

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
 Data File : BG062527.D
 Acq On : 22 Aug 2024 4:12
 Operator : MA/JU
 Sample : P3626-18
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCXZ7

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: BG062527.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	16	20	24	rVB2	12628	15558	0.99%	0.068%
2	3.664	60	64	76	rVB	69895	110006	7.00%	0.481%
3	3.811	84	89	100	rBV	165342	264759	16.85%	1.156%
4	4.151	143	147	150	rBV4	4310	7207	0.46%	0.031%
5	5.050	295	300	304	rBV5	3530	4864	0.31%	0.021%
6	6.019	460	465	473	rVB2	7661	12395	0.79%	0.054%
7	7.100	641	649	658	rBV	249452	437728	27.86%	1.912%
8	7.265	666	677	685	rBV	170455	288918	18.39%	1.262%
9	7.470	705	712	718	rBV	237571	389859	24.82%	1.703%
10	7.523	718	721	731	rVB	22484	36982	2.35%	0.162%
11	7.940	783	792	799	rBV	265365	450148	28.66%	1.966%
12	7.999	799	802	807	rVV5	6271	9151	0.58%	0.040%
13	8.633	898	910	919	rBV	246406	423234	26.94%	1.849%
14	9.086	979	987	996	rBV	222007	383955	24.44%	1.677%
15	9.538	1059	1064	1069	rVB3	2055	3410	0.22%	0.015%
16	9.644	1077	1082	1089	rBV	28303	46968	2.99%	0.205%
17	9.808	1103	1110	1119	rBV	187407	320285	20.39%	1.399%
18	10.349	1194	1202	1213	rBV	307475	537124	34.19%	2.346%
19	10.731	1257	1267	1277	rBV	409815	709211	45.15%	3.098%
20	10.860	1280	1289	1297	rBV	230151	409353	26.06%	1.788%
21	11.259	1350	1357	1362	rBV3	12550	21869	1.39%	0.096%
22	11.465	1385	1392	1398	rBV	69008	126847	8.07%	0.554%
23	12.311	1532	1536	1543	rVB3	4504	8016	0.51%	0.035%
24	13.392	1715	1720	1723	rBV4	3506	4948	0.31%	0.022%
25	13.462	1729	1732	1739	rBV4	5887	9903	0.63%	0.043%
26	13.961	1811	1817	1824	rVB	566956	790276	50.31%	3.452%
27	14.202	1855	1858	1859	rBV2	6738	5910	0.38%	0.026%
28	14.249	1859	1866	1878	rVB2	642374	990486	63.05%	4.327%
29	14.461	1899	1902	1908	rVB2	3421	4455	0.28%	0.019%
30	14.555	1911	1918	1929	rVB	690167	1077193	68.57%	4.705%
31	14.743	1940	1950	1961	rBV3	242884	468051	29.79%	2.044%
32	14.895	1972	1976	1983	rVB6	4086	7405	0.47%	0.032%
33	14.972	1985	1989	1992	rBV4	3955	6007	0.38%	0.026%
34	15.348	2051	2053	2060	rVB5	5292	8444	0.54%	0.037%
35	15.412	2060	2064	2067	rBV2	5577	7376	0.47%	0.032%
36	15.548	2080	2087	2096	rVB	851036	1267532	80.69%	5.537%

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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

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
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37	15.653	2099	2105	2114	rBV	224853	326960	20.81%	1.428%
38	15.788	2124	2128	2132	rBV3	3752	6713	0.43%	0.029%
39	15.924	2145	2151	2154	rVB3	2025	3450	0.22%	0.015%
40	16.106	2177	2182	2189	rBV7	4886	10340	0.66%	0.045%
41	16.270	2207	2210	2214	rBV4	6000	7226	0.46%	0.032%
42	16.699	2279	2283	2288	rBV4	5142	7805	0.50%	0.034%
43	16.858	2304	2310	2313	rBV5	5219	7536	0.48%	0.033%
44	17.016	2333	2337	2339	rVB4	3461	3612	0.23%	0.016%
45	17.063	2342	2345	2347	rBV4	4977	6141	0.39%	0.027%
46	17.192	2363	2367	2374	rBV3	20511	32026	2.04%	0.140%
47	17.257	2374	2378	2379	rVV	11939	14683	0.93%	0.064%
48	17.298	2379	2385	2394	rVV	850141	1266435	80.62%	5.532%
49	17.398	2394	2402	2410	rVV	819756	1274717	81.15%	5.568%
50	17.468	2411	2414	2417	rVB4	3504	4231	0.27%	0.018%
51	17.797	2467	2470	2475	rBV2	4139	6515	0.41%	0.028%
52	17.974	2496	2500	2503	rVB2	9807	11702	0.74%	0.051%
53	18.074	2510	2517	2521	rBV7	8720	18114	1.15%	0.079%
54	18.132	2521	2527	2532	rBV5	12037	21850	1.39%	0.095%
55	18.191	2532	2537	2547	rBV2	179447	268888	17.12%	1.175%
56	18.491	2585	2588	2590	rBV2	8458	10486	0.67%	0.046%
57	18.667	2616	2618	2622	rVB4	4518	4249	0.27%	0.019%
58	18.861	2647	2651	2653	rBV4	4568	5412	0.34%	0.024%
59	18.919	2658	2661	2662	rVB2	4433	3654	0.23%	0.016%
60	19.060	2682	2685	2690	rBV5	3475	5802	0.37%	0.025%
61	19.119	2690	2695	2700	rBV4	9738	18417	1.17%	0.080%
62	19.178	2701	2705	2710	rVV5	5087	9117	0.58%	0.040%
63	19.243	2712	2716	2721	rVB3	12961	15729	1.00%	0.069%
64	19.313	2723	2728	2731	rBV2	23693	38298	2.44%	0.167%
65	19.342	2731	2733	2735	rVV3	11633	13497	0.86%	0.059%
66	19.366	2735	2737	2741	rVV4	12847	19103	1.22%	0.083%
67	19.401	2741	2743	2747	rVB5	8639	8525	0.54%	0.037%
68	19.460	2749	2753	2765	rVV	47099	69477	4.42%	0.303%
69	19.613	2776	2779	2784	rVB6	8535	8930	0.57%	0.039%
70	19.677	2784	2790	2796	rBV	1098631	1540176	98.04%	6.728%
71	19.789	2805	2809	2812	rVB6	5190	7511	0.48%	0.033%
72	19.977	2839	2841	2847	rVB7	6410	8300	0.53%	0.036%
73	20.059	2852	2855	2858	rVV3	16735	20787	1.32%	0.091%
74	20.124	2862	2866	2872	rVB	48759	62577	3.98%	0.273%
75	20.200	2875	2879	2880	rVV3	16692	19077	1.21%	0.083%
76	20.224	2881	2883	2888	rVB3	22073	26292	1.67%	0.115%

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77	20.447	2919	2921	2923	rVB3	5488	5107	0.33%	0.022%
78	20.506	2928	2931	2933	rBV4	6800	8056	0.51%	0.035%
79	20.553	2936	2939	2944	rVV4	17387	22831	1.45%	0.100%
80	20.605	2944	2948	2951	rVV6	10140	17920	1.14%	0.078%
81	20.652	2952	2956	2960	rVV	58190	75868	4.83%	0.331%
82	20.717	2963	2967	2971	rVB5	14754	19382	1.23%	0.085%
83	21.210	3046	3051	3061	rBV	199489	367540	23.40%	1.605%
84	21.440	3087	3090	3096	rVB3	13735	19294	1.23%	0.084%
85	21.534	3100	3106	3119	rVV2	858058	1351328	86.02%	5.903%
86	21.739	3138	3141	3150	rVB4	18901	34702	2.21%	0.152%
87	21.939	3172	3175	3178	rBV	26542	35816	2.28%	0.156%
88	22.350	3236	3245	3262	rBV2	297811	823323	52.41%	3.596%
89	22.990	3350	3354	3358	rVB3	22814	33803	2.15%	0.148%
90	23.319	3405	3410	3417	rVB5	25848	46858	2.98%	0.205%
91	23.754	3478	3484	3490	rBV2	81362	168117	10.70%	0.734%
92	23.830	3491	3497	3507	rVB7	38157	113074	7.20%	0.494%
93	24.394	3583	3593	3609	rVB	595215	1570906	100.00%	6.862%
94	24.612	3619	3630	3644	rBV	557654	1525058	97.08%	6.662%
95	25.152	3717	3722	3728	rVB8	24462	49472	3.15%	0.216%
96	25.728	3814	3820	3831	rVB	64697	169769	10.81%	0.742%
97	25.857	3837	3842	3843	rBV5	20699	32551	2.07%	0.142%
98	29.681	4482	4493	4510	rBV5	52666	271187	17.26%	1.185%
99	30.797	4674	4683	4695	rBV10	57042	251068	15.98%	1.097%
100	31.702	4824	4837	4865	rVB10	157248	1000089	63.66%	4.368%

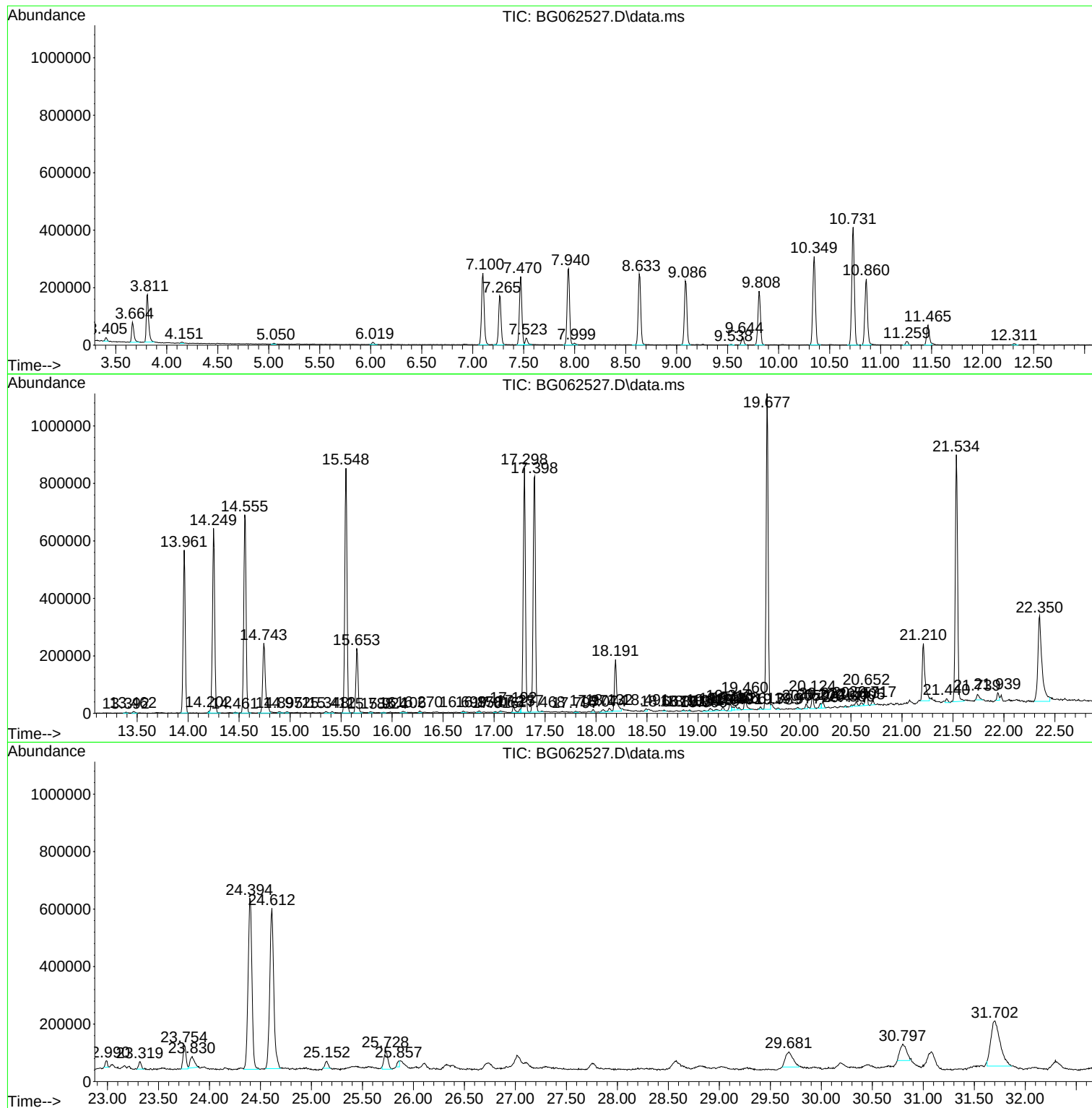
Sum of corrected areas: 22893312

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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



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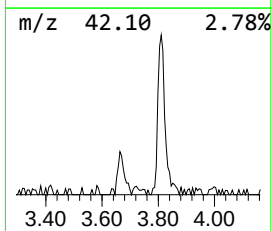
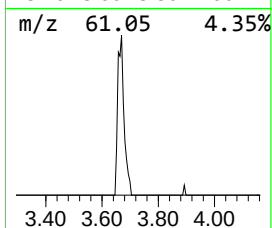
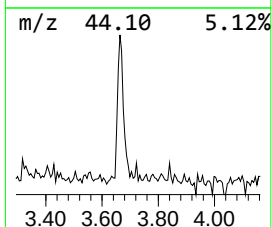
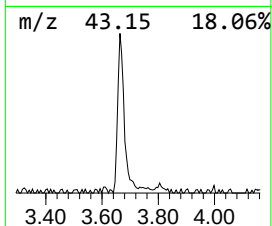
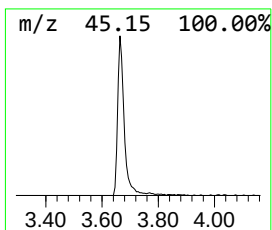
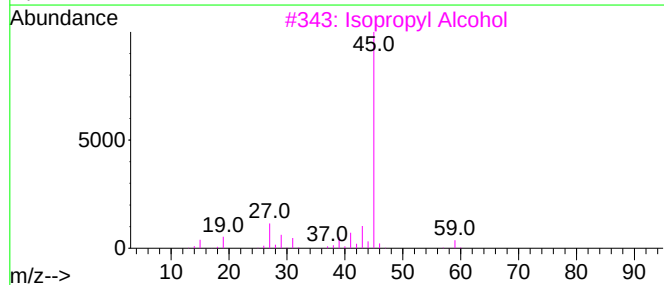
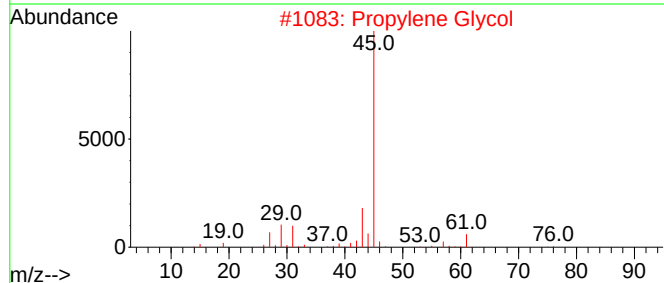
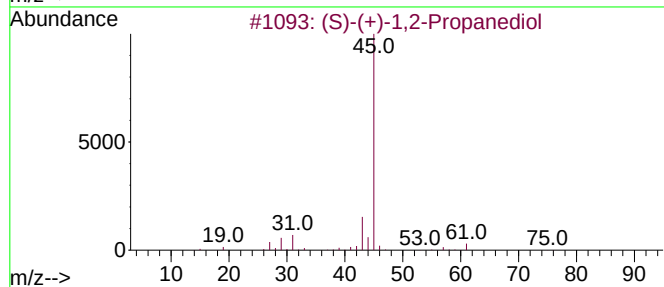
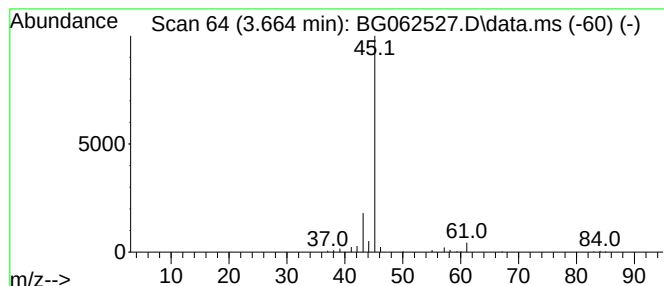
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (S)-(+)-1,2-Propanediol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.664	4.89 ng/ul	110006	1,4-Dichlorobenzene-d4	7.940

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol		76	C3H8O2	004254-15-3	64
2	Propylene Glycol		76	C3H8O2	000057-55-6	43
3	Isopropyl Alcohol		60	C3H8O	000067-63-0	9
4	Propylene Glycol		76	C3H8O2	000057-55-6	9
5	Formamide		45	CH3NO	000075-12-7	7



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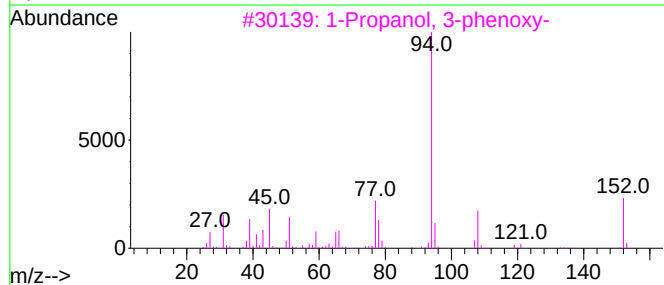
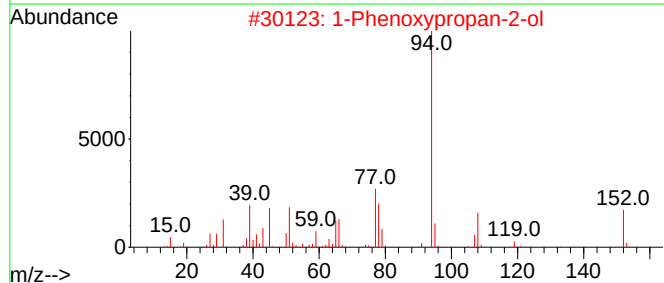
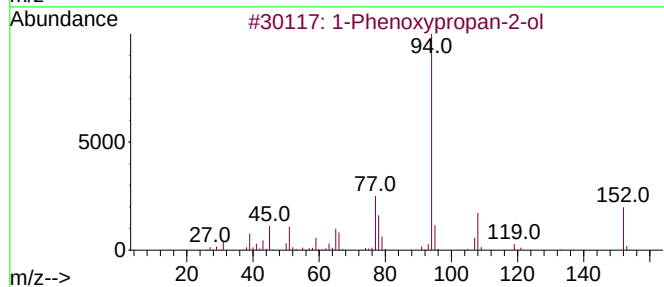
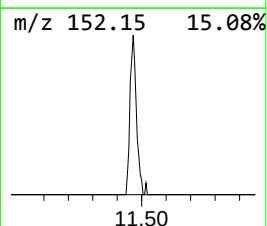
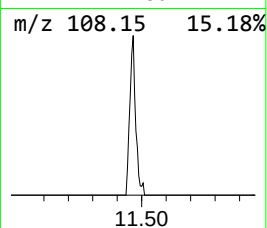
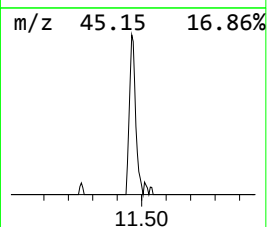
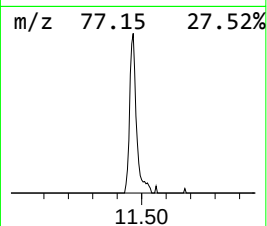
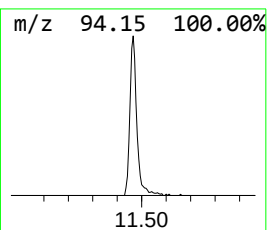
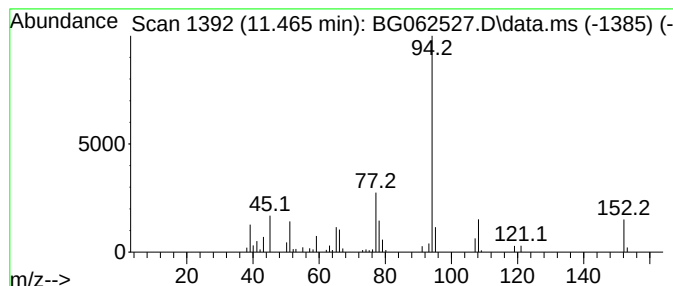
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.465	3.58 ng/ul	126847	Naphthalene-d8	10.730

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



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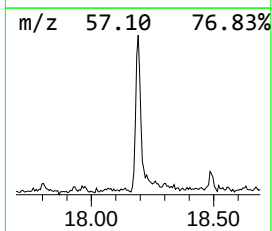
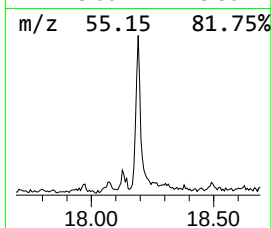
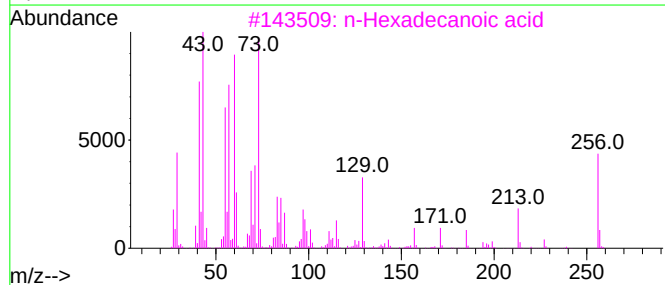
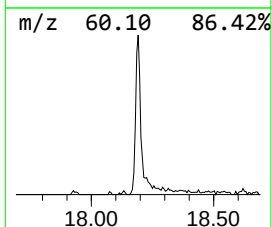
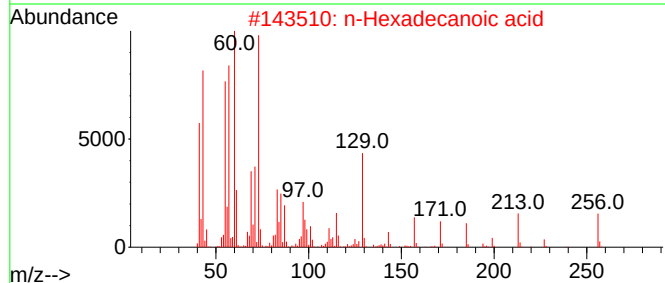
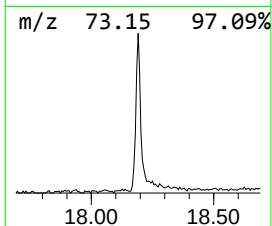
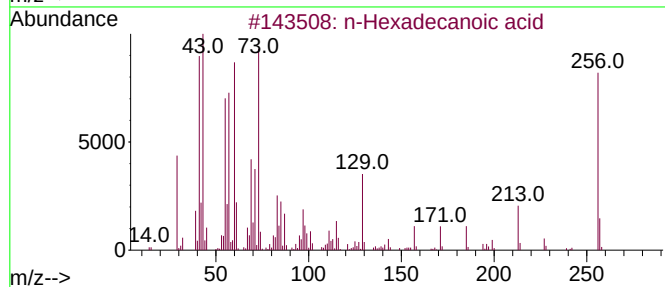
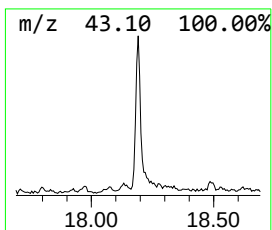
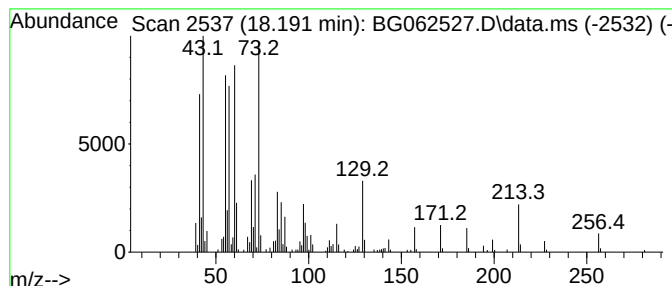
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.191	4.25 ng/ul	268888	Phenanthrene-d10	17.298

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	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4			Tridecanoic acid	214	C13H26O2	000638-53-9	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	93



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Acq On : 22 Aug 2024 4:12
Operator : MA/JU
Sample : P3626-18
Misc :
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Instrument :
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ClientSampleId :
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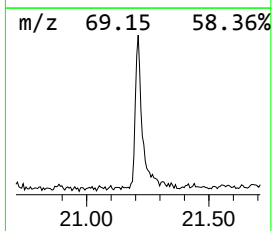
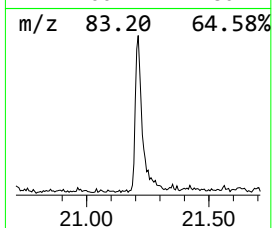
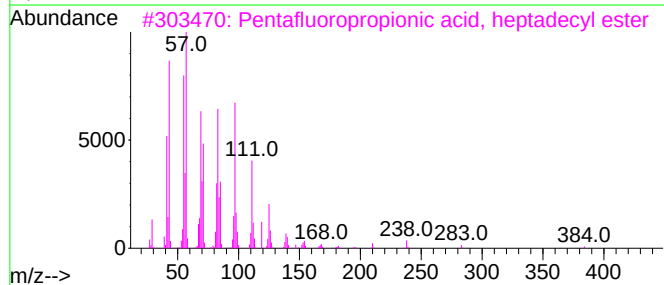
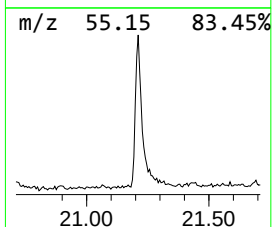
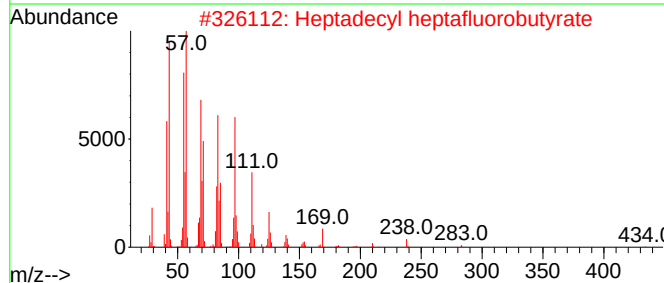
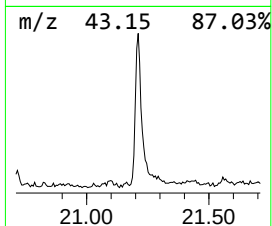
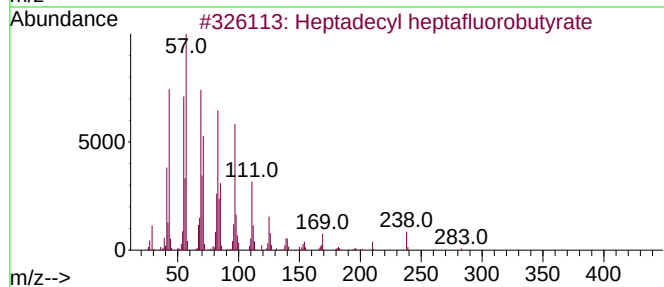
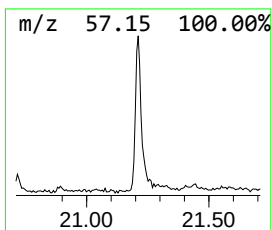
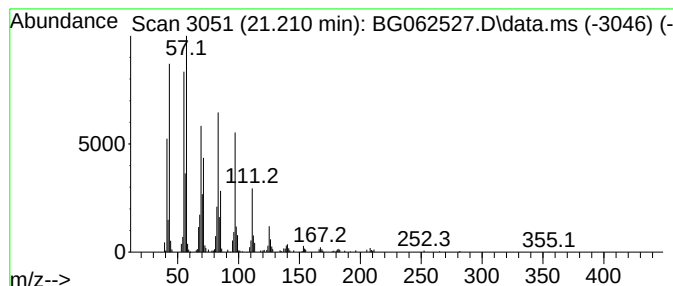
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Heptadecyl heptafluorobutyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.211	5.44 ng/ul	367540	Chrysene-d12	21.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
4			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91
5			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062527.D
Acq On : 22 Aug 2024 4:12
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Sample : P3626-18
Misc :
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Instrument :
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ClientSampleId :
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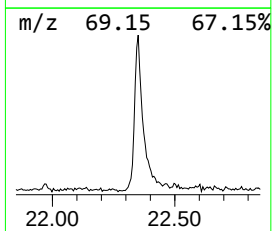
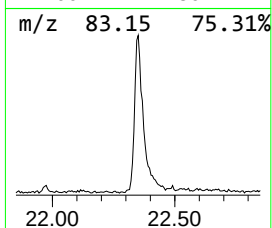
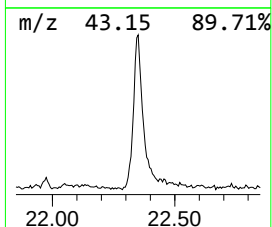
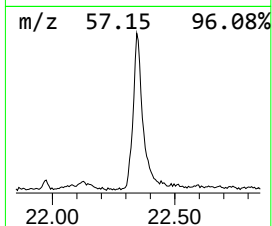
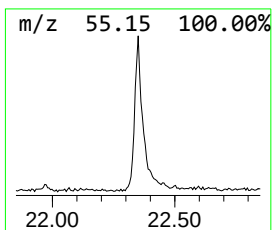
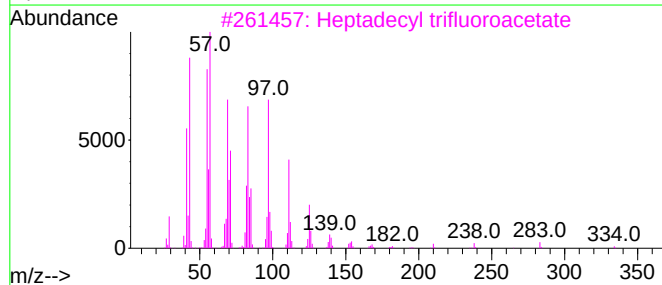
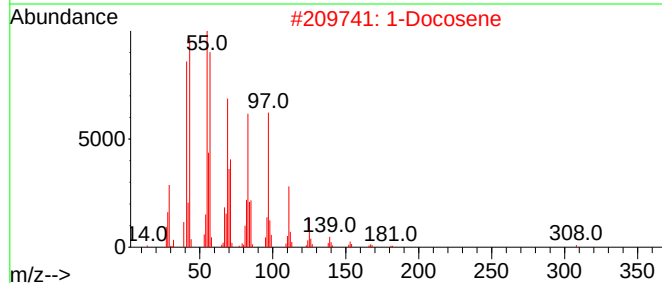
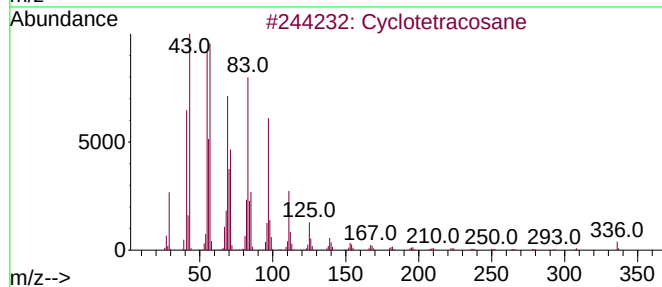
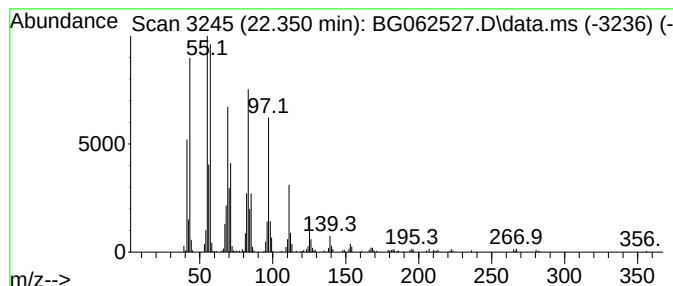
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic22.350 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.350	12.19 ng/ul	823323	Chrysene-d12	21.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	95
2			1-Docosene	308	C22H44	001599-67-3	94
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	94
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
5			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062527.D
Acq On : 22 Aug 2024 4:12
Operator : MA/JU
Sample : P3626-18
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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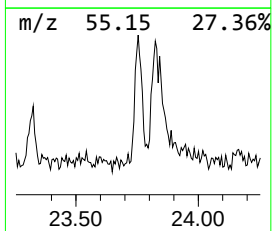
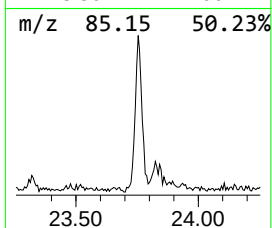
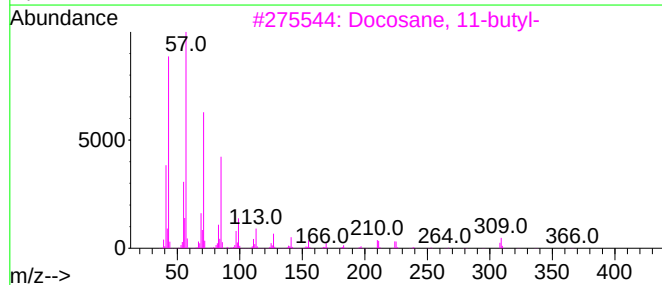
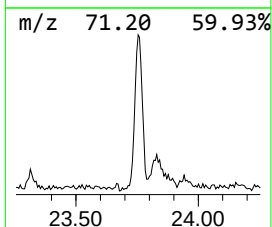
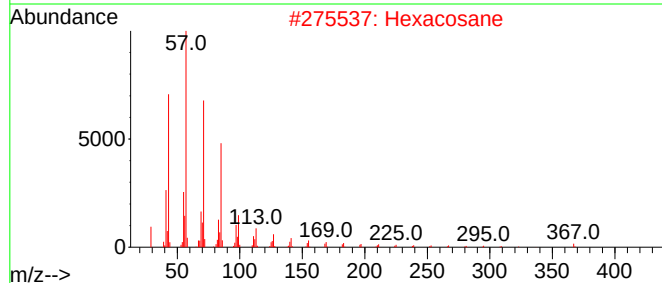
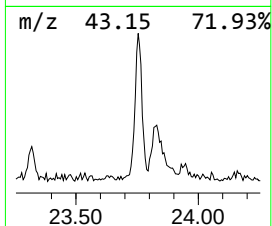
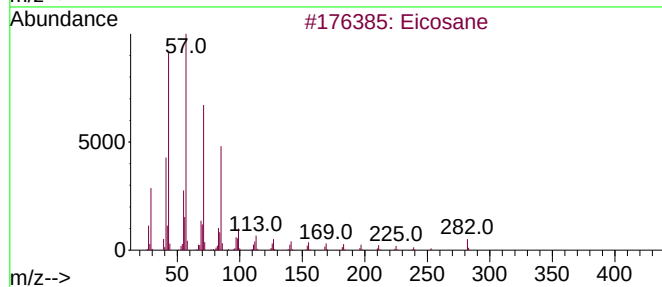
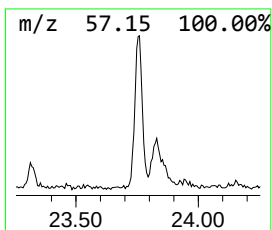
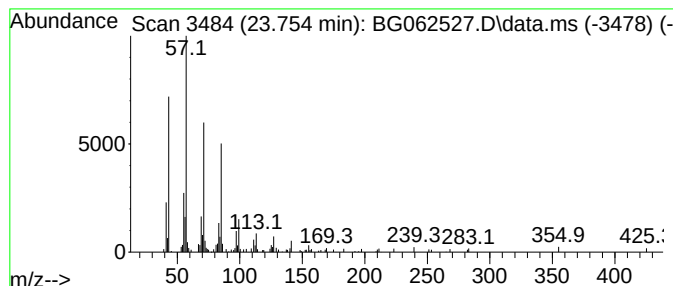
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.754	2.20 ng/ul	168117	Perylene-d12	24.612

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	90
2		Hexacosane	366	C26H54	000630-01-3	87
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	87
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062527.D
Acq On : 22 Aug 2024 4:12
Operator : MA/JU
Sample : P3626-18
Misc :
ALS Vial : 29 Sample Multiplier: 1

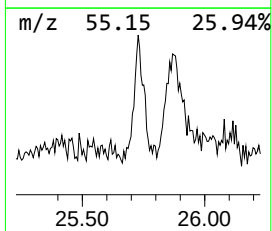
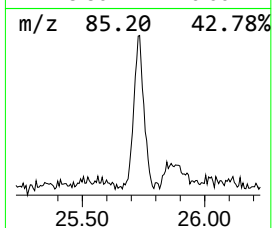
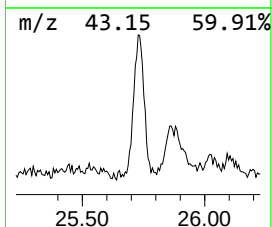
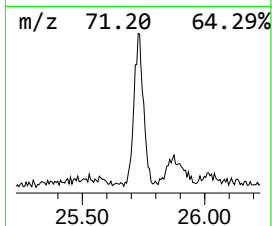
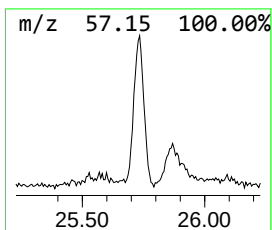
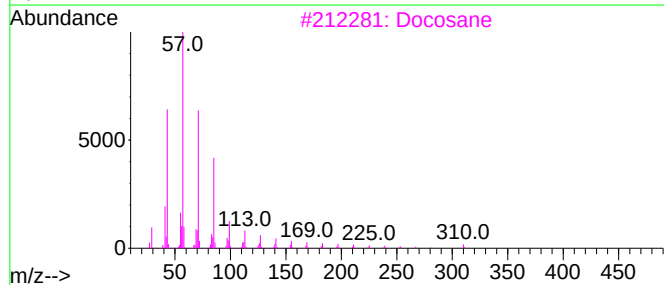
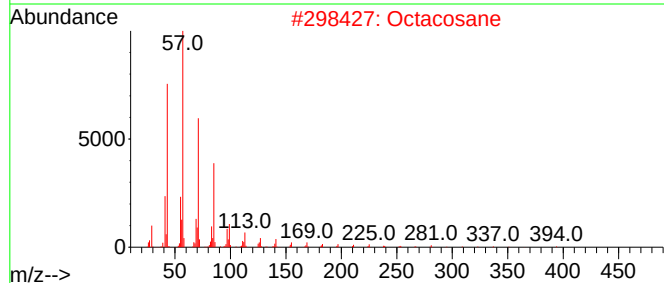
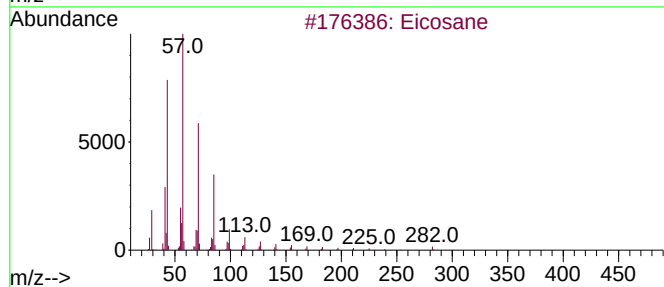
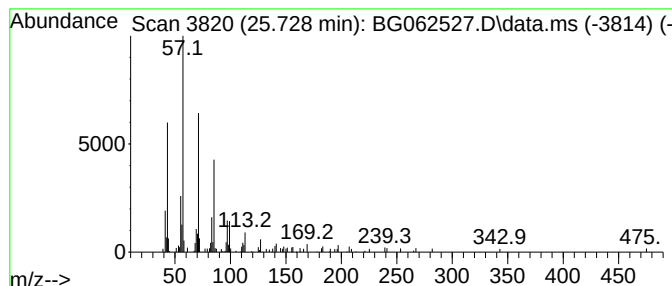
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.728	2.23 ng/ul	169769	Perylene-d12	24.612	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Octacosane	394	C28H58	000630-02-4	90
3	Docosane	310	C22H46	000629-97-0	87
4	Tetracosane	338	C24H50	000646-31-1	87
5	Hexatriacontane	507	C36H74	000630-06-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062527.D
Acq On : 22 Aug 2024 4:12
Operator : MA/JU
Sample : P3626-18
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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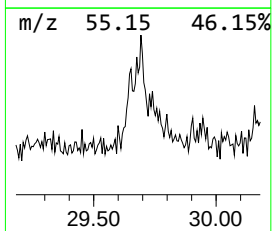
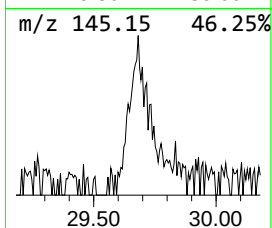
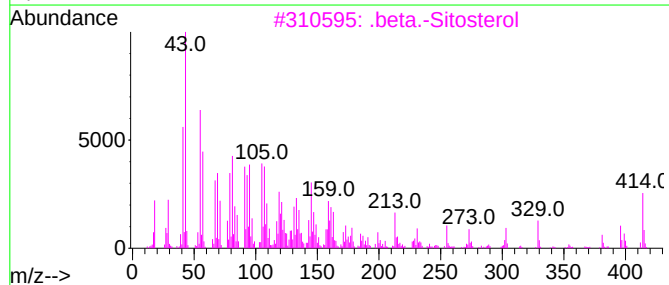
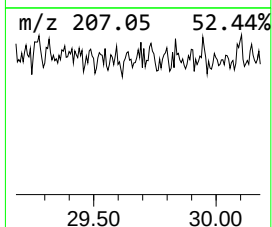
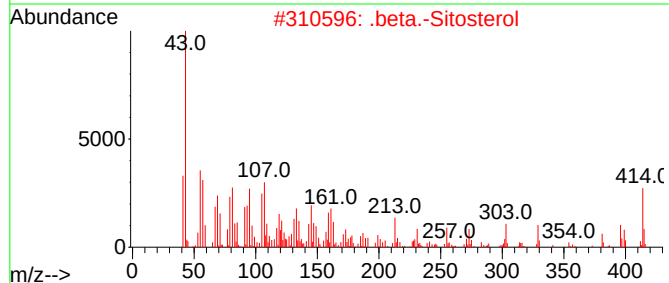
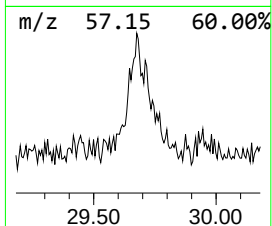
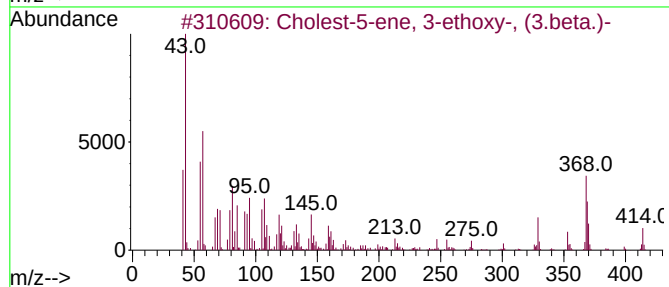
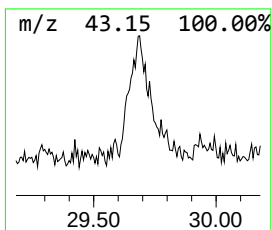
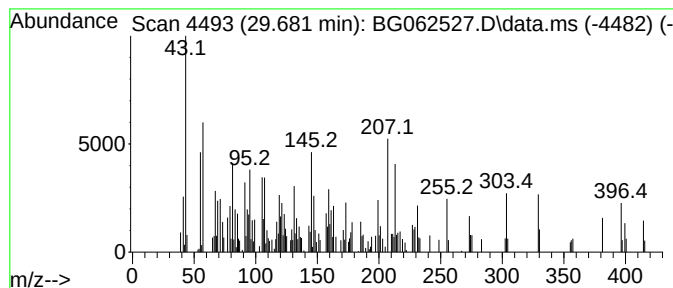
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.681	3.56 ng/ul	271187	Perylene-d12	24.612

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	30
2			.beta.-Sitosterol	414	C29H50O	000083-46-5	22
3			.beta.-Sitosterol	414	C29H50O	000083-46-5	12
4			(3S,8S,9S,10R,13R,14S,17R)-17-((...	414	C29H50O	029365-29-5	11
5			.gamma.-Sitosterol	414	C29H50O	000083-47-6	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062527.D
Acq On : 22 Aug 2024 4:12
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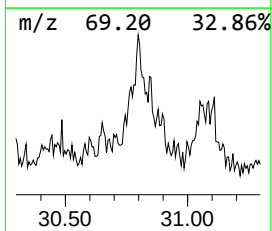
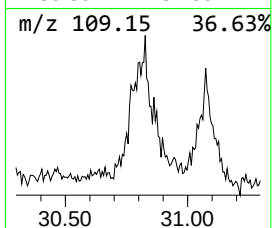
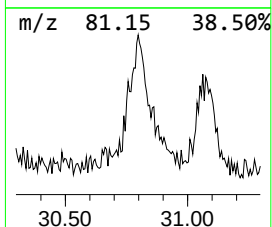
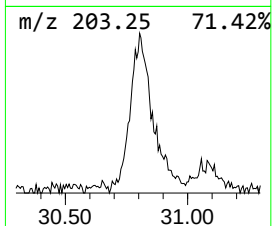
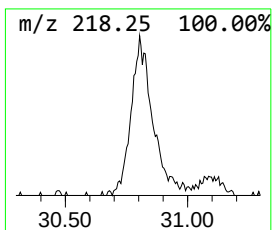
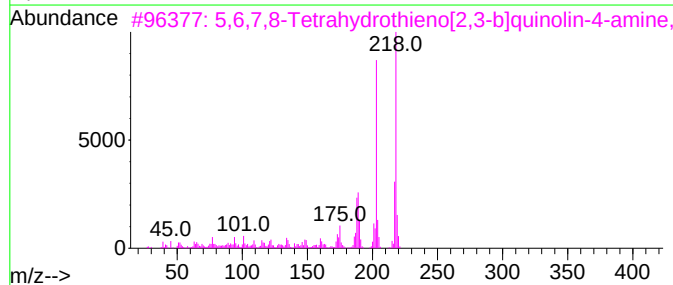
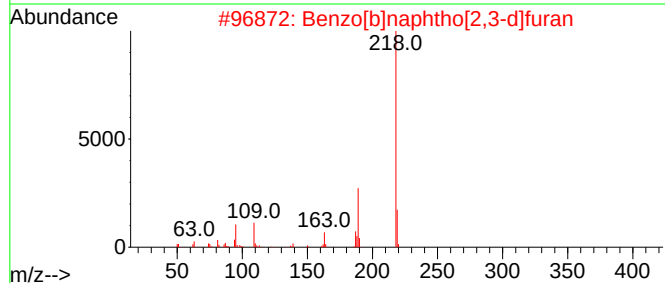
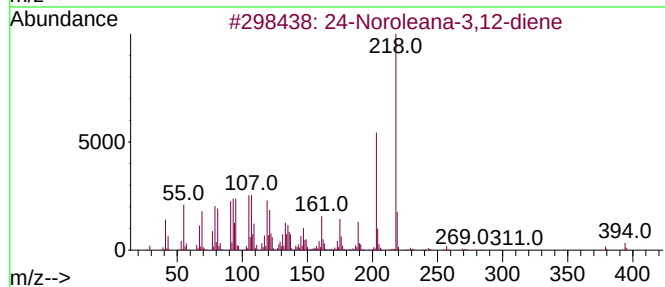
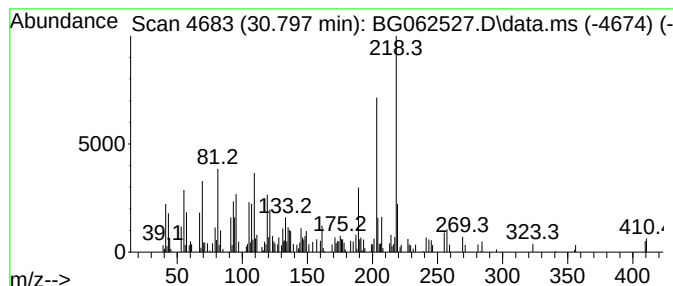
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 24-Noroleana-3,12-diene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.797	3.29 ng/ul	251068	Perylene-d12	24.612

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			24-Noroleana-3,12-diene	394	C29H46	201358-24-9	72
2			Benzo[b]naphtho[2,3-d]furan	218	C16H10O	000243-42-5	53
3			5,6,7,8-Tetrahydrothieno[2,3-b]q...	218	C12H14N2S	134315-44-9	53
4			8-Amino-6,7-dimethoxy-4-methylqu...	218	C12H14N2O2	1000213-07-2	50
5			7-Isopropenyl-1,4a-dimethyl-4,4a...	218	C15H22O	000473-08-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
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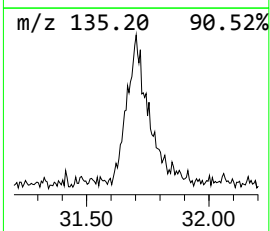
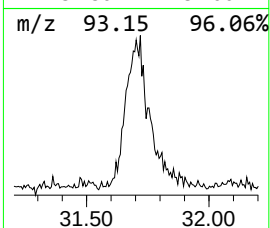
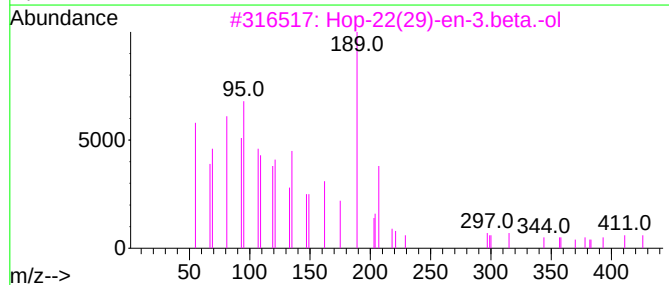
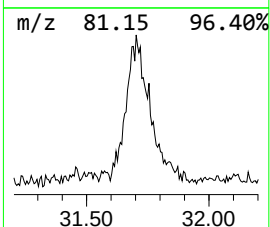
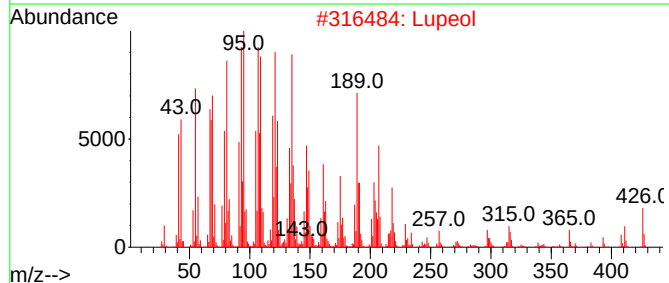
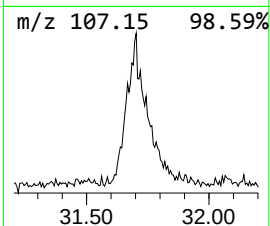
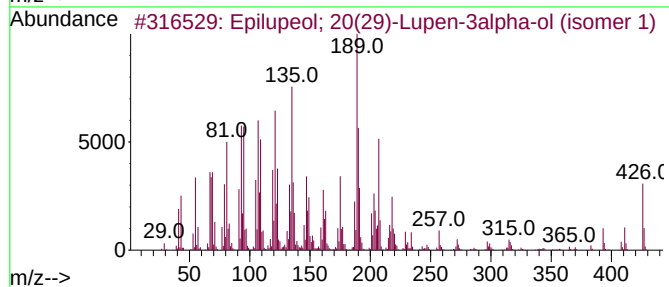
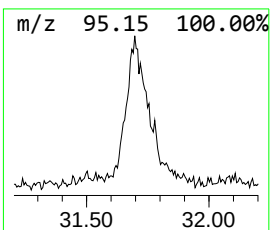
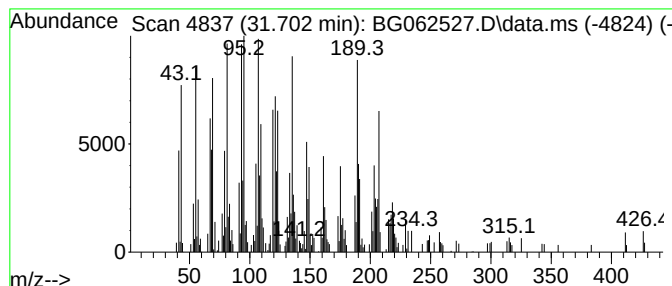
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Epilupeol; 20(29)-Lupen-3a1... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.702	13.12 ng/ul	1000090	Perylene-d12	24.612

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Epilupeol; 20(29)-Lupen-3alpha-o...	426	C30H50O	1000513-01-3	92
2			Lupeol	426	C30H50O	000545-47-1	80
3			Hop-22(29)-en-3.beta.-ol	426	C30H50O	058801-23-3	78
4			Guaia-9,11-diene	204	C15H24	1000374-19-8	70
5			1,3-Benzodioxole-5-methanamine, ...	325	C19H16FN03	1000351-17-0	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062527.D
Acq On : 22 Aug 2024 4:12
Operator : MA/JU
Sample : P3626-18
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCXZ7

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
(S)-(+)-1,2-Pro...	3.664	4.9	ng/ul	110006	1	7.940	450148	20.0
1-Phenoxypropan...	11.465	3.6	ng/ul	126847	2	10.730	709211	20.0
n-Hexadecanoic ...	18.191	4.3	ng/ul	268888	4	17.298	1266440	20.0
Heptadecyl hept...	21.211	5.4	ng/ul	367540	5	21.534	1351330	20.0
(DEL) Alkane: C...	22.350	12.2	ng/ul	823323	5	21.534	1351330	20.0
(DEL) Alkane: S...	23.754	2.2	ng/ul	168117	6	24.612	1525060	20.0
(DEL) Alkane: S...	25.728	2.2	ng/ul	169769	6	24.612	1525060	20.0
unknown-01	29.681	3.6	ng/ul	271187	6	24.612	1525060	20.0
24-Noroleana-3,...	30.797	3.3	ng/ul	251068	6	24.612	1525060	20.0
Epilupeol; 20(2...	31.702	13.1	ng/ul	1000090	6	24.612	1525060	20.0