

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
 Data File : BG062185.D
 Acq On : 23 Jul 2024 14:31
 Operator : MA/JU
 Sample : P3263-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4BA5

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-BG071924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062185.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.737	71	76	83	rBV	34918	57519	4.93%	0.471%
2	3.884	99	101	109	rBV4	13300	20148	1.73%	0.165%
3	3.990	115	119	120	rBV3	4409	5762	0.49%	0.047%
4	4.219	156	158	163	rVB3	3757	4072	0.35%	0.033%
5	4.577	215	219	220	rBV	2734	3385	0.29%	0.028%
6	4.636	226	229	234	rVB2	3272	4433	0.38%	0.036%
7	4.783	249	254	257	rBV4	3211	4071	0.35%	0.033%
8	5.135	307	314	320	rBV2	11222	20563	1.76%	0.168%
9	5.229	327	330	332	rBV2	2077	2839	0.24%	0.023%
10	6.116	476	481	483	rBV2	2478	3618	0.31%	0.030%
11	6.169	487	490	492	rBV	2063	2725	0.23%	0.022%
12	6.927	616	619	623	rBV2	2557	3794	0.33%	0.031%
13	6.974	623	627	630	rVV	1663	2259	0.19%	0.019%
14	7.156	655	658	660	rVB	3368	3401	0.29%	0.028%
15	7.238	664	672	682	rVV	163629	283921	24.35%	2.326%
16	7.385	690	697	705	rBV	98428	162479	13.93%	1.331%
17	7.603	727	734	739	rBV2	137849	243893	20.92%	1.998%
18	7.867	772	779	780	rBV2	1716	2030	0.17%	0.017%
19	8.067	806	813	819	rBV	146252	249021	21.36%	2.040%
20	8.302	850	853	854	rBV	2445	2581	0.22%	0.021%
21	8.548	893	895	898	rBV2	2145	2731	0.23%	0.022%
22	8.783	926	935	942	rBV2	194006	336537	28.86%	2.757%
23	8.889	949	953	958	rBV3	3304	6106	0.52%	0.050%
24	9.236	1004	1012	1020	rBV	153895	272576	23.38%	2.233%
25	9.594	1069	1073	1076	rBV3	1923	2549	0.22%	0.021%
26	9.776	1099	1104	1108	rBV4	6554	11299	0.97%	0.093%
27	9.958	1128	1135	1142	rVV	152673	262474	22.51%	2.150%
28	10.211	1175	1178	1183	rBV	2413	3965	0.34%	0.032%
29	10.311	1193	1195	1198	rVB	2562	2193	0.19%	0.018%
30	10.505	1218	1228	1236	rBV2	227559	408073	35.00%	3.343%
31	10.634	1246	1250	1252	rVB	2486	3145	0.27%	0.026%
32	10.881	1285	1292	1299	rVB	215447	378453	32.46%	3.100%
33	11.016	1309	1315	1324	rBV2	102098	194090	16.64%	1.590%
34	11.110	1329	1331	1333	rBV	2076	2060	0.18%	0.017%
35	11.392	1373	1379	1384	rBV4	8707	15925	1.37%	0.130%
36	11.556	1403	1407	1409	rVB2	1598	2048	0.18%	0.017%

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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-BG071924.MA.M
 Title : SVOA CALIBRATION

37	11.609	1409	1416	1428	rBV2	45646	95493	8.19%	0.782%
38	11.709	1430	1433	1437	rVB	2890	2805	0.24%	0.023%
39	11.750	1437	1440	1442	rBV	1901	2409	0.21%	0.020%
40	12.261	1523	1527	1528	rBV	1922	2606	0.22%	0.021%
41	12.326	1534	1538	1539	rBV	1758	2121	0.18%	0.017%
42	12.449	1555	1559	1564	rBV3	4133	6521	0.56%	0.053%
43	12.743	1606	1609	1611	rBV	2273	2961	0.25%	0.024%
44	12.807	1617	1620	1623	rVB	1795	2207	0.19%	0.018%
45	12.854	1626	1628	1631	rVB	1889	2023	0.17%	0.017%
46	13.031	1655	1658	1661	rBV	2993	4204	0.36%	0.034%
47	13.072	1661	1665	1667	rVV	2519	2038	0.17%	0.017%
48	13.148	1675	1678	1679	rBV	3287	2649	0.23%	0.022%
49	13.195	1683	1686	1688	rVV	1673	2375	0.20%	0.019%
50	13.295	1701	1703	1707	rVB	2769	2373	0.20%	0.019%
51	13.365	1714	1715	1719	rBV2	1928	2172	0.19%	0.018%
52	13.489	1734	1736	1742	rVB3	3826	5632	0.48%	0.046%
53	13.542	1742	1745	1748	rBV	3329	3113	0.27%	0.025%
54	13.577	1748	1751	1753	rBV	1438	2283	0.20%	0.019%
55	13.624	1757	1759	1762	rVB	2491	2299	0.20%	0.019%
56	13.871	1796	1801	1805	rVV2	2760	5521	0.47%	0.045%
57	13.900	1805	1806	1810	rVV	1870	2065	0.18%	0.017%
58	13.953	1812	1815	1816	rBV	3031	3269	0.28%	0.027%
59	14.094	1831	1839	1845	rBV	409654	606444	52.01%	4.968%
60	14.311	1873	1876	1877	rBV	4174	4542	0.39%	0.037%
61	14.393	1882	1890	1901	rVB	438452	697720	59.84%	5.715%
62	14.511	1907	1910	1913	rBV3	2930	4092	0.35%	0.034%
63	14.564	1917	1919	1923	rVV2	3082	4475	0.38%	0.037%
64	14.617	1926	1928	1930	rVV2	2719	2869	0.25%	0.024%
65	14.693	1934	1941	1949	rBV	397954	619320	53.11%	5.073%
66	14.752	1949	1951	1958	rVB3	9147	14493	1.24%	0.119%
67	14.916	1966	1979	1990	rBV2	165333	346222	29.69%	2.836%
68	15.051	2000	2002	2005	rVV4	11220	14944	1.28%	0.122%
69	15.104	2007	2011	2017	rVV5	9228	19162	1.64%	0.157%
70	15.151	2017	2019	2021	rVV2	3299	3728	0.32%	0.031%
71	15.192	2025	2026	2031	rVB3	3827	3125	0.27%	0.026%
72	15.345	2050	2052	2053	rBV	3331	2411	0.21%	0.020%
73	15.404	2061	2062	2064	rBV	2547	2282	0.20%	0.019%
74	15.510	2079	2080	2087	rVB4	8343	10881	0.93%	0.089%
75	15.580	2090	2092	2095	rVB3	2502	2807	0.24%	0.023%
76	15.627	2095	2100	2103	rBV	3214	5705	0.49%	0.047%

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Integrator: RTE

Smoothing : OFF

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Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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 Title : SVOA CALIBRATION

77	15.686	2103	2110	2117	rBV	584707	902123	77.36%	7.389%
78	15.739	2117	2119	2125	rVB3	11702	14868	1.28%	0.122%
79	15.809	2125	2131	2139	rBV	166313	260411	22.33%	2.133%
80	15.897	2144	2146	2147	rBV2	10174	7277	0.62%	0.060%
81	16.179	2191	2194	2195	rBV3	6021	3772	0.32%	0.031%
82	16.314	2212	2217	2219	rBV4	4356	8050	0.69%	0.066%
83	16.373	2225	2227	2228	rBV2	6612	5810	0.50%	0.048%
84	16.502	2246	2249	2250	rBV	4080	3604	0.31%	0.030%
85	16.661	2271	2276	2279	rBV5	6926	10588	0.91%	0.087%
86	16.731	2286	2288	2291	rBV3	4707	4653	0.40%	0.038%
87	17.096	2347	2350	2352	rBV4	7818	10328	0.89%	0.085%
88	17.254	2375	2377	2383	rVB4	12340	17774	1.52%	0.146%
89	17.436	2402	2408	2413	rBV	516169	809799	69.45%	6.633%
90	17.483	2413	2416	2420	rVB	154171	207443	17.79%	1.699%
91	17.536	2420	2425	2431	rBV	633391	973070	83.45%	7.971%
92	18.024	2505	2508	2513	rVB7	13949	23183	1.99%	0.190%
93	18.077	2515	2517	2520	rBV3	13606	15699	1.35%	0.129%
94	18.312	2554	2557	2561	rVB2	23330	28710	2.46%	0.235%
95	18.506	2586	2590	2594	rBV2	37590	53328	4.57%	0.437%
96	19.487	2753	2757	2763	rVB	406212	525856	45.10%	4.307%
97	19.821	2808	2814	2817	rBV	736539	1166067	100.00%	9.552%
98	21.566	3108	3111	3116	rVB	113302	139979	12.00%	1.147%
99	21.701	3128	3134	3139	rBV3	515981	994750	85.31%	8.148%
100	24.715	3642	3647	3655	rVB2	226658	509966	43.73%	4.177%

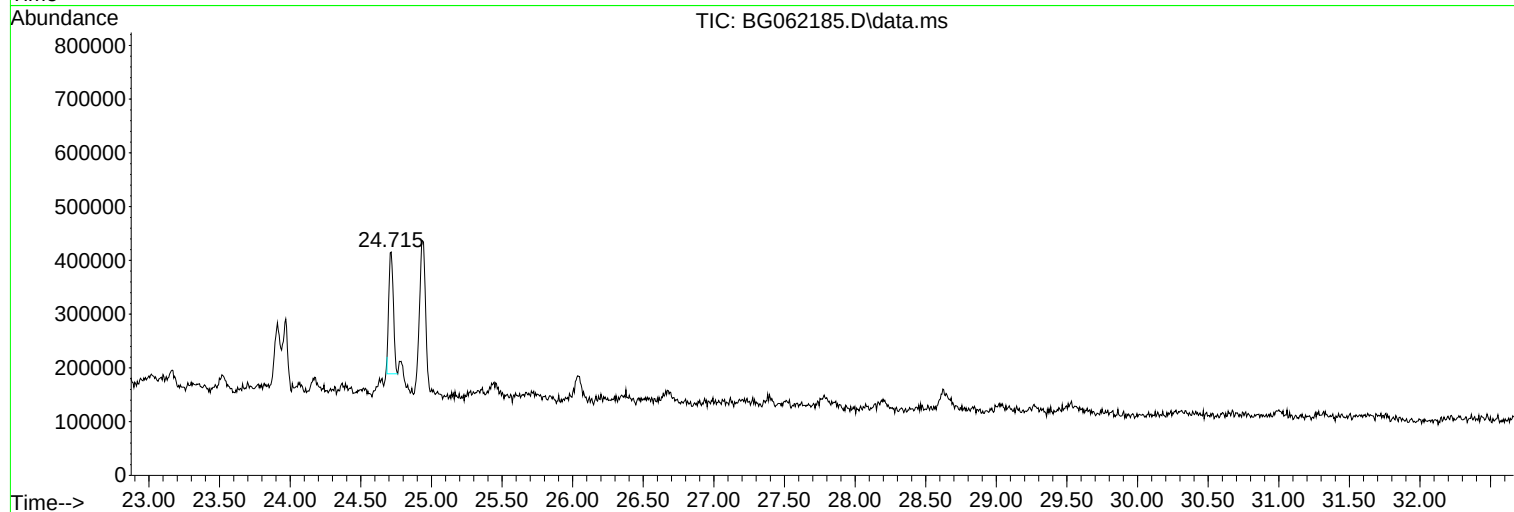
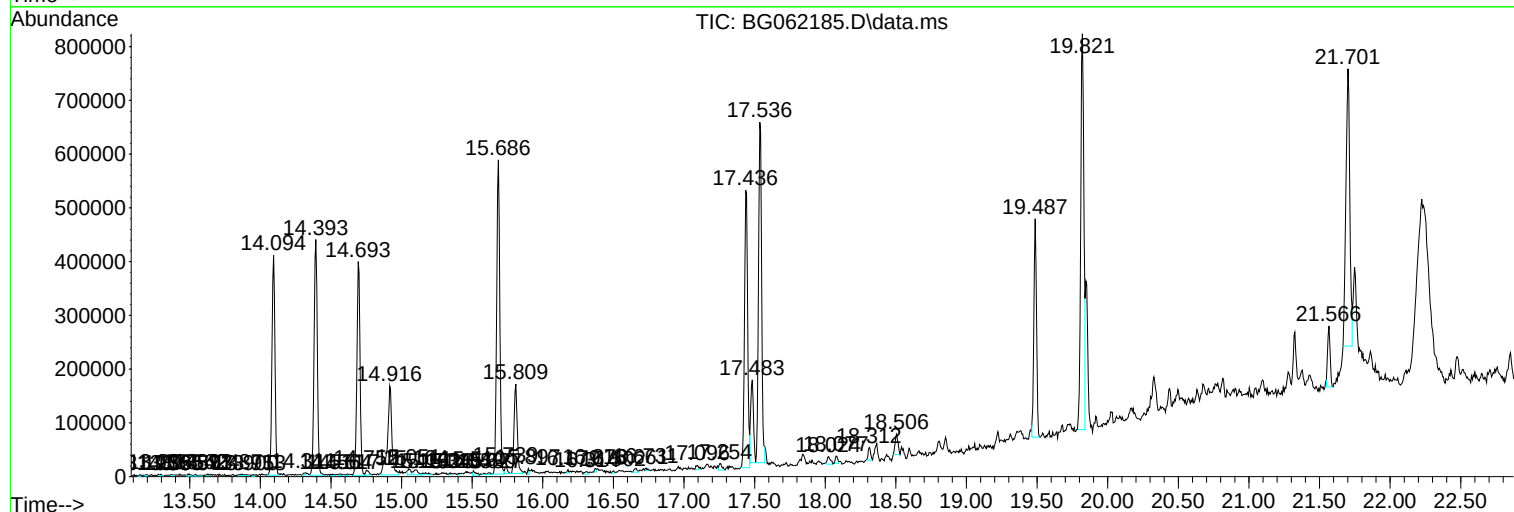
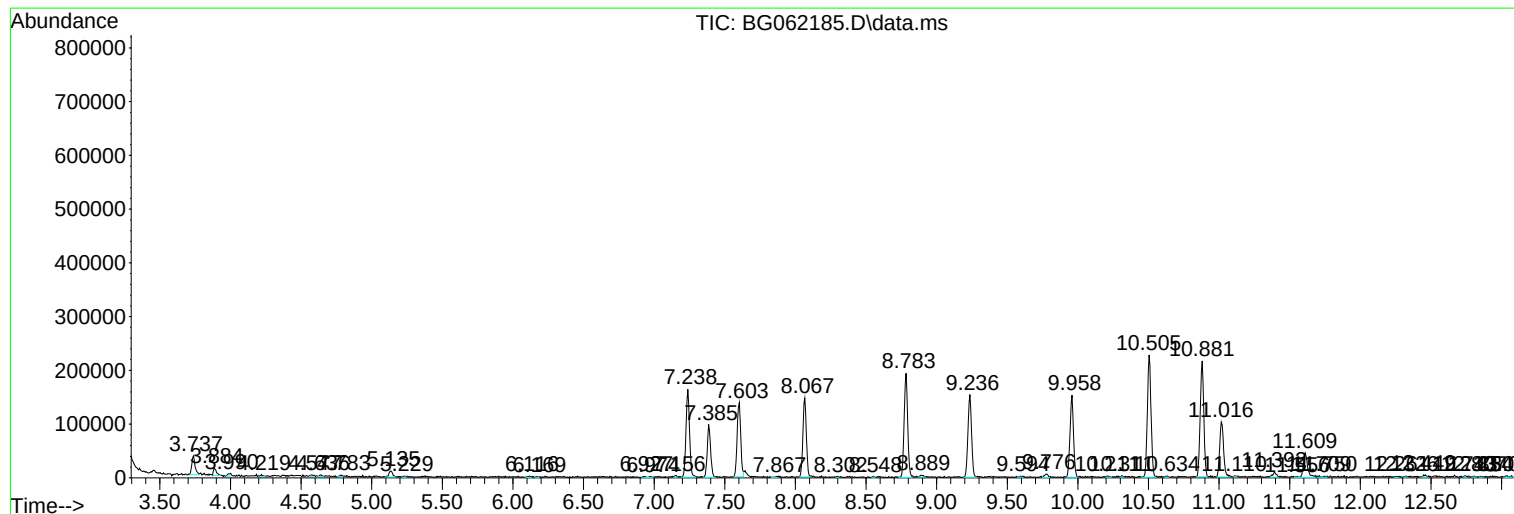
Sum of corrected areas: 12208207

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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



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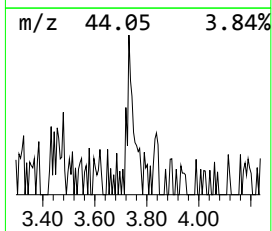
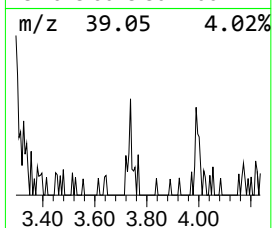
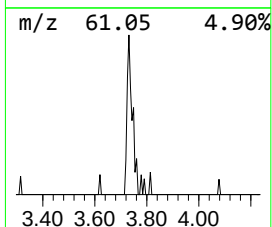
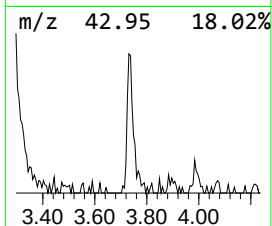
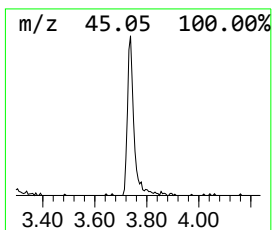
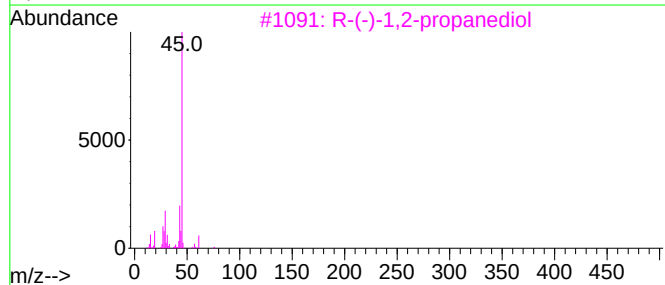
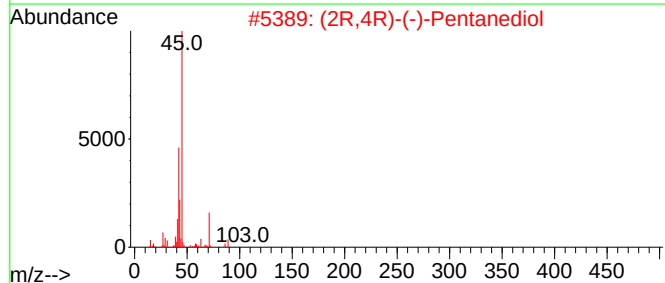
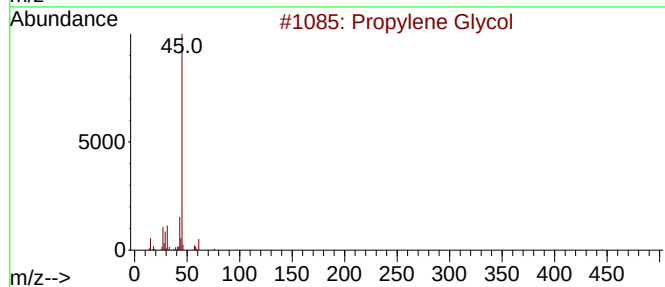
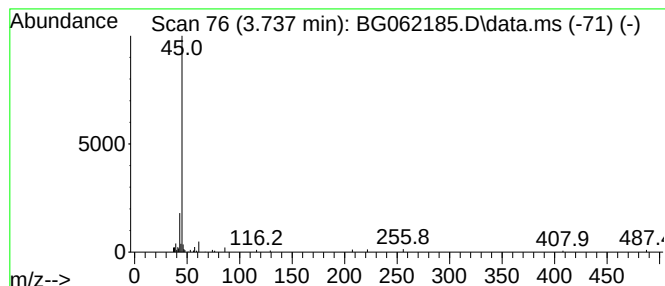
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.737	4.62 ng/ul	57519	1,4-Dichlorobenzene-d4	8.067

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	64
2			(2R,4R)-(-)-Pentanediol	104	C5H12O2	042075-32-1	39
3			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	39
4			Propylene Glycol	76	C3H8O2	000057-55-6	39
5			Propylene Glycol	76	C3H8O2	000057-55-6	9



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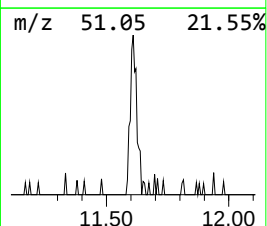
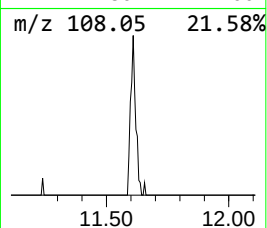
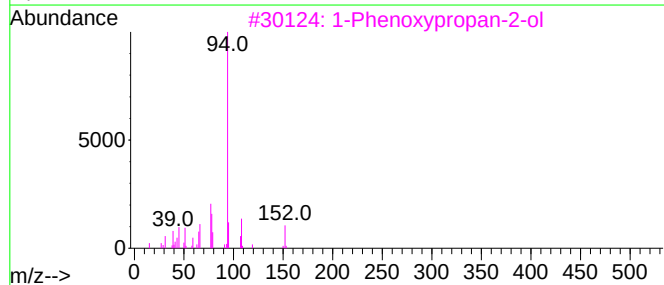
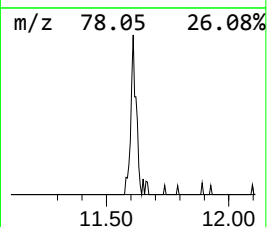
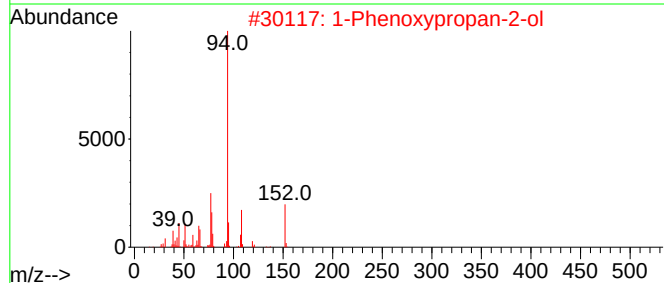
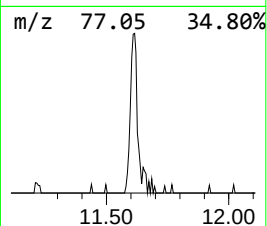
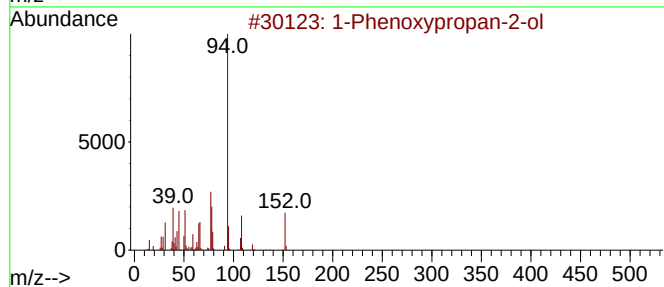
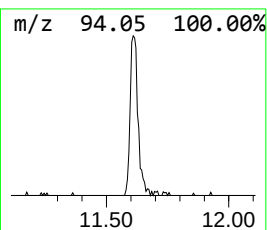
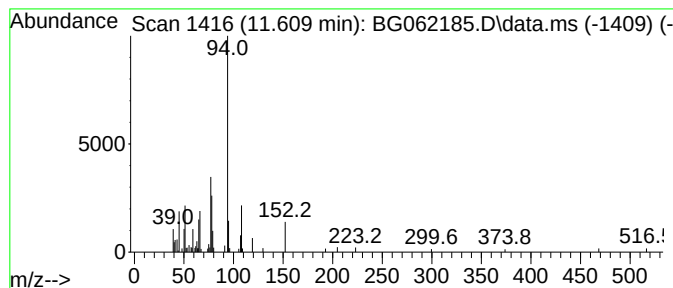
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.609	5.05 ng/ul	95493	Naphthalene-d8	10.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	78
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	59



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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
Propylene Glycol	3.737	4.6	ng/u1	57519	1	8.067	249021	20.0
1-Phenoxypropan...	11.609	5.0	ng/u1	95493	2	10.881	378453	20.0