

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071024\
 Data File : BG062001.D
 Acq On : 10 Jul 2024 16:37
 Operator : MA/JU
 Sample : P3035-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4CX7

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: BG062001.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.484	31	33	38	rVB6	8975	9482	0.62%	0.055%
2	3.754	76	79	85	rVB	19563	31461	2.05%	0.182%
3	3.831	90	92	95	rVB3	3524	3144	0.21%	0.018%
4	3.913	103	106	115	rBV2	17076	28214	1.84%	0.163%
5	4.019	121	124	130	rVB5	9067	13558	0.88%	0.078%
6	4.248	161	163	167	rVB5	4012	5644	0.37%	0.033%
7	4.365	180	183	187	rBV4	1871	3116	0.20%	0.018%
8	4.618	221	226	227	rBV3	2257	3093	0.20%	0.018%
9	4.812	254	259	262	rBV6	5796	8049	0.53%	0.046%
10	5.017	287	294	297	rBV2	2711	3868	0.25%	0.022%
11	5.064	300	302	308	rVB4	3157	5113	0.33%	0.030%
12	5.164	313	319	326	rBV2	19104	28864	1.88%	0.167%
13	5.264	333	336	339	rVB4	4098	4058	0.26%	0.023%
14	5.358	349	352	354	rBV2	3308	3972	0.26%	0.023%
15	6.063	467	472	474	rBV2	2527	4211	0.27%	0.024%
16	6.146	481	486	490	rBV3	7298	14302	0.93%	0.083%
17	6.351	517	521	523	rVB2	2195	3041	0.20%	0.018%
18	6.381	523	526	531	rBV4	3852	5827	0.38%	0.034%
19	6.968	625	626	632	rVV3	2859	3724	0.24%	0.021%
20	7.050	638	640	643	rBV2	4047	4179	0.27%	0.024%
21	7.103	646	649	652	rBV3	3629	3192	0.21%	0.018%
22	7.250	667	674	685	rBV2	236414	427925	27.92%	2.470%
23	7.421	696	703	711	rBV	148080	256437	16.73%	1.480%
24	7.556	723	726	728	rBV	2597	3741	0.24%	0.022%
25	7.626	731	738	760	rVB	224379	409762	26.74%	2.365%
26	8.096	811	818	825	rBV	216007	372221	24.29%	2.148%
27	8.190	831	834	838	rVB2	3743	5429	0.35%	0.031%
28	8.619	903	907	910	rBV4	3529	3128	0.20%	0.018%
29	8.795	930	937	945	rVB	249972	447916	29.23%	2.585%
30	8.919	953	958	967	rVB4	6850	16922	1.10%	0.098%
31	9.113	987	991	995	rBV	2681	3974	0.26%	0.023%
32	9.260	1008	1016	1024	rVV2	231574	431631	28.17%	2.491%
33	9.318	1024	1026	1030	rVB2	2306	2946	0.19%	0.017%
34	9.424	1043	1044	1047	rVB	4208	3082	0.20%	0.018%
35	9.718	1092	1094	1098	rVB4	2298	2950	0.19%	0.017%
36	9.806	1102	1109	1114	rBV2	20694	34697	2.26%	0.200%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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37	9.906	1123	1126	1129	rVB2	2883	3597	0.23%	0.021%
38	9.982	1129	1139	1147	rBV	212372	398484	26.00%	2.300%
39	10.435	1214	1216	1220	rVB	3329	3519	0.23%	0.020%
40	10.523	1223	1231	1240	rVV	312467	556264	36.30%	3.211%
41	10.605	1243	1245	1250	rVB	2399	3266	0.21%	0.019%
42	10.670	1252	1256	1258	rBV	4838	5749	0.38%	0.033%
43	10.911	1289	1297	1303	rBV	346350	601834	39.27%	3.474%
44	11.040	1312	1319	1327	rBV	173986	330895	21.59%	1.910%
45	11.140	1334	1336	1339	rVB2	3516	3771	0.25%	0.022%
46	11.198	1344	1346	1352	rVV4	2816	5401	0.35%	0.031%
47	11.245	1352	1354	1359	rVB3	2305	3353	0.22%	0.019%
48	11.422	1380	1384	1389	rVB4	11663	18464	1.20%	0.107%
49	11.639	1413	1421	1428	rBV3	32108	67765	4.42%	0.391%
50	11.868	1458	1460	1463	rBV	2358	3197	0.21%	0.018%
51	11.898	1463	1465	1470	rVV	1884	3259	0.21%	0.019%
52	12.027	1484	1487	1490	rBV2	3828	4096	0.27%	0.024%
53	12.291	1529	1532	1535	rVB	2488	3689	0.24%	0.021%
54	12.485	1561	1565	1570	rBV4	5529	8384	0.55%	0.048%
55	12.550	1573	1576	1578	rBV2	3653	3906	0.25%	0.023%
56	12.661	1590	1595	1596	rBV3	2529	3049	0.20%	0.018%
57	12.961	1644	1646	1649	rVV2	2250	3148	0.21%	0.018%
58	13.190	1682	1685	1688	rVB	3430	3952	0.26%	0.023%
59	13.320	1703	1707	1708	rBV2	2753	3027	0.20%	0.017%
60	13.531	1738	1743	1747	rVB3	5057	9067	0.59%	0.052%
61	13.625	1756	1759	1761	rBV2	2584	3408	0.22%	0.020%
62	13.778	1784	1785	1789	rVB	2778	3219	0.21%	0.019%
63	13.819	1789	1792	1795	rBV2	3463	4309	0.28%	0.025%
64	13.854	1795	1798	1801	rVB	3866	4102	0.27%	0.024%
65	13.895	1801	1805	1809	rBV5	5300	9136	0.60%	0.053%
66	14.119	1836	1843	1849	rVV	572757	814903	53.18%	4.704%
67	14.348	1877	1882	1884	rVV4	5381	8027	0.52%	0.046%
68	14.412	1886	1893	1903	rVV	572442	923726	60.28%	5.332%
69	14.718	1938	1945	1952	rVB	557831	836311	54.57%	4.827%
70	14.783	1952	1956	1960	rBV4	7023	11272	0.74%	0.065%
71	14.912	1971	1978	1986	rBV3	345607	650231	42.43%	3.753%
72	15.065	2000	2004	2009	rVB4	18481	27136	1.77%	0.157%
73	15.117	2009	2013	2021	rVB10	16481	29282	1.91%	0.169%
74	15.176	2021	2023	2026	rBV4	4483	4947	0.32%	0.029%
75	15.264	2035	2038	2039	rBV3	3154	3983	0.26%	0.023%
76	15.488	2074	2076	2077	rBV2	7129	4075	0.27%	0.024%

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77	15.517	2079	2081	2084	rVB3	5840	6482	0.42%	0.037%
78	15.705	2107	2113	2119	rBV	744420	1146896	74.84%	6.620%
79	15.817	2127	2132	2145	rVB	249520	380948	24.86%	2.199%
80	15.940	2148	2153	2157	rVV7	11059	18903	1.23%	0.109%
81	16.904	2315	2317	2320	rVV3	5846	3882	0.25%	0.022%
82	16.933	2320	2322	2324	rVB3	5860	3117	0.20%	0.018%
83	17.462	2405	2412	2416	rBV	636881	1011805	66.03%	5.840%
84	17.503	2416	2419	2423	rVV	101787	134941	8.81%	0.779%
85	17.556	2423	2428	2437	rVB	784113	1178088	76.88%	6.800%
86	17.720	2454	2456	2457	rBV2	9328	8803	0.57%	0.051%
87	18.055	2508	2513	2518	rVB5	54119	96614	6.30%	0.558%
88	18.108	2519	2522	2524	rVB4	15839	16381	1.07%	0.095%
89	18.161	2528	2531	2532	rBV3	14267	15502	1.01%	0.089%
90	18.331	2557	2560	2565	rBV6	38781	54459	3.55%	0.314%
91	18.513	2587	2591	2597	rBV4	79075	128072	8.36%	0.739%
92	18.572	2598	2601	2605	rVV5	47585	72417	4.73%	0.418%
93	18.619	2605	2609	2614	rVB7	41165	56166	3.67%	0.324%
94	18.801	2636	2640	2642	rBV4	39276	51130	3.34%	0.295%
95	19.330	2727	2730	2732	rBV3	52352	53265	3.48%	0.307%
96	19.501	2755	2759	2764	rVB	199551	288694	18.84%	1.666%
97	19.835	2810	2816	2828	rBV2	866867	1532457	100.00%	8.845%
98	21.716	3130	3136	3142	rBV	619719	1128939	73.67%	6.516%
99	24.736	3645	3650	3659	rVB2	329079	794628	51.85%	4.587%
100	24.965	3682	3689	3701	rVB	411204	1140598	74.43%	6.583%

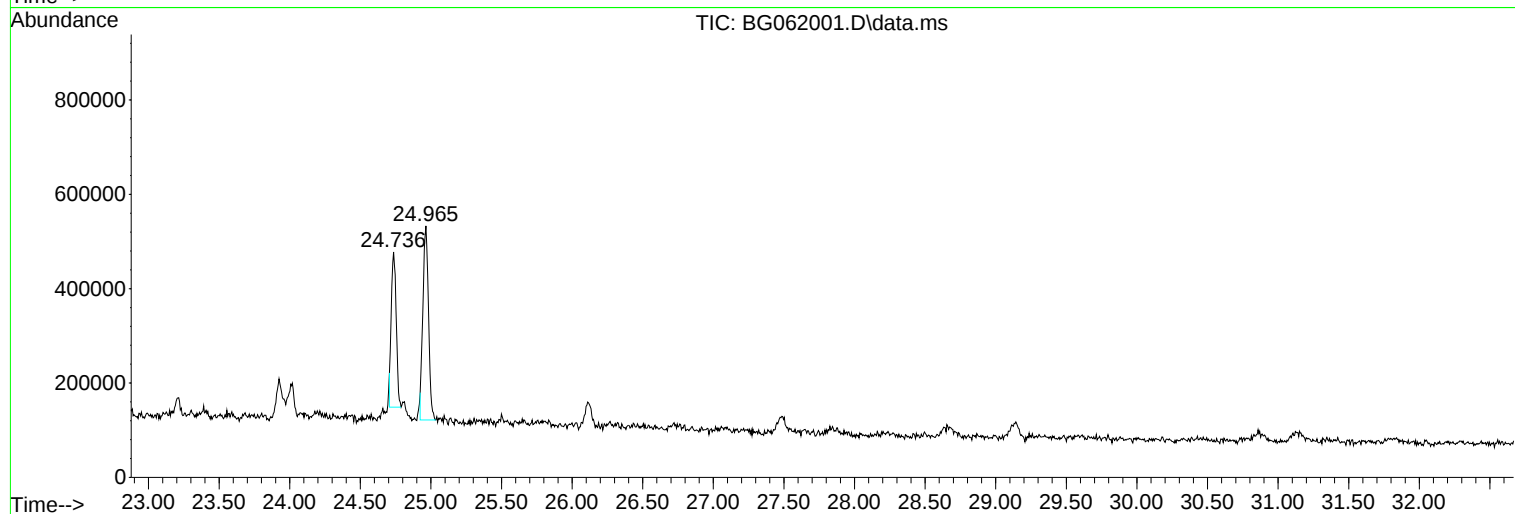
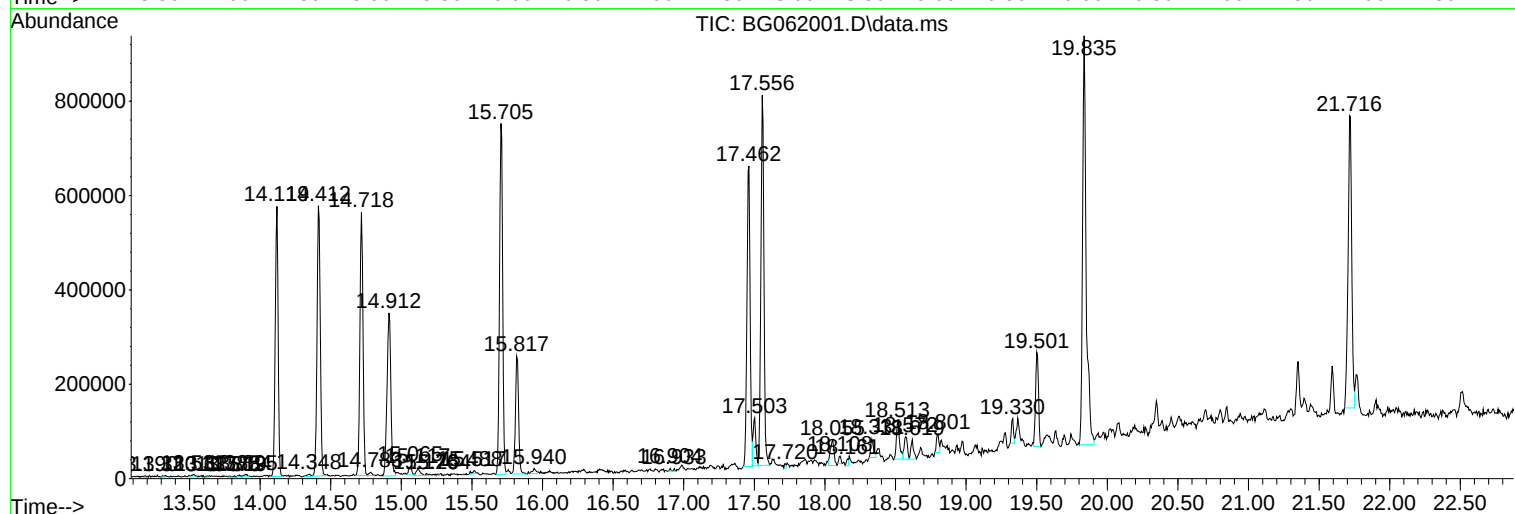
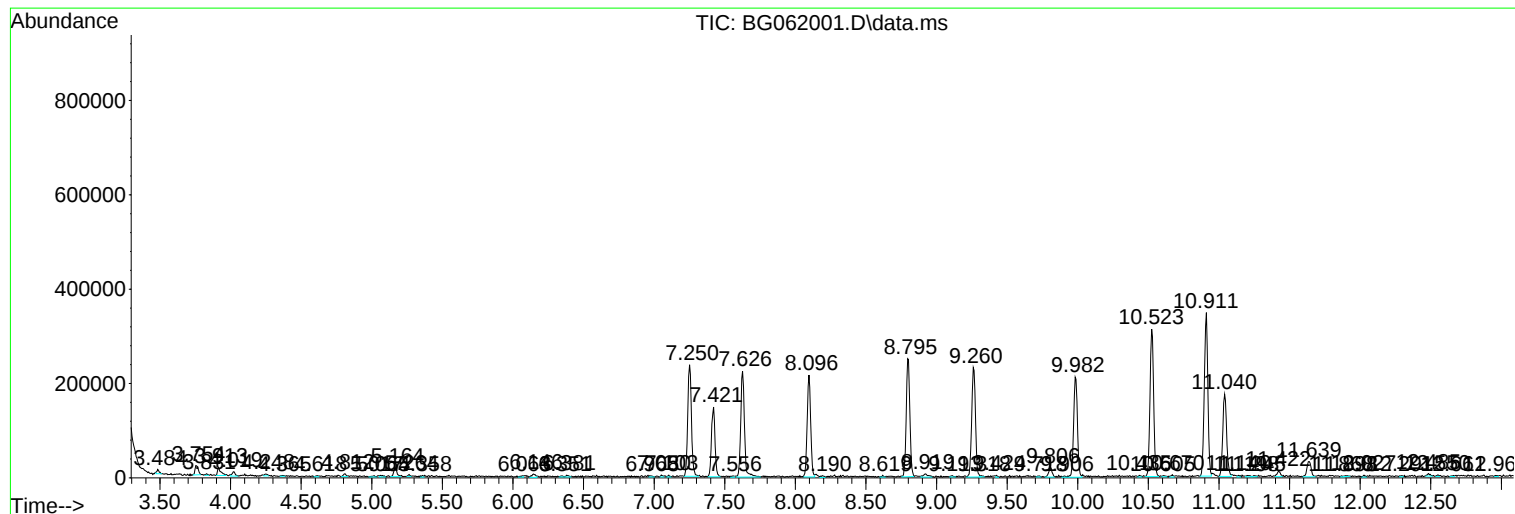
Sum of corrected areas: 17325263

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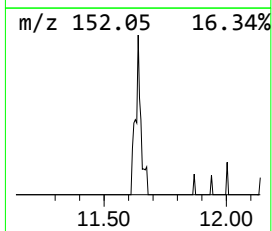
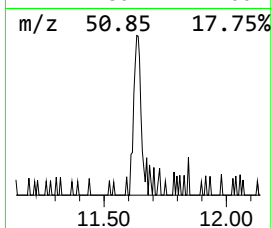
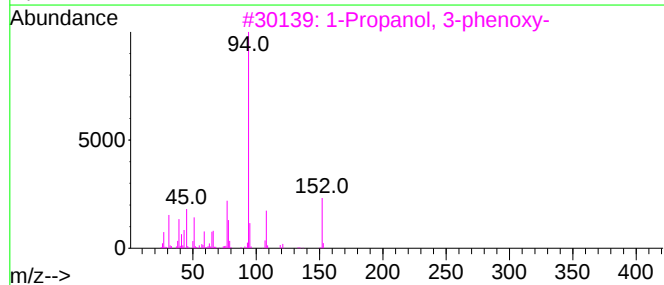
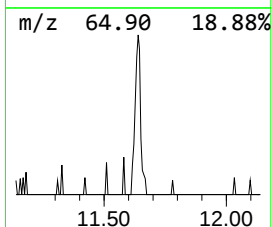
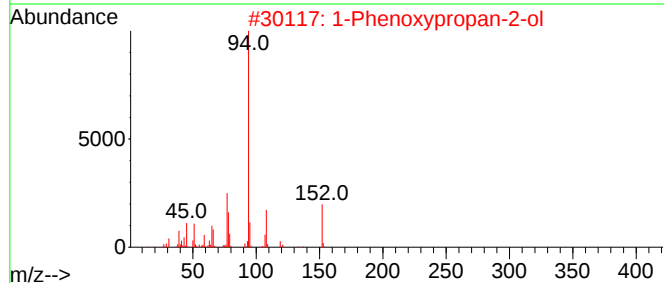
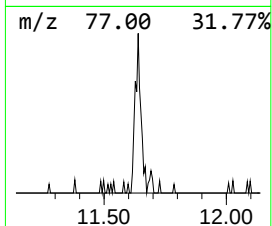
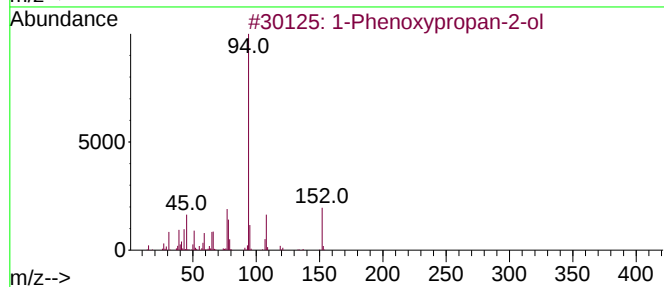
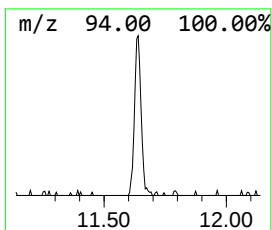
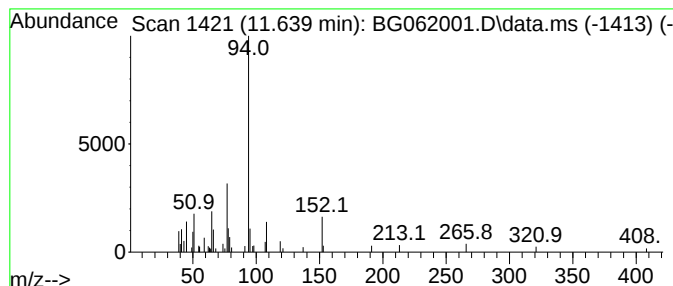
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.639	2.25 ng/ul	67765	Naphthalene-d8	10.911

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	81
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	80
5			3-(.alpha.-Hydroxyethyl)-aniline	137	C8H11NO	002454-37-7	64



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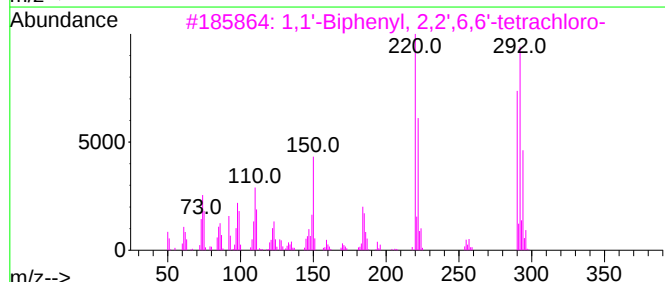
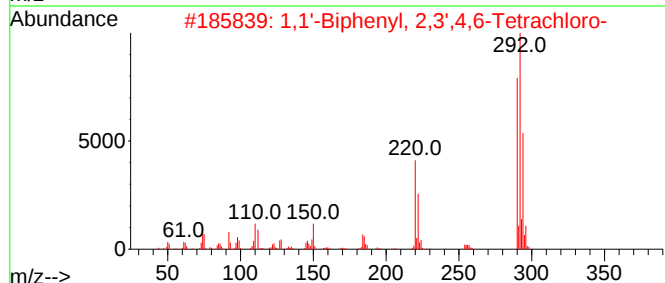
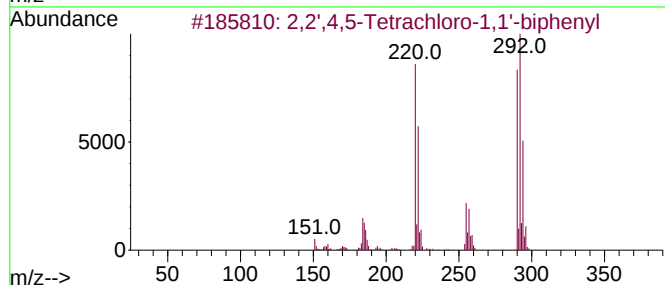
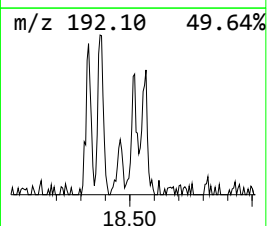
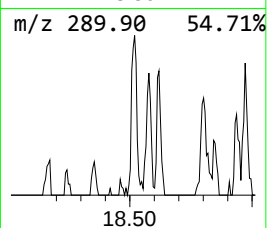
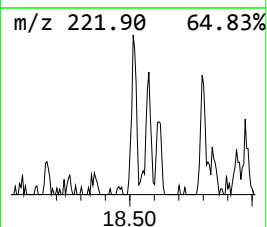
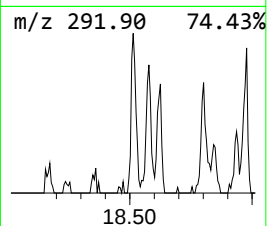
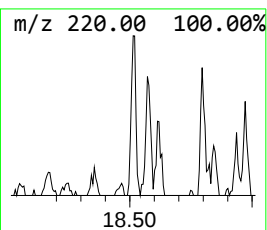
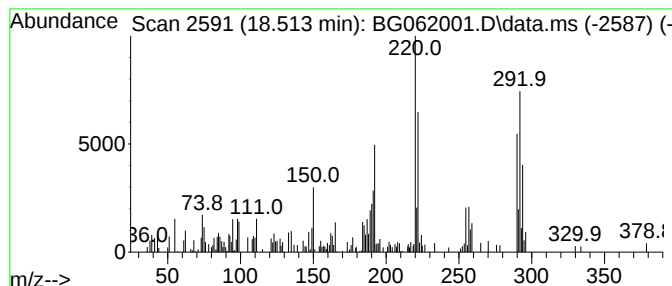
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.513	2.53 ng/ul	128072	Phenanthrene-d10	17.462

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	97		
2	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	96		
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	95		
4	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	94		
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	94		



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TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Phenoxypropan...	11.639	2.3	ng/u1	67765	2	10.911	601834	20.0
2,2',4,5-Tetrac...	18.513	2.5	ng/u1	128072	4	17.462	1011810	20.0