

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
 Data File : BG062095.D
 Acq On : 16 Jul 2024 15:34
 Operator : MA/JU
 Sample : P3177-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4BS1

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-BG070224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062095.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.317	3	5	8	rBV3	3683	3814	0.40%	0.034%
2	3.346	8	10	14	rVV3	4198	5244	0.55%	0.047%
3	3.382	14	16	20	rVB5	3684	4730	0.49%	0.042%
4	3.440	24	26	34	rVB5	8877	12136	1.27%	0.108%
5	3.687	64	68	69	rBV2	2222	2412	0.25%	0.021%
6	3.722	70	74	82	rVV	23727	38045	3.98%	0.339%
7	3.863	95	98	107	rBV	68921	110583	11.57%	0.986%
8	3.957	113	114	116	rBV	3150	2394	0.25%	0.021%
9	4.187	152	153	155	rVV2	3482	2787	0.29%	0.025%
10	4.204	155	156	159	rVV3	3505	3185	0.33%	0.028%
11	4.392	185	188	194	rBV2	1523	2955	0.31%	0.026%
12	4.950	279	283	284	rBV2	2549	3051	0.32%	0.027%
13	5.115	308	311	312	rBV	4069	3464	0.36%	0.031%
14	5.133	312	314	316	rVB3	3392	2828	0.30%	0.025%
15	5.415	359	362	364	rBV	2418	2311	0.24%	0.021%
16	5.456	364	369	370	rBV	1691	2329	0.24%	0.021%
17	5.667	402	405	408	rVB	1865	2521	0.26%	0.022%
18	5.691	408	409	413	rBV2	2449	2515	0.26%	0.022%
19	6.272	505	508	510	rBV	2240	2241	0.23%	0.020%
20	6.378	522	526	528	rBV2	1969	2438	0.25%	0.022%
21	6.960	622	625	626	rBV2	3334	2257	0.24%	0.020%
22	7.060	638	642	646	rVB2	3276	5827	0.61%	0.052%
23	7.230	664	671	679	rVB	193085	331596	34.68%	2.955%
24	7.330	685	688	689	rBV	2753	2342	0.24%	0.021%
25	7.377	689	696	704	rVB	108003	182289	19.07%	1.625%
26	7.588	725	732	738	rBV	182570	307082	32.12%	2.737%
27	7.823	766	772	775	rBV2	1688	2954	0.31%	0.026%
28	8.053	804	811	818	rVV	176040	301177	31.50%	2.684%
29	8.100	818	819	825	rVB3	2992	3860	0.40%	0.034%
30	8.241	841	843	848	rVB3	1958	2485	0.26%	0.022%
31	8.317	852	856	859	rVB2	2499	3688	0.39%	0.033%
32	8.364	859	864	865	rBV2	2185	3155	0.33%	0.028%
33	8.417	870	873	879	rVB4	1865	2912	0.30%	0.026%
34	8.534	889	893	895	rBV4	2942	4113	0.43%	0.037%
35	8.564	897	898	903	rVB	1724	2225	0.23%	0.020%
36	8.775	923	934	943	rBV2	239468	407922	42.66%	3.636%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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37	8.899	953	955	960	rVB	2344	2636	0.28%	0.023%
38	9.222	1001	1010	1018	rBV2	196376	349298	36.53%	3.113%
39	9.275	1018	1019	1024	rVB2	2266	2431	0.25%	0.022%
40	9.322	1024	1027	1031	rBV	2566	3349	0.35%	0.030%
41	9.439	1043	1047	1049	rBV	2291	2943	0.31%	0.026%
42	9.668	1084	1086	1089	rBV	2113	2678	0.28%	0.024%
43	9.762	1098	1102	1111	rVB2	17433	29231	3.06%	0.261%
44	9.850	1116	1117	1121	rVV	1974	2227	0.23%	0.020%
45	9.944	1126	1133	1141	rBV	185797	330982	34.62%	2.950%
46	10.044	1147	1150	1152	rBV	2853	2363	0.25%	0.021%
47	10.185	1167	1174	1177	rBV3	2038	4962	0.52%	0.044%
48	10.356	1201	1203	1206	rBV2	2961	2654	0.28%	0.024%
49	10.497	1218	1227	1234	rBV	239519	429113	44.88%	3.825%
50	10.749	1267	1270	1272	rBV	2152	2556	0.27%	0.023%
51	10.867	1283	1290	1299	rVB	262881	461739	48.29%	4.115%
52	11.002	1306	1313	1320	rBV	197888	352166	36.83%	3.139%
53	11.125	1331	1334	1338	rBV	1371	2220	0.23%	0.020%
54	11.302	1363	1364	1368	rBV	2638	2598	0.27%	0.023%
55	11.384	1373	1378	1385	rVB4	11968	20900	2.19%	0.186%
56	11.549	1405	1406	1409	rBV4	3394	2741	0.29%	0.024%
57	11.601	1409	1415	1423	rVV	66599	120311	12.58%	1.072%
58	11.660	1423	1425	1426	rVV	5265	3963	0.41%	0.035%
59	11.695	1429	1431	1435	rVB3	1647	2289	0.24%	0.020%
60	12.254	1523	1526	1528	rVB	2607	2371	0.25%	0.021%
61	12.318	1534	1537	1538	rBV	3262	2329	0.24%	0.021%
62	12.518	1568	1571	1573	rBV	2586	2912	0.30%	0.026%
63	12.600	1581	1585	1587	rVB3	2157	2237	0.23%	0.020%
64	12.735	1606	1608	1611	rBV2	2089	2536	0.27%	0.023%
65	12.794	1614	1618	1620	rVB	2411	2459	0.26%	0.022%
66	13.082	1665	1667	1670	rBV	2319	2949	0.31%	0.026%
67	13.188	1682	1685	1688	rVB3	4159	4832	0.51%	0.043%
68	13.294	1699	1703	1707	rBV3	3587	6308	0.66%	0.056%
69	13.323	1707	1708	1710	rBV	4185	2477	0.26%	0.022%
70	13.452	1729	1730	1733	rBV	2204	2306	0.24%	0.021%
71	13.482	1733	1735	1737	rBV2	3403	2917	0.31%	0.026%
72	14.010	1824	1825	1829	rVB2	3383	2459	0.26%	0.022%
73	14.087	1831	1838	1845	rVV	478666	691982	72.37%	6.168%
74	14.157	1847	1850	1853	rBV3	2809	2752	0.29%	0.025%
75	14.234	1861	1863	1867	rVB3	2448	2248	0.24%	0.020%
76	14.322	1875	1878	1882	rVV3	6685	9110	0.95%	0.081%

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77	14.386	1882	1889	1896	rVV	468563	764678	79.98%	6.815%
78	14.469	1899	1903	1907	rBV4	3536	4792	0.50%	0.043%
79	14.551	1915	1917	1922	rVB3	3445	3653	0.38%	0.033%
80	14.621	1925	1929	1930	rBV3	4776	4649	0.49%	0.041%
81	14.686	1934	1940	1946	rVV	447299	679425	71.06%	6.056%
82	14.792	1956	1958	1960	rBV2	3457	2221	0.23%	0.020%
83	14.868	1967	1971	1972	rBV2	28563	33183	3.47%	0.296%
84	14.909	1973	1978	1988	rVV	186058	333775	34.91%	2.975%
85	14.974	1988	1989	1992	rVB3	5868	3393	0.35%	0.030%
86	15.091	2003	2009	2012	rBV6	8789	16285	1.70%	0.145%
87	15.450	2068	2070	2072	rBV3	6602	5005	0.52%	0.045%
88	15.509	2076	2080	2082	rBV4	7257	9892	1.03%	0.088%
89	15.567	2089	2090	2093	rVV3	4939	3949	0.41%	0.035%
90	15.679	2103	2109	2114	rBV	597996	892874	93.39%	7.958%
91	15.796	2123	2129	2135	rBV	185697	286689	29.99%	2.555%
92	15.984	2159	2161	2162	rVB2	6717	2960	0.31%	0.026%
93	16.366	2223	2226	2227	rBV3	5110	5374	0.56%	0.048%
94	16.572	2259	2261	2262	rBV2	7013	4097	0.43%	0.037%
95	17.436	2402	2408	2412	rBV	491432	763581	79.86%	6.806%
96	17.530	2418	2424	2431	rBV	603963	926146	96.87%	8.255%
97	18.194	2534	2537	2541	rBV3	47073	60399	6.32%	0.538%
98	18.487	2584	2587	2593	rVB6	64411	88048	9.21%	0.785%
99	19.815	2808	2813	2824	rVB	621931	956108	100.00%	8.522%
100	21.695	3129	3133	3138	rVB	421985	709156	74.17%	6.321%

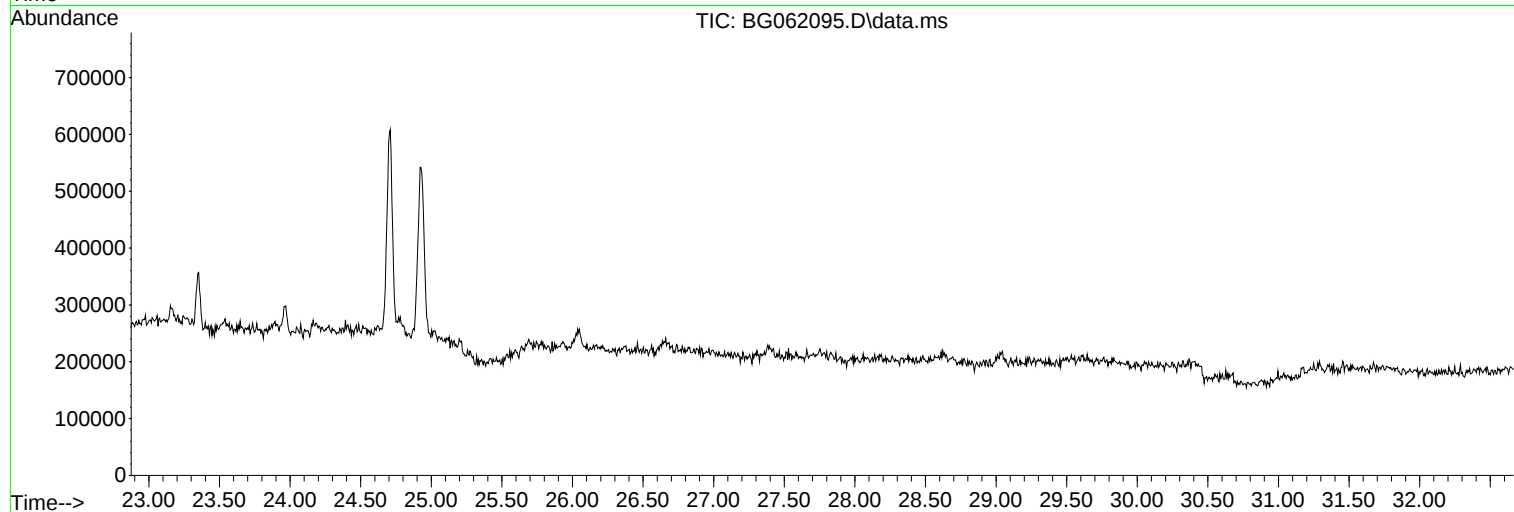
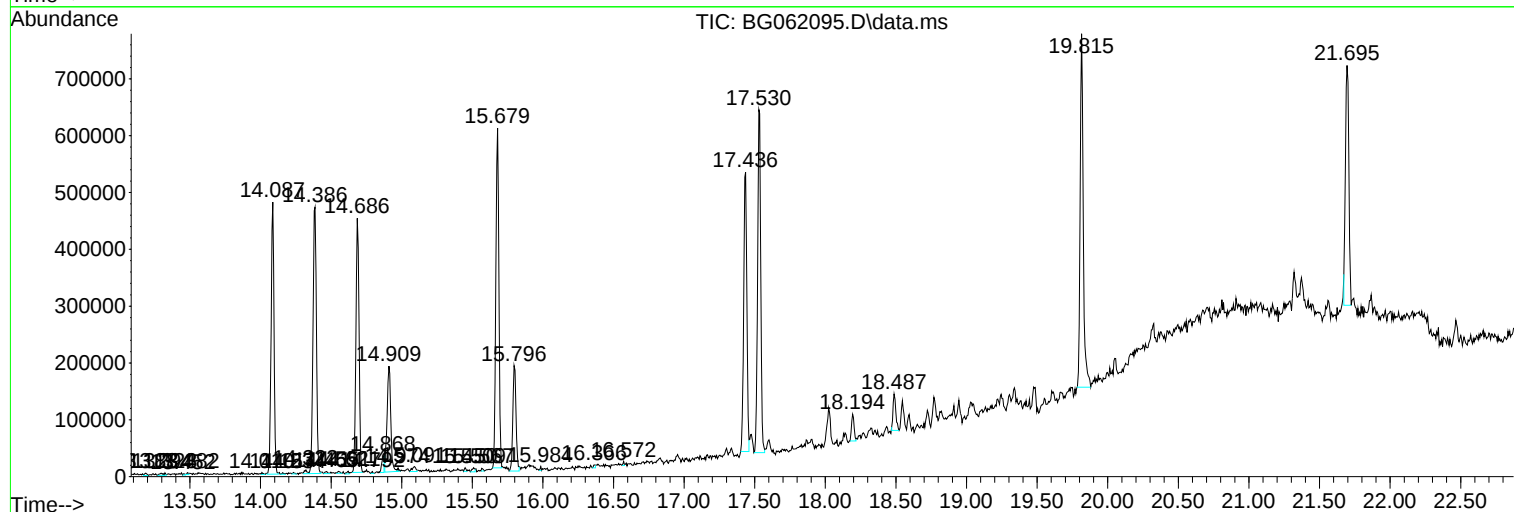
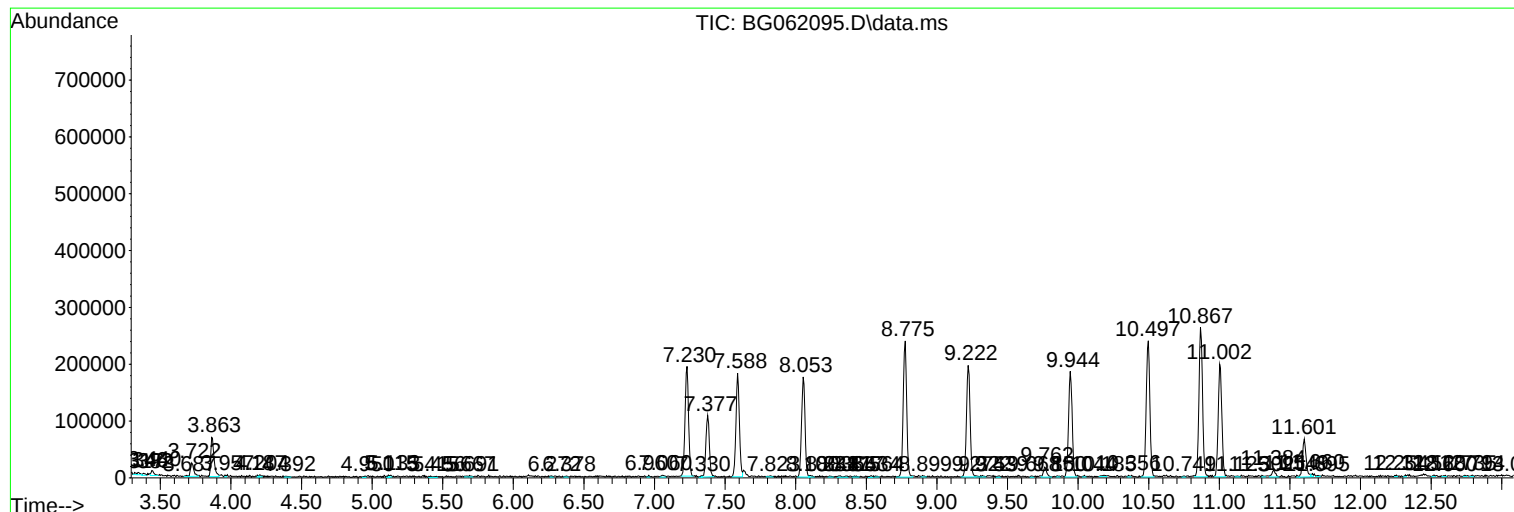
Sum of corrected areas: 11219733

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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



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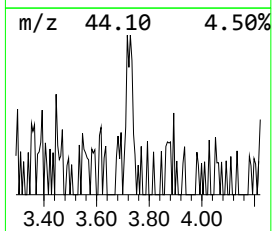
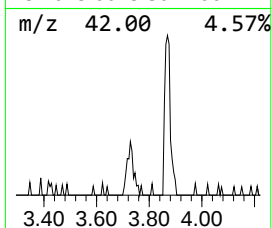
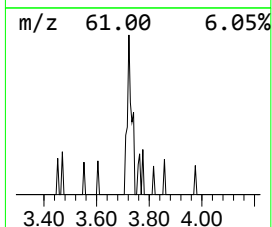
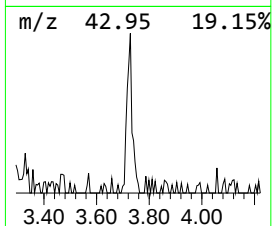
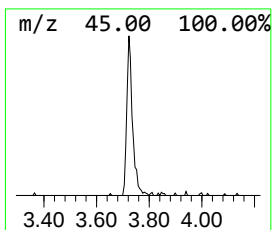
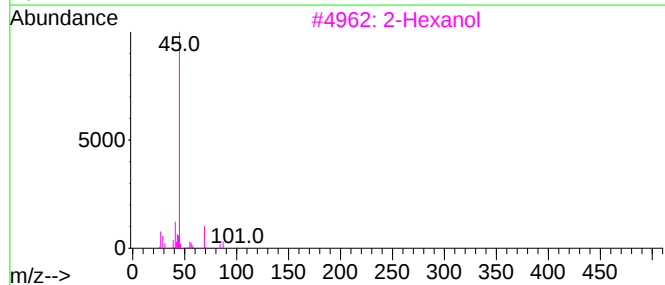
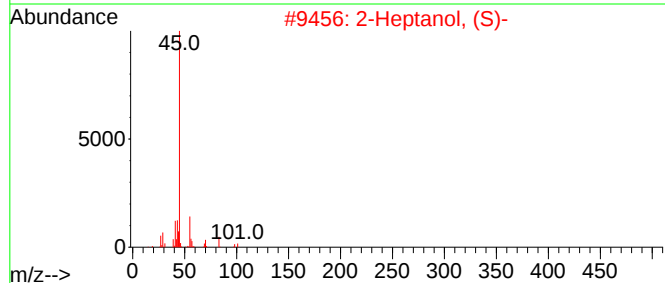
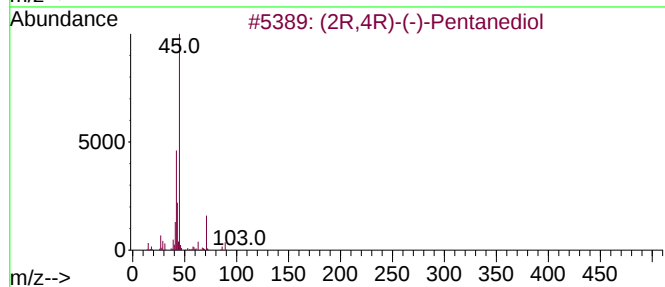
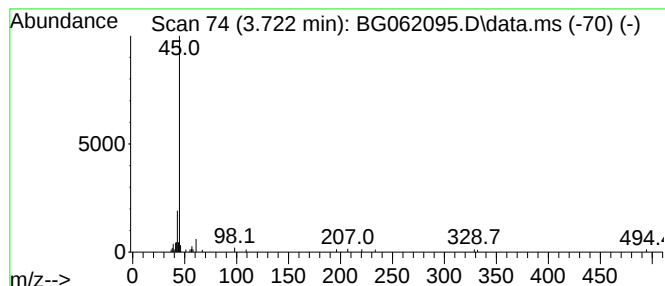
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.722	2.53 ng/ul	38045	1,4-Dichlorobenzene-d4	8.053

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2R,4R)-(-)-Pentanediol			104	C5H12O2	042075-32-1	38
2	2-Heptanol, (S)-			116	C7H16O	006033-23-4	9
3	2-Hexanol			102	C6H14O	000626-93-7	9
4	2,4-Pentanediol			104	C5H12O2	000625-69-4	9
5	Propane, 2-ethoxy-			88	C5H12O	000625-54-7	9



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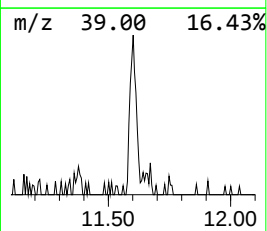
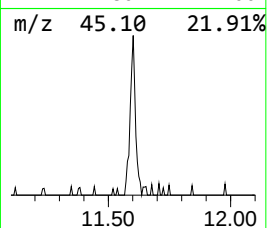
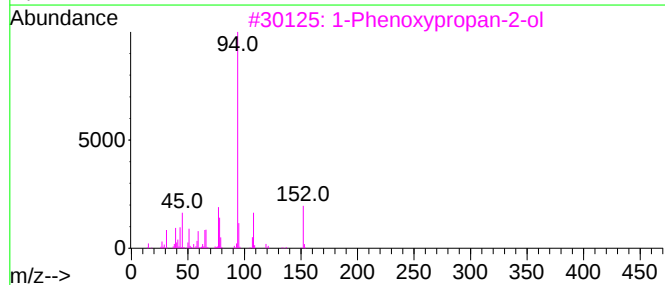
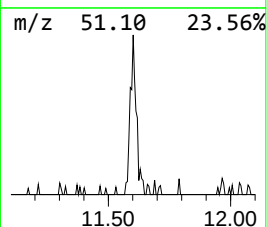
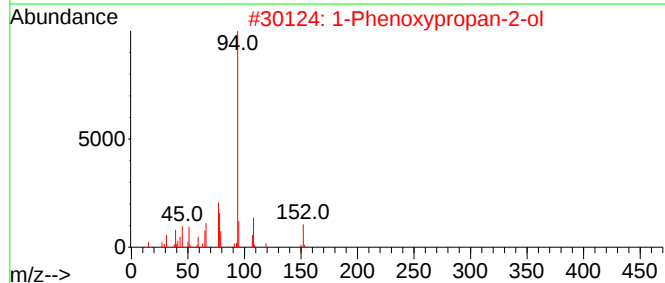
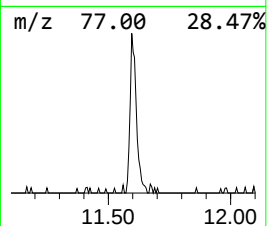
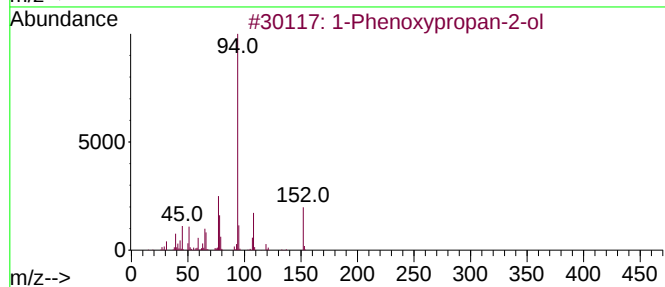
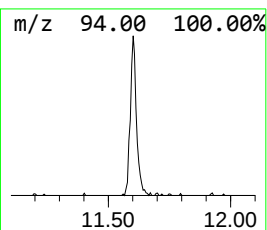
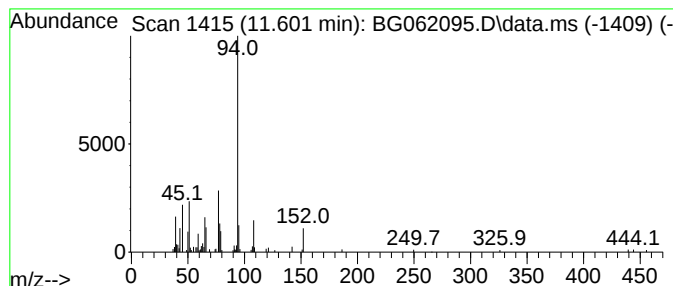
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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.
11.601	5.21 ng/ul	120311	Naphthalene-d8	10.867

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	83
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	80
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	59
5			Diphenyl phosphate	250	C12H11O4P	000838-85-7	53



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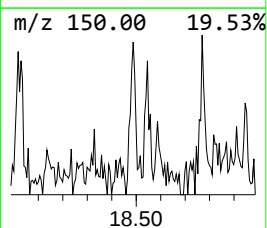
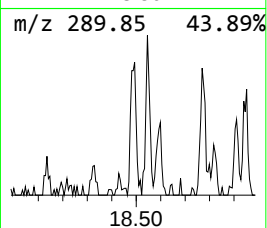
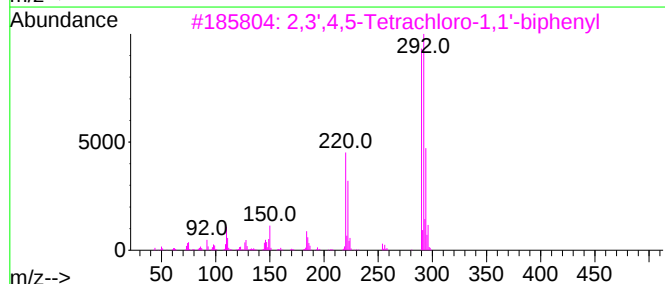
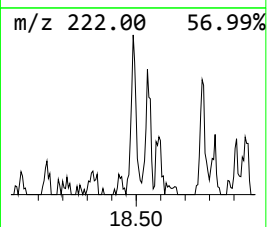
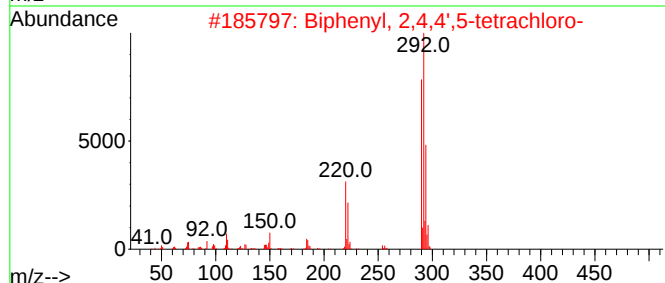
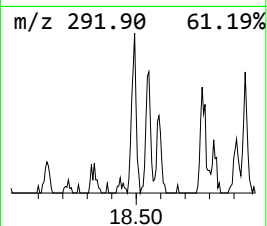
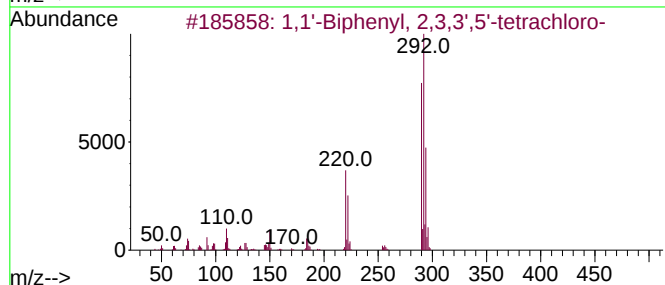
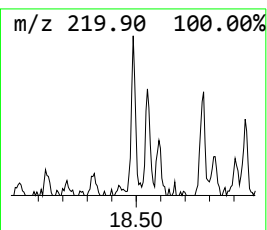
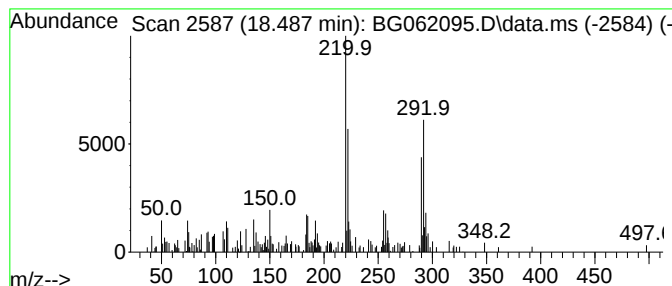
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 2,3,3',5'-te... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.487	2.31 ng/ul	88048	Phenanthrene-d10	17.436

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	97		
2	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	96		
3	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	93		
4	1,1'-Biphenyl, 2,3,4,4'-tetrachl...	290	C12H6Cl4	033025-41-1	93		
5	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	93		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062095.D
Acq On : 16 Jul 2024 15:34
Operator : MA/JU
Sample : P3177-03
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BS1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.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
unknown-01	3.722	2.5	ng/ul	38045	1	8.053	301177	20.0
1-Phenoxypropan...	11.601	5.2	ng/ul	120311	2	10.867	461739	20.0
1,1'-Biphenyl, ...	18.487	2.3	ng/ul	88048	4	17.436	763581	20.0