

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082422\
 Data File : BP011566.D
 Acq On : 25 Aug 2022 17:17
 Operator : CG/JU
 Sample : N4349-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGN62

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-BP081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011566.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	11	rBV	185721	246432	8.78%	1.229%
2	3.064	11	13	17	rVB	23339	22424	0.80%	0.112%
3	3.358	61	63	68	rBV4	2385	3218	0.11%	0.016%
4	3.493	84	86	99	rBV	22130	39255	1.40%	0.196%
5	3.687	114	119	126	rVB	14098	22120	0.79%	0.110%
6	3.775	130	134	139	rVB2	10077	13318	0.47%	0.066%
7	4.205	204	207	213	rVB2	3812	5228	0.19%	0.026%
8	4.287	215	221	235	rBV	2069463	2806634	100.00%	14.002%
9	6.734	631	637	649	rBV	73363	119707	4.27%	0.597%
10	6.899	659	665	680	rVB	228871	353722	12.60%	1.765%
11	7.081	690	696	706	rBV	251077	391020	13.93%	1.951%
12	7.546	768	775	786	rBV	288818	433794	15.46%	2.164%
13	8.269	890	898	909	rBV	204809	325451	11.60%	1.624%
14	8.710	966	973	988	rBV	308315	497888	17.74%	2.484%
15	8.846	991	996	1000	rBV6	1986	3178	0.11%	0.016%
16	9.110	1038	1041	1045	rBV6	2666	4039	0.14%	0.020%
17	9.428	1087	1095	1106	rBV	246813	406245	14.47%	2.027%
18	9.963	1179	1186	1198	rBV	329945	551385	19.65%	2.751%
19	10.334	1241	1249	1261	rBV	374330	615155	21.92%	3.069%
20	10.487	1268	1275	1290	rBV	240462	431330	15.37%	2.152%
21	11.616	1462	1467	1472	rBV7	2338	5049	0.18%	0.025%
22	11.945	1516	1523	1528	rVB3	4760	8376	0.30%	0.042%
23	13.187	1730	1734	1740	rBV8	2005	3841	0.14%	0.019%
24	13.645	1805	1812	1823	rBV	713400	966706	34.44%	4.823%
25	13.910	1850	1857	1869	rBV	697269	978325	34.86%	4.881%
26	14.222	1903	1910	1920	rBV	526020	770734	27.46%	3.845%
27	14.457	1944	1950	1958	rBV3	37030	85714	3.05%	0.428%
28	14.669	1977	1986	1994	rVB3	4787	13074	0.47%	0.065%
29	15.222	2074	2080	2094	rBV	927629	1305651	46.52%	6.514%
30	15.363	2096	2104	2116	rBV	377362	524041	18.67%	2.614%
31	15.469	2119	2122	2126	rVB6	3187	4504	0.16%	0.022%
32	15.922	2194	2199	2201	rBV6	2188	3601	0.13%	0.018%
33	15.957	2201	2205	2212	rVV3	13692	24171	0.86%	0.121%
34	16.016	2212	2215	2221	rVB8	2940	4333	0.15%	0.022%
35	16.769	2339	2343	2347	rBV6	2776	4374	0.16%	0.022%
36	16.992	2375	2381	2390	rBV	621258	866309	30.87%	4.322%

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37	17.092	2391	2398	2411	rVV	1187811	1615762	57.57%	8.061%
38	17.298	2429	2433	2436	rBV5	3821	4665	0.17%	0.023%
39	17.692	2494	2500	2503	rBV7	7509	9571	0.34%	0.048%
40	17.910	2531	2537	2546	rBV2	52363	81118	2.89%	0.405%
41	17.980	2547	2549	2554	rVB3	4601	6745	0.24%	0.034%
42	18.927	2706	2710	2715	rBV4	14227	20538	0.73%	0.102%
43	18.998	2718	2722	2728	rVB4	12005	19325	0.69%	0.096%
44	19.163	2746	2750	2754	rBV3	15137	18622	0.66%	0.093%
45	19.351	2777	2782	2791	rBV	1304239	1714548	61.09%	8.554%
46	20.857	3034	3038	3046	rBV3	81115	169780	6.05%	0.847%
47	21.121	3078	3083	3089	rBV	736132	910105	32.43%	4.540%
48	23.263	3440	3447	3454	rBV2	942986	1706560	60.80%	8.514%
49	23.410	3466	3472	3479	rVB2	483927	907180	32.32%	4.526%

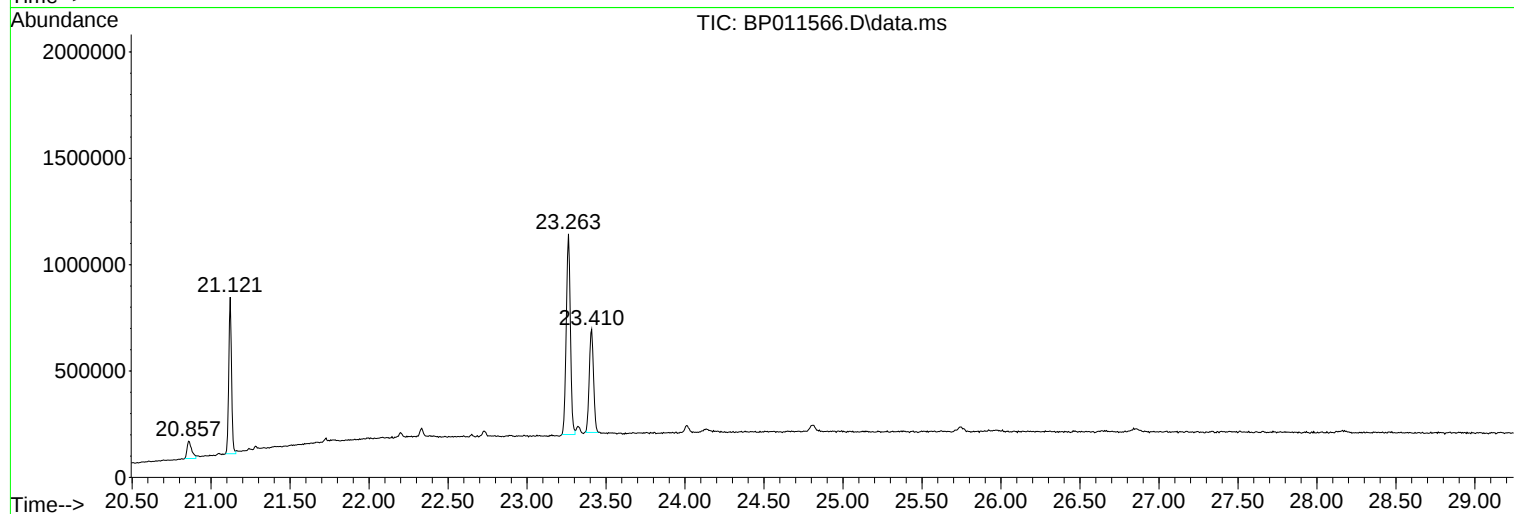
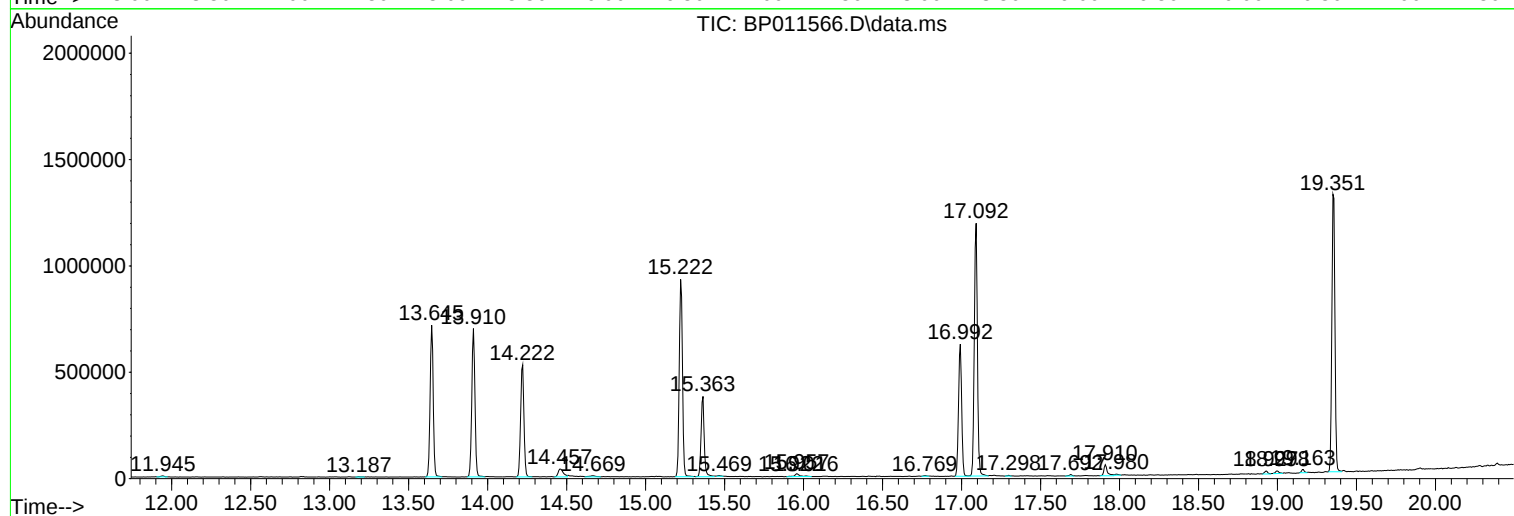
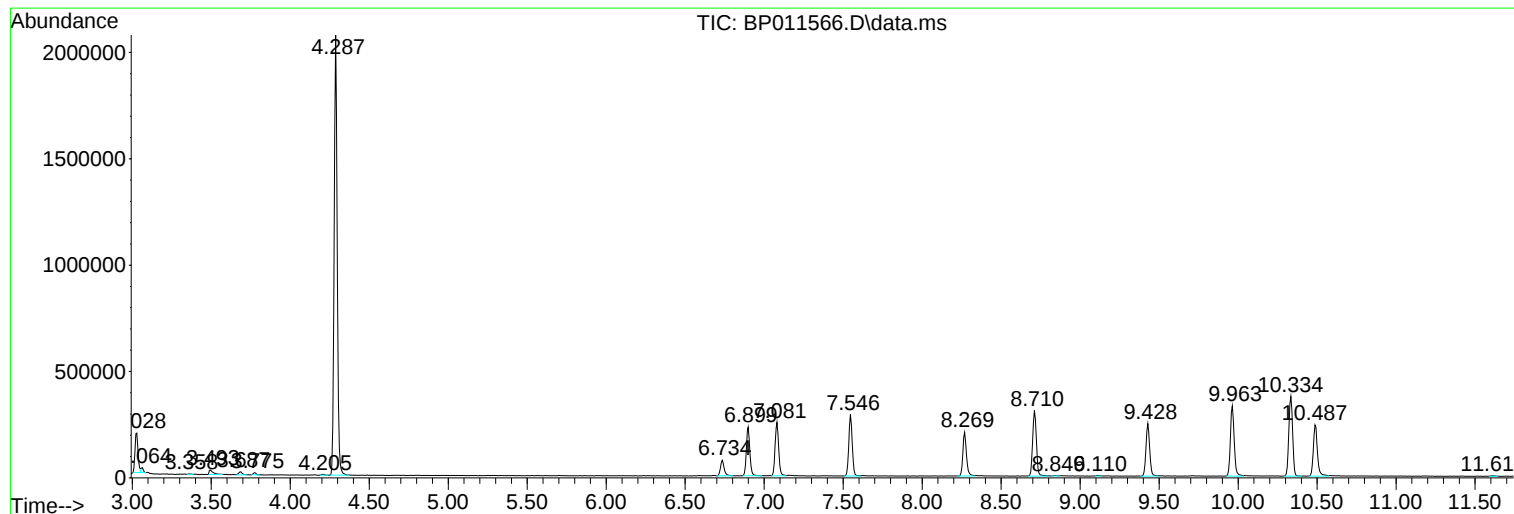
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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



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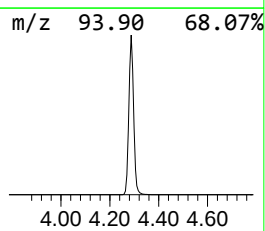
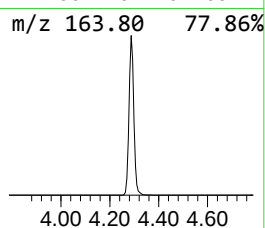
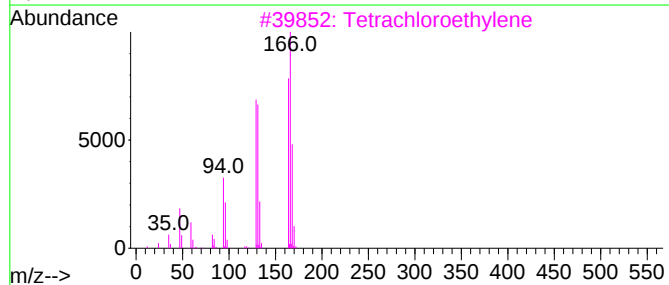
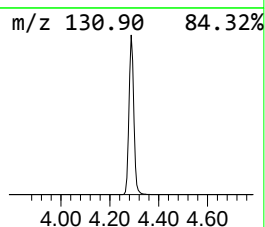
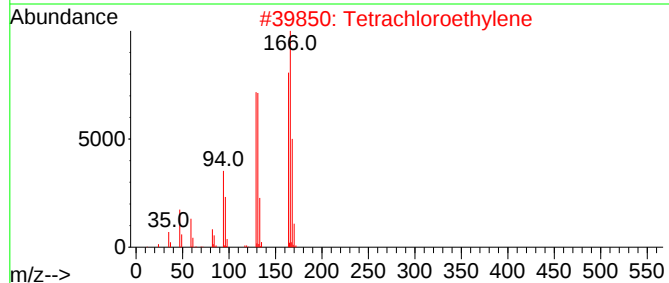
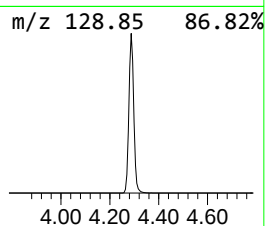
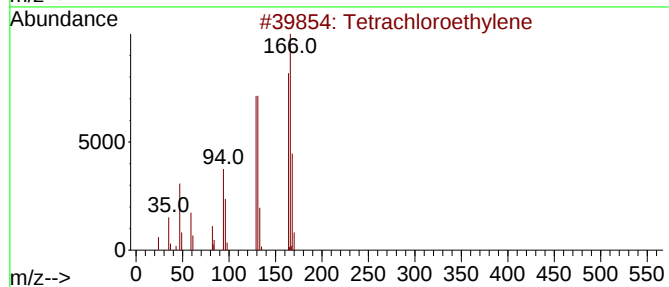
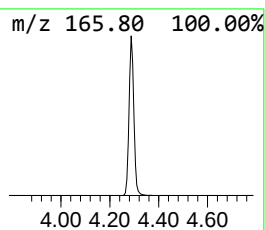
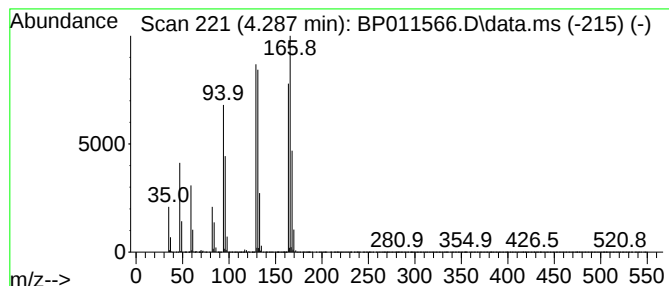
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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.
4.287	129.40 ng/ul	2806630	1,4-Dichlorobenzene-d4	7.546

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	98
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	27



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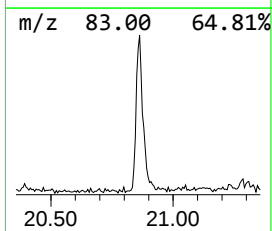
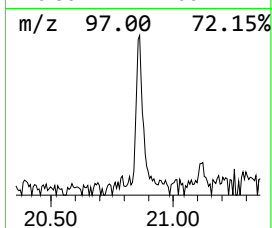
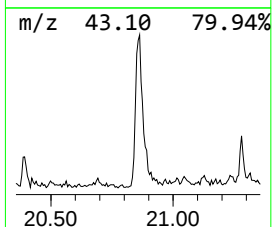
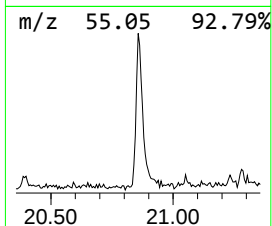
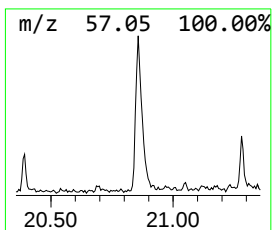
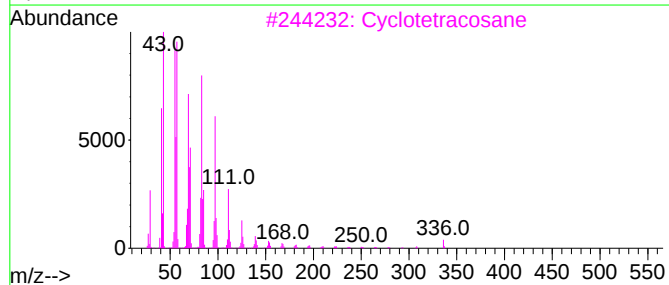
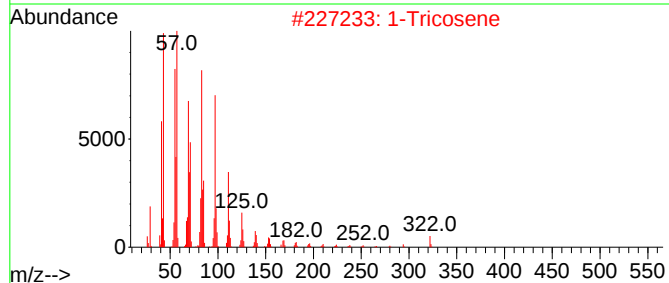
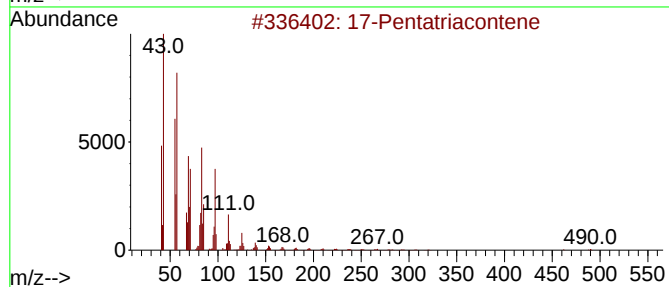
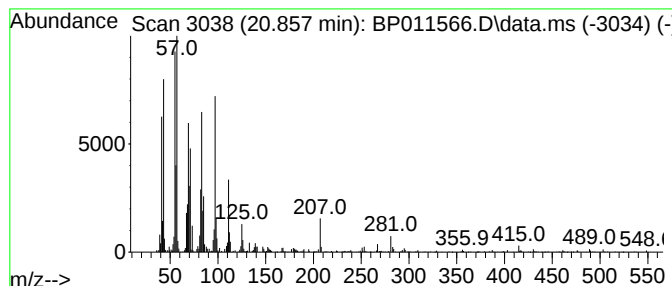
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.857	3.73 ng/ul	169780	Chrysene-d12	21.121	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Pentatriacontene	491	C35H70	006971-40-0	95
2	1-Tricosene	322	C23H46	018835-32-0	91
3	Cyclotetracosane	336	C24H48	000297-03-0	91
4	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
5	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
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
Tetrachloroethy...	4.287	129.4	ng/u1	2806630	1	7.546	433794	20.0
(DEL) Alkane: S...	20.857	3.7	ng/u1	169780	5	21.121	910105	20.0