

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013215.D
 Acq On : 21 Dec 2022 19:12
 Operator : CG/JU
 Sample : N6032-02
 Misc :
 ALS Vial : 97 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0L85

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013215.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.299	14	19	20	rBV	31621209	37842456	100.00%	28.841%
2	3.781	99	101	109	rBV	81272	122495	0.32%	0.093%
3	4.093	151	154	160	rVB	36403	49619	0.13%	0.038%
4	4.199	167	172	180	rVB	294854	400880	1.06%	0.306%
5	4.628	240	245	255	rVB	502528	751001	1.98%	0.572%
6	5.040	309	315	320	rBV2	29840	55544	0.15%	0.042%
7	7.110	661	667	676	rVB	405506	641211	1.69%	0.489%
8	7.275	688	695	705	rBV	1499303	2328242	6.15%	1.774%
9	7.475	722	729	740	rBV	1169783	1886716	4.99%	1.438%
10	7.952	803	810	817	rBV	1758619	2759714	7.29%	2.103%
11	8.657	921	930	946	rBV	934486	1513621	4.00%	1.154%
12	9.116	1000	1008	1022	rBV	1863277	3047794	8.05%	2.323%
13	9.840	1124	1131	1145	rBV	1265509	2135342	5.64%	1.627%
14	10.381	1214	1223	1235	rBV	1719305	2859382	7.56%	2.179%
15	10.763	1280	1288	1297	rBV	2377416	3876637	10.24%	2.955%
16	10.904	1305	1312	1323	rBV	682633	1187939	3.14%	0.905%
17	12.345	1550	1557	1564	rBV	38750	68726	0.18%	0.052%
18	13.998	1831	1838	1848	rBV	3772384	5095372	13.46%	3.883%
19	14.287	1880	1887	1907	rBV	4412587	6379469	16.86%	4.862%
20	14.592	1933	1939	1951	rVB	3186717	4599609	12.15%	3.506%
21	14.792	1968	1973	1989	rBV	135257	289911	0.77%	0.221%
22	15.586	2100	2108	2122	rBV	5555824	7788325	20.58%	5.936%
23	15.704	2122	2128	2141	rVV	1117111	1547346	4.09%	1.179%
24	15.828	2144	2149	2155	rVB3	25763	47891	0.13%	0.036%
25	17.345	2401	2407	2418	rBV	3245079	4702486	12.43%	3.584%
26	17.445	2418	2424	2437	rVV	5442754	7725988	20.42%	5.888%
27	18.216	2551	2555	2562	rBV6	65053	96005	0.25%	0.073%
28	19.257	2728	2732	2736	rBV2	65331	87976	0.23%	0.067%
29	19.322	2740	2743	2747	rVB	66733	85243	0.23%	0.065%
30	19.680	2798	2804	2815	rBV	6942331	9314378	24.61%	7.099%
31	21.433	3097	3102	3109	rBV	4162448	5363876	14.17%	4.088%
32	23.751	3489	3496	3507	rVB2	5037290	10444258	27.60%	7.960%
33	23.910	3517	3523	3535	rVB2	2972651	6115231	16.16%	4.661%

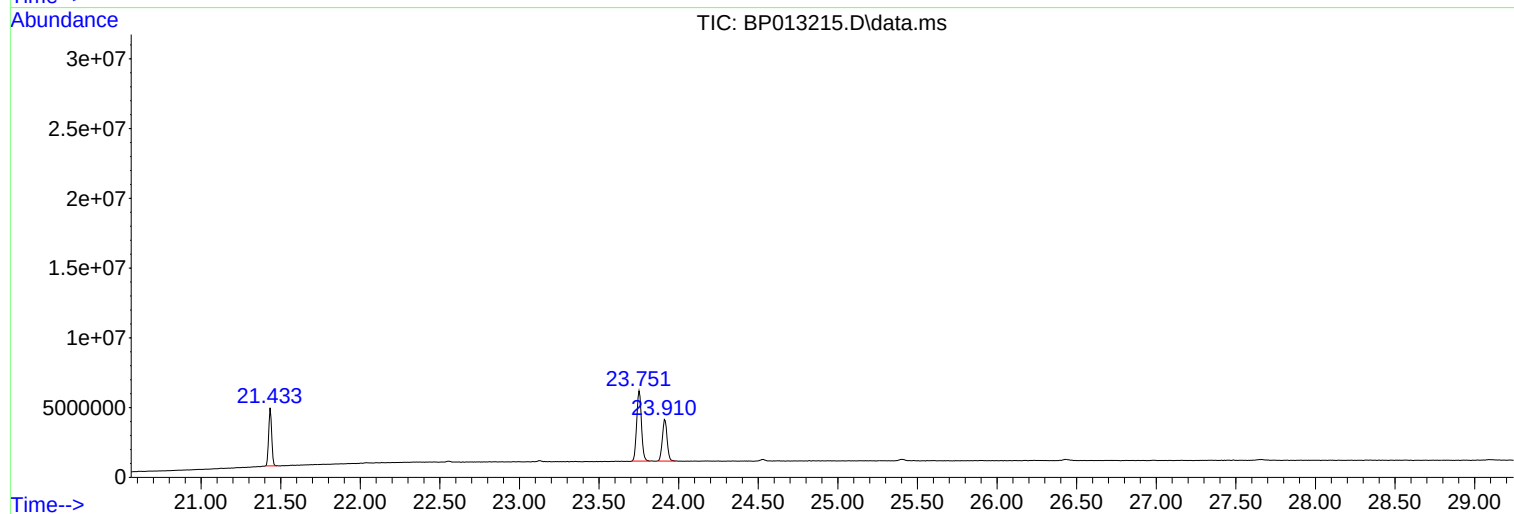
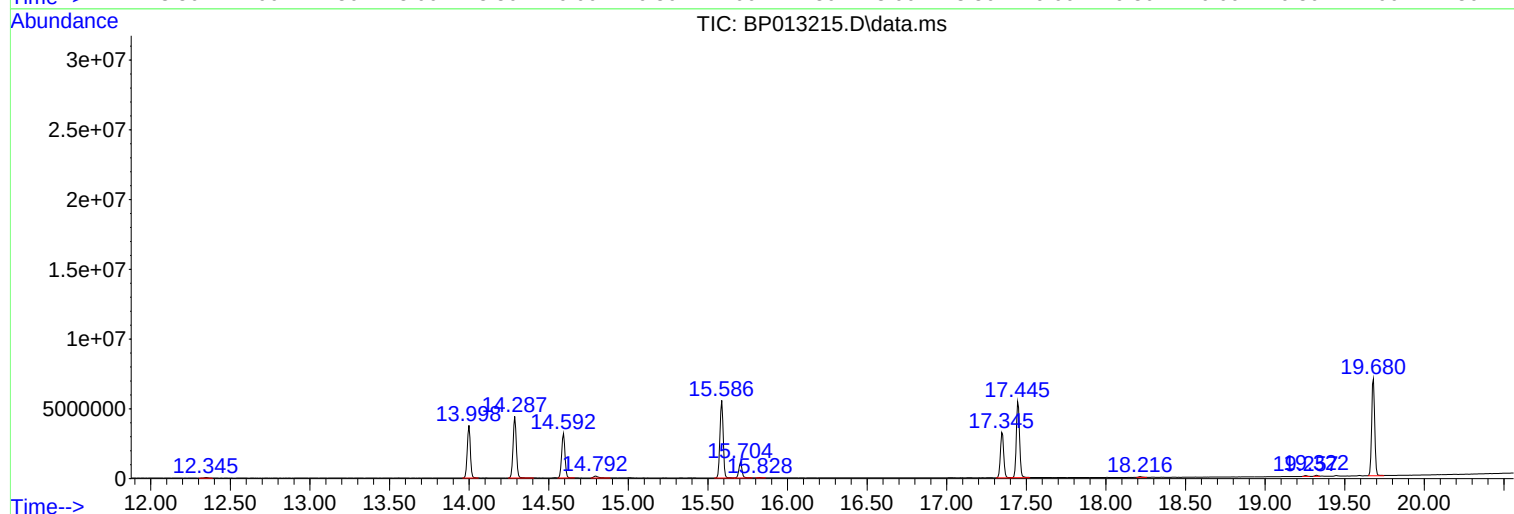
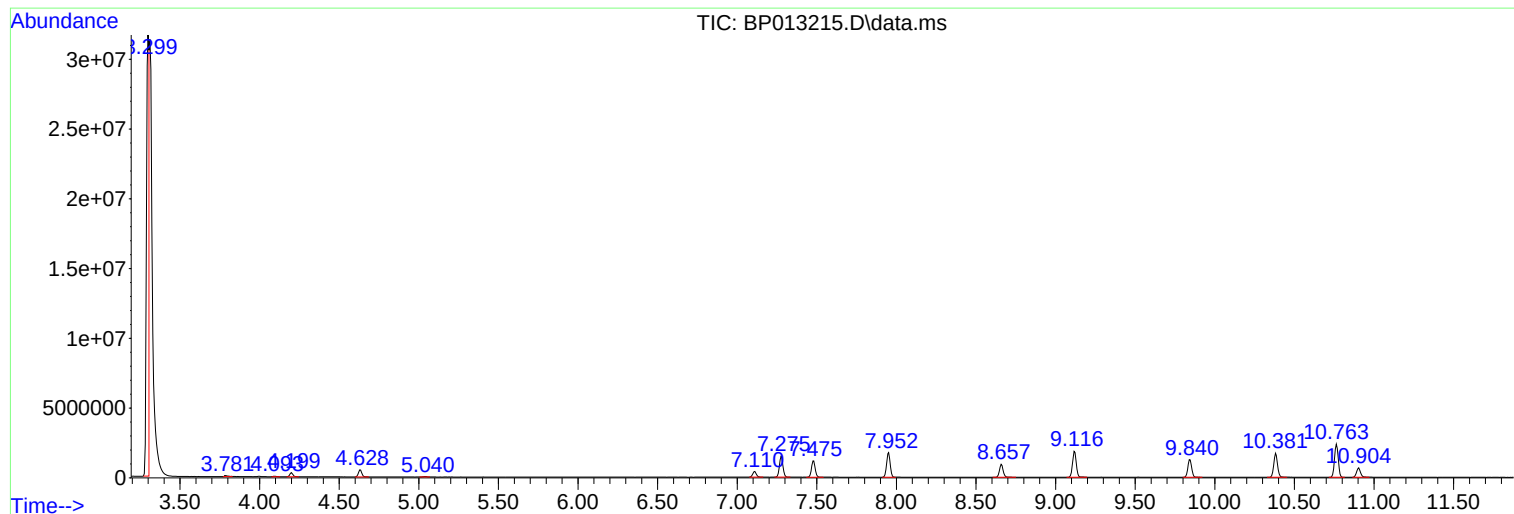
Sum of corrected areas: 131210683

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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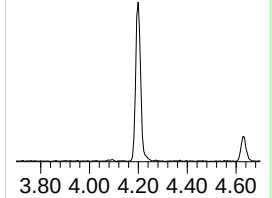
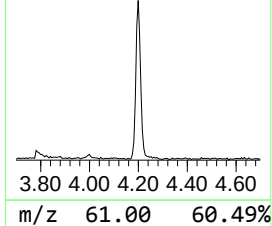
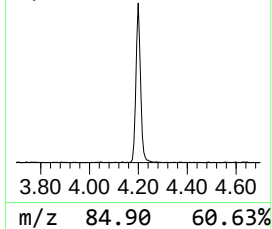
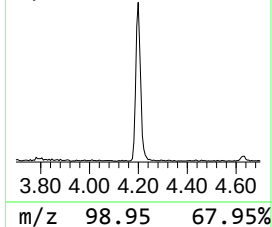
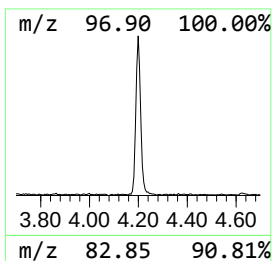
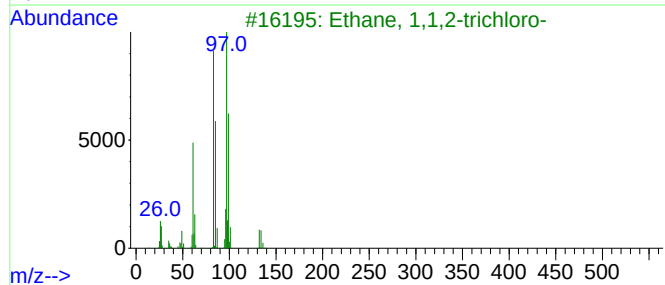
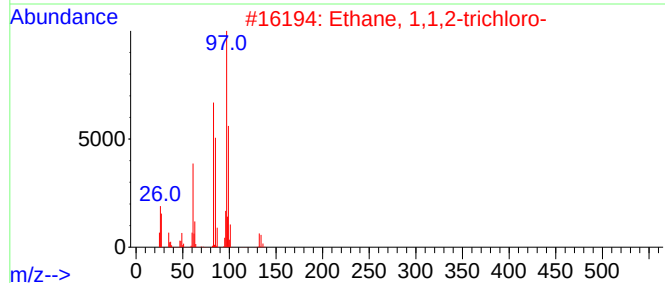
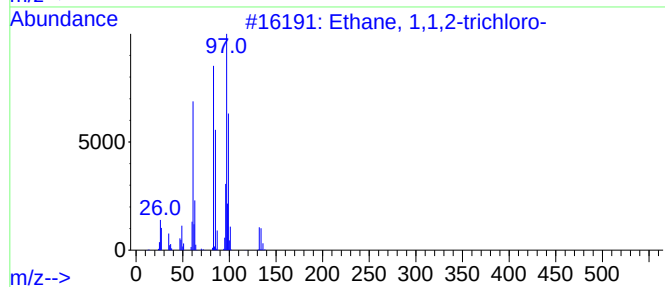
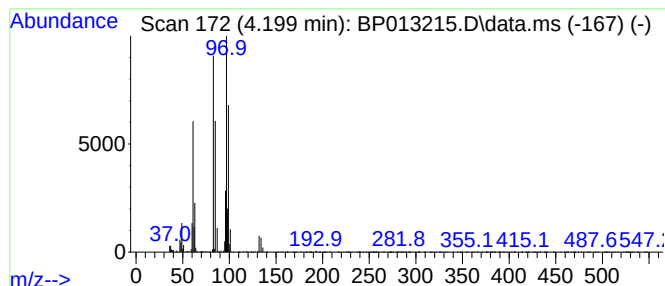
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethane, 1,1,2-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.199	2.91 ng/ul	400880	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	97
2			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	95
3			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	81
4			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	76
5			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	64



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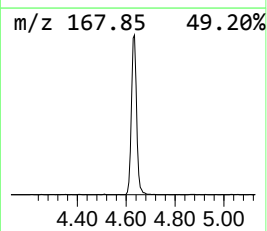
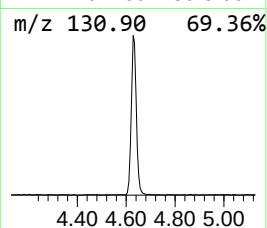
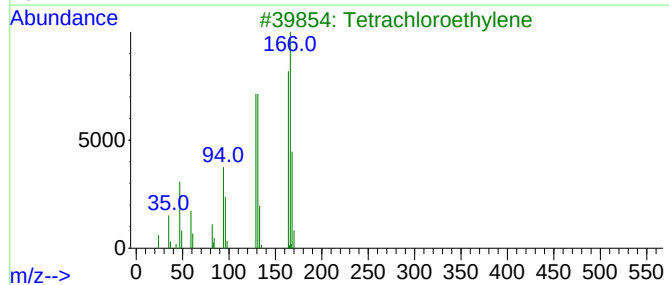
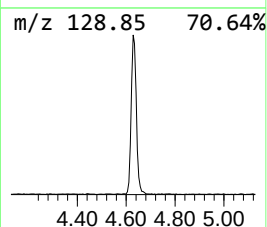
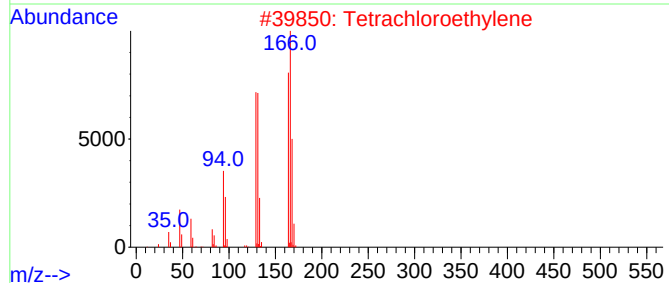
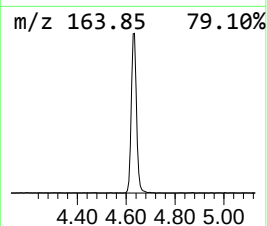
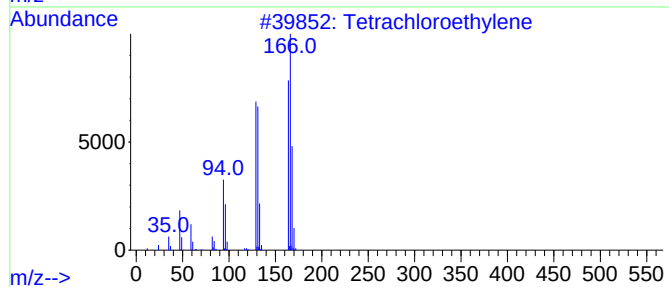
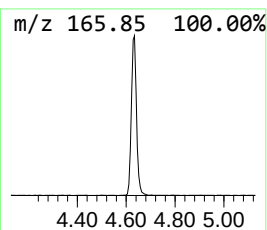
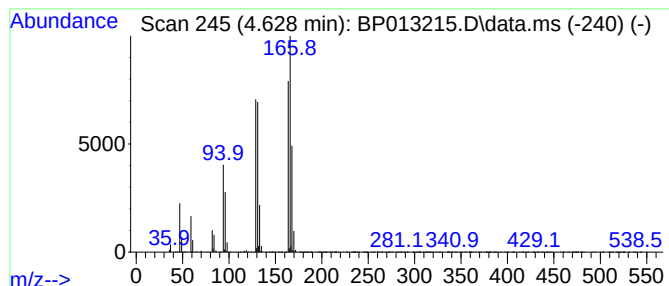
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Peak Number 2 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.628	5.44 ng/ul	751001	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,1,2-t...	4.199	2.9	ng/ul	400880	1	7.952	2759710	20.0
Tetrachloroethy...	4.628	5.4	ng/ul	751001	1	7.952	2759710	20.0