

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
 Data File : BG062384.D
 Acq On : 13 Aug 2024 14:10
 Operator : MA/JU
 Sample : P3502-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCXU7

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-BG080524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062384.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.426	18	23	30	rVB2	27901	49903	4.32%	0.281%
2	3.679	61	66	78	rVB	40847	71789	6.21%	0.404%
3	3.820	85	90	106	rVB	109654	194257	16.80%	1.093%
4	3.931	106	109	113	rBV4	2944	4068	0.35%	0.023%
5	4.160	144	148	151	rBV5	3183	4880	0.42%	0.027%
6	4.207	152	156	161	rVB	9270	13698	1.18%	0.077%
7	4.331	170	177	183	rBV4	5181	11096	0.96%	0.062%
8	4.519	207	209	213	rVB4	2302	2415	0.21%	0.014%
9	5.288	334	340	345	rBV4	5015	9878	0.85%	0.056%
10	5.999	458	461	464	rBV4	3353	4176	0.36%	0.024%
11	6.034	464	467	473	rVV	9878	13929	1.20%	0.078%
12	6.510	546	548	555	rVB4	1952	2681	0.23%	0.015%
13	6.657	567	573	579	rVB2	17701	30204	2.61%	0.170%
14	6.898	610	614	617	rBV4	2874	5039	0.44%	0.028%
15	6.921	617	618	629	rVB5	3590	7038	0.61%	0.040%
16	7.109	643	650	661	rVB	182560	311235	26.92%	1.751%
17	7.280	672	679	687	rBV	117029	191184	16.53%	1.076%
18	7.403	694	700	704	rBV4	6447	9107	0.79%	0.051%
19	7.485	707	714	720	rVV	162932	271145	23.45%	1.526%
20	7.544	720	724	734	rVB	13892	23732	2.05%	0.134%
21	7.750	757	759	764	rVB3	4096	6008	0.52%	0.034%
22	7.855	771	777	783	rVB6	2326	5671	0.49%	0.032%
23	7.955	787	794	800	rVV	204440	342424	29.61%	1.927%
24	8.014	800	804	812	rVV5	7638	15782	1.36%	0.089%
25	8.096	812	818	823	rVB4	1849	3467	0.30%	0.020%
26	8.649	906	912	923	rBV2	195181	338917	29.31%	1.907%
27	9.101	982	989	1000	rVV	153205	265530	22.96%	1.494%
28	9.189	1000	1004	1007	rVV3	2021	3613	0.31%	0.020%
29	9.218	1007	1009	1014	rVV5	1810	2299	0.20%	0.013%
30	9.277	1014	1019	1022	rVB4	2281	3694	0.32%	0.021%
31	9.553	1064	1066	1071	rVB3	2806	2784	0.24%	0.016%
32	9.665	1078	1085	1092	rBV2	27761	44490	3.85%	0.250%
33	9.823	1104	1112	1123	rBV	126343	227191	19.65%	1.279%
34	10.053	1145	1151	1156	rVB4	6343	11933	1.03%	0.067%
35	10.364	1197	1204	1215	rVB	219828	392476	33.94%	2.209%
36	10.522	1227	1231	1236	rVB5	2178	4146	0.36%	0.023%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
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37	10.746	1261	1269	1277	rBV	319601	573320	49.58%	3.226%
38	10.875	1282	1291	1298	rBV	174775	321437	27.80%	1.809%
39	11.139	1331	1336	1341	rVB3	5614	9378	0.81%	0.053%
40	11.227	1345	1351	1355	rBV2	3678	5754	0.50%	0.032%
41	11.280	1355	1360	1365	rVB2	12033	19746	1.71%	0.111%
42	11.480	1388	1394	1406	rBV2	54042	108299	9.37%	0.609%
43	11.973	1473	1478	1483	rBV3	1296	2548	0.22%	0.014%
44	12.220	1516	1520	1527	rVB5	4007	6742	0.58%	0.038%
45	12.326	1535	1538	1544	rVB2	4051	6166	0.53%	0.035%
46	13.330	1702	1709	1710	rBV4	1956	3109	0.27%	0.017%
47	13.407	1718	1722	1728	rBV4	3111	5483	0.47%	0.031%
48	13.471	1728	1733	1734	rBV4	2591	3722	0.32%	0.021%
49	13.577	1746	1751	1757	rVB	33968	45794	3.96%	0.258%
50	13.853	1796	1798	1802	rBV4	2419	2769	0.24%	0.016%
51	13.912	1802	1808	1813	rBV4	23278	35021	3.03%	0.197%
52	13.977	1813	1819	1827	rVV	400508	580430	50.20%	3.266%
53	14.264	1856	1868	1875	rBV	474782	761991	65.90%	4.288%
54	14.482	1901	1905	1910	rBV3	3750	5289	0.46%	0.030%
55	14.570	1912	1920	1930	rVV	544850	862422	74.58%	4.853%
56	14.646	1930	1933	1938	rVB5	3100	4828	0.42%	0.027%
57	14.752	1944	1951	1957	rBV	191887	331096	28.63%	1.863%
58	14.905	1973	1977	1982	rVB6	2574	4216	0.36%	0.024%
59	14.987	1987	1991	1995	rBV3	6123	8942	0.77%	0.050%
60	15.163	2018	2021	2023	rBV3	1744	2113	0.18%	0.012%
61	15.310	2043	2046	2050	rVB2	8176	10615	0.92%	0.060%
62	15.369	2052	2056	2059	rBV5	5460	7548	0.65%	0.042%
63	15.433	2063	2067	2071	rBV5	3564	5931	0.51%	0.033%
64	15.480	2074	2075	2079	rBV4	2209	2887	0.25%	0.016%
65	15.563	2083	2089	2098	rVB	624746	940288	81.32%	5.292%
66	15.668	2100	2107	2115	rBV2	144524	222430	19.24%	1.252%
67	15.839	2134	2136	2141	rVB5	3654	4166	0.36%	0.023%
68	15.939	2148	2153	2158	rBB3	15503	20465	1.77%	0.115%
69	15.992	2161	2162	2166	rVB4	1935	2359	0.20%	0.013%
70	16.109	2180	2182	2183	rVV2	2599	2058	0.18%	0.012%
71	16.185	2191	2195	2201	rVB	32049	47595	4.12%	0.268%
72	16.291	2210	2213	2218	rBV6	5637	6791	0.59%	0.038%
73	16.420	2231	2235	2242	rBV7	9038	19520	1.69%	0.110%
74	16.655	2273	2275	2278	rBV4	1596	2429	0.21%	0.014%
75	16.714	2281	2285	2290	rBV3	24129	36115	3.12%	0.203%
76	16.879	2309	2313	2315	rBV4	5155	7440	0.64%	0.042%

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 Title : SVOA CALIBRATION

77	16.961	2322	2327	2334	rVB3	27735	40934	3.54%	0.230%
78	17.025	2334	2338	2343	rBV7	4492	7928	0.69%	0.045%
79	17.078	2343	2347	2351	rBV6	8765	14754	1.28%	0.083%
80	17.213	2365	2370	2377	rBV	101458	145935	12.62%	0.821%
81	17.313	2377	2387	2393	rVV	671166	1098632	95.01%	6.183%
82	17.407	2397	2403	2412	rVB2	645446	984795	85.17%	5.542%
83	17.489	2413	2417	2420	rBV5	13855	19319	1.67%	0.109%
84	17.948	2492	2495	2499	rBV4	20571	29505	2.55%	0.166%
85	18.095	2514	2520	2525	rBV	52120	92805	8.03%	0.522%
86	18.153	2525	2530	2535	rBV	156383	224447	19.41%	1.263%
87	18.212	2535	2540	2549	rVV	612429	888509	76.84%	5.000%
88	18.418	2573	2575	2581	rVB4	27575	36163	3.13%	0.204%
89	18.512	2587	2591	2596	rBV3	52062	92172	7.97%	0.519%
90	18.647	2610	2614	2618	rBV4	23411	40776	3.53%	0.229%
91	19.334	2729	2731	2733	rBV3	53933	55938	4.84%	0.315%
92	19.363	2734	2736	2738	rVV2	67380	66668	5.77%	0.375%
93	19.481	2752	2756	2766	rBV	326541	466412	40.34%	2.625%
94	19.692	2787	2792	2800	rBV	800524	1144486	98.98%	6.441%
95	21.044	3018	3022	3033	rBV	152973	379241	32.80%	2.134%
96	21.190	3043	3047	3050	rBV	301704	417727	36.13%	2.351%
97	21.226	3050	3053	3058	rVB3	206034	289946	25.07%	1.632%
98	21.555	3104	3109	3120	rVB	670849	1098214	94.97%	6.180%
99	24.427	3591	3598	3612	rVB	425593	1156330	100.00%	6.507%
100	24.639	3627	3634	3643	rVB	392240	1026903	88.81%	5.779%

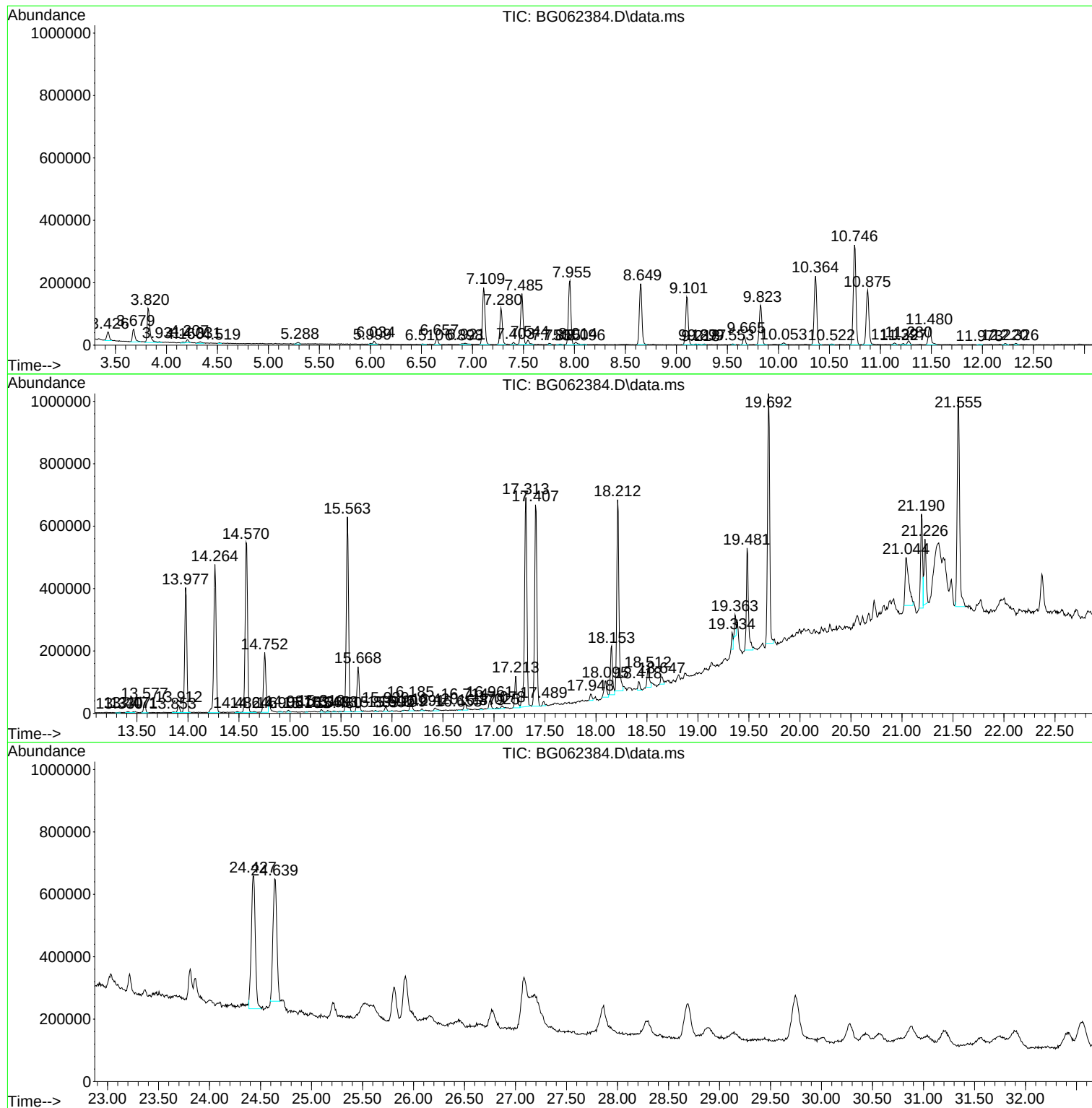
Sum of corrected areas: 17769670

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
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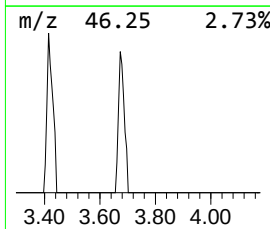
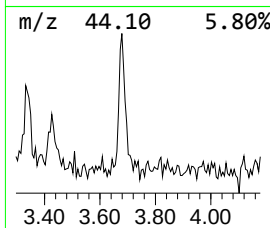
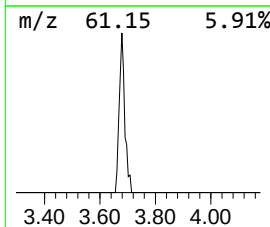
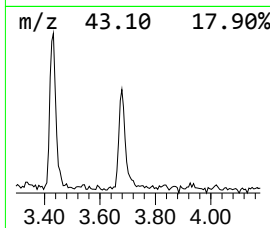
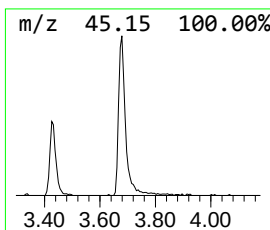
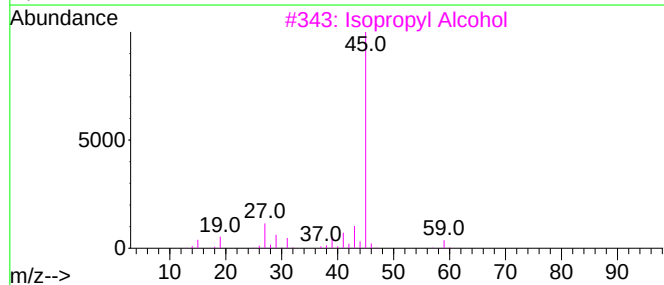
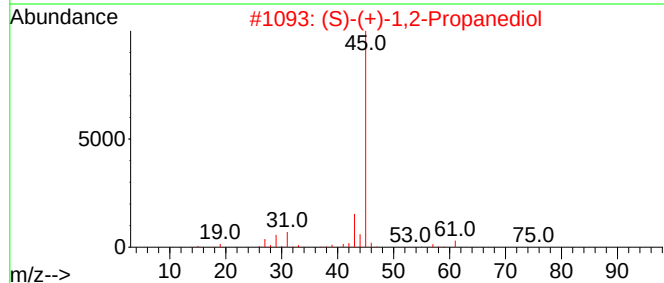
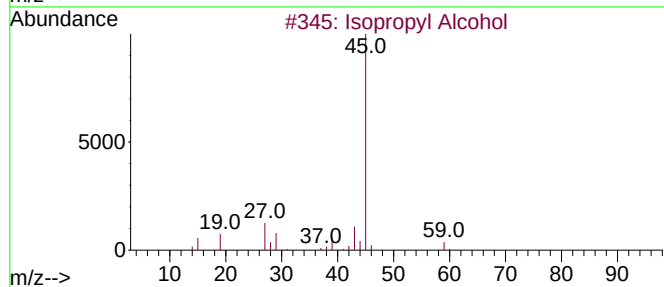
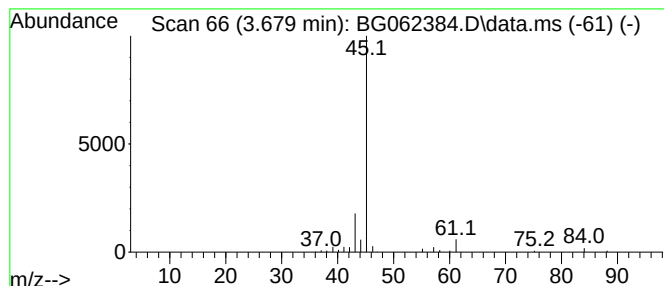
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Isopropyl Alcohol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.679	4.19 ng/ul	71789	1,4-Dichlorobenzene-d4	7.955

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	80
2			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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ALS Vial : 8 Sample Multiplier: 1

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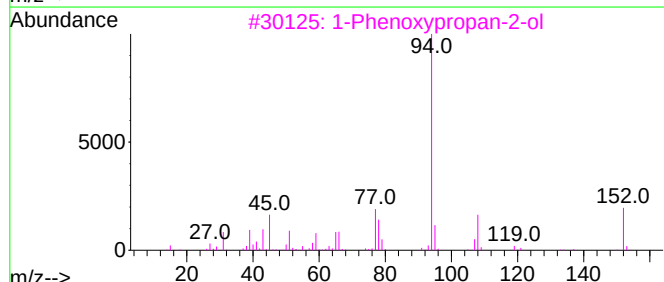
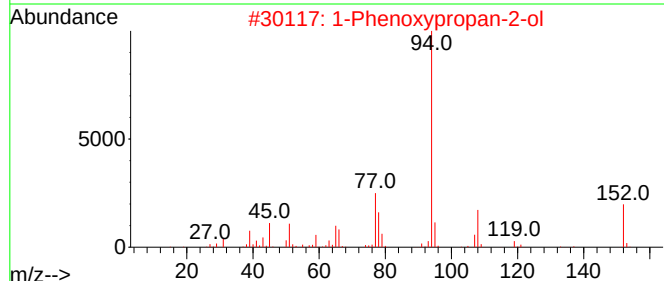
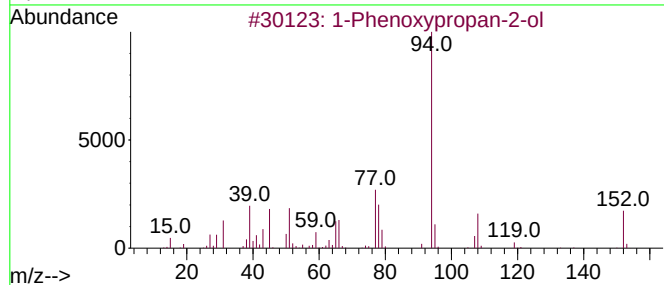
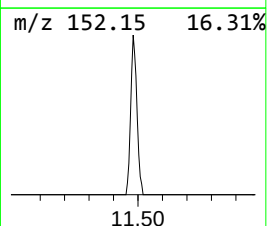
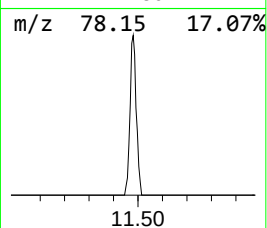
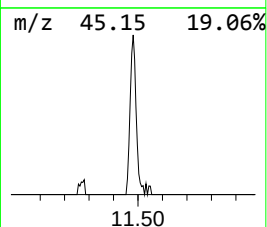
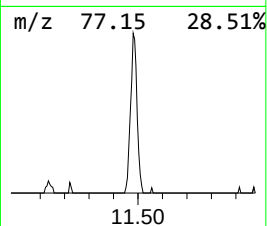
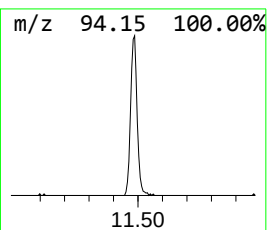
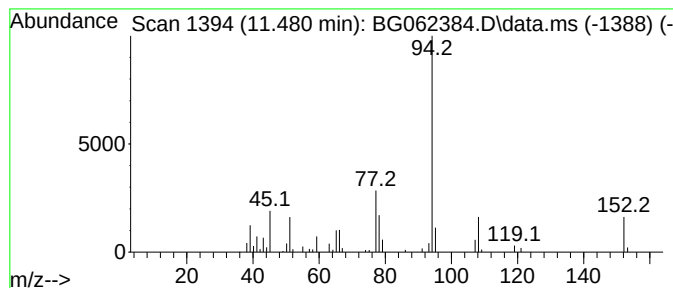
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.480	3.78 ng/ul	108299	Naphthalene-d8	10.746

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-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	94
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93



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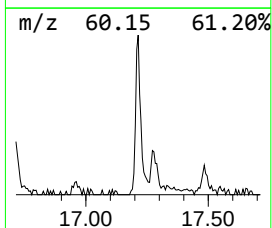
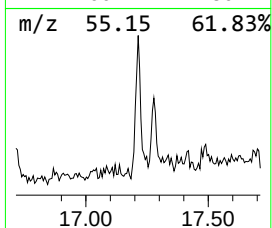
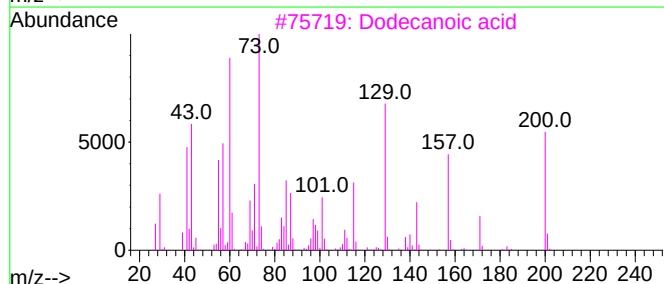
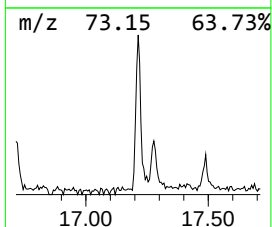
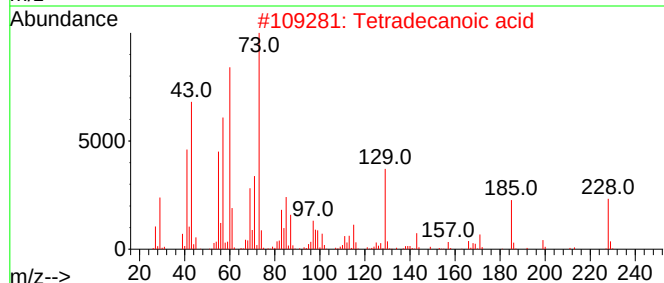
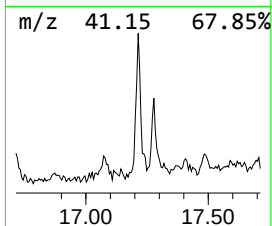
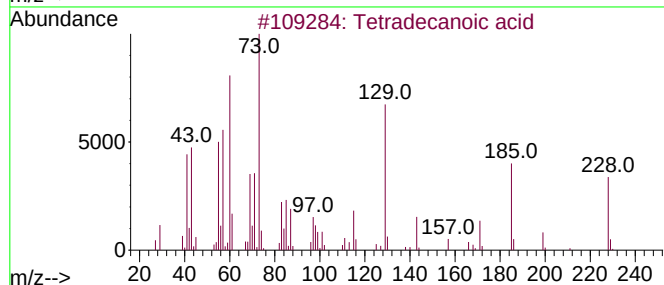
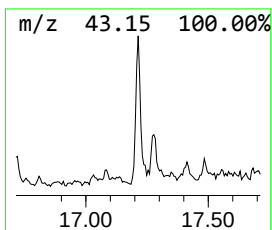
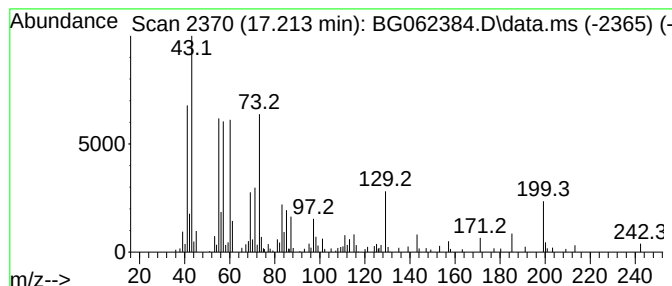
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Tetradecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.213	2.66 ng/ul	145935	Phenanthrene-d10	17.313

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	80
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3			Dodecanoic acid	200	C12H24O2	000143-07-7	64
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	58
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
Data File : BG062384.D
Acq On : 13 Aug 2024 14:10
Operator : MA/JU
Sample : P3502-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCXU7

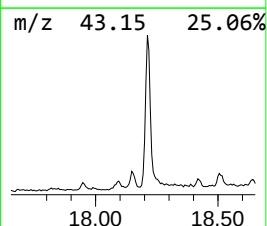
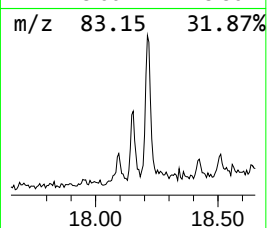
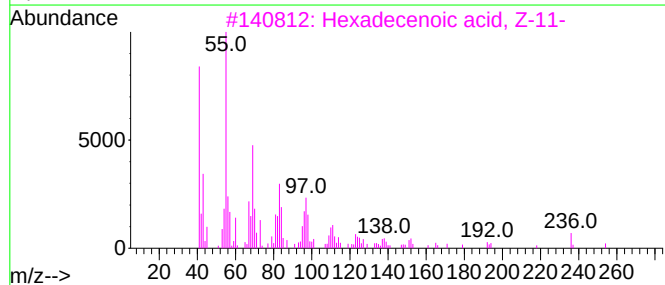
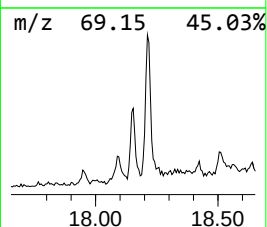
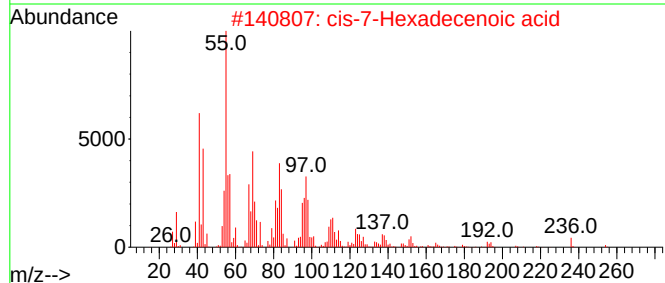
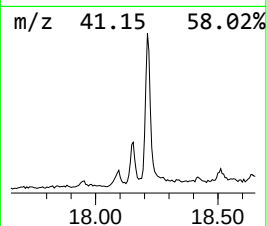
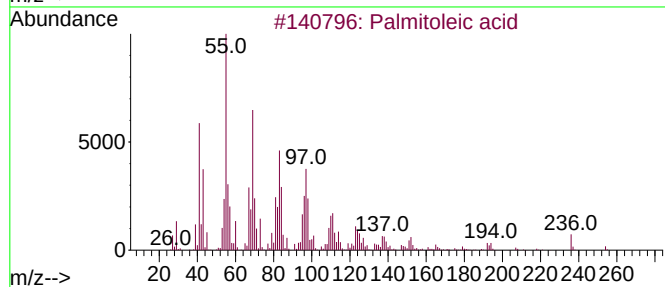
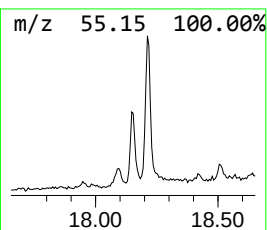
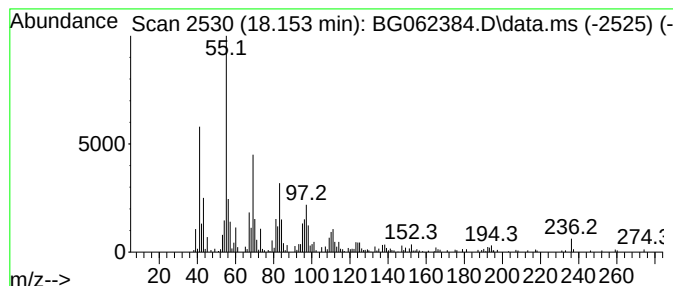
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Palmitoleic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.153	4.09 ng/ul	224447	Phenanthrene-d10	17.313

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	91
2			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	91
3			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91
4			cis-Vaccenic acid	282	C18H34O2	000506-17-2	91
5			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
Data File : BG062384.D
Acq On : 13 Aug 2024 14:10
Operator : MA/JU
Sample : P3502-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCXU7

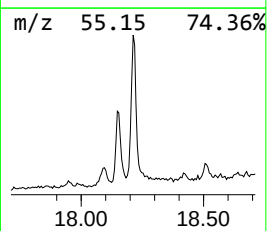
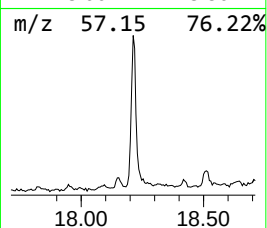
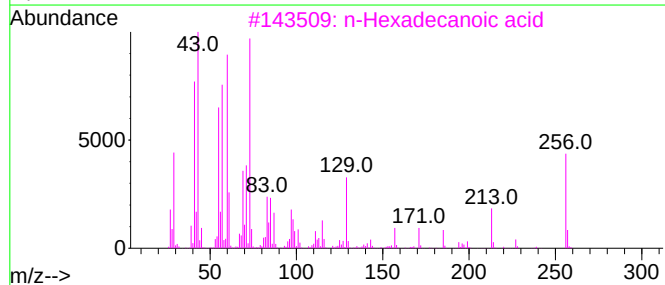
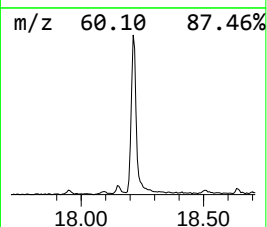
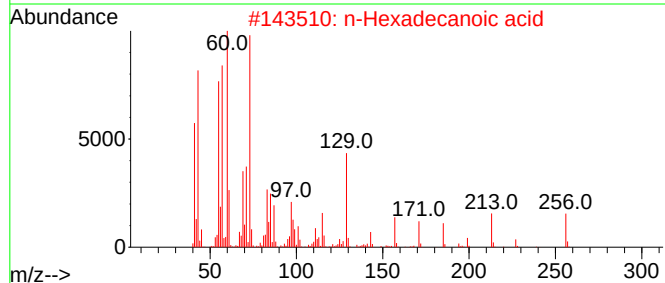
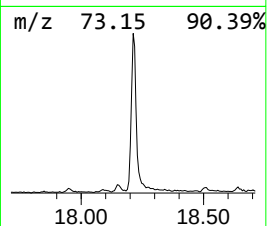
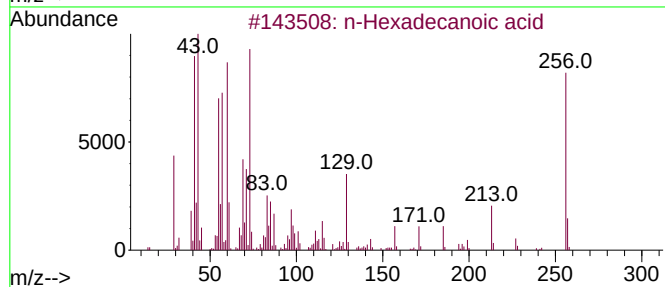
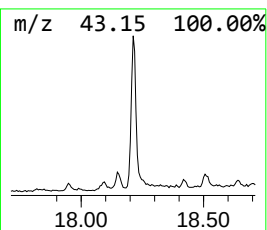
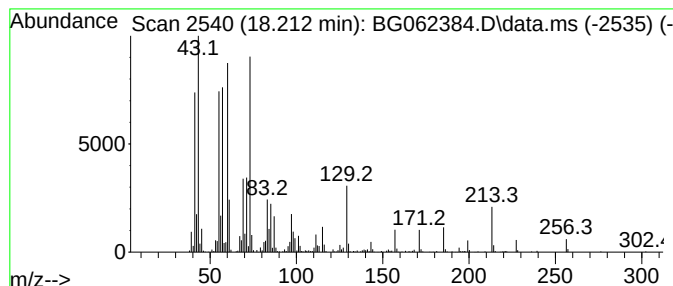
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.212	16.17 ng/ul	888509	Phenanthrene-d10	17.313

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	94
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
Data File : BG062384.D
Acq On : 13 Aug 2024 14:10
Operator : MA/JU
Sample : P3502-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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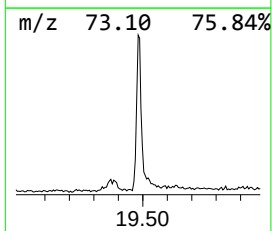
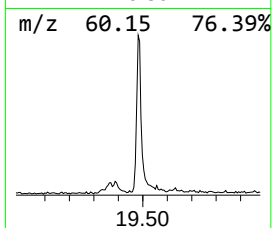
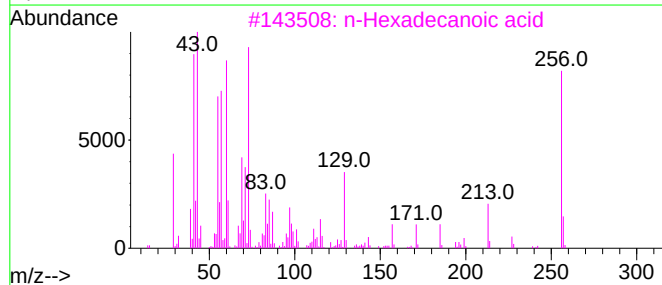
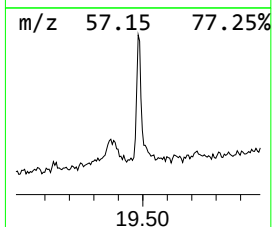
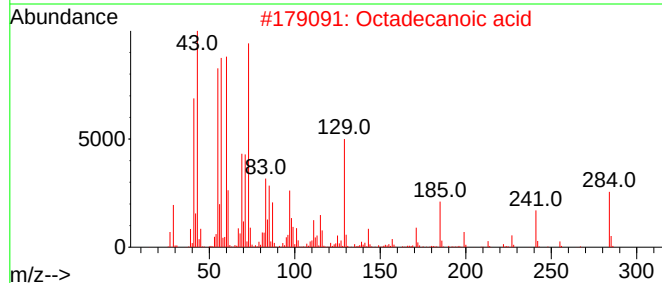
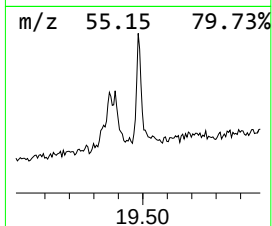
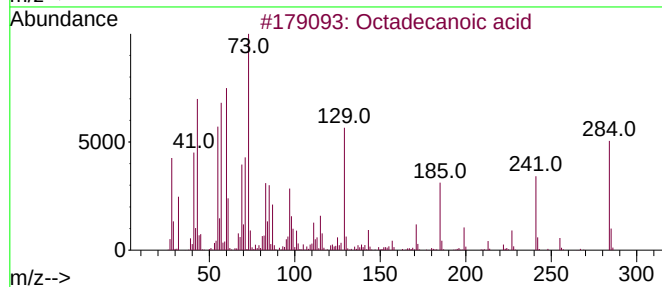
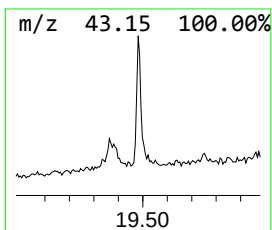
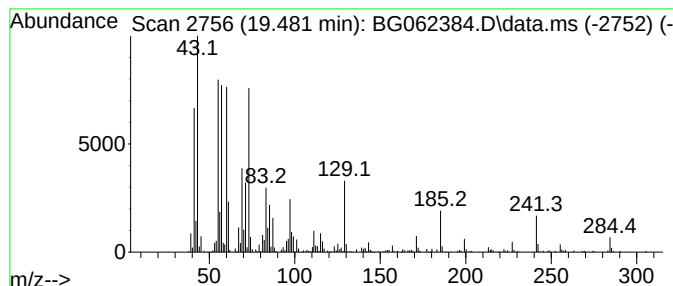
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.481	8.49 ng/ul	466412	Chrysene-d12	21.555

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4			Octadecanoic acid	284	C18H36O2	000057-11-4	90
5			Octadecanoic acid	284	C18H36O2	000057-11-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
Data File : BG062384.D
Acq On : 13 Aug 2024 14:10
Operator : MA/JU
Sample : P3502-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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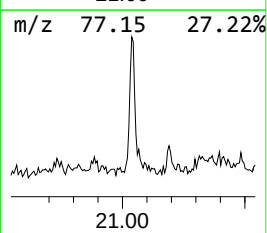
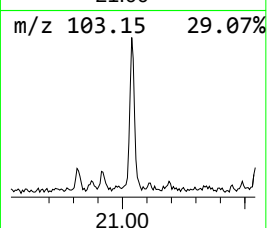
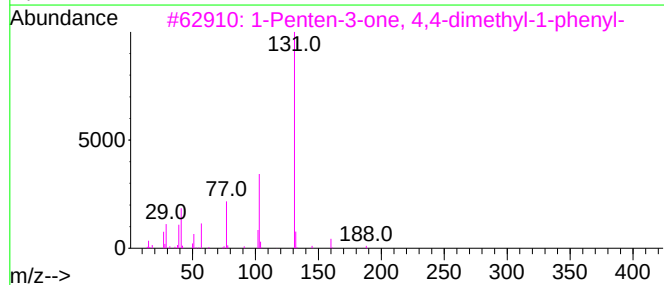
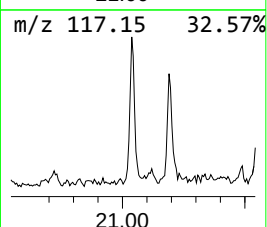
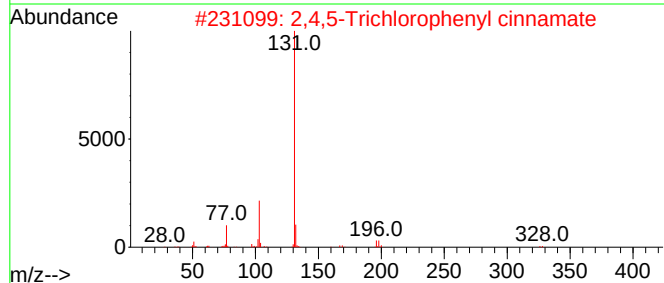
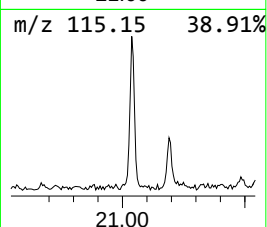
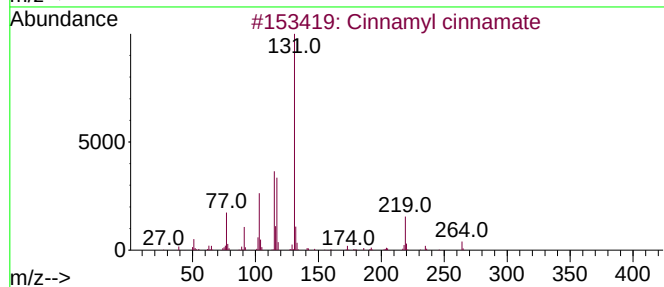
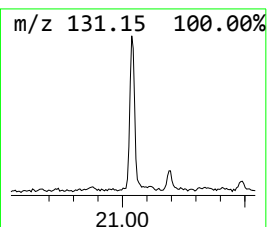
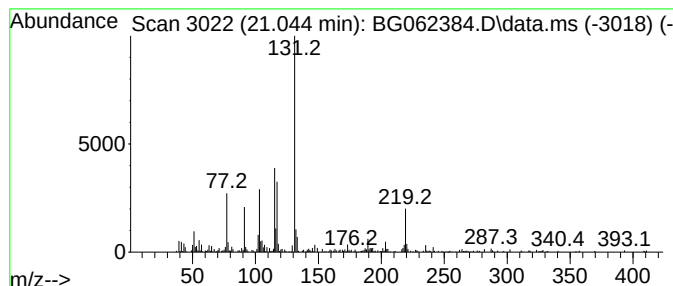
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cinnamyl cinnamate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.044	6.91 ng/ul	379241	Chrysene-d12	21.555

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2			2,4,5-Trichlorophenyl cinnamate	326	C15H9Cl3O2	1010242-39-6	50
3			1-Penten-3-one, 4,4-dimethyl-1-p...	188	C13H16O	000538-44-3	49
4			Propanedioic acid, nitrile, hydr...	229	C12H11N3O2	294878-35-6	46
5			Oxalic acid, monoamide monohydra...	323	C18H17N3O3	1000294-01-4	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
Data File : BG062384.D
Acq On : 13 Aug 2024 14:10
Operator : MA/JU
Sample : P3502-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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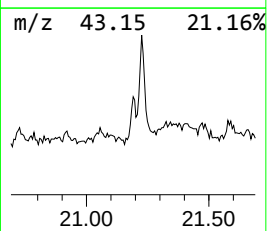
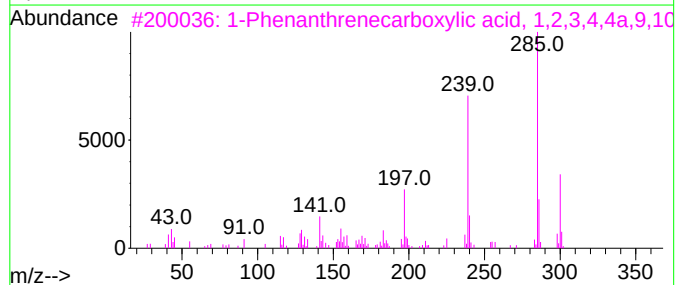
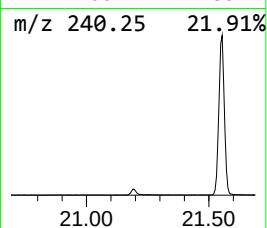
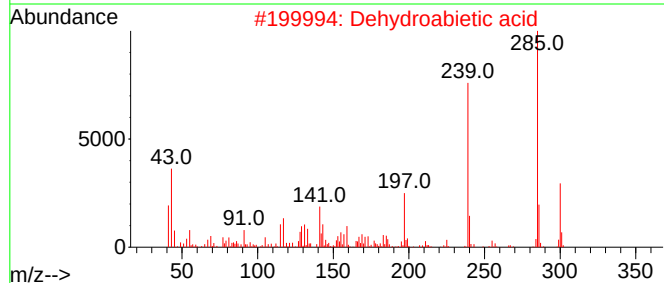
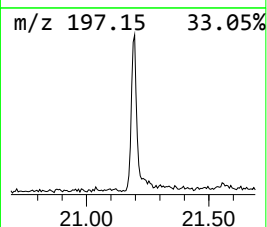
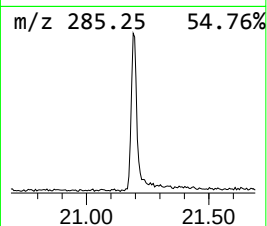
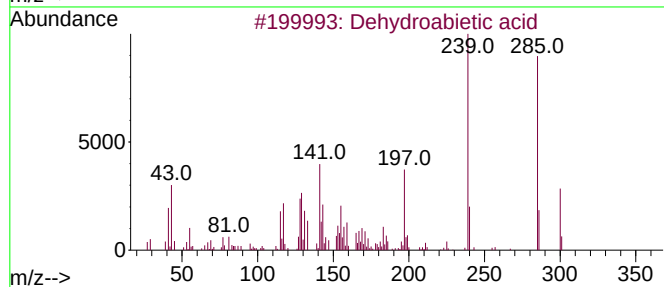
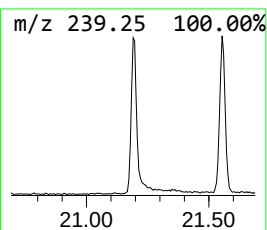
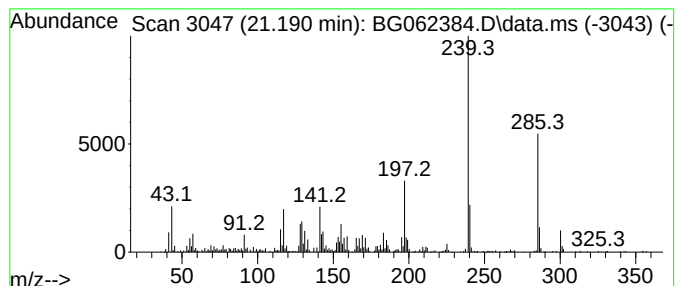
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Dehydroabietic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.190	7.61 ng/ul	417727	Chrysene-d12	21.555

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabietic acid	300	C20H28O2	001740-19-8	96
2			Dehydroabietic acid	300	C20H28O2	001740-19-8	90
3			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	83
4			Butanoic acid, 2-(cyano)(2,4,6-t...	300	C17H20N2O3	1000267-71-7	76
5			3-Quinolinecarboxylic acid, 4-hy...	285	C13H10F3NO3	023851-84-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
Data File : BG062384.D
Acq On : 13 Aug 2024 14:10
Operator : MA/JU
Sample : P3502-07
Misc :
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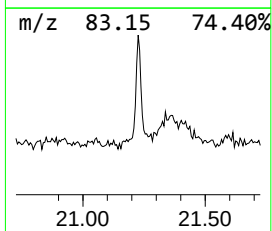
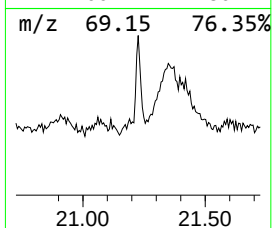
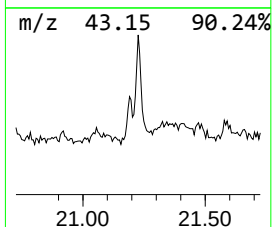
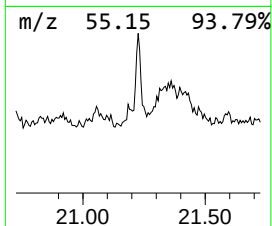
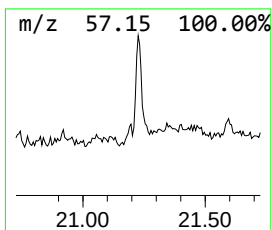
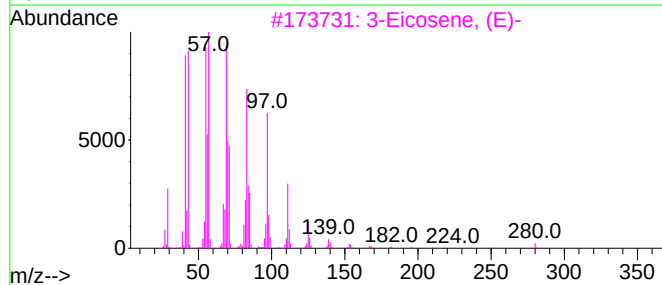
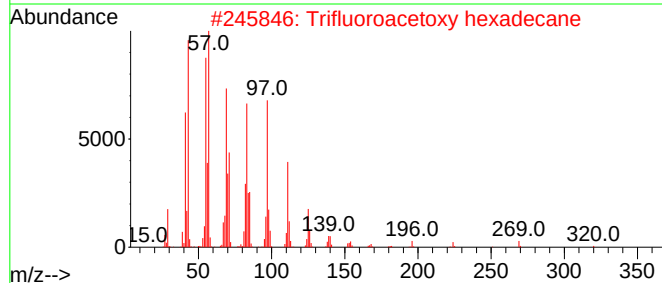
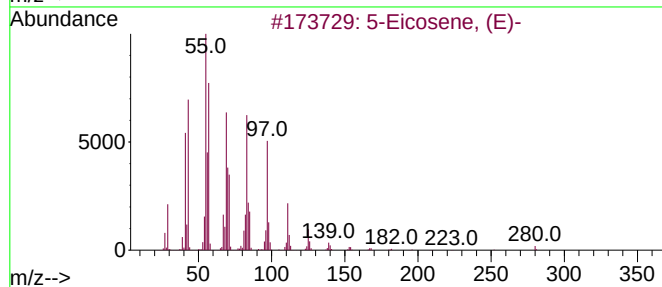
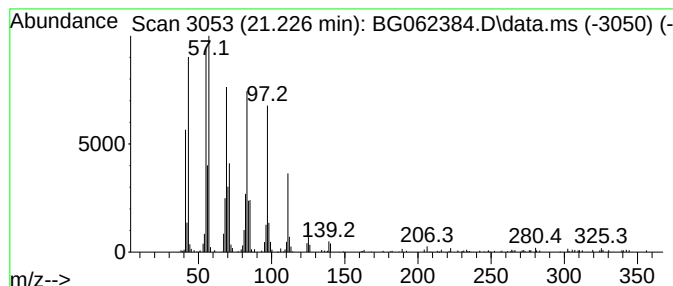
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.226	5.28 ng/ul	289946	Chrysene-d12		21.555	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	95
2	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	94
3	3-Eicosene, (E)-		280	C20H40	074685-33-9	94
4	Behenic alcohol		326	C22H46O	000661-19-8	91
5	n-Tetracosanol-1		354	C24H50O	000506-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081324\
Data File : BG062384.D
Acq On : 13 Aug 2024 14:10
Operator : MA/JU
Sample : P3502-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCXU7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.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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1-Phenoxypropan...	11.480	3.8	ng/ul	108299	2	10.746	573320	20.0
Tetradecanoic acid	17.213	2.7	ng/ul	145935	4	17.313	1098630	20.0
Palmitoleic acid	18.153	4.1	ng/ul	224447	4	17.313	1098630	20.0
n-Hexadecanoic ...	18.212	16.2	ng/ul	888509	4	17.313	1098630	20.0
Octadecanoic acid	19.481	8.5	ng/ul	466412	5	21.555	1098210	20.0
Cinnamyl cinnamate	21.044	6.9	ng/ul	379241	5	21.555	1098210	20.0
Dehydroabietic ...	21.190	7.6	ng/ul	417727	5	21.555	1098210	20.0
(DEL) Alkane: S...	21.226	5.3	ng/ul	289946	5	21.555	1098210	20.0