

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063296.D
 Acq On : 11 Nov 2024 19:39
 Operator : RC/JU
 Sample : P4644-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0CE9

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-BG110624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063296.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.301	33	36	43	rVB5	23905	38987	0.81%	0.097%
2	3.713	102	106	112	rBV	62108	99736	2.07%	0.249%
3	4.036	160	161	166	rVB4	8807	10436	0.22%	0.026%
4	4.130	175	177	181	rVB3	8327	10134	0.21%	0.025%
5	4.171	181	184	187	rBV3	6118	10208	0.21%	0.026%
6	4.570	244	252	262	rBV	400005	592857	12.29%	1.481%
7	5.140	343	349	353	rBV8	12919	28445	0.59%	0.071%
8	7.009	658	667	676	rBV2	127909	223366	4.63%	0.558%
9	7.167	686	694	701	rBV	314461	517684	10.73%	1.294%
10	7.367	722	728	737	rBV	357177	624510	12.95%	1.561%
11	7.837	800	808	816	rBV	348011	582816	12.09%	1.456%
12	8.542	922	928	936	rBV2	334181	576592	11.96%	1.441%
13	8.989	996	1004	1012	rBV	427682	779117	16.16%	1.947%
14	9.218	1038	1043	1046	rBV5	10050	15528	0.32%	0.039%
15	9.424	1075	1078	1082	rVB4	9117	11398	0.24%	0.028%
16	9.553	1095	1100	1102	rBV2	10099	12201	0.25%	0.030%
17	9.711	1119	1127	1137	rBV	416030	740204	15.35%	1.850%
18	10.252	1211	1219	1232	rBV	628294	1157782	24.01%	2.893%
19	10.622	1276	1282	1290	rBV	569868	988253	20.49%	2.469%
20	10.757	1296	1305	1316	rBV	278944	547534	11.35%	1.368%
21	11.451	1417	1423	1429	rVB7	12704	27321	0.57%	0.068%
22	12.214	1548	1553	1558	rBV4	8983	16903	0.35%	0.042%
23	12.567	1608	1613	1616	rBV	4759	7878	0.16%	0.020%
24	12.667	1627	1630	1638	rVB4	5066	11083	0.23%	0.028%
25	13.072	1693	1699	1700	rBV3	5612	7107	0.15%	0.018%
26	13.877	1829	1836	1843	rBV	1447009	2092537	43.39%	5.229%
27	14.165	1877	1885	1894	rVB2	1396253	2192484	45.46%	5.479%
28	14.477	1931	1938	1945	rBV	1104872	1742304	36.13%	4.354%
29	14.688	1968	1974	1983	rBV2	113666	235802	4.89%	0.589%
30	14.900	2007	2010	2014	rBV4	14690	20228	0.42%	0.051%
31	15.469	2100	2107	2119	rBV	2186854	3264673	67.70%	8.158%
32	15.593	2122	2128	2137	rVV	735429	1080056	22.40%	2.699%
33	15.704	2145	2147	2152	rVB4	12878	18663	0.39%	0.047%
34	15.834	2164	2169	2175	rBV	172028	251360	5.21%	0.628%
35	16.169	2224	2226	2230	rVB3	5377	6835	0.14%	0.017%
36	16.656	2305	2309	2312	rBV3	5787	8555	0.18%	0.021%

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37	16.762	2323	2327	2331	rBV3	14496	22095	0.46%	0.055%
38	16.991	2362	2366	2367	rBV3	4963	7345	0.15%	0.018%
39	17.097	2382	2384	2389	rVB4	16544	19451	0.40%	0.049%
40	17.226	2400	2406	2414	rBV	1590724	2280031	47.28%	5.697%
41	17.326	2416	2423	2431	rVV	2593513	3859345	80.03%	9.644%
42	17.508	2451	2454	2456	rBV3	7353	7977	0.17%	0.020%
43	17.579	2463	2466	2470	rBV6	9614	16934	0.35%	0.042%
44	17.890	2513	2519	2525	rBV7	10156	16721	0.35%	0.042%
45	18.125	2555	2559	2566	rBV4	51157	88429	1.83%	0.221%
46	18.460	2612	2616	2620	rBV5	8040	12684	0.26%	0.032%
47	18.613	2636	2642	2647	rBV8	14423	31422	0.65%	0.079%
48	18.924	2693	2695	2699	rVB4	8330	10063	0.21%	0.025%
49	18.965	2699	2702	2706	rBV5	11406	15429	0.32%	0.039%
50	19.183	2735	2739	2745	rBV2	46092	71915	1.49%	0.180%
51	19.259	2747	2752	2757	rVB2	37578	59841	1.24%	0.150%
52	19.623	2808	2814	2822	rBV	3515445	4822606	100.00%	12.051%
53	20.851	3021	3023	3025	rBV3	14529	14159	0.29%	0.035%
54	21.480	3121	3130	3138	rBV	1686477	2680413	55.58%	6.698%
55	24.312	3602	3612	3625	rVB	1790381	4633906	96.09%	11.579%
56	24.518	3637	3647	3660	rVB2	1055901	2778036	57.60%	6.942%
57	24.764	3687	3689	3690	rBV2	17887	9202	0.19%	0.023%
58	28.748	4365	4367	4368	rBV2	15910	9437	0.20%	0.024%

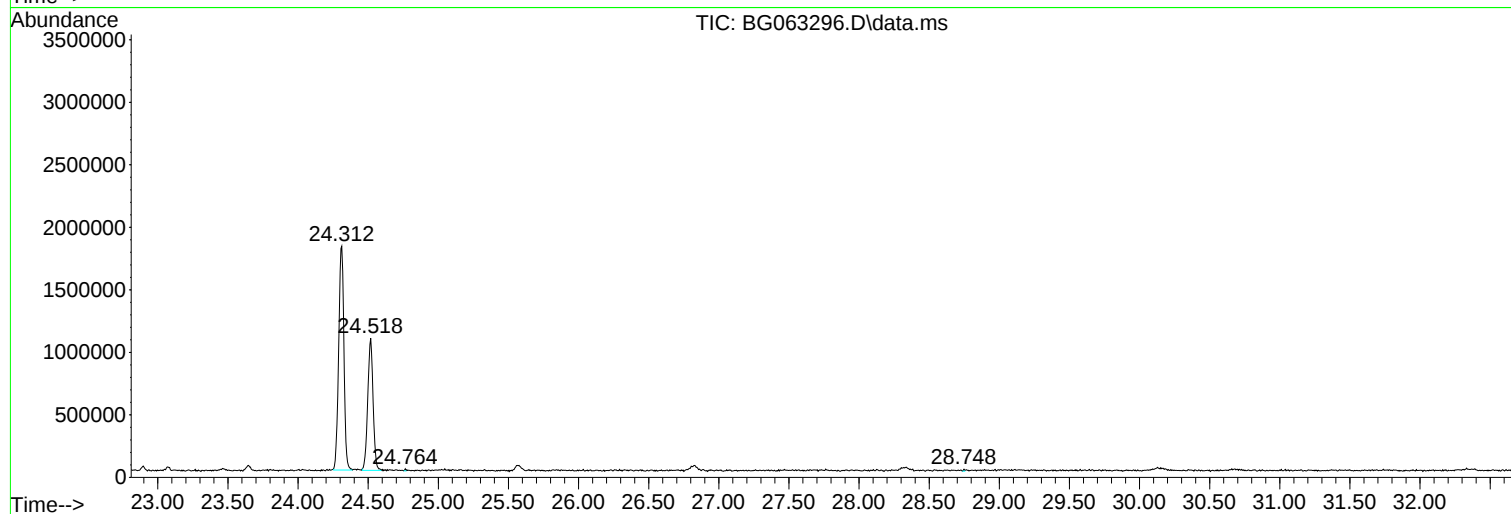
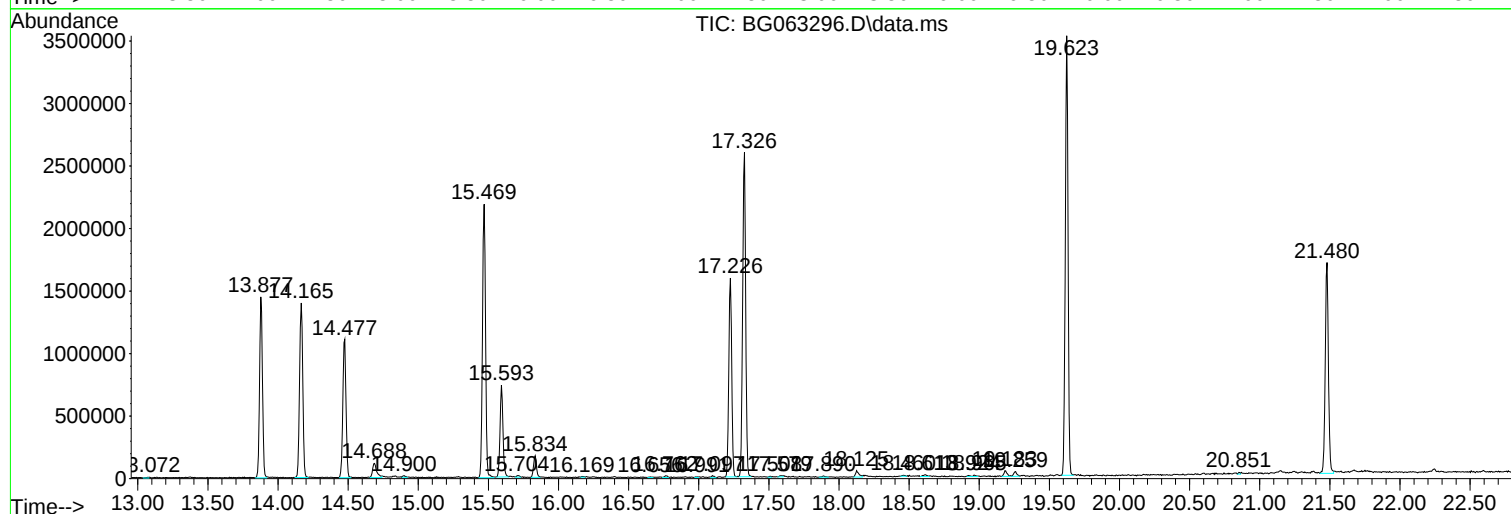
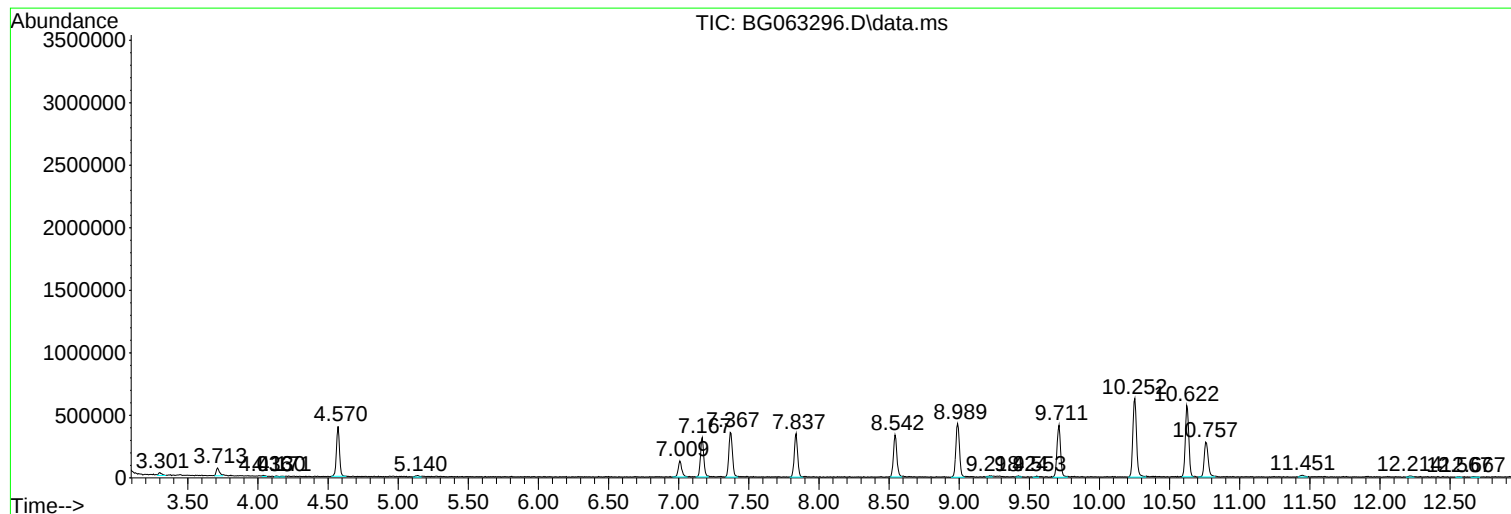
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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



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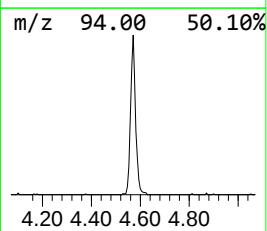
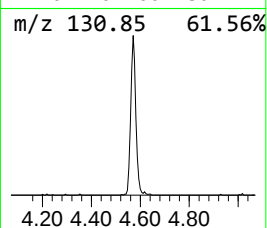
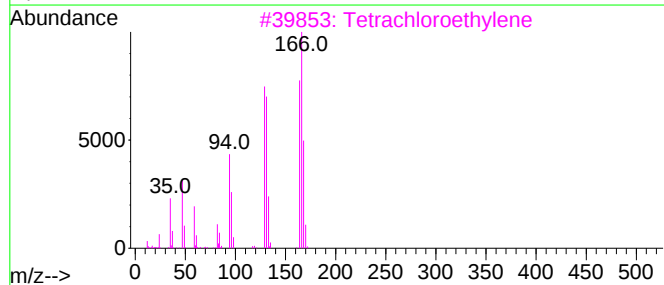
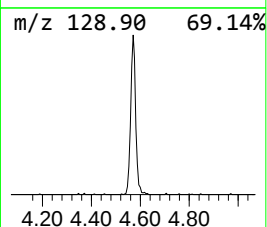
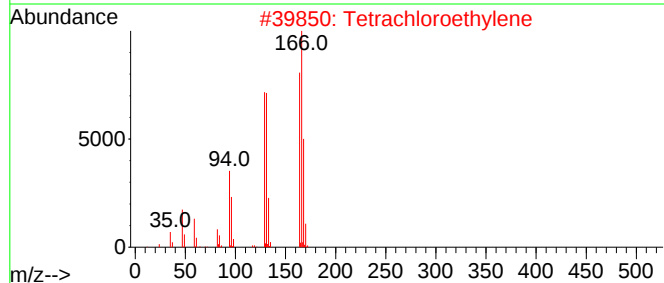
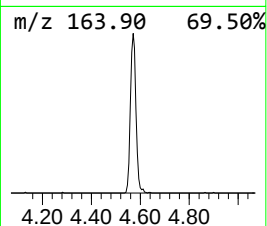
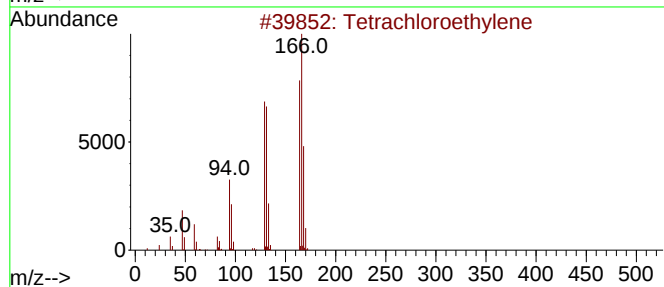
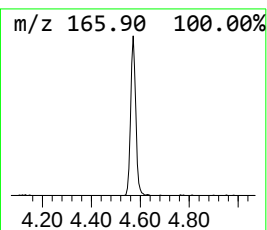
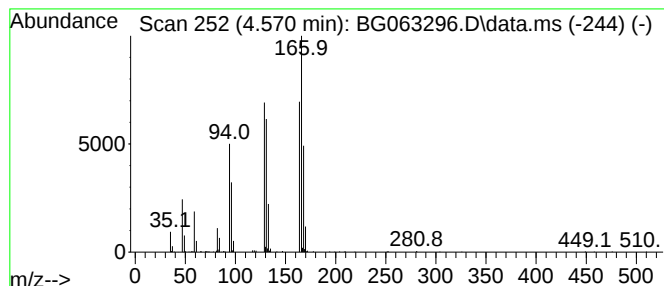
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.570	20.34 ng/ul	592857	1,4-Dichlorobenzene-d4	7.837

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	93
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	43



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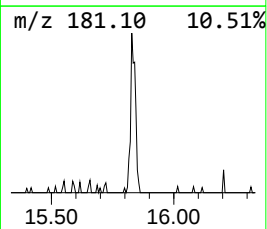
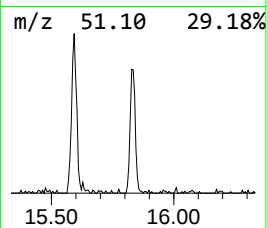
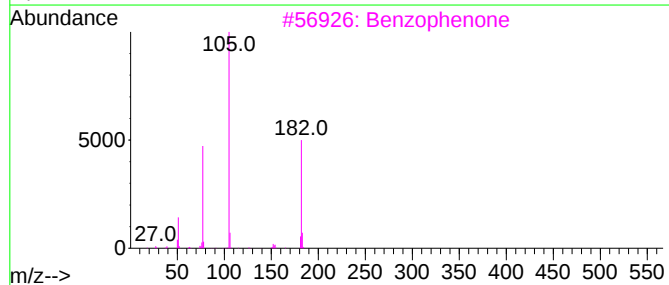
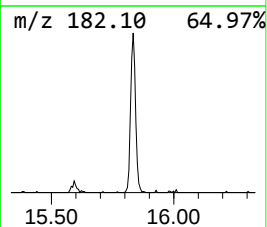
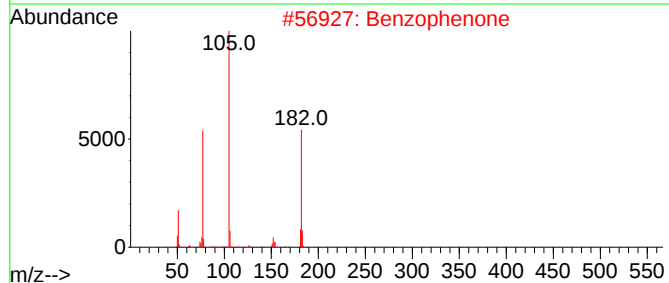
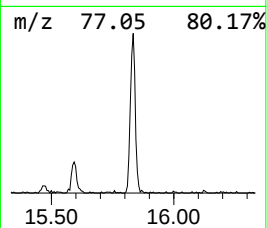
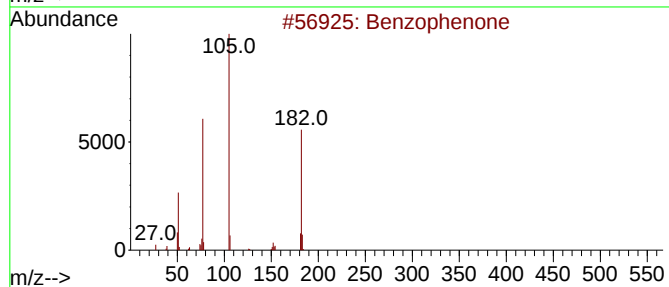
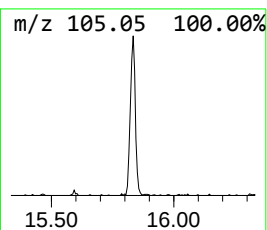
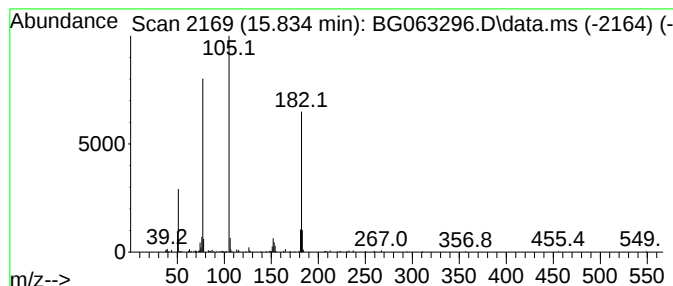
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.834	2.89 ng/ul	251360	Acenaphthene-d10	14.477

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	91
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	68



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.570	20.3	ng/u1	592857	1	7.837	582816	20.0
Benzophenone	15.834	2.9	ng/u1	251360	3	14.477	1742300	20.0