

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013226.D
 Acq On : 22 Dec 2022 02:01
 Operator : CG/JU
 Sample : N6032-09
 Misc :
 ALS Vial : 109 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COLA1

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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: BP013226.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	32603607	40150371	82.83%	20.681%
2	3.317	20	22	50	rVB	31924601	48474069	100.00%	24.969%
3	3.781	98	101	110	rBV	117055	195730	0.40%	0.101%
4	3.993	134	137	143	rVB	37798	54650	0.11%	0.028%
5	4.093	150	154	161	rVB	43144	62813	0.13%	0.032%
6	4.199	167	172	182	rVB	375202	514328	1.06%	0.265%
7	4.628	239	245	259	rVB	894855	1290641	2.66%	0.665%
8	5.040	310	315	323	rVB2	56855	89462	0.18%	0.046%
9	7.105	660	666	674	rBV	483748	748602	1.54%	0.386%
10	7.275	686	695	706	rBV	1818815	2781739	5.74%	1.433%
11	7.475	721	729	739	rBV	1542987	2389202	4.93%	1.231%
12	7.946	802	809	818	rBV	1793605	2855629	5.89%	1.471%
13	8.657	922	930	945	rBV	1187300	1899711	3.92%	0.979%
14	9.116	1000	1008	1015	rBV	2127535	3398618	7.01%	1.751%
15	9.840	1124	1131	1145	rBV	1544300	2446549	5.05%	1.260%
16	10.381	1215	1223	1235	rBV	2104681	3449165	7.12%	1.777%
17	10.763	1280	1288	1295	rBV	2377580	3904141	8.05%	2.011%
18	10.899	1303	1311	1330	rBV	919493	1623306	3.35%	0.836%
19	12.346	1551	1557	1564	rBV2	41675	75964	0.16%	0.039%
20	13.545	1755	1761	1770	rBV9	23089	50008	0.10%	0.026%
21	13.998	1830	1838	1850	rBV	3789883	5396987	11.13%	2.780%
22	14.281	1879	1886	1900	rBV	4572264	6862674	14.16%	3.535%
23	14.592	1932	1939	1947	rVB	3130460	4586421	9.46%	2.362%
24	14.792	1965	1973	1983	rBV	153988	335211	0.69%	0.173%
25	15.581	2099	2107	2122	rBV	5902546	8439407	17.41%	4.347%
26	15.704	2122	2128	2136	rBV	1220720	1710268	3.53%	0.881%
27	15.822	2142	2148	2155	rVB4	30151	56660	0.12%	0.029%
28	17.345	2399	2407	2417	rBV	3499205	4885312	10.08%	2.516%
29	17.445	2417	2424	2441	rBV	6331093	8769344	18.09%	4.517%
30	18.210	2550	2554	2559	rBV4	42579	67870	0.14%	0.035%
31	19.257	2728	2732	2735	rVB2	72945	95817	0.20%	0.049%
32	19.322	2739	2743	2747	rVV	66820	96783	0.20%	0.050%
33	19.675	2797	2803	2813	rBV	8349545	11059804	22.82%	5.697%
34	21.433	3097	3102	3108	rVB	4376906	5862848	12.09%	3.020%
35	23.745	3488	3495	3508	rBV2	6136571	12764640	26.33%	6.575%
36	23.910	3516	3523	3534	rVB2	3340217	6693344	13.81%	3.448%

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BNA_P
ClientSampleId :
C0LA1

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

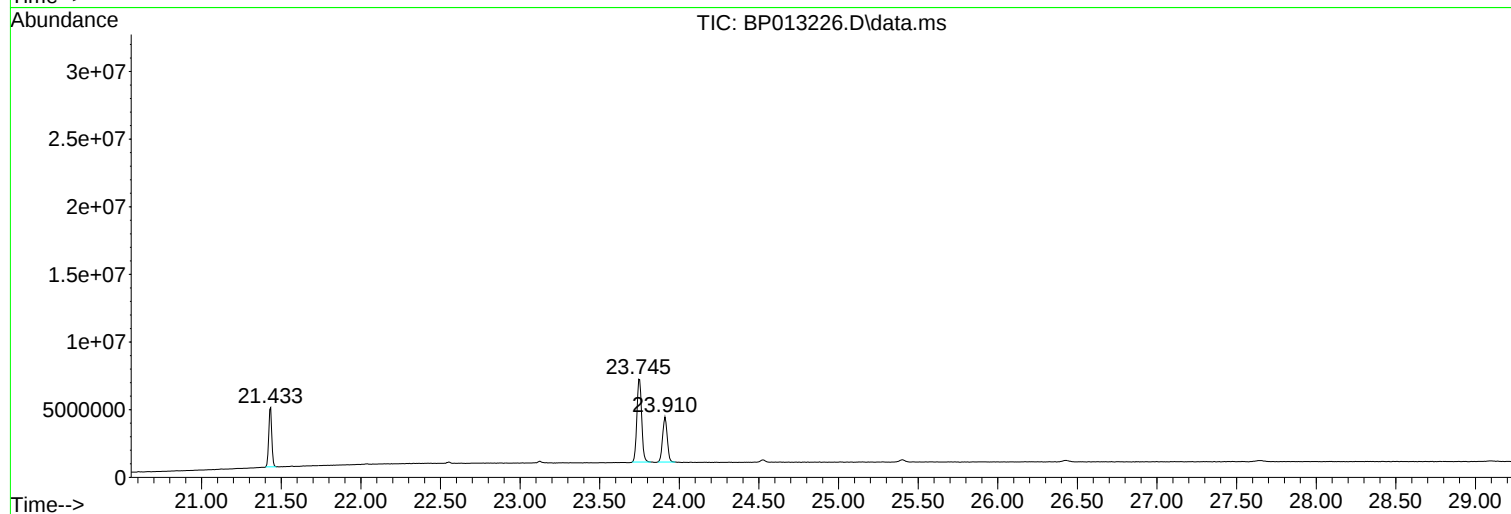
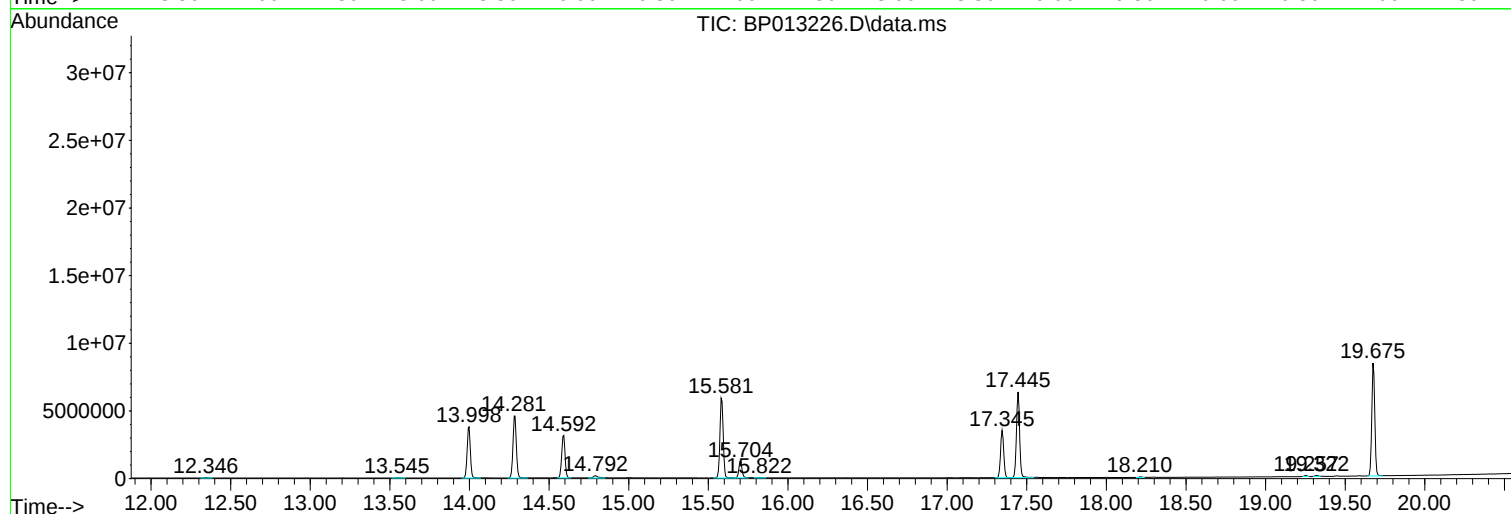
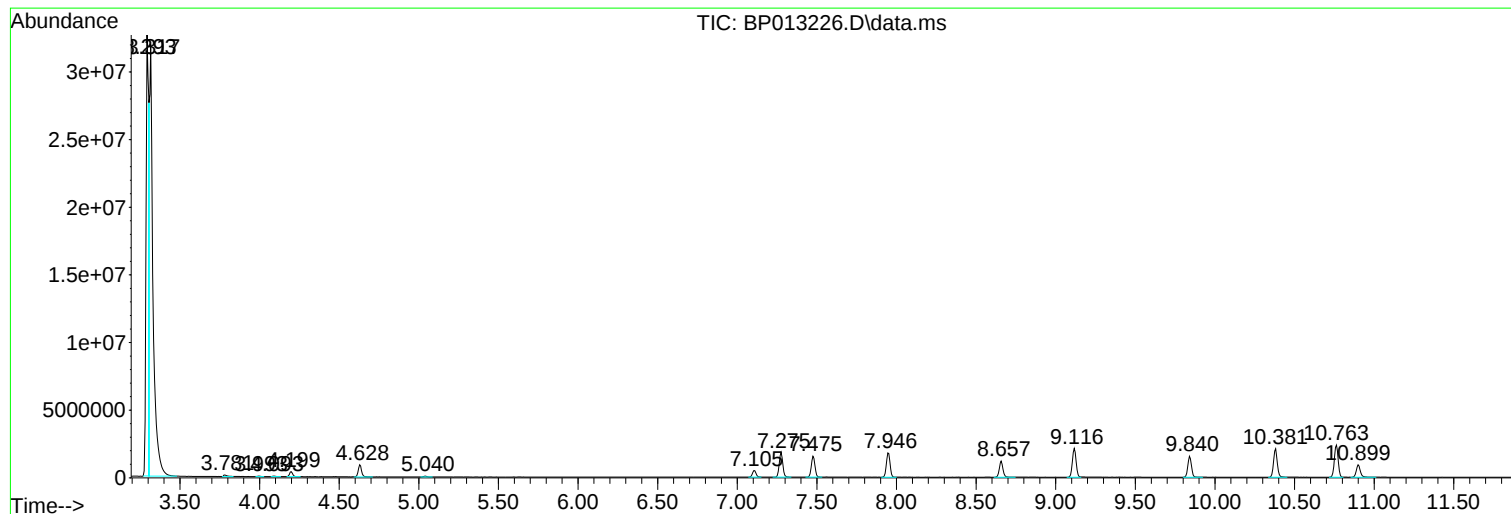
Sum of corrected areas: 194138088

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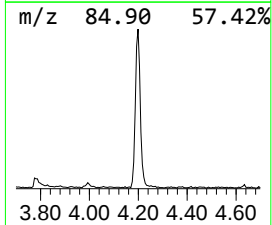
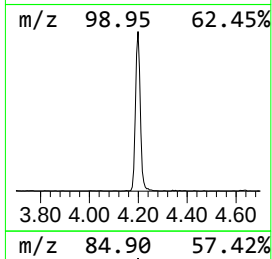
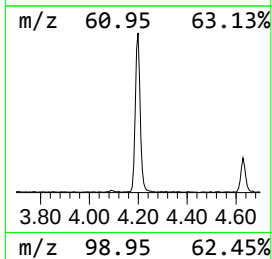
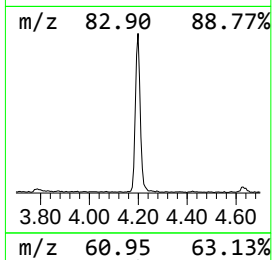
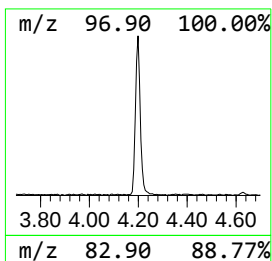
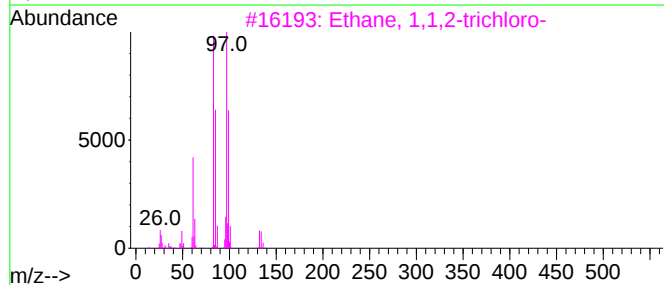
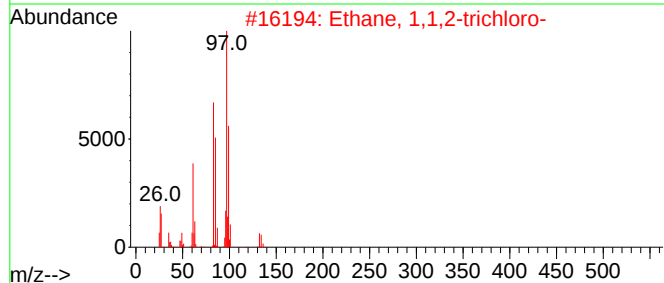
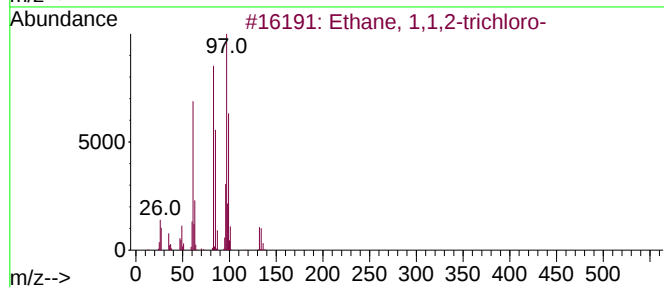
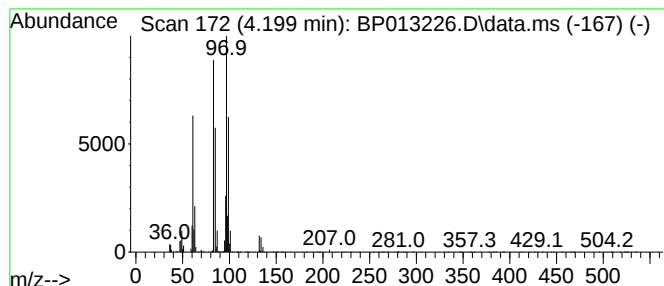
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.199	3.60 ng/ul	514328	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	96
2			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	95
3			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	93
4			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	87
5			Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	76



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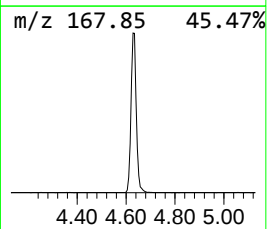
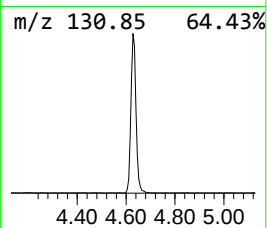
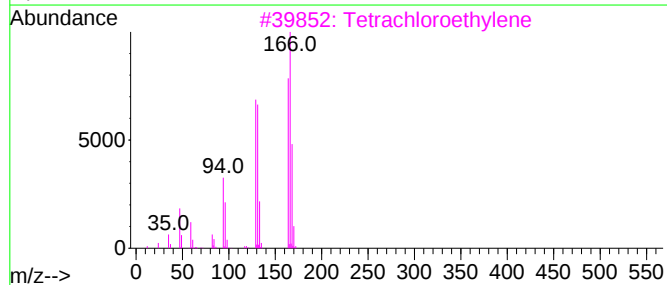
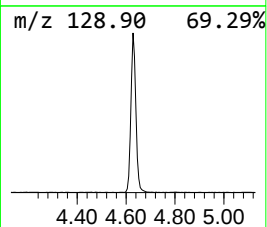
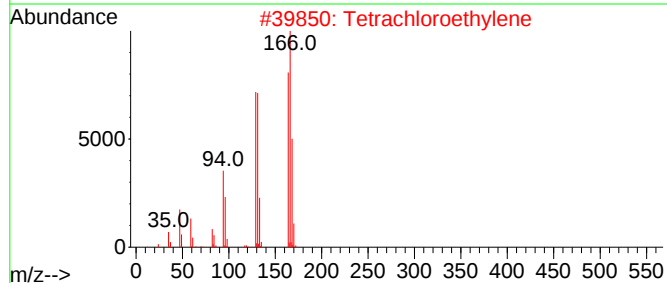
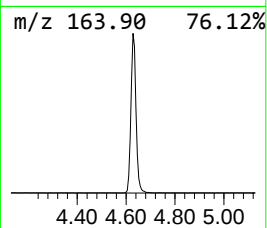
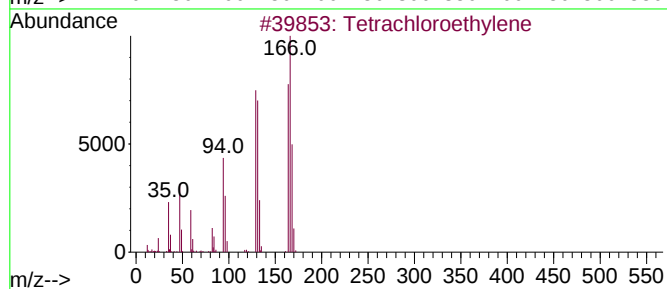
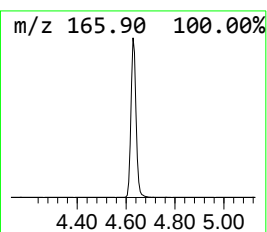
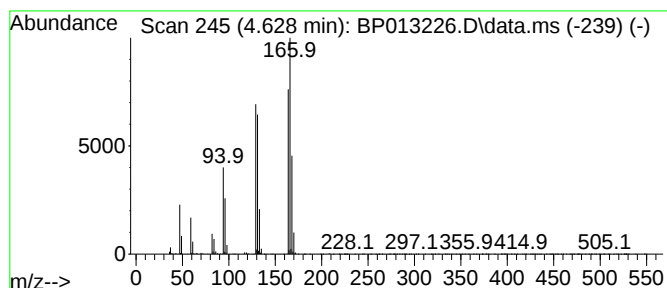
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.628	9.04 ng/ul	1290640	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	98
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HCl2FN2	002927-71-1	38



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,1,2-t...	4.199	3.6	ng/ul	514328	1	7.952	2855630	20.0
Tetrachloroethy...	4.628	9.0	ng/ul	1290640	1	7.952	2855630	20.0