

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017152.D
 Acq On : 13 Sep 2023 16:51
 Operator : MA/JU
 Sample : 04324-16
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBKG8

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-BP091223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017152.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	26	30	35	rBV	35111	53165	0.19%	0.064%
2	3.317	35	39	44	rVB	49709	62963	0.22%	0.075%
3	3.752	110	113	126	rBV	269917	418867	1.46%	0.501%
4	3.946	140	146	151	rBV2	46897	71977	0.25%	0.086%
5	4.587	248	255	273	rBV	17110861	28731037	100.00%	34.381%
6	7.105	675	683	693	rBV	247893	395466	1.38%	0.473%
7	7.293	709	715	725	rVB	849936	1276508	4.44%	1.528%
8	7.469	738	745	756	rBV	807778	1298521	4.52%	1.554%
9	7.952	817	827	839	rBV	685498	1066330	3.71%	1.276%
10	8.663	939	948	958	rBV	661966	1094457	3.81%	1.310%
11	9.152	1020	1031	1042	rBV	956168	1535828	5.35%	1.838%
12	9.863	1145	1152	1162	rBV	763635	1211934	4.22%	1.450%
13	10.387	1232	1241	1258	rBV	1147484	1897196	6.60%	2.270%
14	10.775	1300	1307	1315	rBV	938241	1555752	5.41%	1.862%
15	10.940	1328	1335	1344	rBV	584526	968656	3.37%	1.159%
16	11.775	1468	1477	1491	rBV	140248	296729	1.03%	0.355%
17	12.357	1570	1576	1583	rBV2	19741	33970	0.12%	0.041%
18	14.028	1852	1860	1869	rBV	2463980	3306160	11.51%	3.956%
19	14.310	1901	1908	1924	rBV2	2517683	3648617	12.70%	4.366%
20	14.610	1953	1959	1967	rBV	1495627	2116167	7.37%	2.532%
21	14.834	1991	1997	2011	rBV	152013	315535	1.10%	0.378%
22	15.610	2121	2129	2143	rBV	3546898	5034029	17.52%	6.024%
23	15.734	2144	2150	2165	rVV	1154628	1508669	5.25%	1.805%
24	17.381	2423	2430	2440	rBV	1907141	2596328	9.04%	3.107%
25	17.481	2441	2447	2463	rVV2	4132057	5751381	20.02%	6.882%
26	18.222	2569	2573	2580	rBV4	49752	72060	0.25%	0.086%
27	19.292	2751	2755	2759	rBV3	46217	60789	0.21%	0.073%
28	19.363	2762	2767	2772	rVB	45324	76081	0.26%	0.091%
29	19.457	2780	2783	2793	rVB6	38523	56087	0.20%	0.067%
30	19.722	2822	2828	2838	rBV	5265781	6792035	23.64%	8.128%
31	21.480	3122	3127	3135	rBV	2031460	2550073	8.88%	3.052%
32	23.839	3521	3528	3539	rVB2	2673746	5481305	19.08%	6.559%
33	24.004	3550	3556	3566	rVB	1132152	2232689	7.77%	2.672%

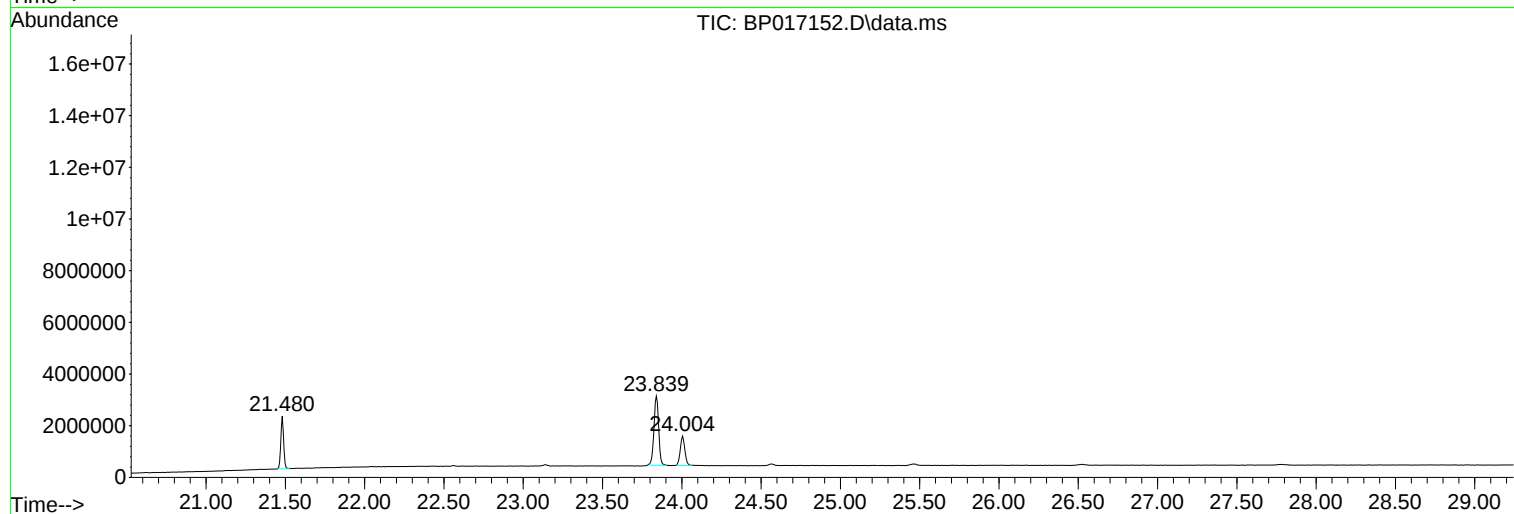
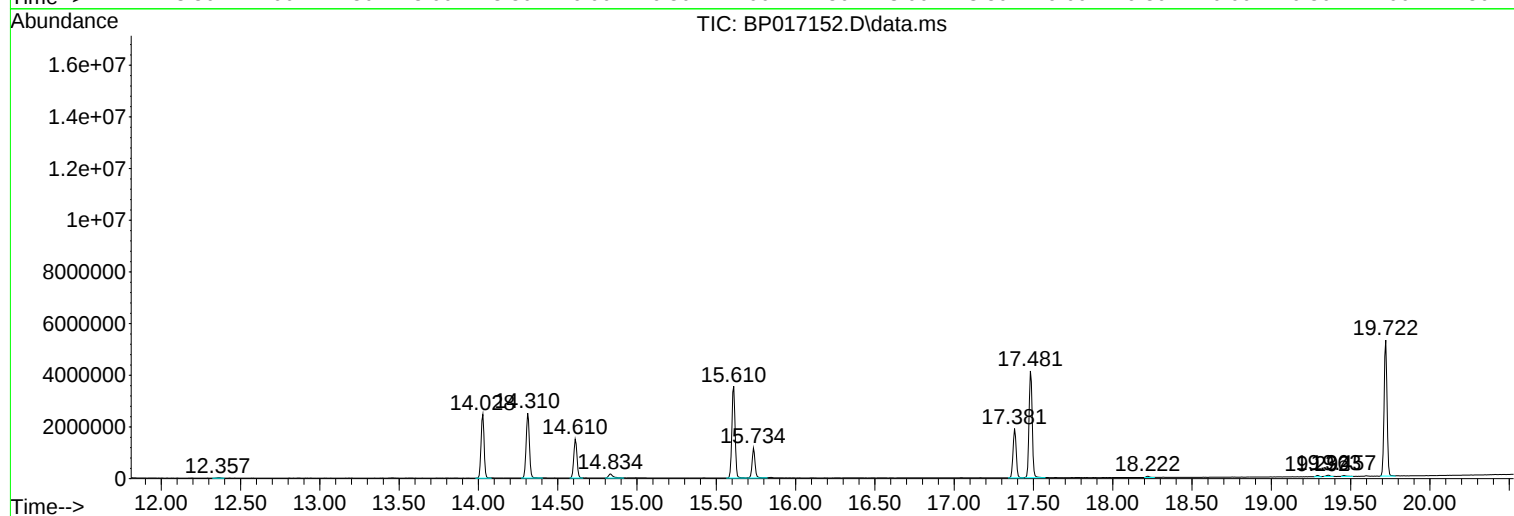
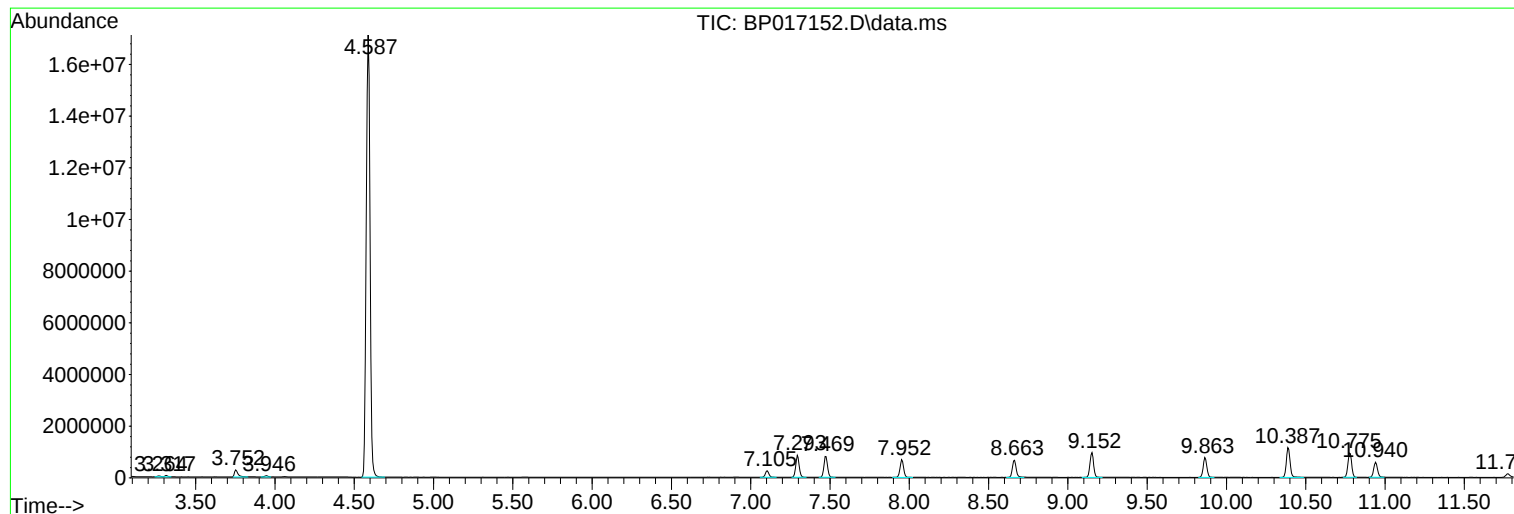
Sum of corrected areas: 83567361

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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



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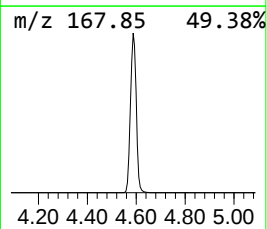
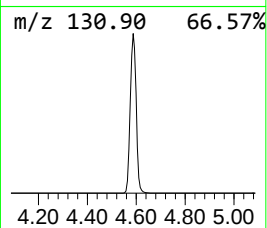
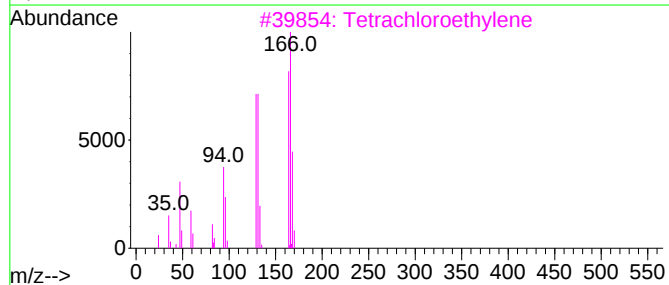
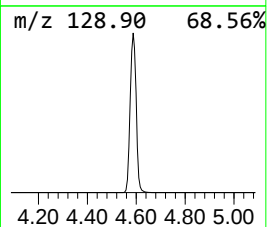
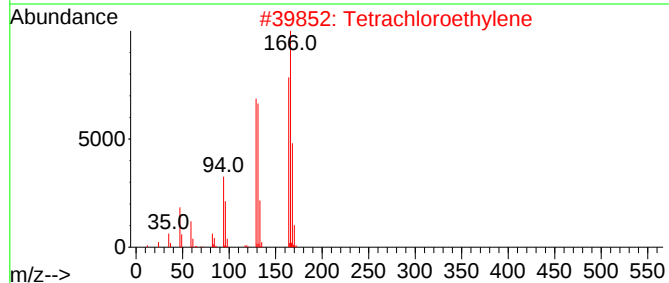
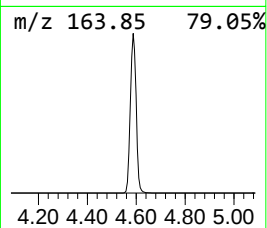
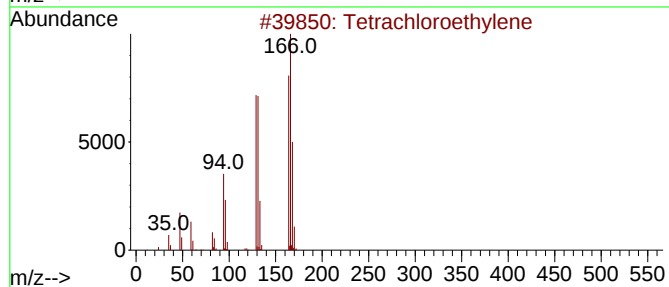
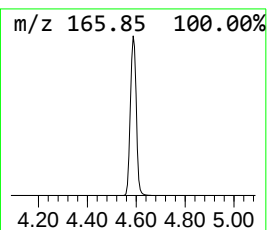
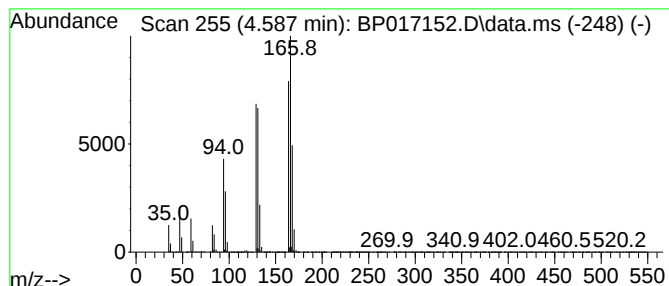
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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.
4.587	538.88 ng/ul	28731000	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	87
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.587	538.9	ng/ul	28731000	1	7.952	1066330	20.0