

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017461.D
 Acq On : 30 Sep 2023 00:07
 Operator : MA/JU
 Sample : 04497-21
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ05

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-BP092023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017461.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	22	28	34	rBV	963293	1325269	3.14%	1.328%
2	3.746	109	112	122	rBV	57574	97751	0.23%	0.098%
3	3.923	138	142	150	rVB	36968	55246	0.13%	0.055%
4	4.570	244	252	255	rBV	22484416	42233639	100.00%	42.313%
5	4.593	255	256	278	rVB	18583856	10221627	24.20%	10.241%
6	7.105	677	683	695	rVB	206208	342480	0.81%	0.343%
7	7.293	708	715	725	rBV	654898	991795	2.35%	0.994%
8	7.469	738	745	756	rBV	698793	1097260	2.60%	1.099%
9	7.952	820	827	837	rBV	740350	1155119	2.74%	1.157%
10	8.669	940	949	959	rBV	530205	869382	2.06%	0.871%
11	9.158	1024	1032	1044	rBV	749767	1245514	2.95%	1.248%
12	9.875	1146	1154	1165	rBV	621937	1005338	2.38%	1.007%
13	10.405	1236	1244	1261	rBV	909244	1541180	3.65%	1.544%
14	10.787	1300	1309	1318	rBV	944778	1577924	3.74%	1.581%
15	10.957	1330	1338	1358	rBV	746617	1269523	3.01%	1.272%
16	11.804	1475	1482	1489	rBV	54974	121734	0.29%	0.122%
17	14.057	1857	1865	1876	rBV	1706738	2263634	5.36%	2.268%
18	14.340	1906	1913	1924	rBV	1896435	2783559	6.59%	2.789%
19	14.646	1957	1965	1973	rBV	1399576	2039688	4.83%	2.044%
20	14.887	1998	2006	2021	rBV	106294	224634	0.53%	0.225%
21	15.645	2128	2135	2143	rBV	2580208	3513968	8.32%	3.521%
22	15.787	2153	2159	2168	rBV	692656	904398	2.14%	0.906%
23	17.428	2432	2438	2449	rBV	1845456	2509318	5.94%	2.514%
24	17.534	2449	2456	2466	rBV2	2759598	4045644	9.58%	4.053%
25	18.275	2578	2582	2587	rBV2	87334	120783	0.29%	0.121%
26	19.351	2761	2765	2769	rBV	34901	46672	0.11%	0.047%
27	19.422	2773	2777	2783	rVB	31468	46743	0.11%	0.047%
28	19.516	2790	2793	2799	rBV3	38985	49997	0.12%	0.050%
29	19.781	2832	2838	2844	rBV	3721380	4909108	11.62%	4.918%
30	21.557	3135	3140	3149	rBV	2233846	2910311	6.89%	2.916%
31	23.963	3542	3549	3564	rVB2	2446035	5182022	12.27%	5.192%
32	24.133	3571	3578	3588	rVB	1465391	3110077	7.36%	3.116%

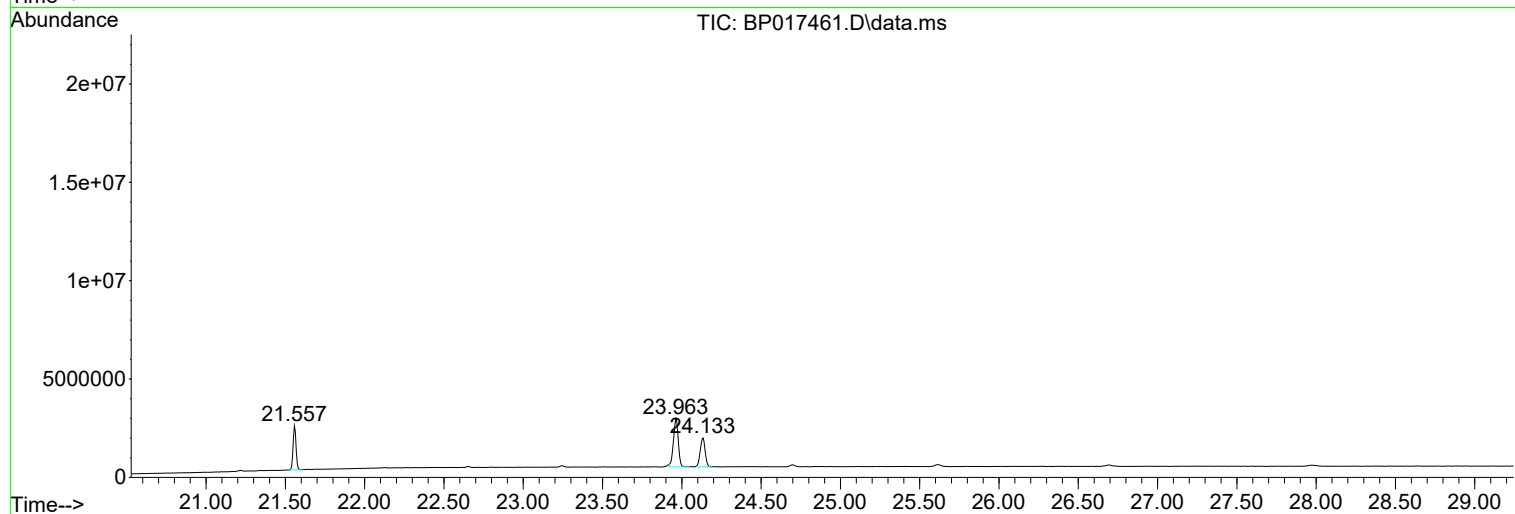
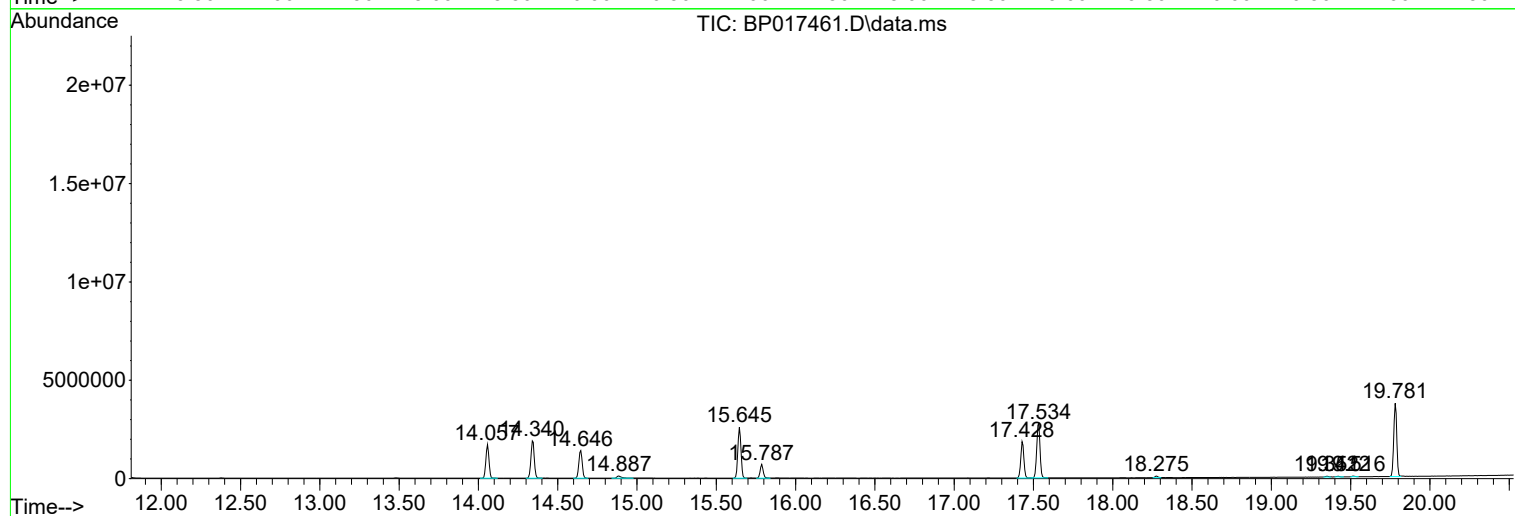
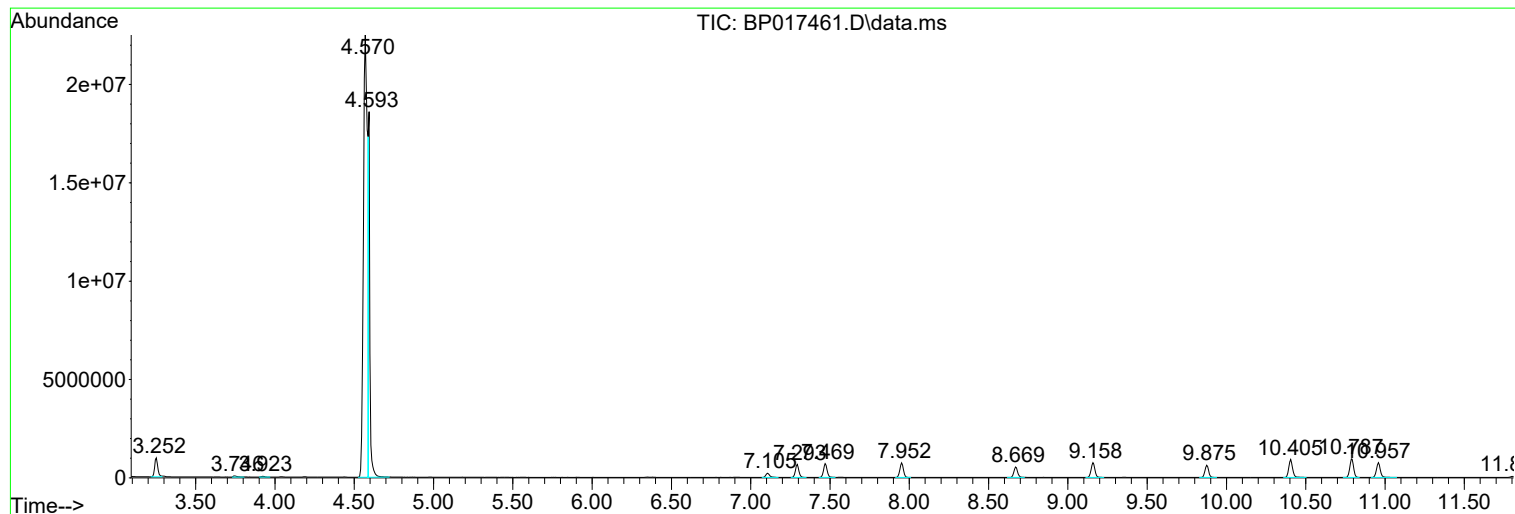
Sum of corrected areas: 99811337

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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



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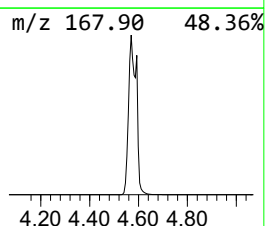
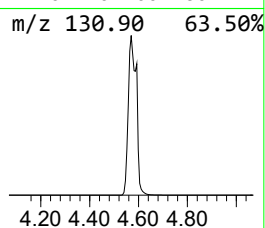
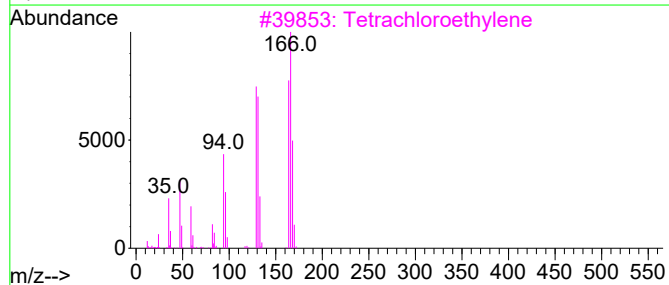
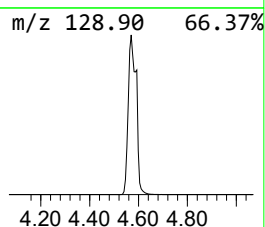
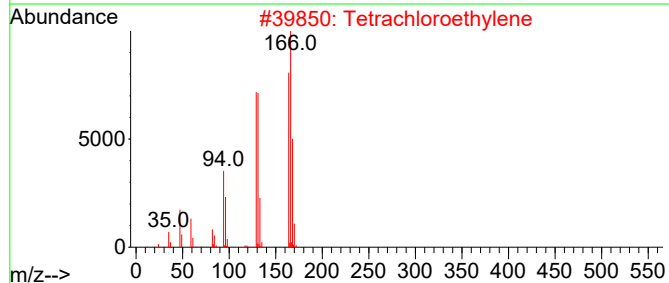
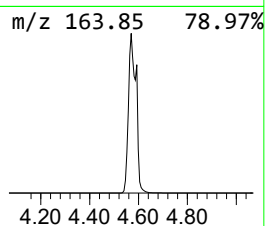
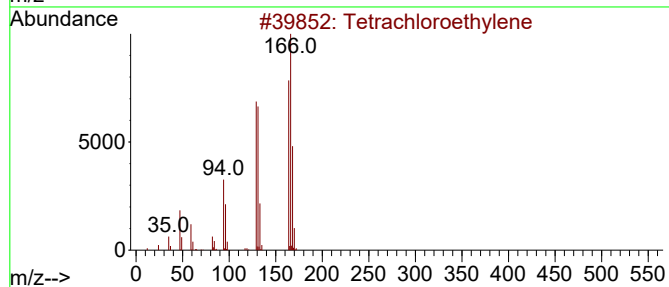
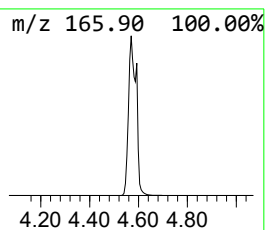
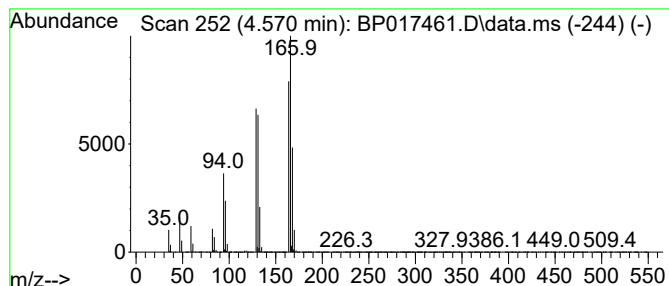
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.570	731.24 ng/ul	42233600	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	97
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
5			2,4-Dichloro-6-fluoropyrimidine	166	C4HC12FN2	003833-57-6	38



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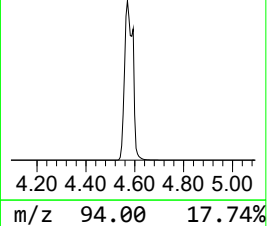
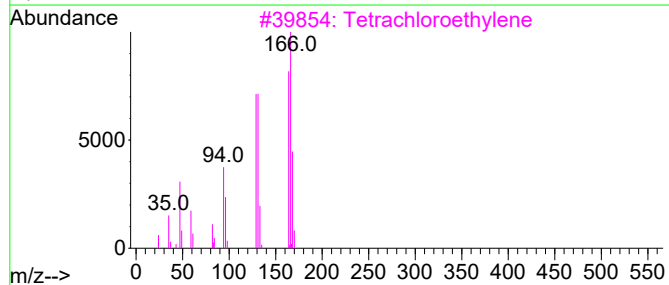
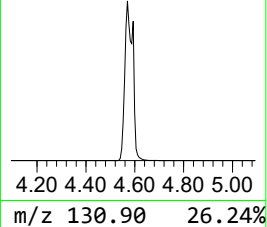
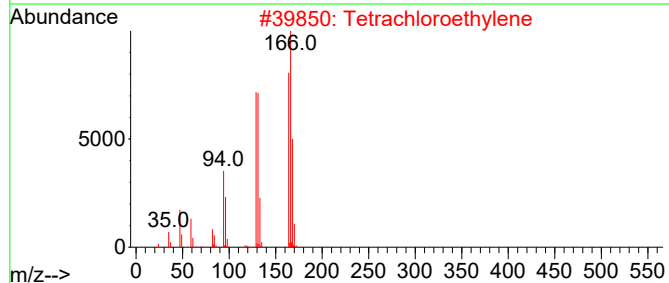
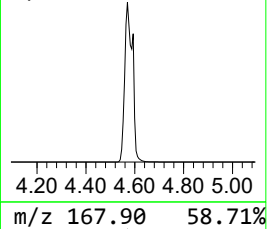
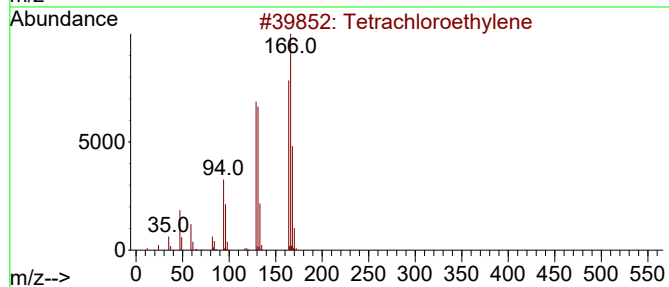
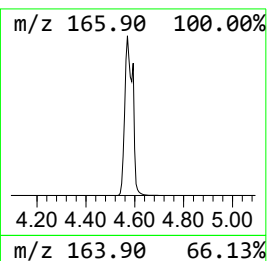
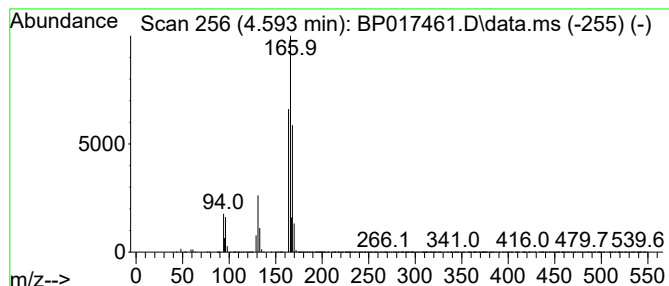
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Peak Number 2 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.593	176.98 ng/ul	10221600	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	56
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	56
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	50
4			Norketamine	223	C12H14ClNO	035211-10-0	33
5			1H-Purine, 8-(methylthio)-	166	C6H6N4S	033426-53-8	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.570	731.2	ng/ul	42233600	1	7.952	1155120	20.0
unknown-02	4.593	177.0	ng/ul	10221600	1	7.952	1155120	20.0