

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082422\
 Data File : BP011570.D
 Acq On : 25 Aug 2022 19:37
 Operator : CG/JU
 Sample : N4349-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGN82

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: BP011570.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	10	rBV	85974	110479	6.34%	0.629%
2	3.064	10	13	16	rVV	23401	28286	1.62%	0.161%
3	3.099	16	19	23	rVB	10952	13370	0.77%	0.076%
4	3.481	81	84	97	rBV	68002	112115	6.43%	0.638%
5	3.687	114	119	126	rVB2	12967	23238	1.33%	0.132%
6	3.775	131	134	139	rVB2	7228	9206	0.53%	0.052%
7	4.158	195	199	200	rBV4	1884	2414	0.14%	0.014%
8	4.205	204	207	212	rVB3	4410	6111	0.35%	0.035%
9	4.287	215	221	235	rVB	733017	994541	57.06%	5.659%
10	5.964	503	506	510	rVB5	1543	2338	0.13%	0.013%
11	6.734	630	637	646	rBV	69419	113346	6.50%	0.645%
12	6.899	658	665	674	rBV	225980	350056	20.08%	1.992%
13	7.081	689	696	706	rBV	244241	384758	22.08%	2.189%
14	7.158	707	709	715	rVB5	2435	3448	0.20%	0.020%
15	7.546	768	775	785	rBV	241860	376759	21.62%	2.144%
16	8.269	890	898	909	rBV	188110	311705	17.88%	1.774%
17	8.710	966	973	980	rBV	310001	490804	28.16%	2.793%
18	8.852	992	997	999	rBV6	1565	2899	0.17%	0.016%
19	9.105	1035	1040	1045	rVB6	2077	4417	0.25%	0.025%
20	9.428	1084	1095	1104	rBV	242022	395409	22.69%	2.250%
21	9.957	1178	1185	1200	rBV	313137	540776	31.03%	3.077%
22	10.104	1208	1210	1217	rVB8	1217	1842	0.11%	0.010%
23	10.334	1241	1249	1256	rBV	322346	527895	30.29%	3.004%
24	10.487	1267	1275	1288	rBV	241694	431167	24.74%	2.453%
25	11.616	1461	1467	1470	rBV7	2167	3248	0.19%	0.018%
26	11.940	1518	1522	1530	rVB4	4550	8838	0.51%	0.050%
27	12.575	1625	1630	1632	rBV6	1138	1930	0.11%	0.011%
28	12.822	1666	1672	1674	rBV7	1593	2370	0.14%	0.013%
29	12.934	1686	1691	1696	rBV8	1432	2531	0.15%	0.014%
30	13.122	1720	1723	1728	rBV7	1056	1916	0.11%	0.011%
31	13.645	1805	1812	1821	rBV	666804	918390	52.69%	5.226%
32	13.904	1849	1856	1867	rBV	690006	991914	56.91%	5.644%
33	14.216	1899	1909	1918	rBV	478348	674373	38.69%	3.837%
34	14.281	1918	1920	1925	rVB6	1385	2072	0.12%	0.012%
35	14.457	1944	1950	1966	rBV2	31894	82559	4.74%	0.470%
36	14.663	1982	1985	1990	rVB6	1887	2960	0.17%	0.017%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
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37	15.222	2071	2080	2096	rBV	923670	1274358	73.12%	7.252%
38	15.357	2096	2103	2116	rBV	371865	506219	29.04%	2.881%
39	15.463	2116	2121	2126	rVB7	4279	8763	0.50%	0.050%
40	15.957	2202	2205	2209	rBV5	3750	5893	0.34%	0.034%
41	16.016	2213	2215	2219	rVB5	2170	1812	0.10%	0.010%
42	16.469	2289	2292	2295	rBV4	1815	2545	0.15%	0.014%
43	16.663	2317	2325	2329	rBV4	1746	3593	0.21%	0.020%
44	16.763	2339	2342	2346	rBV4	2636	3636	0.21%	0.021%
45	16.986	2374	2380	2390	rBV	572927	779217	44.71%	4.434%
46	17.086	2391	2397	2412	rBV	1175764	1627130	93.36%	9.259%
47	17.298	2428	2433	2435	rBV6	2792	3748	0.22%	0.021%
48	17.439	2456	2457	2461	rVB4	1731	1782	0.10%	0.010%
49	17.686	2496	2499	2503	rVB5	2885	2939	0.17%	0.017%
50	17.910	2532	2537	2543	rBV	31141	48465	2.78%	0.276%
51	17.975	2546	2548	2554	rVB2	4259	6872	0.39%	0.039%
52	18.928	2706	2710	2714	rBV2	12190	16237	0.93%	0.092%
53	18.992	2717	2721	2725	rBV2	13624	18996	1.09%	0.108%
54	19.157	2746	2749	2753	rBV5	11001	14246	0.82%	0.081%
55	19.351	2776	2782	2792	rBV	1408809	1742896	100.00%	9.918%
56	20.857	3033	3038	3047	rBV2	94496	205344	11.78%	1.168%
57	21.116	3077	3082	3089	rBV	680412	829846	47.61%	4.722%
58	23.257	3439	3446	3453	rBV2	943787	1735691	99.59%	9.877%
59	23.404	3465	3471	3480	rVB2	419895	802843	46.06%	4.568%

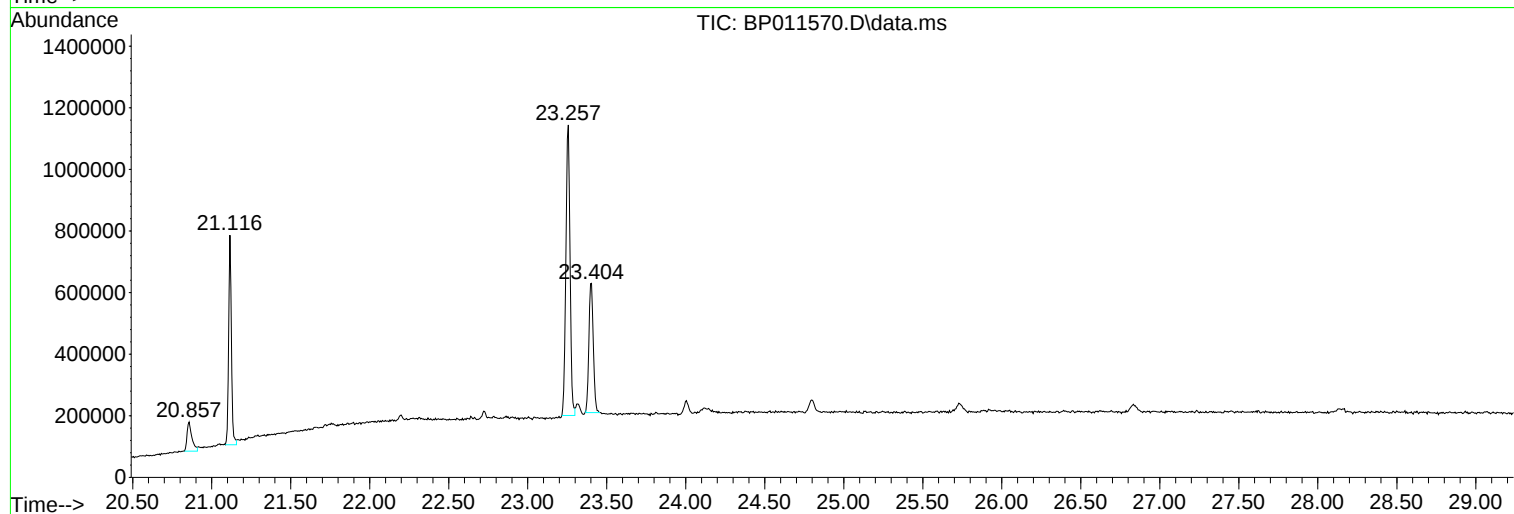
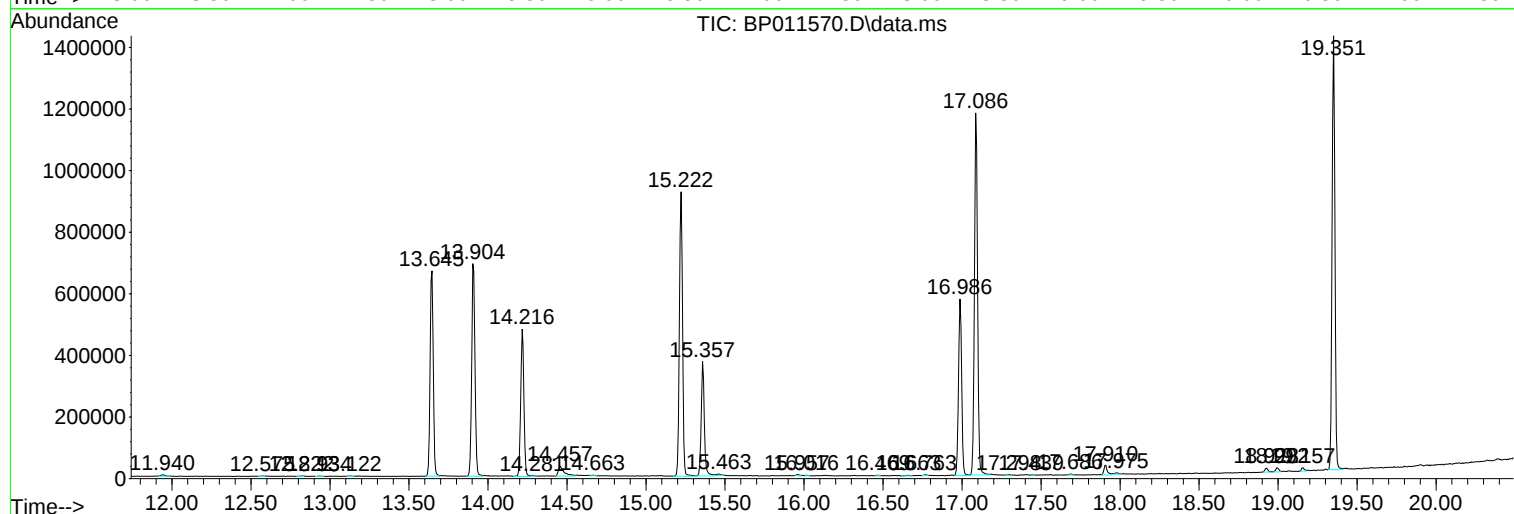
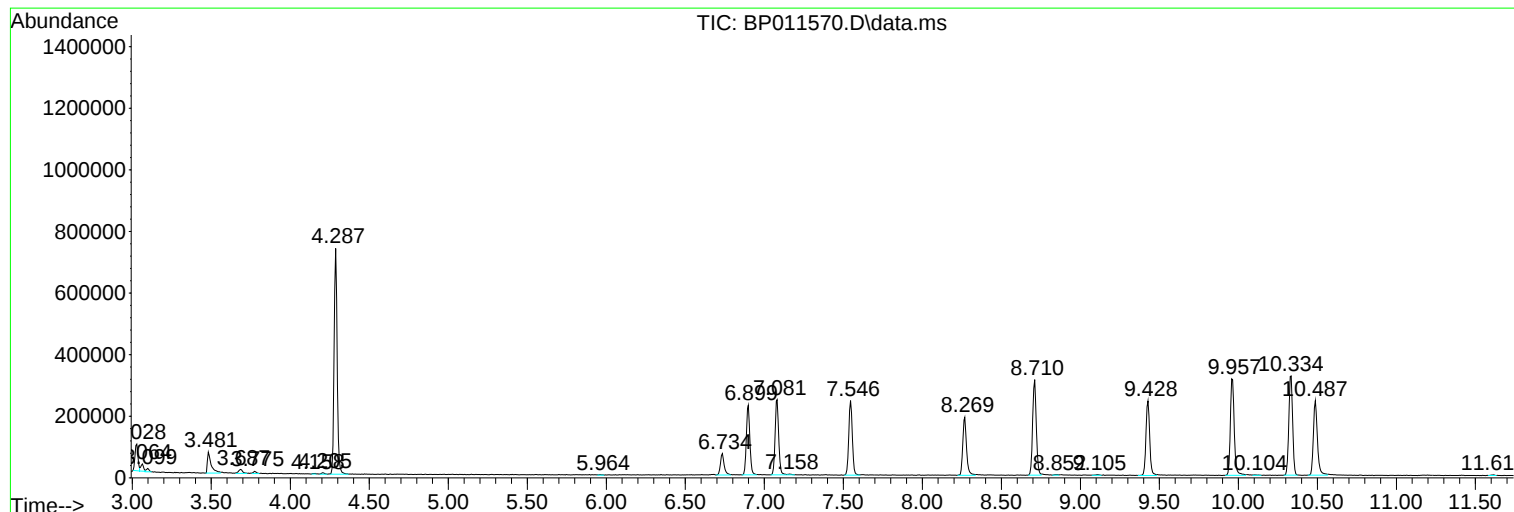
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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



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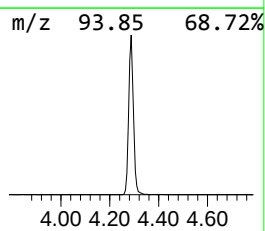
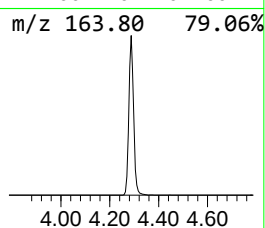
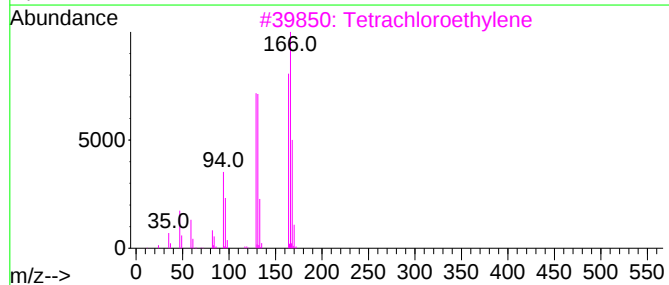
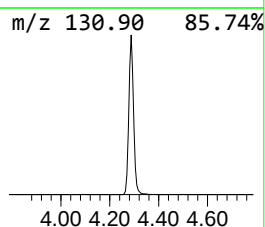
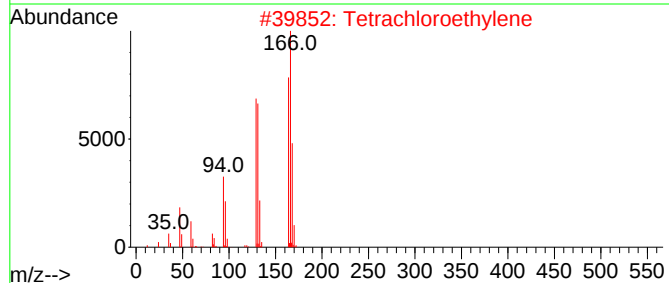
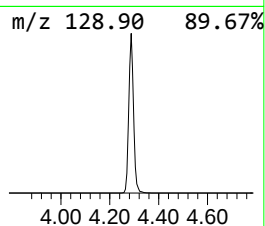
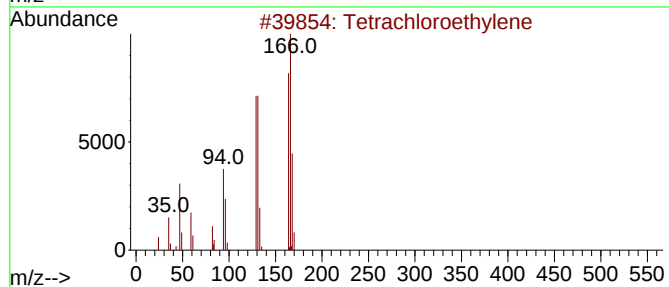
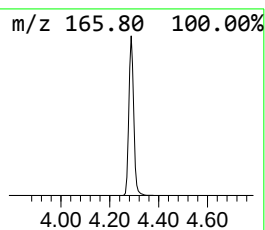
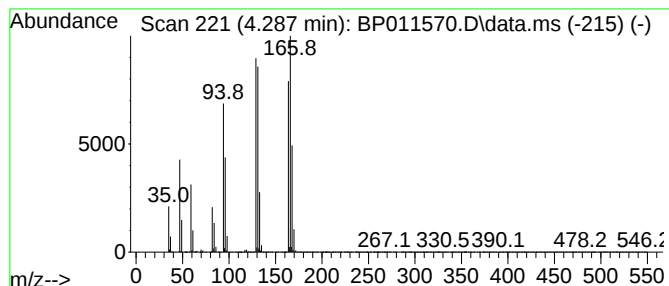
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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	52.79 ng/ul	994541	1,4-Dichlorobenzene-d4	7.546

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	95
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	27



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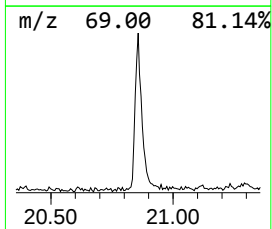
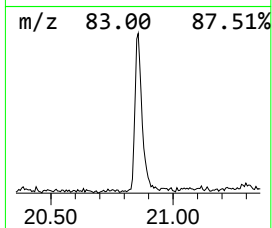
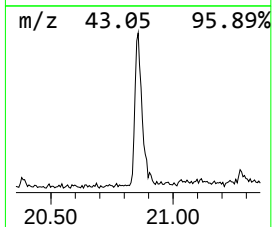
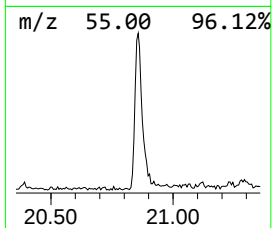
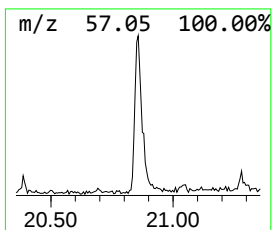
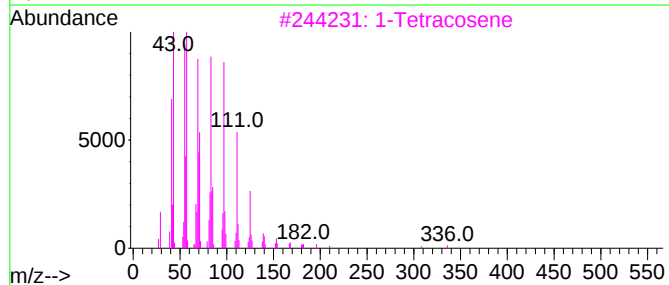
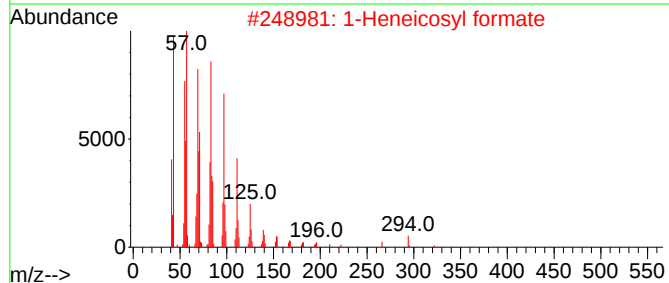
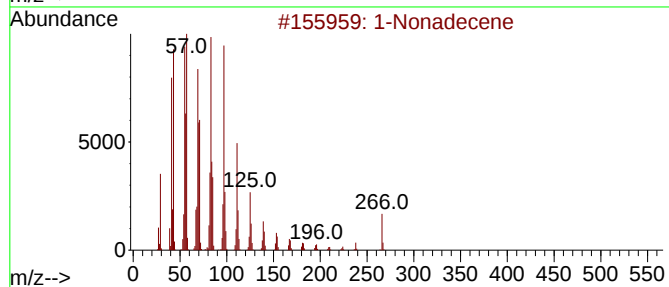
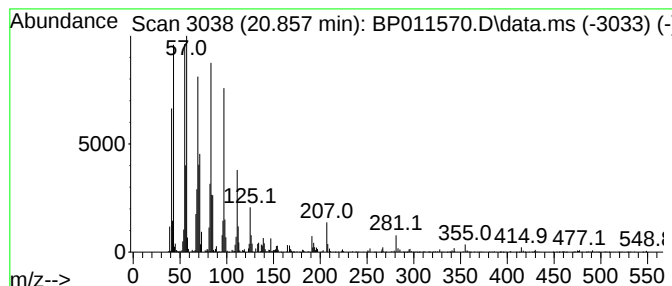
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.857	4.95 ng/ul	205344	Chrysene-d12	21.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	98
2	1-Heneicosyl formate			340	C22H44O2	077899-03-7	96
3	1-Tetracosene			336	C24H48	010192-32-2	95
4	Behenic alcohol			326	C22H46O	000661-19-8	95
5	1-Nonadecene			266	C19H38	018435-45-5	94



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.287	52.8	ng/u1	994541	1	7.546	376759	20.0
(DEL) Alkane: S...	20.857	5.0	ng/u1	205344	5	21.116	829846	20.0