

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021425\
 Data File : BP024062.D
 Acq On : 14 Feb 2025 15:24
 Operator : RC/JU
 Sample : Q1275-06DL 100X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCZT3DL

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-BP012925.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP024062.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.911	134	140	147	rBV	121999	207766	0.42%	0.225%
2	5.364	381	387	394	rBV2	31885	58326	0.12%	0.063%
3	6.946	650	656	668	rVB2	38497	79189	0.16%	0.086%
4	7.093	675	681	692	rBV	130256	220390	0.44%	0.239%
5	7.305	710	717	728	rBV	140356	223734	0.45%	0.242%
6	7.758	787	794	801	rBV	1601404	2550651	5.14%	2.761%
7	7.822	801	805	817	rVB	102222	174473	0.35%	0.189%
8	8.463	906	914	922	rBV	106645	189137	0.38%	0.205%
9	8.905	977	989	996	rBV	145311	250141	0.50%	0.271%
10	9.040	1002	1012	1024	rBV	106845	194810	0.39%	0.211%
11	9.616	1104	1110	1119	rBV	93054	178004	0.36%	0.193%
12	10.163	1195	1203	1219	rBV2	156704	342454	0.69%	0.371%
13	10.534	1257	1266	1276	rBV	2334425	3937344	7.94%	4.262%
14	10.669	1282	1289	1304	rVB	163905	316599	0.64%	0.343%
15	12.634	1614	1623	1629	rBV4	31149	68959	0.14%	0.075%
16	13.310	1728	1738	1750	rBV	142685	310043	0.63%	0.336%
17	13.781	1812	1818	1832	rVB	374914	532279	1.07%	0.576%
18	14.063	1860	1866	1875	rBV	382339	599147	1.21%	0.649%
19	14.375	1911	1919	1935	rBV	3452900	5215855	10.52%	5.646%
20	14.687	1965	1972	1978	rBV	52127	101504	0.20%	0.110%
21	14.922	2007	2012	2016	rBV2	52398	74300	0.15%	0.080%
22	15.004	2016	2026	2045	rVV	3856455	5468235	11.03%	5.919%
23	15.375	2083	2089	2103	rBV	515792	848197	1.71%	0.918%
24	15.492	2103	2109	2122	rVB3	87410	156358	0.32%	0.169%
25	16.216	2227	2232	2237	rBV2	62412	81817	0.17%	0.089%
26	16.816	2325	2334	2353	rBV	32708579	49576785	100.00%	53.663%
27	17.157	2386	2392	2403	rBV	4531984	6061349	12.23%	6.561%
28	17.263	2403	2410	2425	rVB2	554673	949849	1.92%	1.028%
29	17.528	2450	2455	2460	rBV	85432	128988	0.26%	0.140%
30	19.627	2806	2812	2819	rBV	745814	1053529	2.13%	1.140%
31	21.592	3140	3146	3155	rBV	3708021	5924107	11.95%	6.412%
32	24.710	3671	3676	3689	rVB2	258786	679895	1.37%	0.736%
33	24.945	3707	3716	3735	rBV	2222348	5631355	11.36%	6.095%

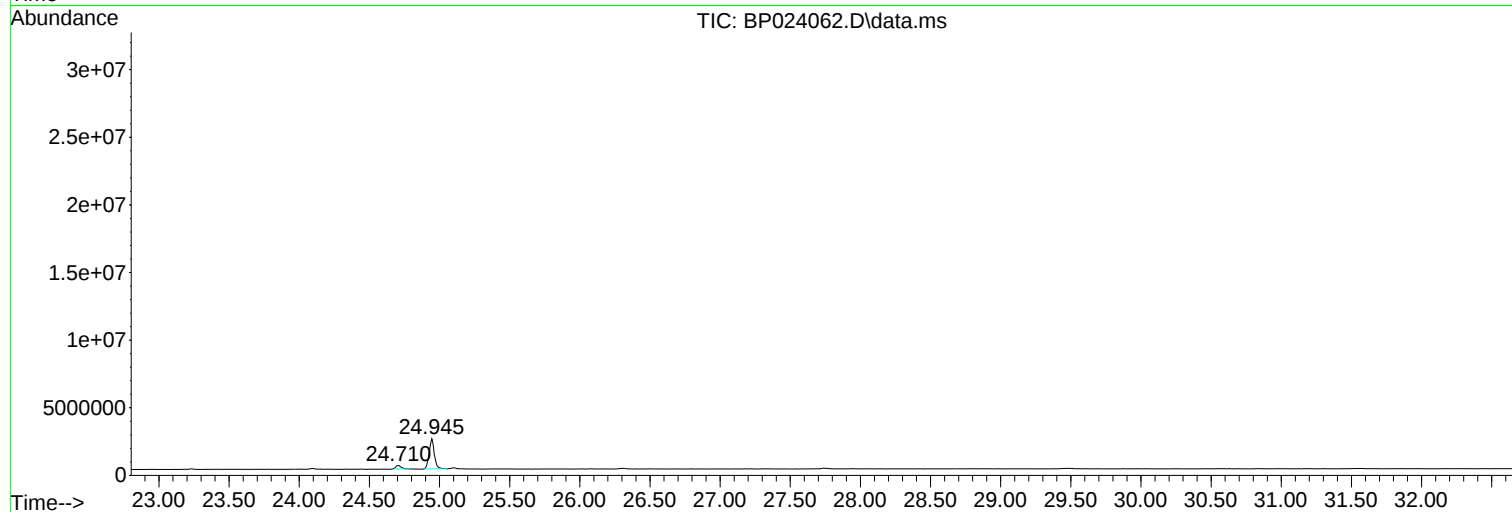
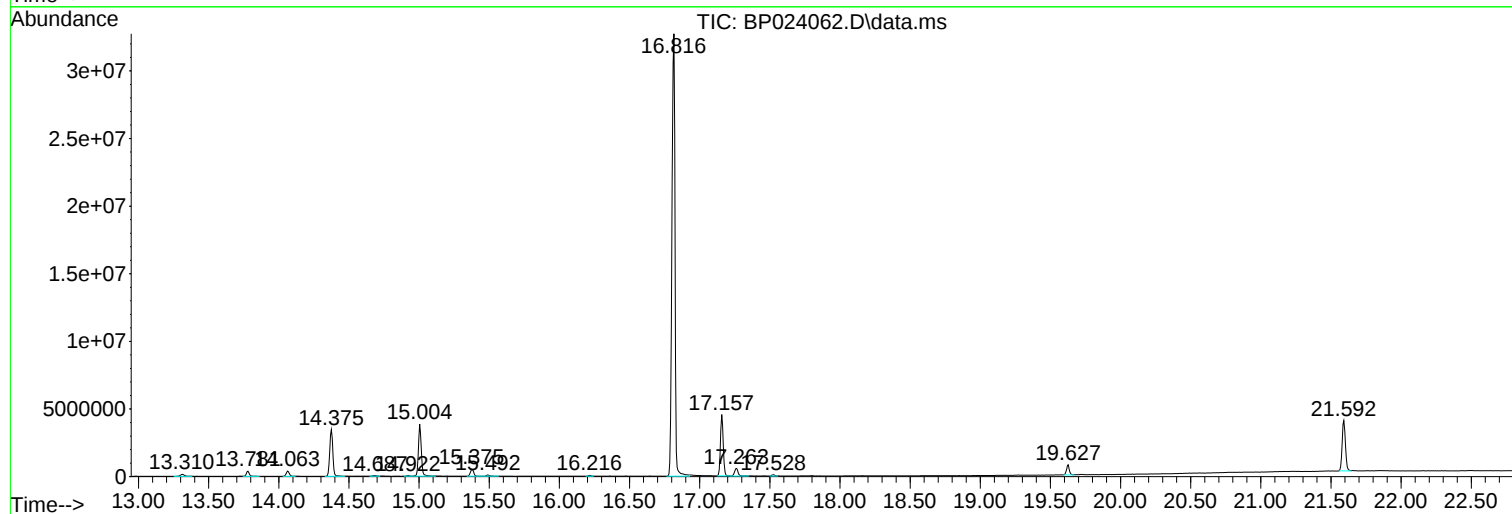
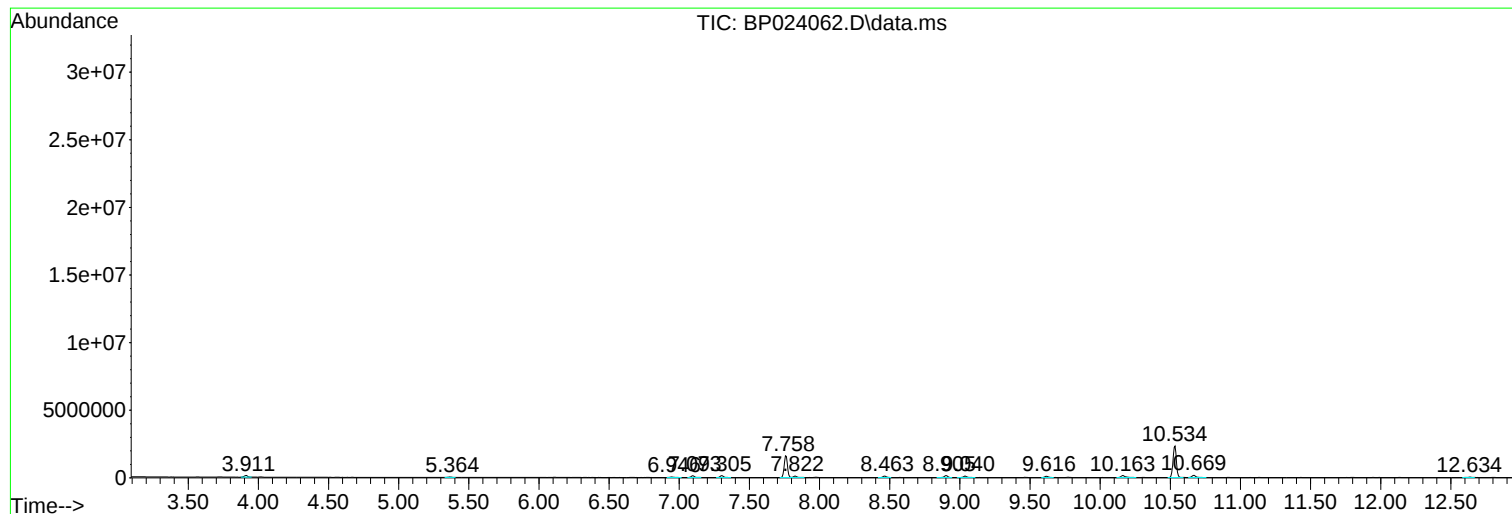
Sum of corrected areas: 92385569

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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



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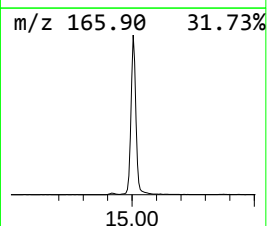
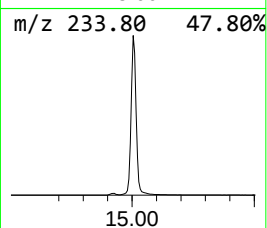
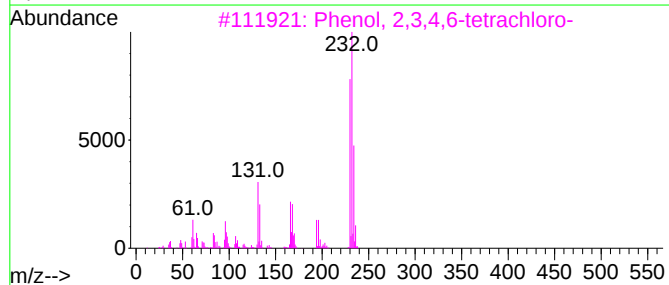
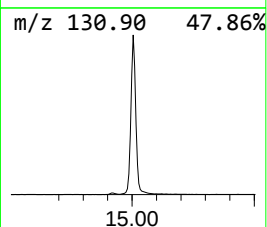
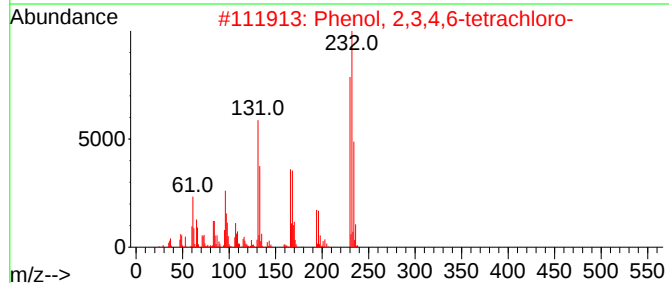
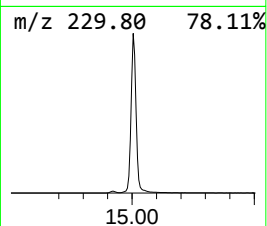
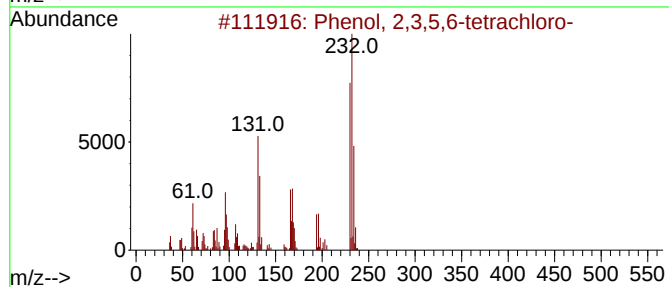
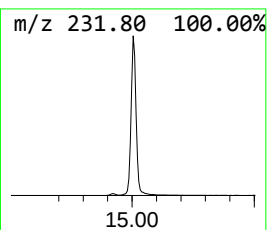
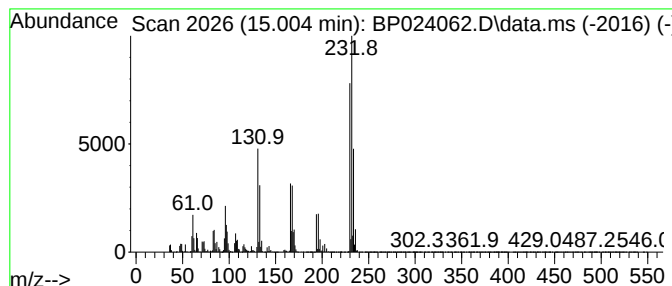
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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.
15.004	20.97 ng/ul	5468240	Acenaphthene-d10	14.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
2			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
3			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
4			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
5			Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	98



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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,3,5,6...	15.004	21.0	ng/ul	5468240	3	14.375	5215860	20.0