

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
 Data File : BP016240.D
 Acq On : 14 Jul 2023 23:04
 Operator : MA/JU
 Sample : 03457-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0C60

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_P\Methods\SFAM-EPA-BP071223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016240.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	14	17	25	rVB3	18339	31747	0.20%	0.043%
2	3.752	93	96	110	rBV	137173	340188	2.18%	0.461%
3	3.946	126	129	135	rVB	19125	32086	0.21%	0.043%
4	7.193	676	681	693	rBV2	66507	213431	1.37%	0.289%
5	7.293	693	698	718	rVB	401886	967772	6.19%	1.311%
6	7.505	728	734	757	rBV	298420	947329	6.06%	1.283%
7	7.963	804	812	830	rBV	429353	1161729	7.43%	1.574%
8	8.722	934	941	969	rBV	223931	834931	5.34%	1.131%
9	9.163	1009	1016	1042	rBV	489442	1344554	8.60%	1.821%
10	9.893	1132	1140	1167	rBV	309213	1065770	6.82%	1.444%
11	10.469	1228	1238	1282	rBV	295989	1544453	9.88%	2.092%
12	10.781	1284	1291	1318	rVB	882495	1942046	12.42%	2.630%
13	10.981	1318	1325	1366	rBV	307470	1432625	9.17%	1.940%
14	12.440	1567	1573	1581	rBV5	10203	28864	0.18%	0.039%
15	13.493	1747	1752	1757	rBV	24085	42438	0.27%	0.057%
16	14.028	1837	1843	1861	rBV	1623464	3097007	19.81%	4.195%
17	14.310	1884	1891	1910	rBV2	2017566	3312402	21.19%	4.487%
18	14.610	1936	1942	1963	rBV	1631255	2710909	17.34%	3.672%
19	15.081	2011	2022	2023	rBV8	18079	45956	0.29%	0.062%
20	15.610	2105	2112	2129	rBV	2706968	4614702	29.52%	6.251%
21	15.810	2140	2146	2172	rBV2	258979	980198	6.27%	1.328%
22	16.116	2191	2198	2227	rBV	8676424	15631317	100.00%	21.173%
23	17.375	2406	2412	2424	rBV	2218091	3411412	21.82%	4.621%
24	17.481	2424	2430	2457	rVB2	2550626	5522485	35.33%	7.480%
25	18.245	2555	2560	2567	rBV2	53021	104204	0.67%	0.141%
26	18.322	2570	2573	2578	rVB2	14302	20659	0.13%	0.028%
27	19.292	2734	2738	2743	rBV	46313	67348	0.43%	0.091%
28	19.369	2747	2751	2757	rBV	58908	93693	0.60%	0.127%
29	19.469	2764	2768	2775	rBV3	54013	97725	0.63%	0.132%
30	19.598	2787	2790	2796	rVB6	20238	30871	0.20%	0.042%
31	19.710	2804	2809	2831	rBV	3867117	6943704	44.42%	9.405%
32	21.475	3104	3109	3122	rBV	2111557	4121962	26.37%	5.583%
33	23.792	3496	3503	3523	rVB2	3034097	6967289	44.57%	9.437%
34	23.963	3526	3532	3556	rVB	1415893	4124388	26.39%	5.586%

Sum of corrected areas: 73828194

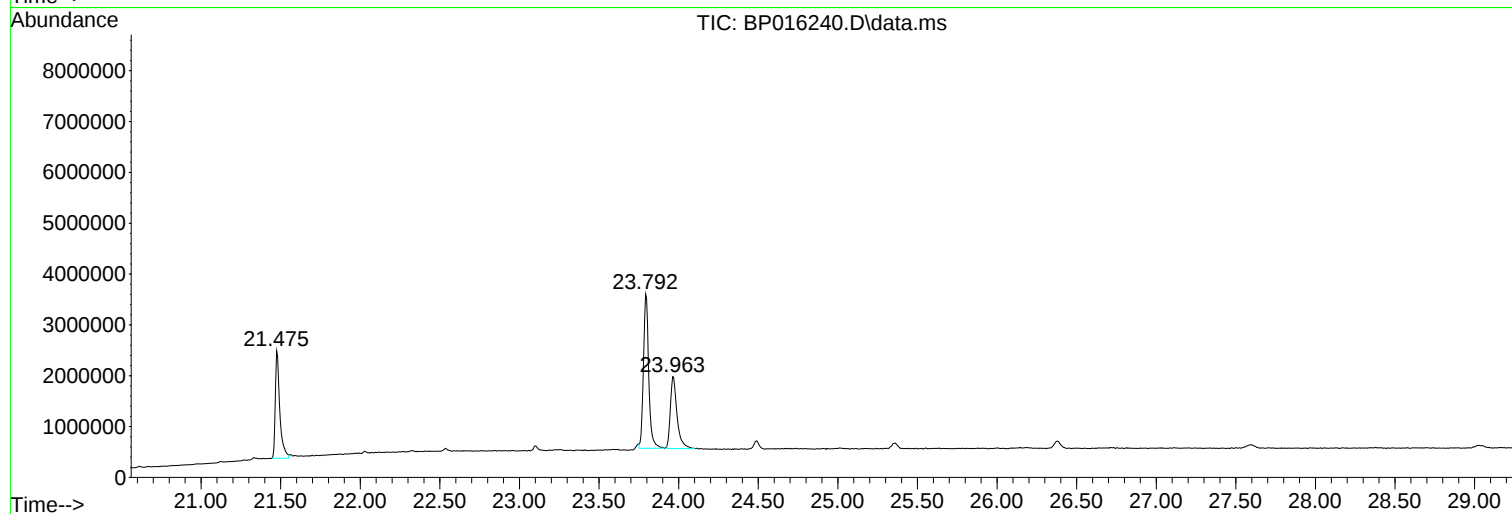
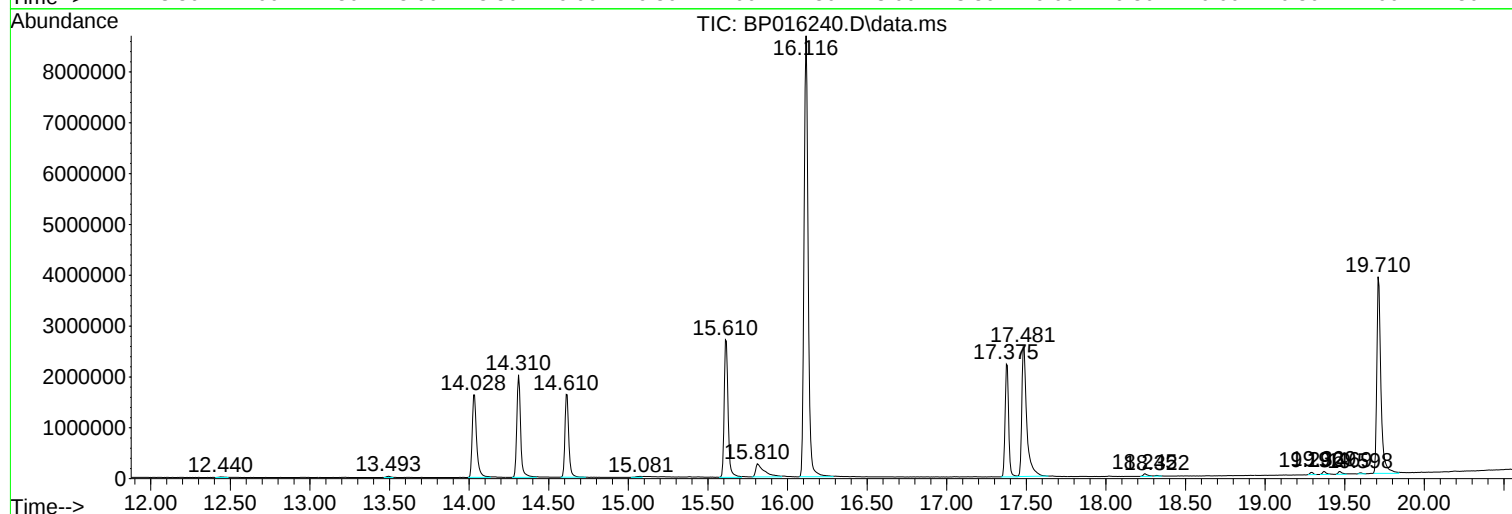
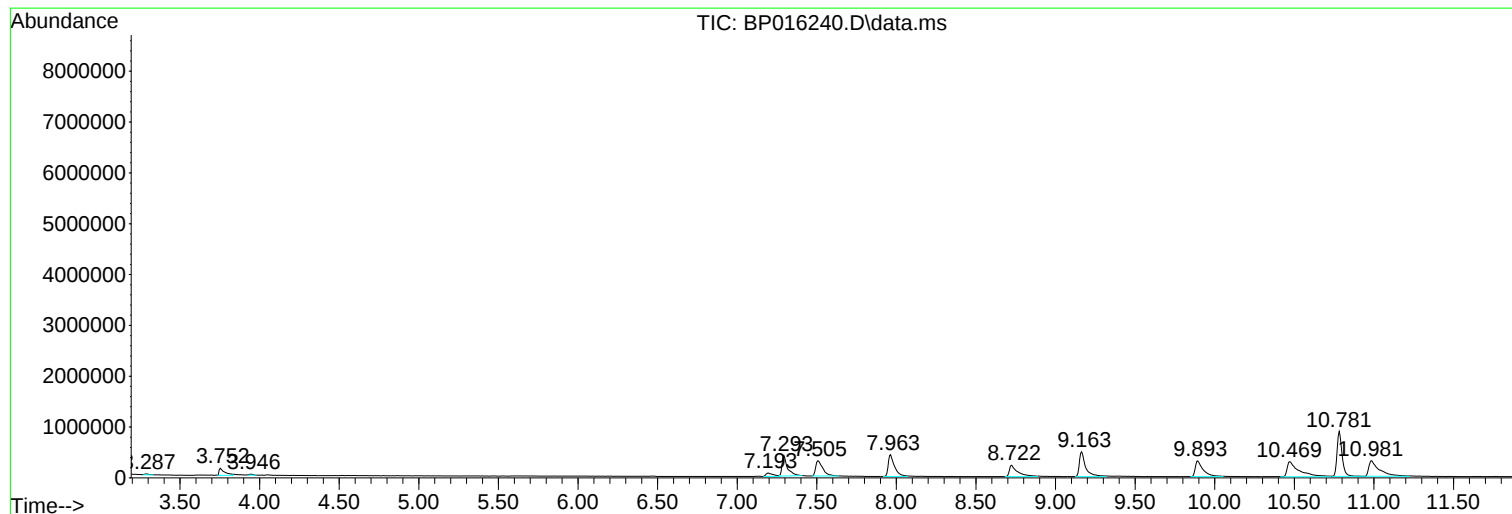
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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



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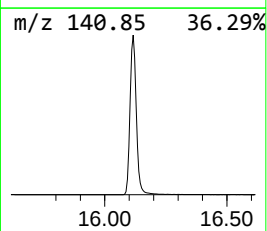
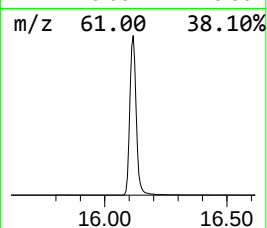
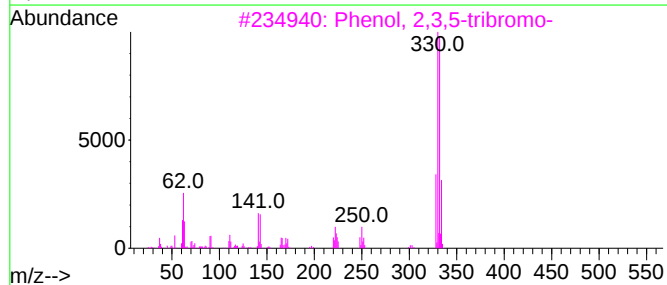
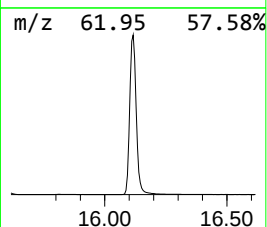
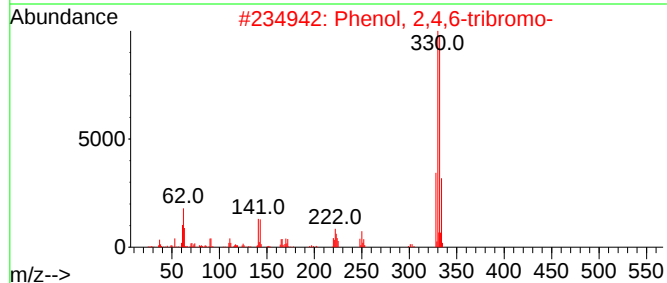
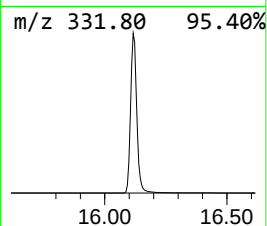
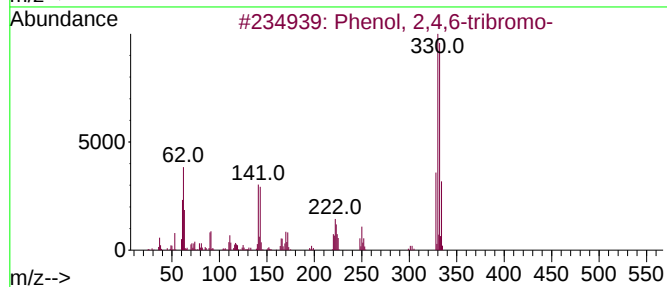
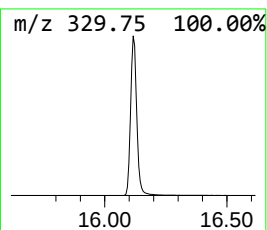
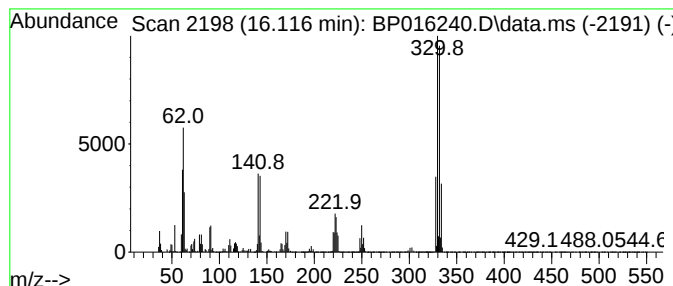
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 2,3,5-tribromo- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.116	91.64 ng/ul	15631300	Phenanthrene-d10	17.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,4,6-tribromo-	328	C6H3Br3O	000118-79-6	99
2			Phenol, 2,4,6-tribromo-	328	C6H3Br3O	000118-79-6	99
3			Phenol, 2,3,5-tribromo-	328	C6H3Br3O	057383-81-0	99
4			2,4,5-Tribromophenol	328	C6H3Br3O	014401-61-7	99
5			2,3,4-Tribromophenol	328	C6H3Br3O	138507-65-0	98



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,3,5-t...	16.116	91.6	ng/ul	15631300	4	17.375	3411410	20.0