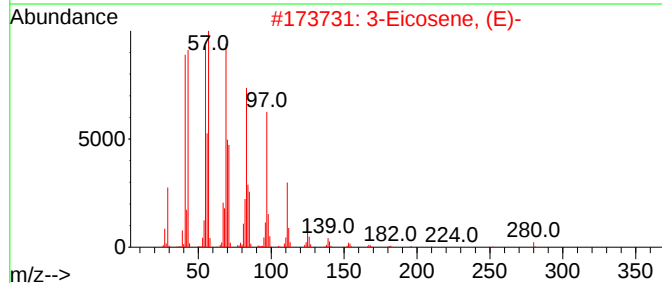
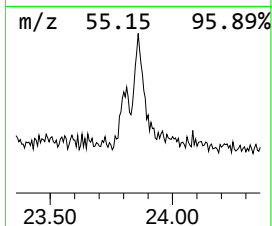
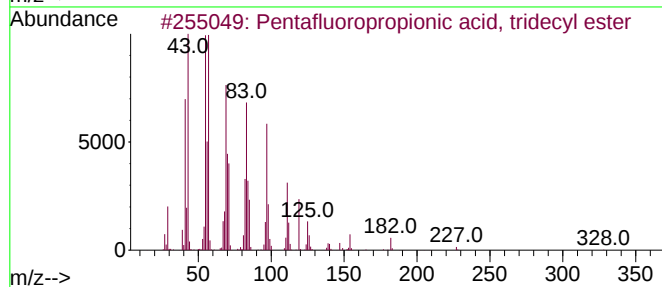
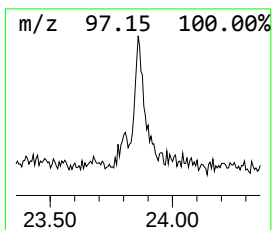
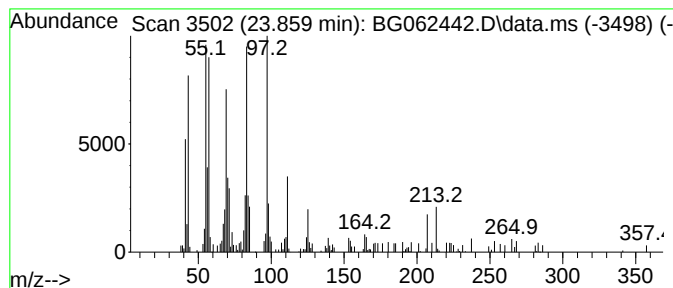
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Pentafluoropropionic acid, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.859	3.17 ng/ul	215375	Perylene-d12	24.653

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, tride...	346	C16H27F5O2	959261-22-4	58
2			3-Eicosene, (E)-	280	C20H40	074685-33-9	52
3			2-Heptafluorobutyroxytetradecane	410	C18H29F7O2	1010245-49-7	49
4			Pentadecafluorooctanoic acid, do...	582	C20H25F15O2	1000406-04-2	46
5			Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
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Sample : P3555-05
Misc :
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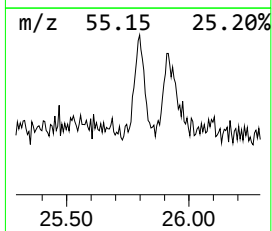
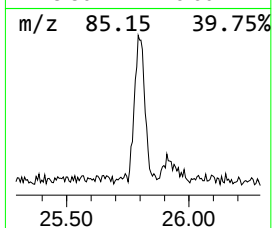
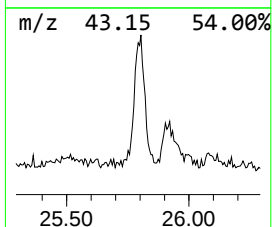
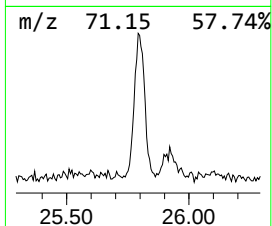
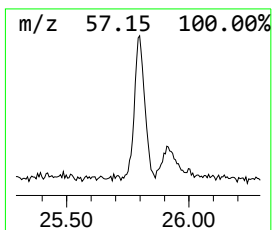
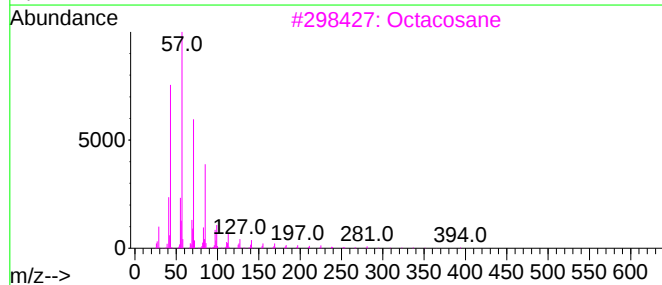
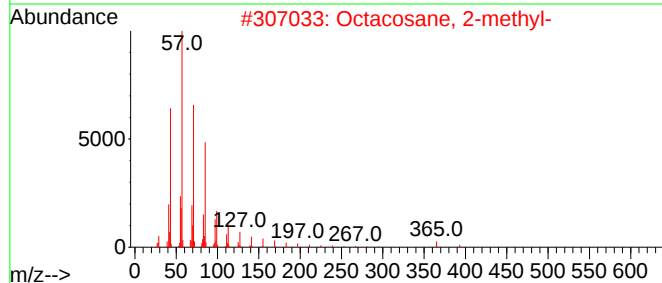
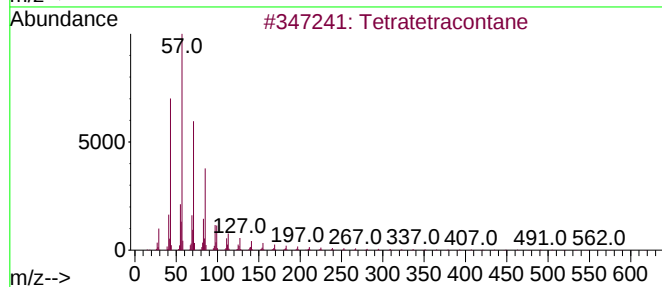
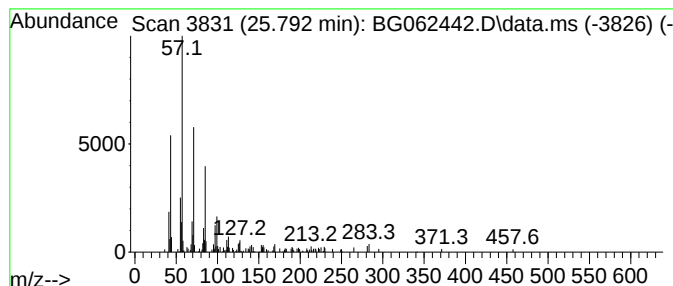
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
25.792	2.77 ng/ul	187916	Perylene-d12		24.653	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetratetracontane		619	C44H90	007098-22-8	87
2	Octacosane, 2-methyl-		408	C29H60	001560-98-1	87
3	Octacosane		394	C28H58	000630-02-4	83
4	Hexacosane		366	C26H54	000630-01-3	83
5	Tetracosane		338	C24H50	000646-31-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
Data File : BG062442.D
Acq On : 16 Aug 2024 13:29
Operator : MA/JU
Sample : P3555-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCXV2

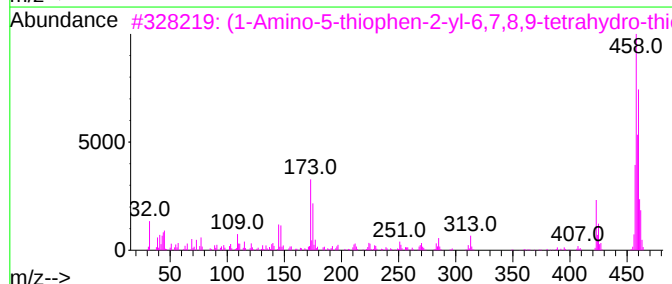
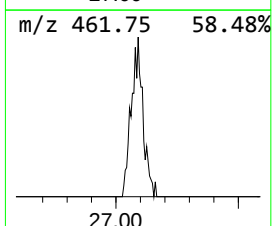
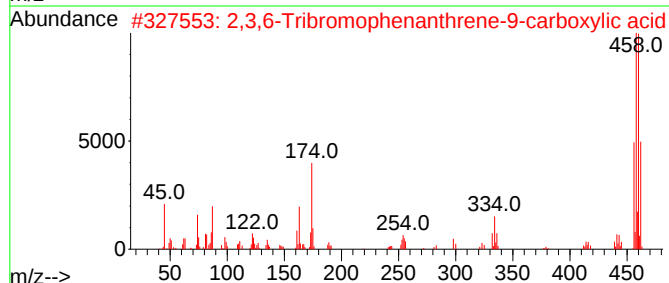
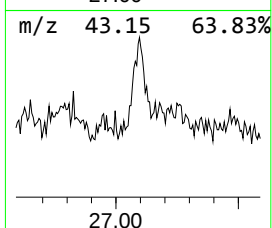
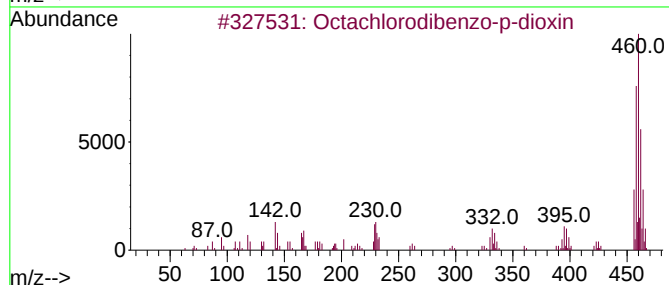
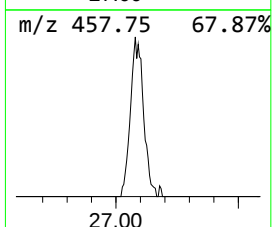
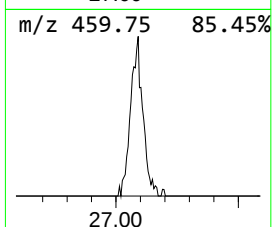
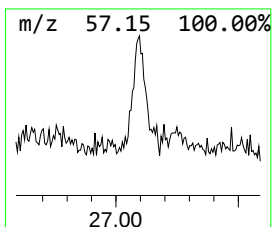
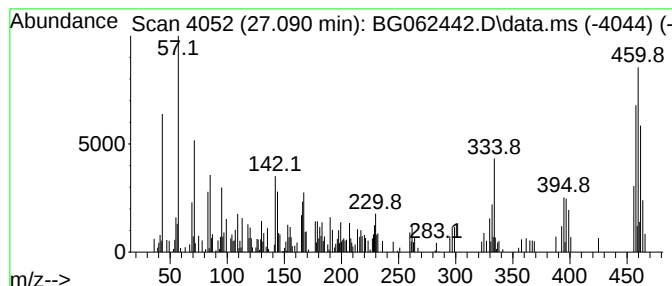
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octachlorodibenzo-p-dioxin Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.090	2.83 ng/ul	192620	Perylene-d12	24.653

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	95
2			2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	43
3			(1-Amino-5-thiophen-2-yl-6,7,8,9...	458	C22H16Cl2N2O5S2	1000318-00-6	22
4			Vanillic acid, 0-heptafluorobuty...	460	C15H10F10O5	1010447-21-1	10
5			Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081524\
Data File : BG062442.D
Acq On : 16 Aug 2024 13:29
Operator : MA/JU
Sample : P3555-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCXV2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.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
(R)-(-)-3-Methy...	3.681	2.9	ng/ul	59107	1	7.957	405805	20.0
(DEL) Alkane: S...	4.380	3.0	ng/ul	61115	1	7.957	405805	20.0
1-Phenoxypropan...	11.482	3.6	ng/ul	117287	2	10.754	645803	20.0
n-Hexadecanoic ...	18.214	6.8	ng/ul	430751	4	17.315	1270400	20.0
Octadecanoic acid	19.483	2.1	ng/ul	135847	5	21.557	1319590	20.0
Pentafluoroprop...	21.228	5.8	ng/ul	381481	5	21.557	1319590	20.0
(DEL) Alkane: C...	22.373	8.4	ng/ul	554641	5	21.557	1319590	20.0
Pentafluoroprop...	23.859	3.2	ng/ul	215375	6	24.653	1359130	20.0
(DEL) Alkane: S...	25.792	2.8	ng/ul	187916	6	24.653	1359130	20.0
Octachlorodiben...	27.090	2.8	ng/ul	192620	6	24.653	1359130	20.0