

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013225.D
 Acq On : 22 Dec 2022 01:27
 Operator : CG/JU
 Sample : N6032-08
 Misc :
 ALS Vial : 108 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COLA0

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: BP013225.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.293	14	18	20	rBV	32444673	40578347	64.19%	19.285%
2	3.776	99	100	111	rBV	126848	217005	0.34%	0.103%
3	4.634	239	246	255	rBV	33213288	63218183	100.00%	30.045%
4	4.717	255	260	272	rVB	2015134	2843020	4.50%	1.351%
5	5.040	308	315	319	rVB2	41521	65832	0.10%	0.031%
6	5.311	356	361	366	rVV4	48371	75281	0.12%	0.036%
7	7.105	656	666	678	rBV	516145	807163	1.28%	0.384%
8	7.275	688	695	707	rVB	1829107	2809104	4.44%	1.335%
9	7.475	722	729	738	rBV	1534890	2390953	3.78%	1.136%
10	7.946	801	809	817	rBV	1838460	2874027	4.55%	1.366%
11	8.658	922	930	943	rBV	1227106	1961547	3.10%	0.932%
12	9.116	1001	1008	1015	rBV	2175084	3510882	5.55%	1.669%
13	9.840	1122	1131	1143	rBV	1520461	2478116	3.92%	1.178%
14	10.381	1213	1223	1236	rBV	2078042	3561198	5.63%	1.693%
15	10.763	1279	1288	1297	rBV	2435754	4025546	6.37%	1.913%
16	10.899	1304	1311	1329	rBV	607010	1103258	1.75%	0.524%
17	12.346	1551	1557	1565	rVB2	48475	89682	0.14%	0.043%
18	13.999	1830	1838	1852	rBV	4263639	5860563	9.27%	2.785%
19	14.287	1879	1887	1903	rBV	4908757	7230228	11.44%	3.436%
20	14.587	1932	1938	1949	rVB	3293526	4803566	7.60%	2.283%
21	14.793	1966	1973	1988	rBV	180985	389907	0.62%	0.185%
22	15.587	2100	2108	2119	rBV	6031975	8892031	14.07%	4.226%
23	15.698	2121	2127	2141	rVV	1275979	1842575	2.91%	0.876%
24	17.345	2400	2407	2417	rBV	3488300	4857472	7.68%	2.309%
25	17.445	2418	2424	2438	rVV	6356150	8825505	13.96%	4.194%
26	19.251	2728	2731	2736	rVB	75306	93286	0.15%	0.044%
27	19.322	2739	2743	2747	rBV2	72412	102398	0.16%	0.049%
28	19.675	2797	2803	2817	rBV	7974181	10591733	16.75%	5.034%
29	21.433	3097	3102	3110	rBV	4122495	5478073	8.67%	2.604%
30	23.751	3488	3496	3513	rVB2	5952442	12450693	19.69%	5.917%
31	23.910	3516	3523	3537	rVB2	3117791	6383301	10.10%	3.034%

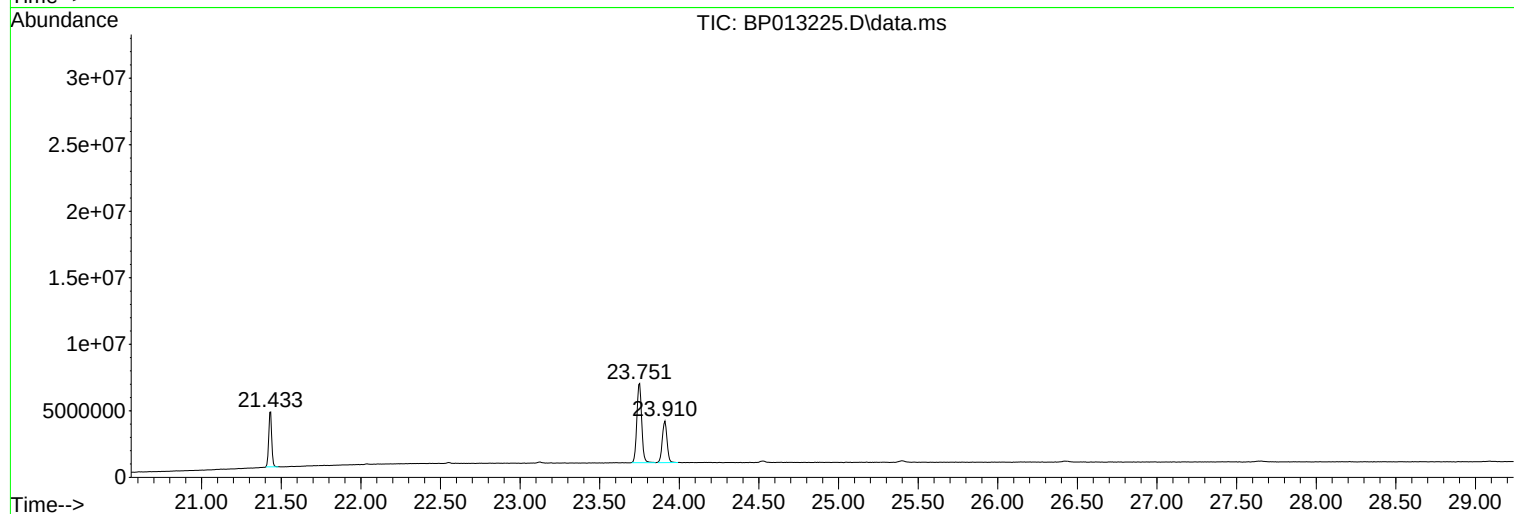
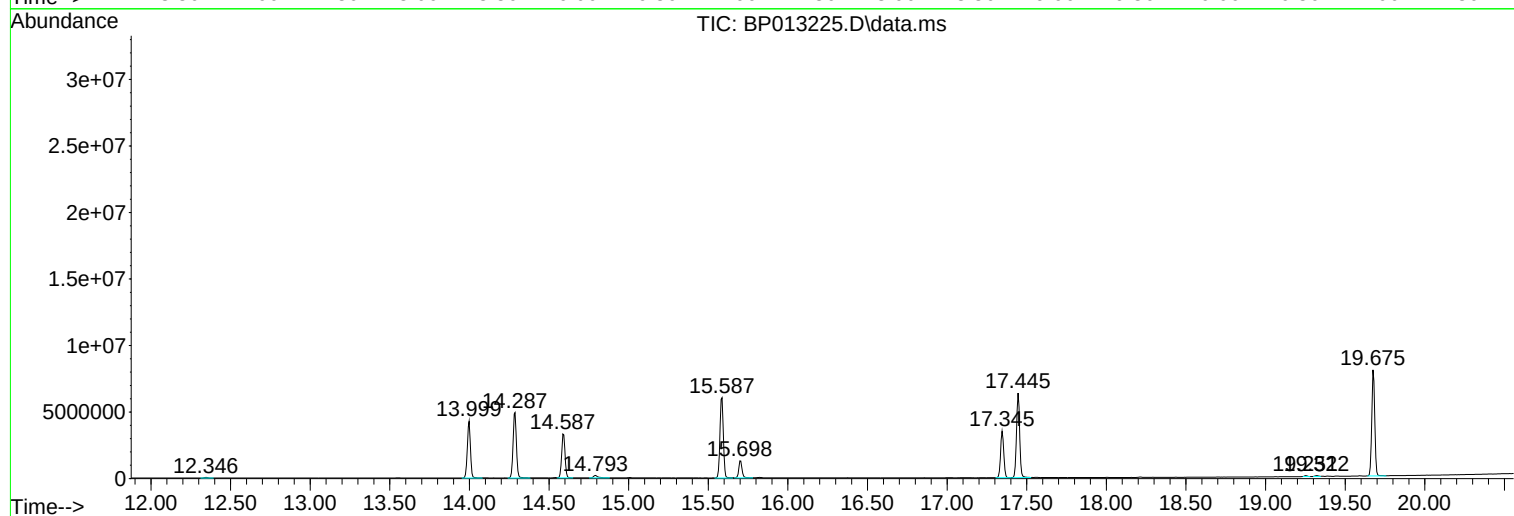
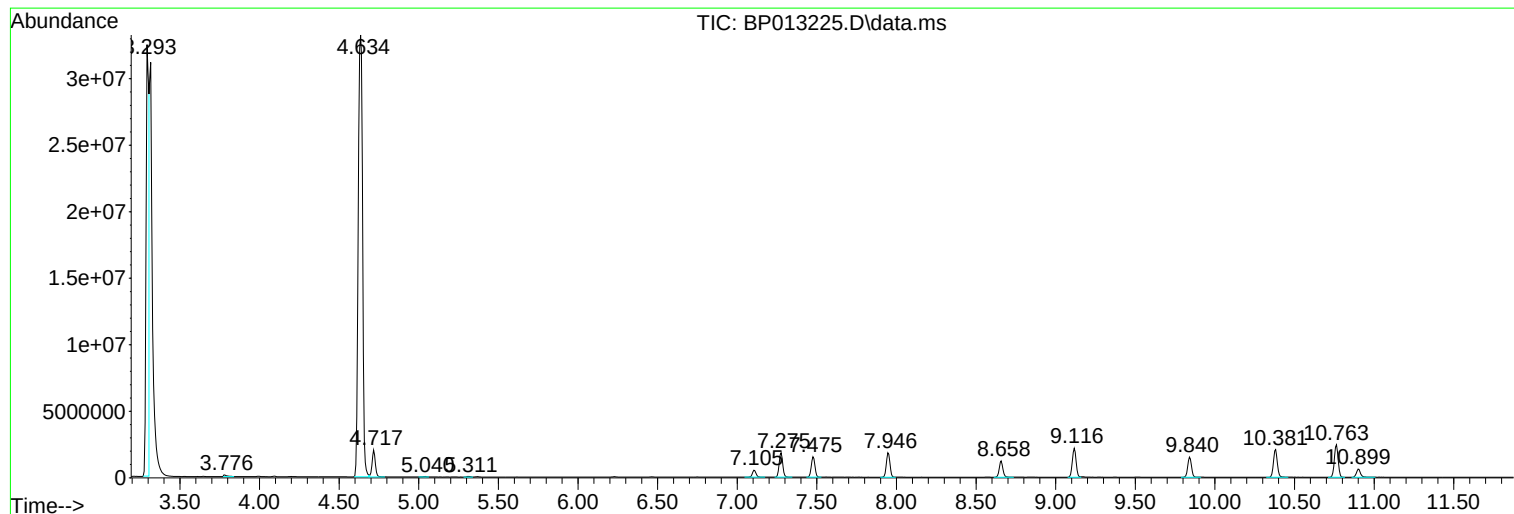
Sum of corrected areas: 210410475

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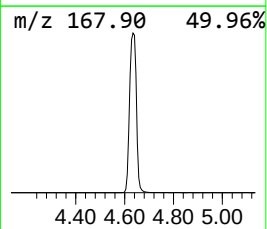
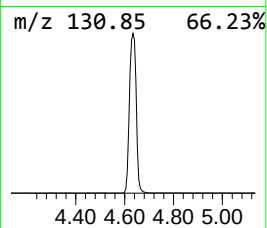
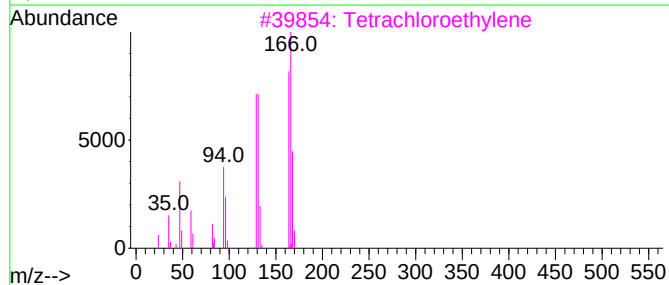
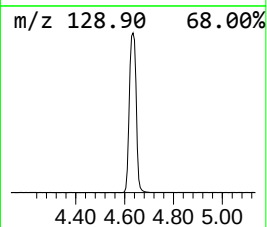
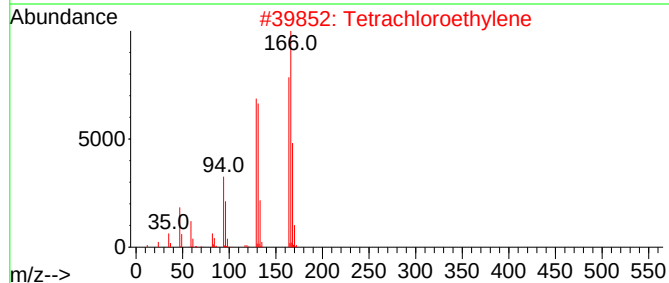
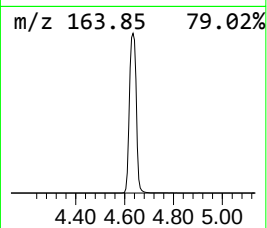
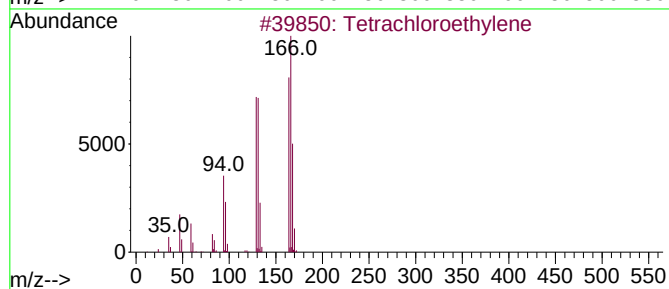
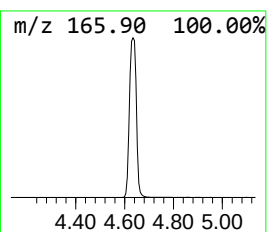
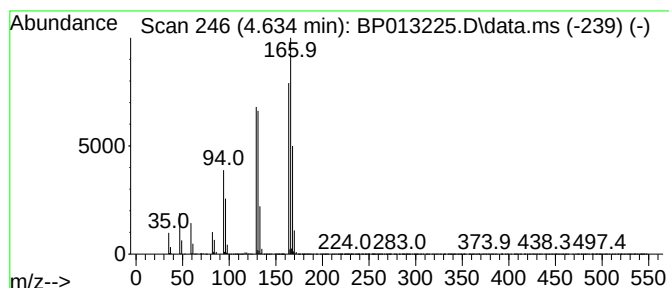
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.634	439.93 ng/ul	63218200	1,4-Dichlorobenzene-d4	7.946

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	98
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	38



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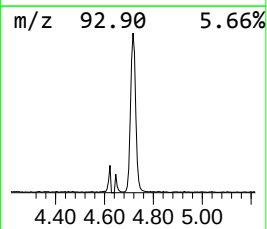
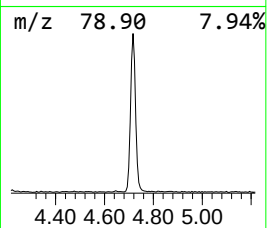
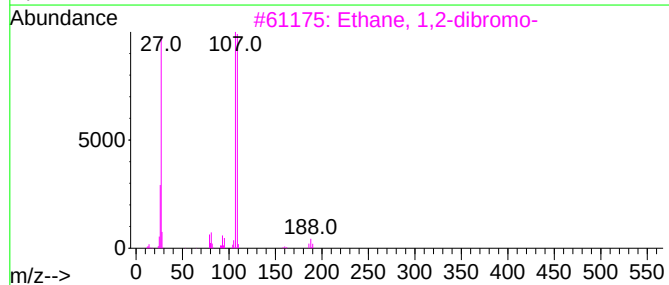
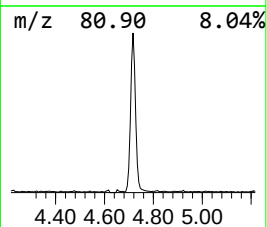
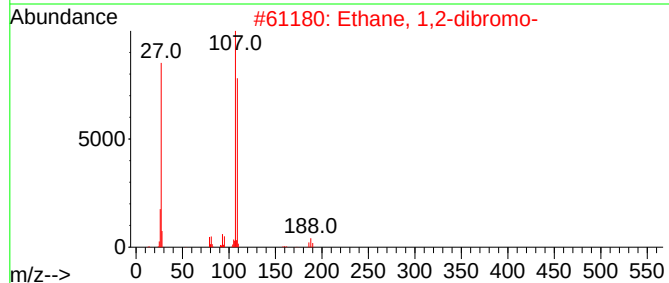
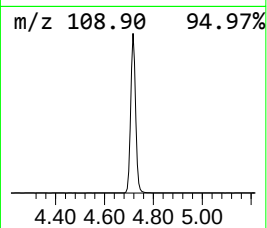
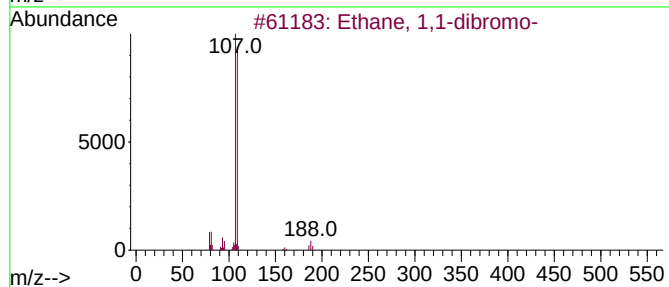
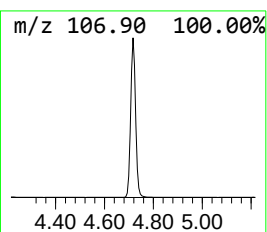
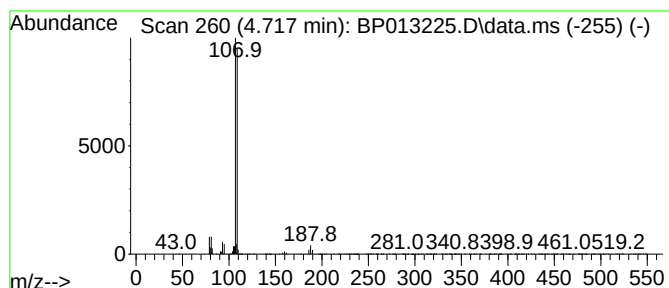
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.717	19.78 ng/ul	2843020	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,1-dibromo-	186	C2H4Br2	000557-91-5	97
2			Ethane, 1,2-dibromo-	186	C2H4Br2	000106-93-4	94
3			Ethane, 1,2-dibromo-	186	C2H4Br2	000106-93-4	93
4			Ethane, 1,1-dibromo-	186	C2H4Br2	000557-91-5	91
5			Ethane, 1,1-dibromo-	186	C2H4Br2	000557-91-5	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.634	439.9	ng/ul	63218200	1	7.946	2874030	20.0
Ethane, 1,1-dib...	4.717	19.8	ng/ul	2843020	1	7.946	2874030	20.0