

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082024\
 Data File : BG062494.D
 Acq On : 21 Aug 2024 2:46
 Operator : MA/JU
 Sample : P3504-23
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCYF1

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: BG062494.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.406	16	20	25	rVB2	14942	19379	1.04%	0.074%
2	3.664	60	64	76	rBV	80549	129631	6.95%	0.492%
3	3.811	84	89	102	rBV	96017	154807	8.30%	0.588%
4	5.051	297	300	305	rVB3	3916	5440	0.29%	0.021%
5	6.026	461	466	472	rVB2	9719	14848	0.80%	0.056%
6	6.637	564	570	577	rVB	113093	172792	9.27%	0.656%
7	6.925	614	619	621	rBV3	4738	7119	0.38%	0.027%
8	7.101	641	649	660	rBV	324513	569489	30.54%	2.162%
9	7.271	670	678	687	rBV	215901	347847	18.65%	1.321%
10	7.389	693	698	705	rVB2	12121	19360	1.04%	0.074%
11	7.471	705	712	718	rVV	290135	488609	26.20%	1.855%
12	7.530	718	722	735	rVB2	23034	39192	2.10%	0.149%
13	7.941	785	792	799	rVV	344950	567978	30.46%	2.157%
14	8.000	799	802	809	rVV5	7477	13281	0.71%	0.050%
15	8.640	904	911	921	rBV	397591	669863	35.92%	2.543%
16	9.092	980	988	999	rBV	280768	499741	26.80%	1.898%
17	9.204	1003	1007	1012	rVB3	2969	4781	0.26%	0.018%
18	9.533	1055	1063	1067	rBV4	6294	10518	0.56%	0.040%
19	9.650	1074	1083	1090	rBV2	29890	48963	2.63%	0.186%
20	9.815	1103	1111	1118	rBV	234979	410747	22.03%	1.560%
21	10.044	1144	1150	1157	rVB5	11580	20234	1.09%	0.077%
22	10.350	1195	1202	1215	rVB	393022	694011	37.22%	2.635%
23	10.508	1226	1229	1238	rVB4	3473	5913	0.32%	0.022%
24	10.731	1260	1267	1275	rBV	510069	910526	48.83%	3.457%
25	10.861	1281	1289	1299	rBV	322861	572479	30.70%	2.174%
26	11.260	1351	1357	1364	rVB2	13477	20491	1.10%	0.078%
27	11.466	1385	1392	1408	rVB	98789	183419	9.84%	0.696%
28	12.218	1515	1520	1524	rVB2	3192	5561	0.30%	0.021%
29	12.312	1527	1536	1542	rBV4	7435	14746	0.79%	0.056%
30	12.541	1572	1575	1580	rVB2	5528	7564	0.41%	0.029%
31	13.175	1678	1683	1688	rBV3	2673	4942	0.27%	0.019%
32	13.392	1716	1720	1725	rBV3	5927	8499	0.46%	0.032%
33	13.968	1811	1818	1825	rVV	658840	933086	50.04%	3.543%
34	14.209	1854	1859	1860	rBV3	15765	18862	1.01%	0.072%
35	14.256	1860	1867	1874	rVB	759264	1208795	64.82%	4.590%
36	14.462	1900	1902	1907	rVB3	4278	5636	0.30%	0.021%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
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37	14.561	1911	1919	1928	rVB	864144	1321215	70.85%	5.017%
38	14.749	1944	1951	1961	rBV2	306630	568472	30.48%	2.159%
39	14.896	1971	1976	1982	rVB5	7383	10741	0.58%	0.041%
40	14.973	1985	1989	1995	rVB3	9998	12150	0.65%	0.046%
41	15.360	2051	2055	2061	rBV3	10602	17052	0.91%	0.065%
42	15.419	2061	2065	2068	rBV4	5196	5678	0.30%	0.022%
43	15.554	2081	2088	2095	rBV	979659	1506386	80.78%	5.720%
44	15.660	2099	2106	2113	rBV	260068	384542	20.62%	1.460%
45	15.789	2123	2128	2132	rBV5	3230	5464	0.29%	0.021%
46	16.101	2175	2181	2184	rBV4	7062	11958	0.64%	0.045%
47	16.277	2207	2211	2214	rBV3	6462	7286	0.39%	0.028%
48	16.441	2235	2239	2244	rVB5	4132	6613	0.35%	0.025%
49	16.700	2280	2283	2285	rBV3	5574	6275	0.34%	0.024%
50	16.858	2305	2310	2313	rBV6	7875	12398	0.66%	0.047%
51	17.017	2334	2337	2341	rVB3	5795	7053	0.38%	0.027%
52	17.064	2341	2345	2352	rBV5	5333	10701	0.57%	0.041%
53	17.193	2361	2367	2370	rBV5	3911	7549	0.40%	0.029%
54	17.223	2370	2372	2376	rVV4	3739	4711	0.25%	0.018%
55	17.299	2378	2385	2395	rVB	1002260	1534095	82.27%	5.825%
56	17.399	2395	2402	2410	rBV	976315	1441532	77.30%	5.474%
57	17.552	2424	2428	2431	rVB4	4006	5945	0.32%	0.023%
58	17.587	2431	2434	2436	rBV2	3980	4687	0.25%	0.018%
59	17.681	2446	2450	2454	rBV7	5567	8739	0.47%	0.033%
60	17.804	2468	2471	2473	rBV3	4740	4446	0.24%	0.017%
61	17.869	2480	2482	2488	rVB6	6001	9814	0.53%	0.037%
62	17.974	2496	2500	2505	rVB2	17356	23816	1.28%	0.090%
63	18.068	2511	2516	2519	rBV5	6807	14458	0.78%	0.055%
64	18.198	2531	2538	2553	rBV	267913	478634	25.67%	1.817%
65	18.315	2553	2558	2565	rVV5	31888	58912	3.16%	0.224%
66	18.644	2610	2614	2615	rVV3	4953	8109	0.43%	0.031%
67	18.674	2615	2619	2624	rVB6	12188	21638	1.16%	0.082%
68	18.744	2628	2631	2636	rBV6	7203	9579	0.51%	0.036%
69	18.785	2636	2638	2641	rBV4	3493	4767	0.26%	0.018%
70	18.862	2646	2651	2652	rVV4	5572	7817	0.42%	0.030%
71	18.914	2656	2660	2667	rVB7	11474	19353	1.04%	0.073%
72	18.985	2667	2672	2679	rVB7	7006	18945	1.02%	0.072%
73	19.079	2681	2688	2692	rBV9	21468	38519	2.07%	0.146%
74	19.114	2692	2694	2696	rVV3	17315	20777	1.11%	0.079%
75	19.138	2696	2698	2705	rVB6	16432	22641	1.21%	0.086%
76	19.249	2713	2717	2721	rBV3	14998	22221	1.19%	0.084%

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77	19.320	2723	2729	2731	rBV4	38380	61314	3.29%	0.233%
78	19.343	2731	2733	2737	rVB4	14311	18864	1.01%	0.072%
79	19.467	2749	2754	2764	rBV	142449	232876	12.49%	0.884%
80	19.613	2773	2779	2783	rVB7	17847	32022	1.72%	0.122%
81	19.684	2784	2791	2796	rBV	1225153	1779123	95.41%	6.755%
82	19.919	2827	2831	2835	rVB7	8530	13248	0.71%	0.050%
83	20.072	2854	2857	2861	rVB4	13856	17534	0.94%	0.067%
84	20.130	2865	2867	2869	rVB3	6881	5172	0.28%	0.020%
85	20.195	2875	2878	2881	rBV4	11094	14034	0.75%	0.053%
86	20.230	2882	2884	2886	rVV2	7019	5702	0.31%	0.022%
87	20.377	2907	2909	2917	rVB9	16554	23633	1.27%	0.090%
88	20.547	2930	2938	2942	rBV4	20721	64739	3.47%	0.246%
89	20.659	2952	2957	2962	rBV	88521	114099	6.12%	0.433%
90	20.712	2962	2966	2971	rBV5	39993	63461	3.40%	0.241%
91	21.182	3037	3046	3048	rBV	348498	580059	31.11%	2.203%
92	21.205	3048	3050	3061	rVB	260376	438596	23.52%	1.665%
93	21.540	3101	3107	3122	rVB2	1017793	1697056	91.01%	6.444%
94	21.658	3124	3127	3133	rVB3	24939	32785	1.76%	0.124%
95	23.761	3483	3485	3490	rVB3	28645	39530	2.12%	0.150%
96	24.407	3585	3595	3607	rVB	699659	1846992	99.05%	7.013%
97	24.618	3621	3631	3648	rVB	662464	1864766	100.00%	7.081%
98	25.858	3834	3842	3859	rVB3	156300	494406	26.51%	1.877%
99	28.595	4298	4308	4321	rVB3	54900	218801	11.73%	0.831%
100	29.670	4484	4491	4507	rVB3	41266	182642	9.79%	0.693%

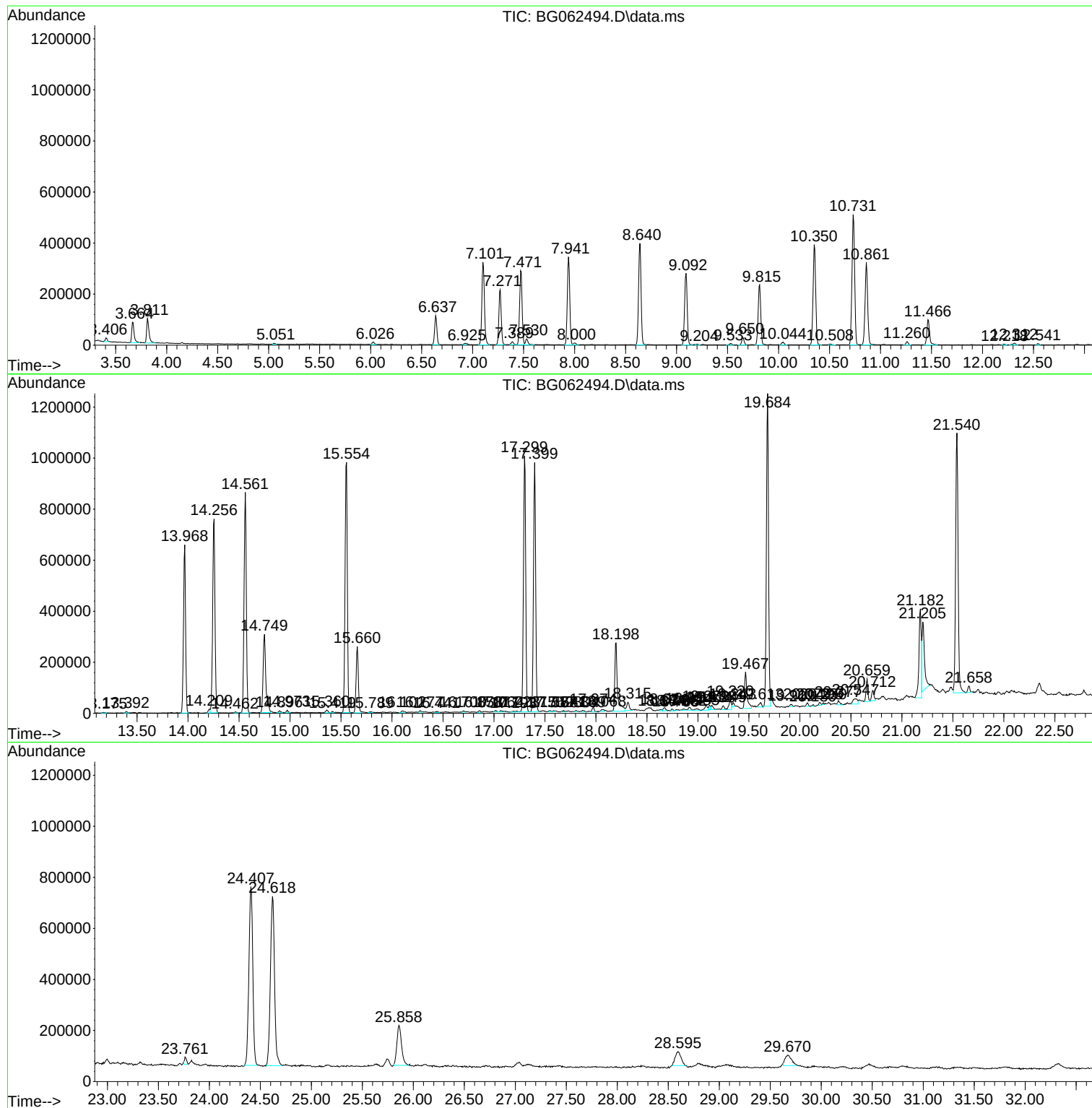
Sum of corrected areas: 26336291

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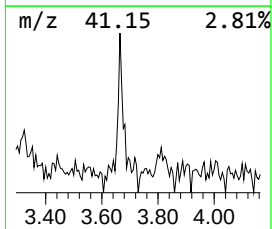
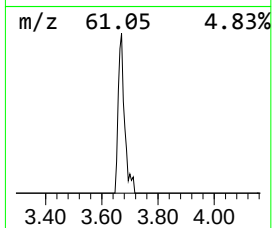
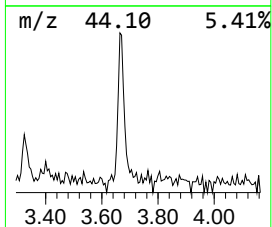
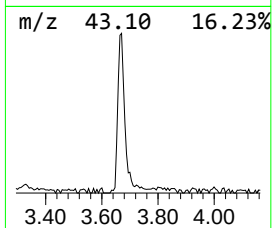
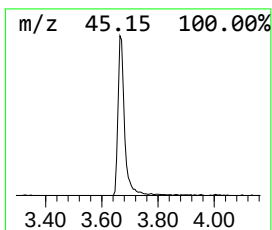
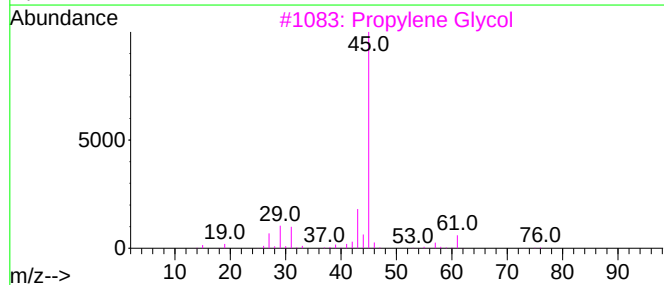
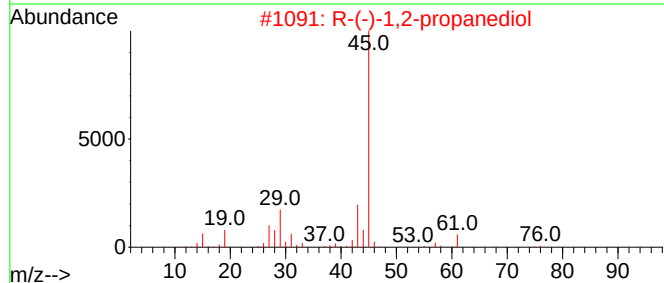
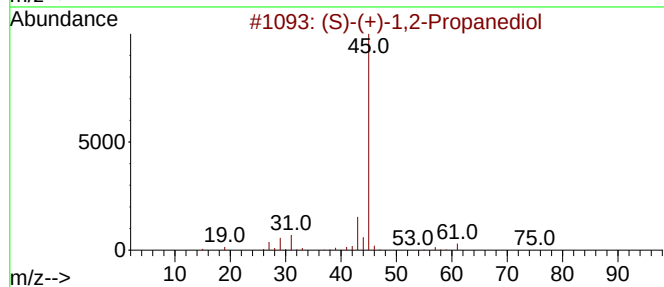
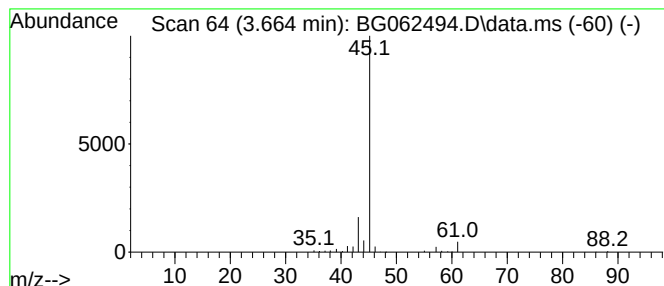
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (S)-(+)-1,2-Propanediol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.664	4.56 ng/ul	129631	1,4-Dichlorobenzene-d4	7.941

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	64
2	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	43
3	Propylene Glycol			76	C3H8O2	000057-55-6	43
4	2-Butanol, 3-methyl-			88	C5H12O	000598-75-4	9
5	2-Butanol, 3-methyl-			88	C5H12O	000598-75-4	9



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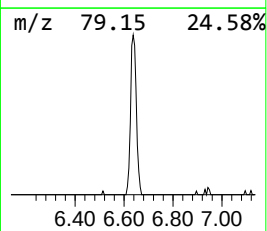
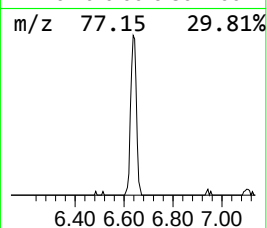
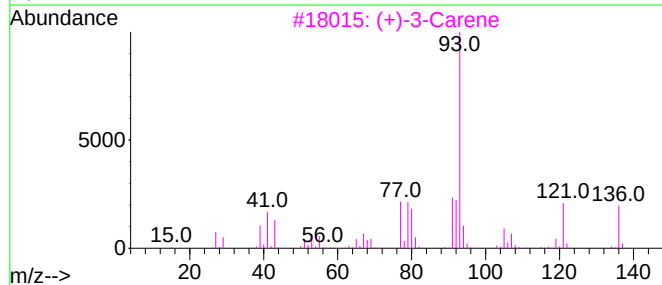
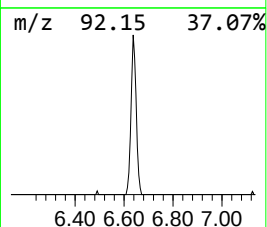
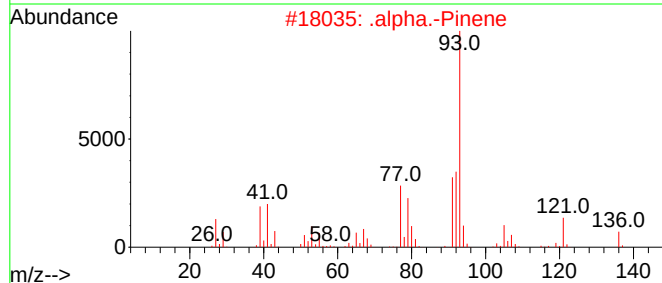
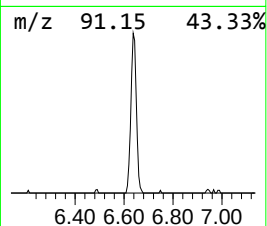
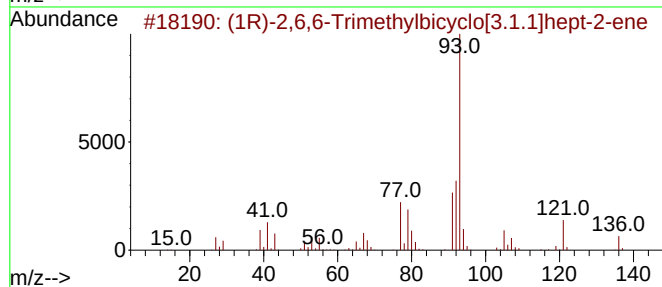
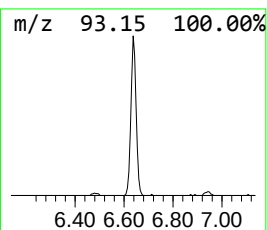
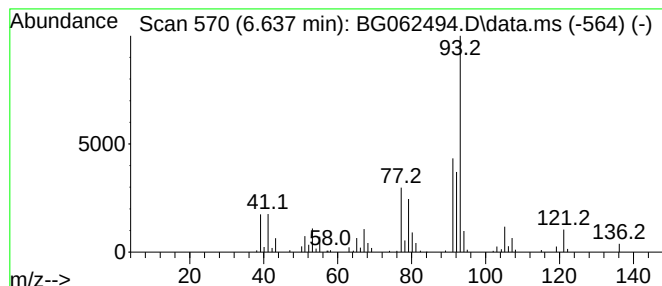
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.637	6.08 ng/ul	172792	1,4-Dichlorobenzene-d4	7.941

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	94
2			.alpha.-Pinene	136	C10H16	000080-56-8	94
3			(+)-3-Carene	136	C10H16	000498-15-7	91
4			.alpha.-Pinene	136	C10H16	000080-56-8	91
5			Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	91



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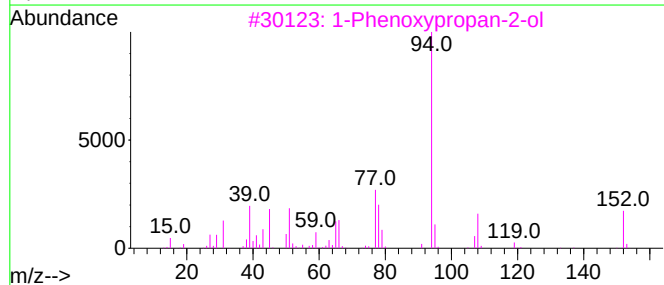
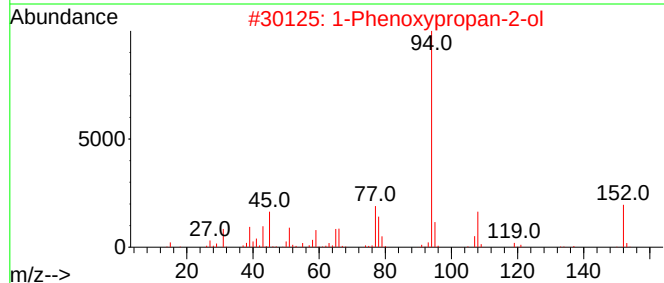
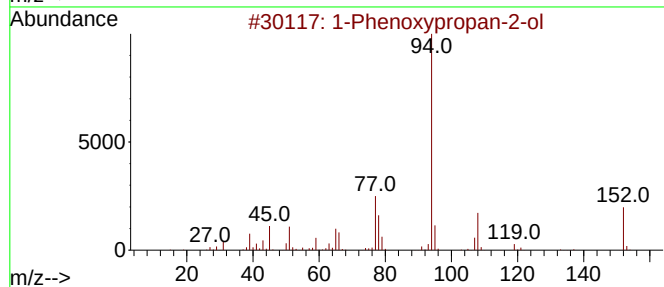
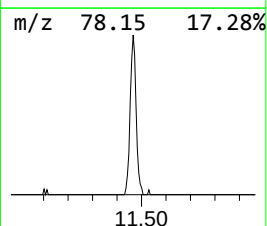
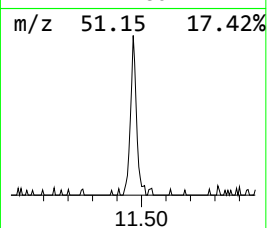
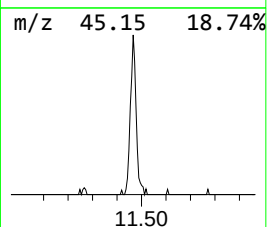
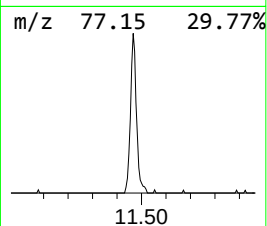
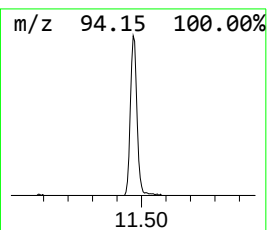
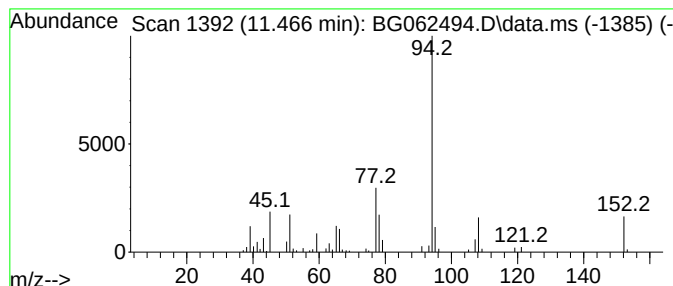
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.466	4.03 ng/ul	183419	Naphthalene-d8	10.731

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	93
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082024\
Data File : BG062494.D
Acq On : 21 Aug 2024 2:46
Operator : MA/JU
Sample : P3504-23
Misc :
ALS Vial : 15 Sample Multiplier: 1

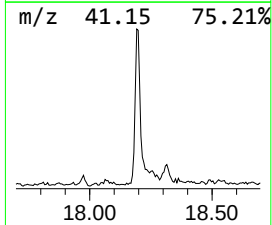
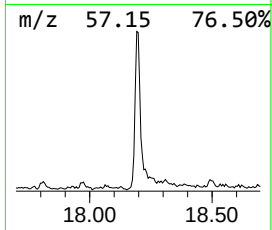
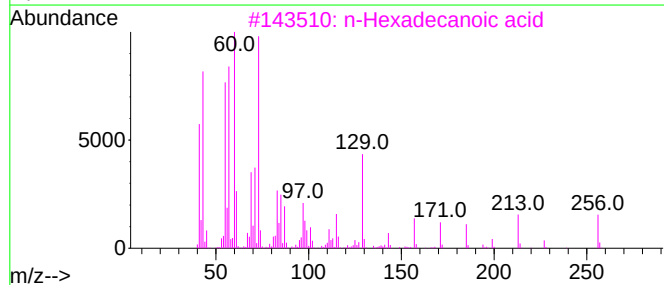
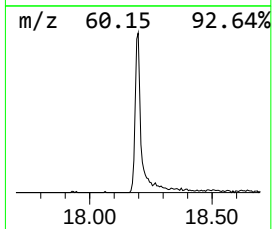
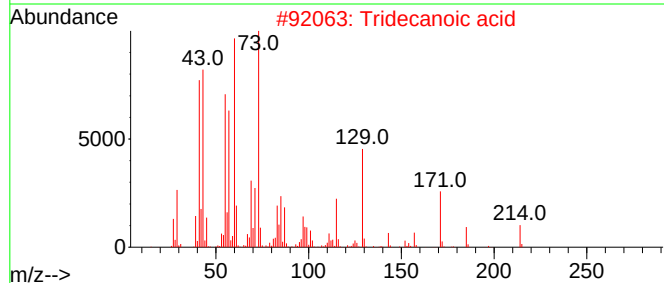
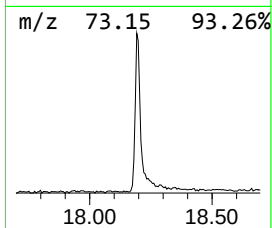
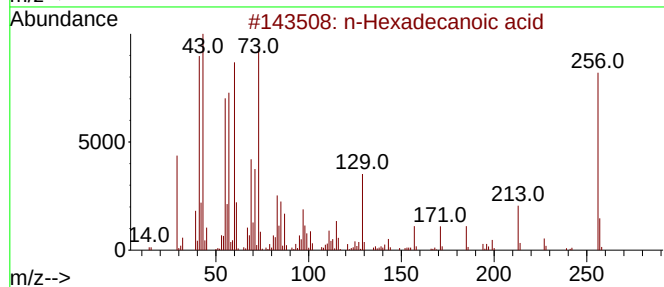
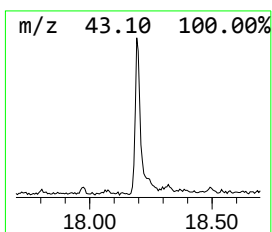
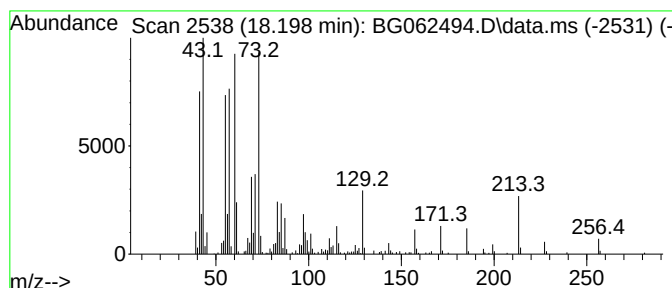
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BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.198	6.24 ng/ul	478634	Phenanthrene-d10	17.299	
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	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082024\
Data File : BG062494.D
Acq On : 21 Aug 2024 2:46
Operator : MA/JU
Sample : P3504-23
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYF1

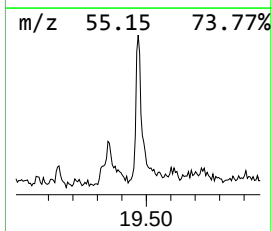
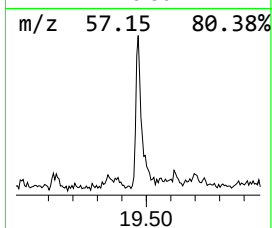
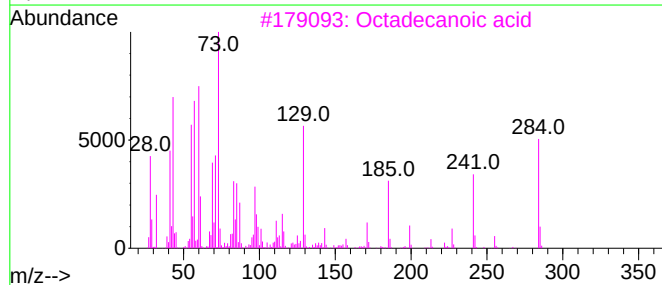
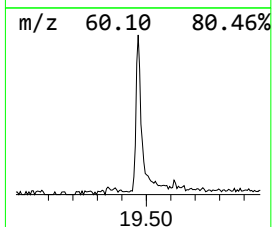
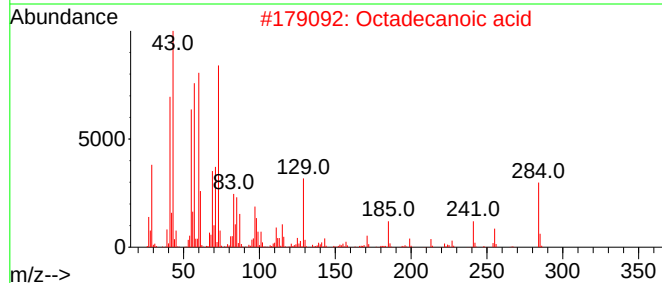
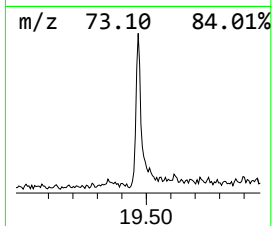
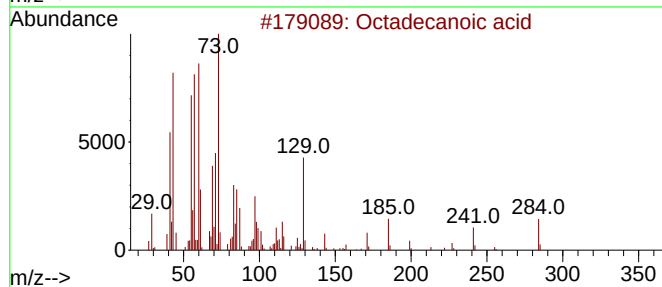
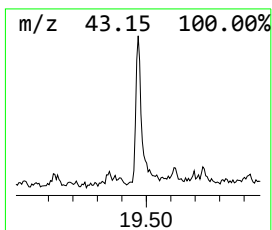
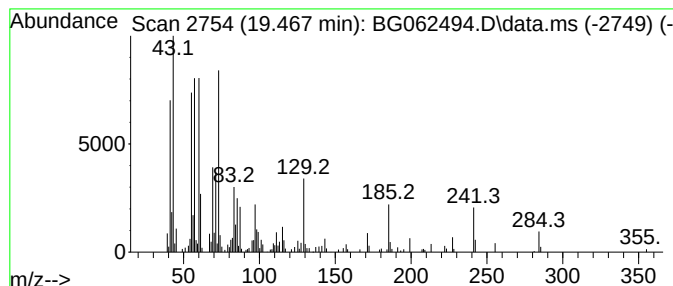
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.467	2.74 ng/ul	232876	Chrysene-d12	21.540

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	95
3			Octadecanoic acid	284	C18H36O2	000057-11-4	93
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082024\
Data File : BG062494.D
Acq On : 21 Aug 2024 2:46
Operator : MA/JU
Sample : P3504-23
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYF1

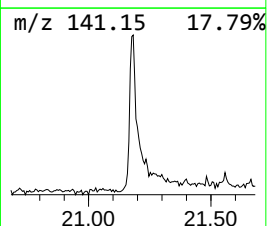
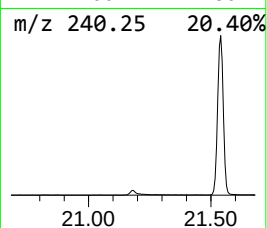
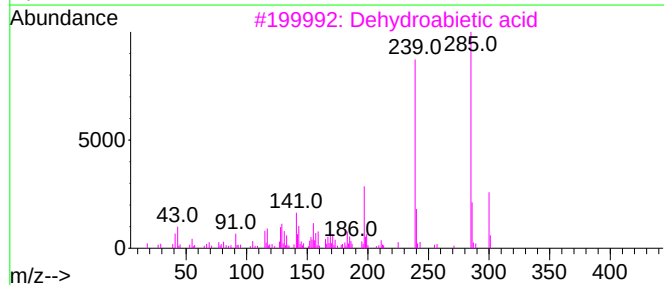
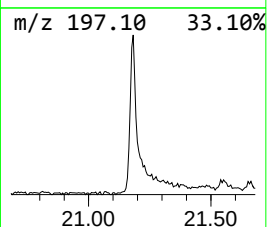
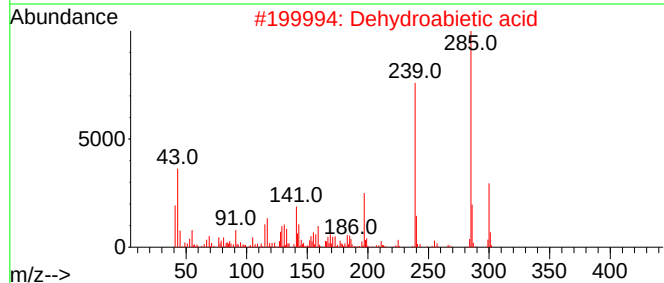
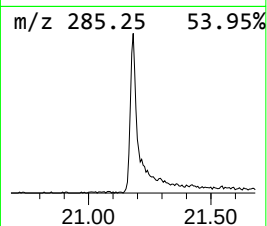
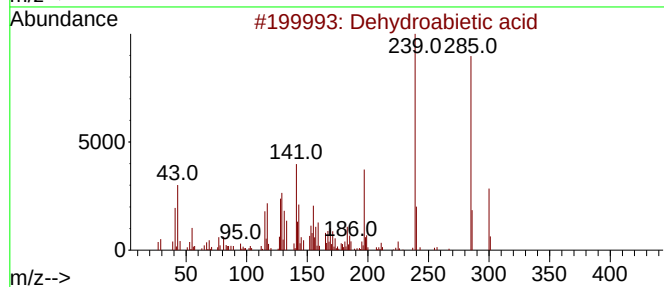
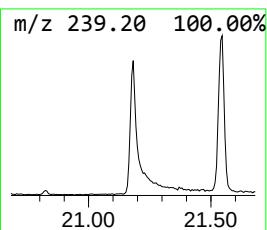
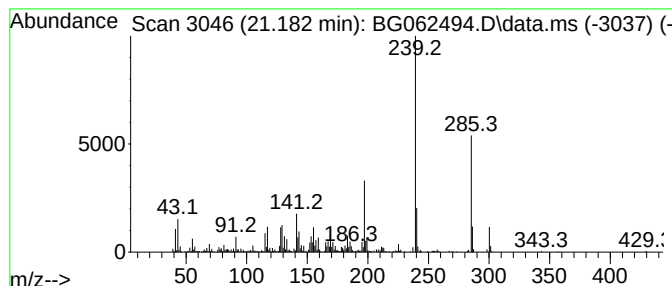
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Dehydroabietic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.182	6.84 ng/ul	580059	Chrysene-d12	21.540

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabietic acid	300	C20H28O2	001740-19-8	99
2			Dehydroabietic acid	300	C20H28O2	001740-19-8	94
3			Dehydroabietic acid	300	C20H28O2	001740-19-8	91
4			Butanoic acid, 2-(cyano)(2,4,6-t...	300	C17H20N2O3	1000267-71-7	74
5			Ethyl 4-hydroxy-7-trifluoromethy...	285	C13H10F3NO3	000391-02-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082024\
Data File : BG062494.D
Acq On : 21 Aug 2024 2:46
Operator : MA/JU
Sample : P3504-23
Misc :
ALS Vial : 15 Sample Multiplier: 1

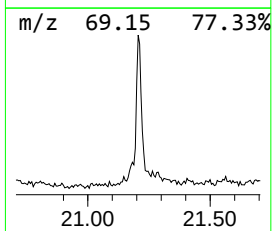
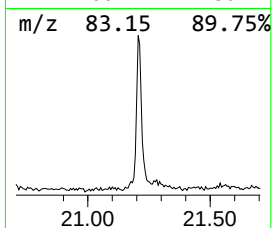
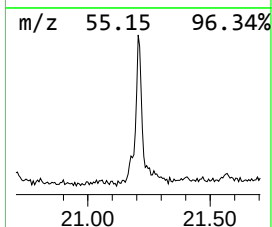
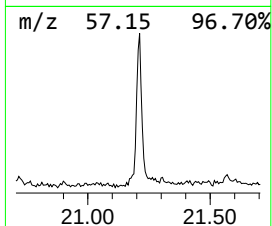
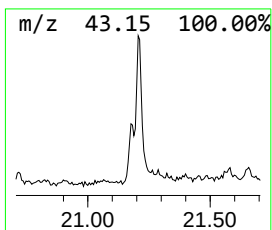
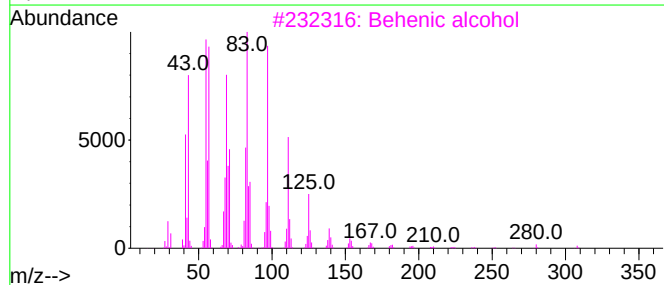
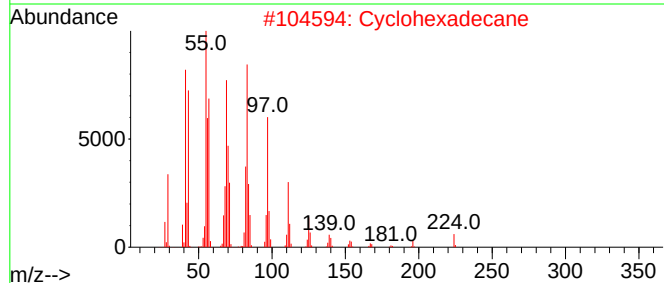
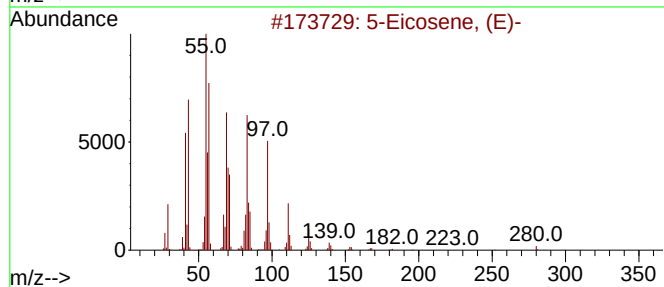
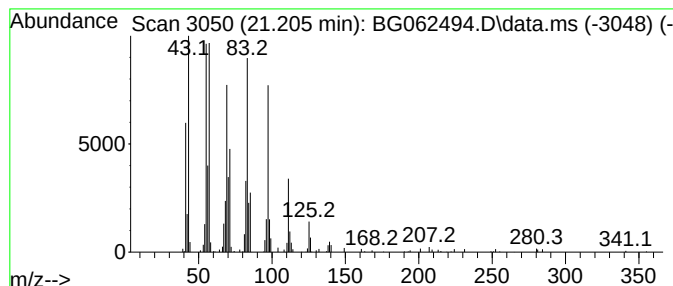
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.205	5.17 ng/ul	438596	Chrysene-d12		21.540	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	95
2	Cyclohexadecane		224	C16H32	000295-65-8	95
3	Behenic alcohol		326	C22H46O	000661-19-8	94
4	1-Hexacosene		364	C26H52	018835-33-1	91
5	1-Heneicosanol		312	C21H44O	015594-90-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082024\
Data File : BG062494.D
Acq On : 21 Aug 2024 2:46
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Sample : P3504-23
Misc :
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Instrument :
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ClientSampleId :
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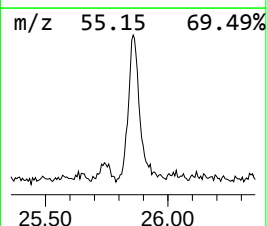
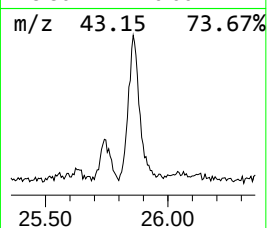
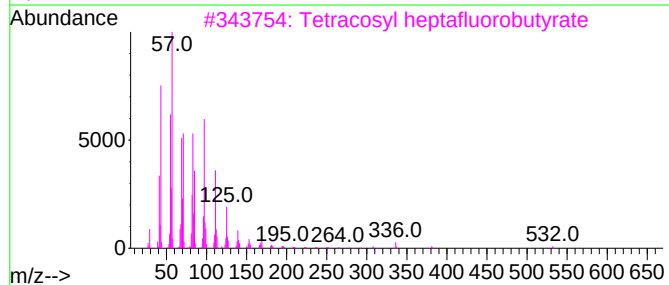
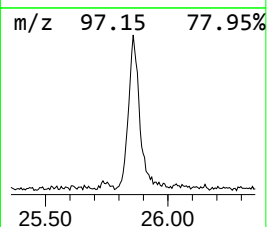
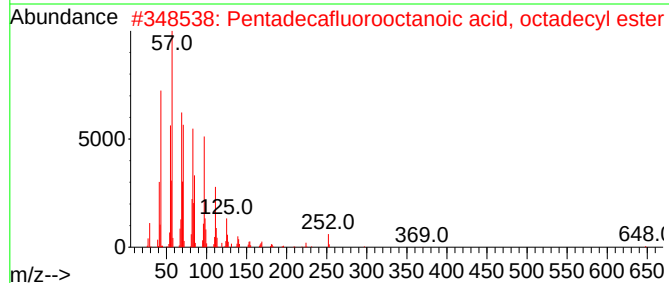
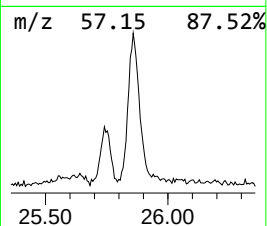
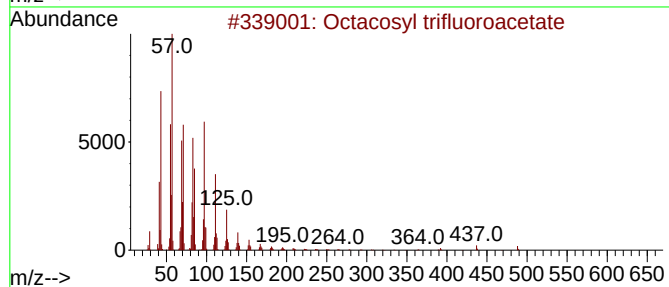
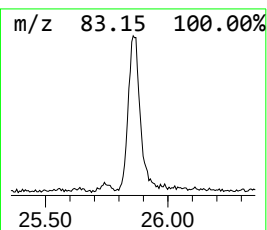
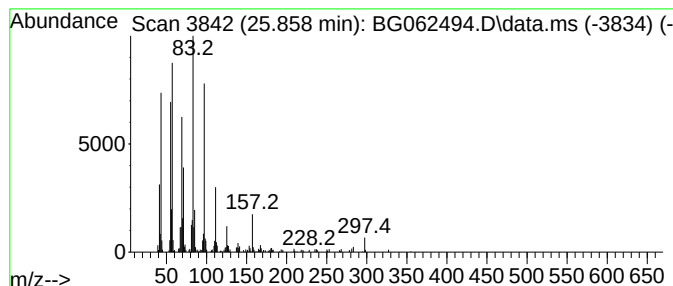
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.858	5.30 ng/ul	494406	Perylene-d12	24.618

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	46
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	41
3			Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	41
4			E-2-Octenyl tiglate	210	C13H22O2	084271-97-6	38
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082024\
Data File : BG062494.D
Acq On : 21 Aug 2024 2:46
Operator : MA/JU
Sample : P3504-23
Misc :
ALS Vial : 15 Sample Multiplier: 1

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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(1R)-2,6,6-Trim...	6.637	6.1	ng/ul	172792	1	7.941	567978	20.0
1-Phenoxypropan...	11.466	4.0	ng/ul	183419	2	10.731	910526	20.0
n-Hexadecanoic ...	18.198	6.2	ng/ul	478634	4	17.299	1534100	20.0
Octadecanoic acid	19.467	2.7	ng/ul	232876	5	21.540	1697060	20.0
Dehydroabietic ...	21.182	6.8	ng/ul	580059	5	21.540	1697060	20.0
(DEL) Alkane: S...	21.205	5.2	ng/ul	438596	5	21.540	1697060	20.0
unknown-01	25.858	5.3	ng/ul	494406	6	24.618	1864770	20.0
unknown-02	28.595	2.4	ng/ul	218801	6	24.618	1864770	20.0