

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020122\
 Data File : BG052241.D
 Acq On : 1 Feb 2022 22:51
 Operator : CG/JU
 Sample : N1307-19
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q62

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-BG013122.M
 Title : SVOA CALIBRATION

Signal : TIC: BG052241.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.996	81	86	101	rBV	37405	75276	0.16%	0.121%
2	7.538	682	689	696	rBV	107778	181797	0.38%	0.292%
3	7.762	720	727	735	rBV	99093	174165	0.36%	0.279%
4	8.232	800	807	814	rBV	230442	398038	0.82%	0.639%
5	8.937	921	927	937	rBV2	77603	134021	0.28%	0.215%
6	9.406	1000	1007	1014	rBV	138313	255350	0.53%	0.410%
7	10.135	1123	1131	1139	rVB	116656	215094	0.44%	0.345%
8	10.687	1218	1225	1233	rBV	162290	287036	0.59%	0.461%
9	11.075	1281	1291	1298	rBV	318783	587906	1.21%	0.943%
10	11.198	1305	1312	1326	rVB	112308	221318	0.46%	0.355%
11	13.395	1680	1686	1696	rBV5	22042	52340	0.11%	0.084%
12	14.265	1828	1834	1840	rBV	387987	578865	1.19%	0.929%
13	14.570	1880	1886	1894	rVB	423921	673953	1.39%	1.081%
14	14.876	1931	1938	1945	rBV	508439	839482	1.73%	1.347%
15	15.064	1963	1970	1976	rBV4	25128	51431	0.11%	0.083%
16	15.416	2023	2030	2036	rBV	173231	278128	0.57%	0.446%
17	15.498	2036	2044	2059	rVB	534480	887875	1.83%	1.425%
18	15.868	2100	2107	2113	rVV	513372	798811	1.65%	1.282%
19	15.986	2120	2127	2144	rBV3	327662	738110	1.52%	1.184%
20	16.714	2243	2251	2254	rBV7	23444	63111	0.13%	0.101%
21	17.337	2340	2357	2361	rBV	17497902	48450580	100.00%	77.747%
22	17.637	2399	2408	2418	rVB	674358	1127791	2.33%	1.810%
23	17.736	2418	2425	2431	rBV	575511	882158	1.82%	1.416%
24	19.264	2680	2685	2689	rBV4	34326	57283	0.12%	0.092%
25	20.004	2805	2811	2817	rVB	719375	1009450	2.08%	1.620%
26	21.942	3133	3141	3148	rBV	640036	1072633	2.21%	1.721%
27	25.173	3680	3691	3704	rVB3	346437	1022218	2.11%	1.640%
28	25.420	3720	3733	3747	rVB2	378099	1153335	2.38%	1.851%
29	28.175	4192	4202	4211	rBV7	14871	51046	0.11%	0.082%

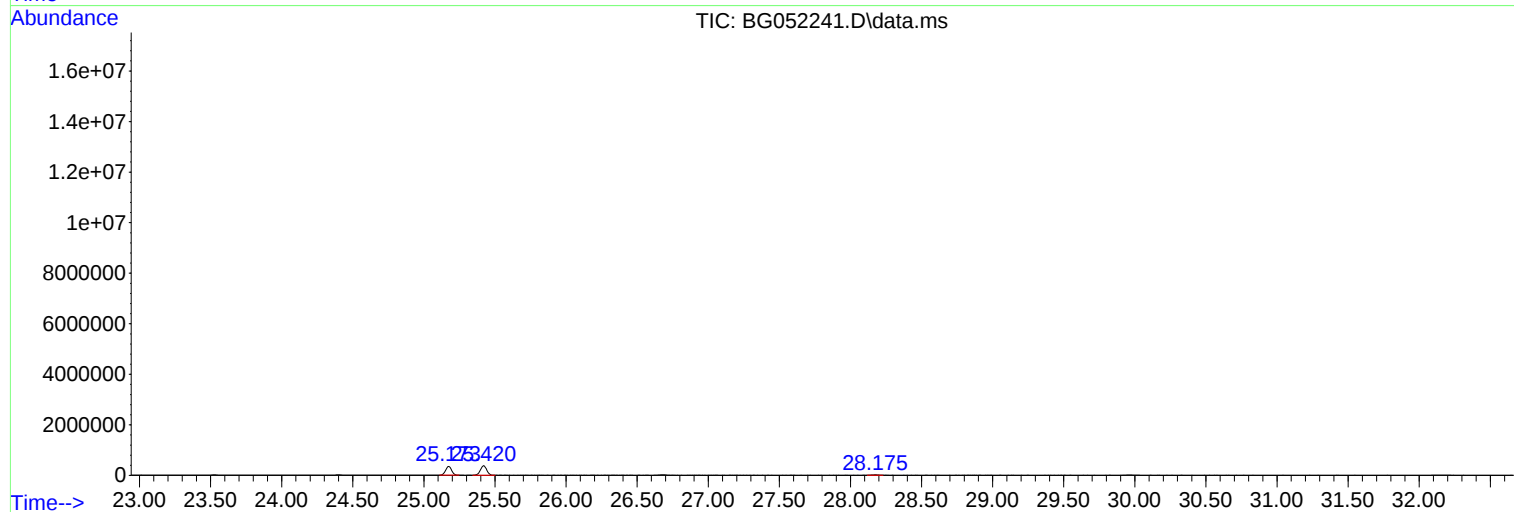
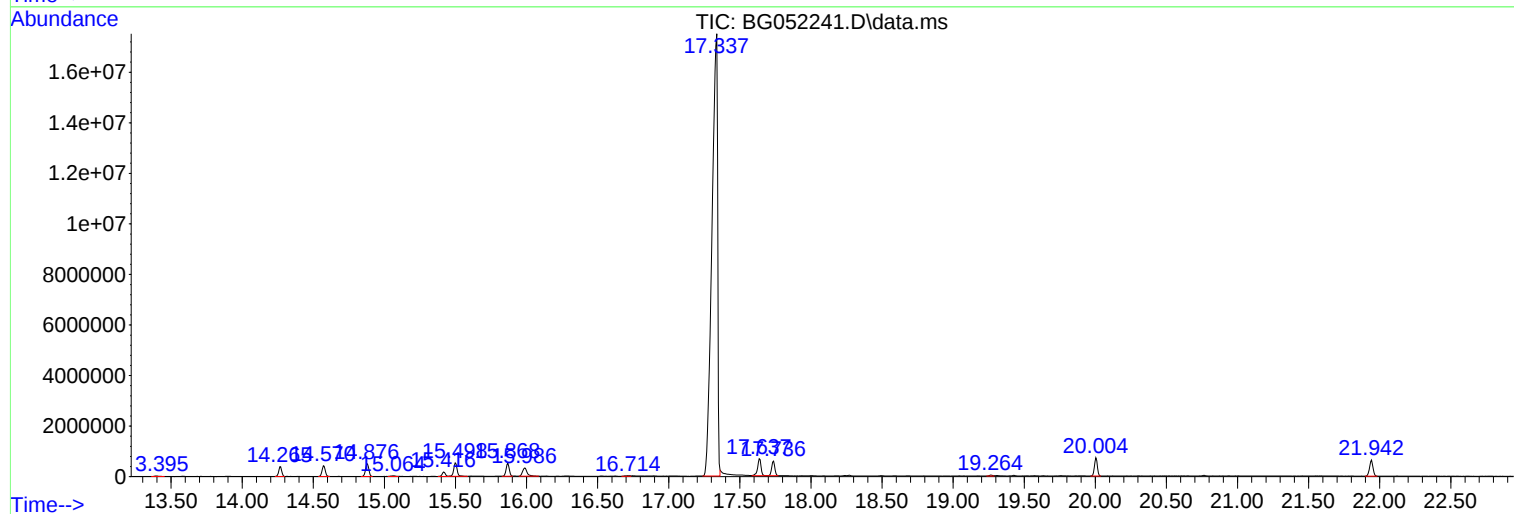
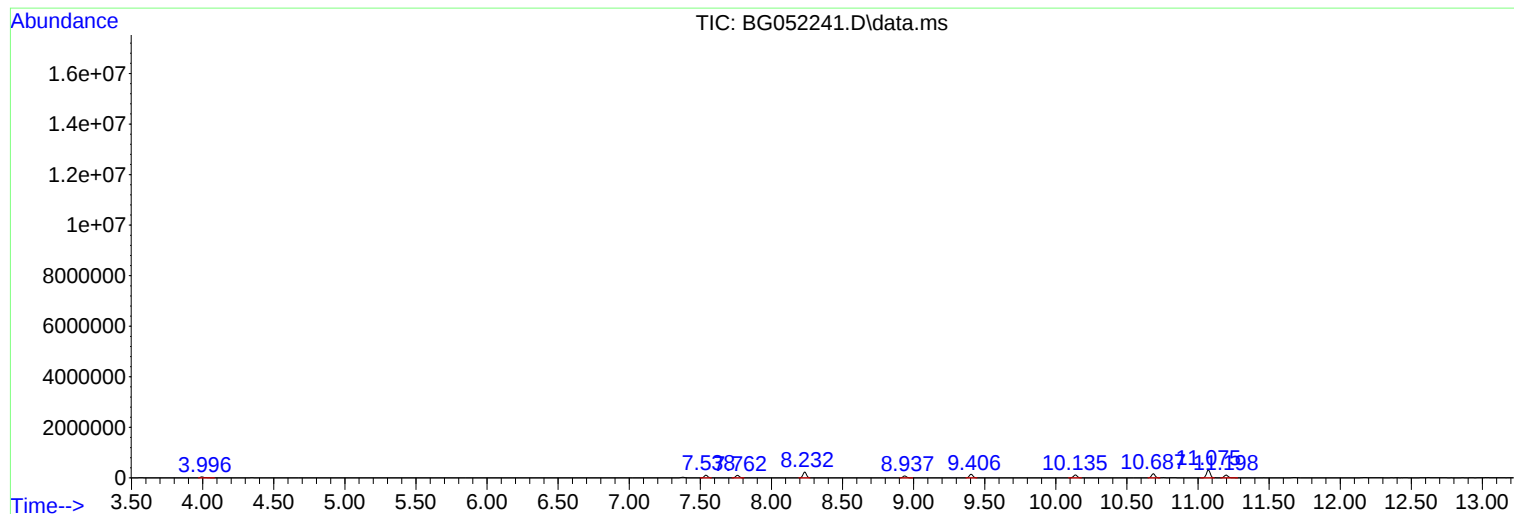
Sum of corrected areas: 62318601

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION

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



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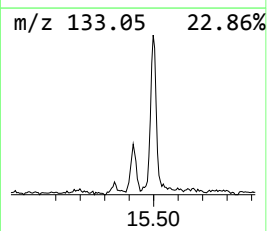
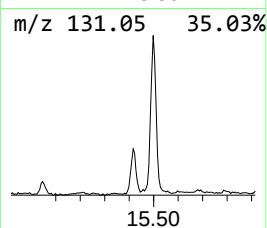
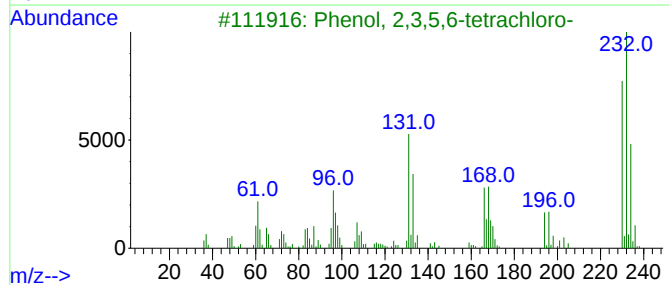
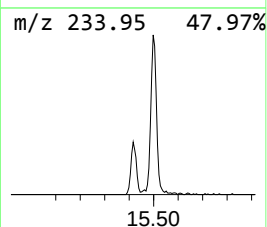
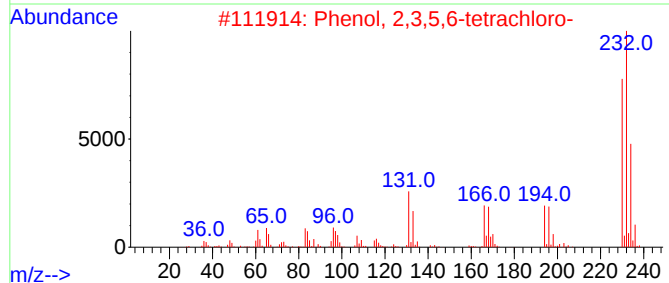
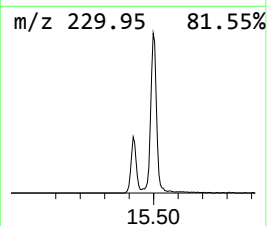
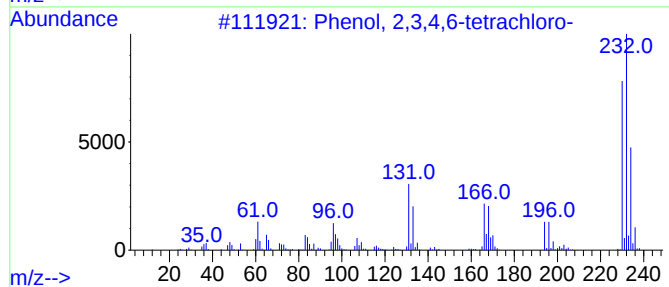
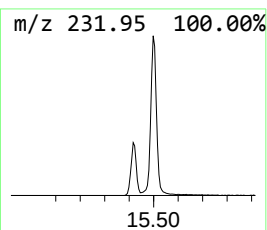
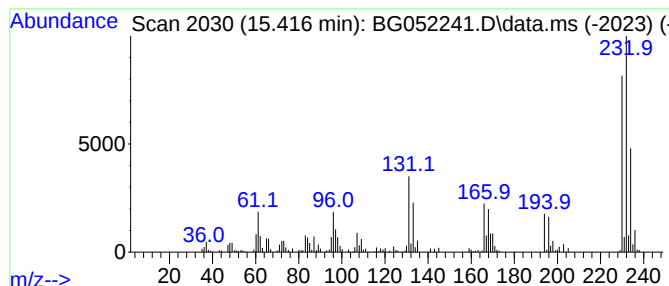
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.416	6.63 ng/ul	278128	Acenaphthene-d10	14.876

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
2			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
3			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
4			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98
5			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,3,5,6...	15.416	6.6	ng/ul	278128	3	14.876	839482	20.0