

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072722\
 Data File : BP011148.D
 Acq On : 28 Jul 2022 02:20
 Operator : CG/JU
 Sample : N3871-16
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJ15

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : 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-BP071422.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011148.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	10	rBV	307411	387550	4.45%	0.442%
2	3.570	80	82	94	rBV	522031	757146	8.69%	0.863%
3	4.393	218	222	228	rVB	430403	593144	6.81%	0.676%
4	6.834	632	637	650	rVB	365871	582291	6.68%	0.664%
5	6.999	659	665	677	rVB	1637375	2538128	29.13%	2.893%
6	7.187	691	697	710	rBV	1497021	2314295	26.56%	2.638%
7	7.658	770	777	785	rBV	1522059	2338234	26.84%	2.665%
8	8.369	892	898	913	rBV	1019815	1662649	19.08%	1.895%
9	8.816	967	974	984	rBV	1981124	3183892	36.54%	3.629%
10	9.234	1040	1045	1051	rVB	58753	94270	1.08%	0.107%
11	9.540	1089	1097	1113	rBV	1398241	2341567	26.87%	2.669%
12	10.069	1180	1187	1202	rBV	1889451	3133589	35.96%	3.572%
13	10.257	1213	1219	1225	rBV3	23749	53147	0.61%	0.061%
14	10.446	1244	1251	1261	rVB	1984226	3271140	37.54%	3.729%
15	10.593	1269	1276	1294	rBV	1672469	2980519	34.21%	3.397%
16	11.010	1343	1347	1353	rBV5	7290	15615	0.18%	0.018%
17	11.075	1354	1358	1361	rVV4	8186	17534	0.20%	0.020%
18	11.116	1361	1365	1373	rVB2	20608	41613	0.48%	0.047%
19	11.246	1383	1387	1389	rBV3	14127	19393	0.22%	0.022%
20	11.734	1463	1470	1477	rBV4	20531	40044	0.46%	0.046%
21	12.046	1518	1523	1528	rBV2	32360	52822	0.61%	0.060%
22	12.675	1624	1630	1634	rBV5	14720	29371	0.34%	0.033%
23	12.928	1667	1673	1678	rVB3	19824	31022	0.36%	0.035%
24	13.034	1687	1691	1696	rBV7	5667	9969	0.11%	0.011%
25	13.222	1720	1723	1729	rVV4	11049	15445	0.18%	0.018%
26	13.287	1729	1734	1742	rVB7	6958	14373	0.16%	0.016%
27	13.646	1790	1795	1799	rVB6	5406	10903	0.13%	0.012%
28	13.740	1804	1811	1823	rVV	3885987	5193503	59.61%	5.920%
29	13.846	1823	1829	1839	rVB6	17641	47044	0.54%	0.054%
30	14.010	1847	1857	1869	rBV	4234309	6007840	68.95%	6.848%
31	14.128	1875	1877	1883	rVB7	6295	9295	0.11%	0.011%
32	14.234	1891	1895	1901	rVB7	5656	9753	0.11%	0.011%
33	14.316	1901	1909	1921	rBV	2478696	3633608	41.70%	4.142%
34	14.540	1941	1947	1966	rBV	120413	275644	3.16%	0.314%
35	14.757	1981	1984	1990	rVB6	10856	17445	0.20%	0.020%
36	15.134	2043	2048	2056	rBV2	45257	75563	0.87%	0.086%

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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-BP071422.M
 Title : SVOA CALIBRATION

37	15.322	2072	2080	2086	rBV	5191475	7233437	83.02%	8.245%
38	15.445	2094	2101	2113	rBV	1456676	1965136	22.55%	2.240%
39	15.557	2116	2120	2131	rVB	59506	95365	1.09%	0.109%
40	15.816	2160	2164	2170	rVB9	4843	8996	0.10%	0.010%
41	16.022	2192	2199	2203	rBV9	17439	39944	0.46%	0.046%
42	16.057	2203	2205	2209	rVV4	12487	16299	0.19%	0.019%
43	16.110	2210	2214	2218	rVB4	18193	29514	0.34%	0.034%
44	16.234	2230	2235	2239	rBV7	6883	11485	0.13%	0.013%
45	16.569	2288	2292	2298	rVB7	6608	12810	0.15%	0.015%
46	16.757	2320	2324	2330	rVB9	10628	20090	0.23%	0.023%
47	16.863	2338	2342	2346	rBV5	16898	26196	0.30%	0.030%
48	16.904	2346	2349	2355	rVB7	9369	15475	0.18%	0.018%
49	17.022	2366	2369	2373	rVB6	8826	11102	0.13%	0.013%
50	17.087	2373	2380	2390	rBV	2624562	3640903	41.79%	4.150%
51	17.187	2390	2397	2410	rBV	5392654	7411887	85.07%	8.449%
52	17.387	2428	2431	2437	rVB	18078	24965	0.29%	0.028%
53	17.716	2484	2487	2492	rVV5	14516	14674	0.17%	0.017%
54	17.781	2493	2498	2508	rVB	57252	75512	0.87%	0.086%
55	17.969	2523	2530	2531	rBV7	6210	10814	0.12%	0.012%
56	17.998	2531	2535	2540	rVV	66885	97499	1.12%	0.111%
57	18.069	2544	2547	2551	rVB	23888	30827	0.35%	0.035%
58	19.016	2704	2708	2712	rBV2	62241	80755	0.93%	0.092%
59	19.081	2716	2719	2723	rVV	53906	71037	0.82%	0.081%
60	19.239	2744	2746	2750	rBV4	13208	14678	0.17%	0.017%
61	19.392	2768	2772	2774	rBV3	23452	30358	0.35%	0.035%
62	19.439	2774	2780	2788	rBV	6453728	8122945	93.23%	9.259%
63	20.939	3031	3035	3042	rBV2	160808	303557	3.48%	0.346%
64	21.204	3075	3080	3086	rBV	3034679	3757125	43.12%	4.283%
65	23.392	3445	3452	3459	rBV2	4577269	8712911	100.00%	9.932%
66	23.539	3471	3477	3485	rVB2	2131431	4044178	46.42%	4.610%

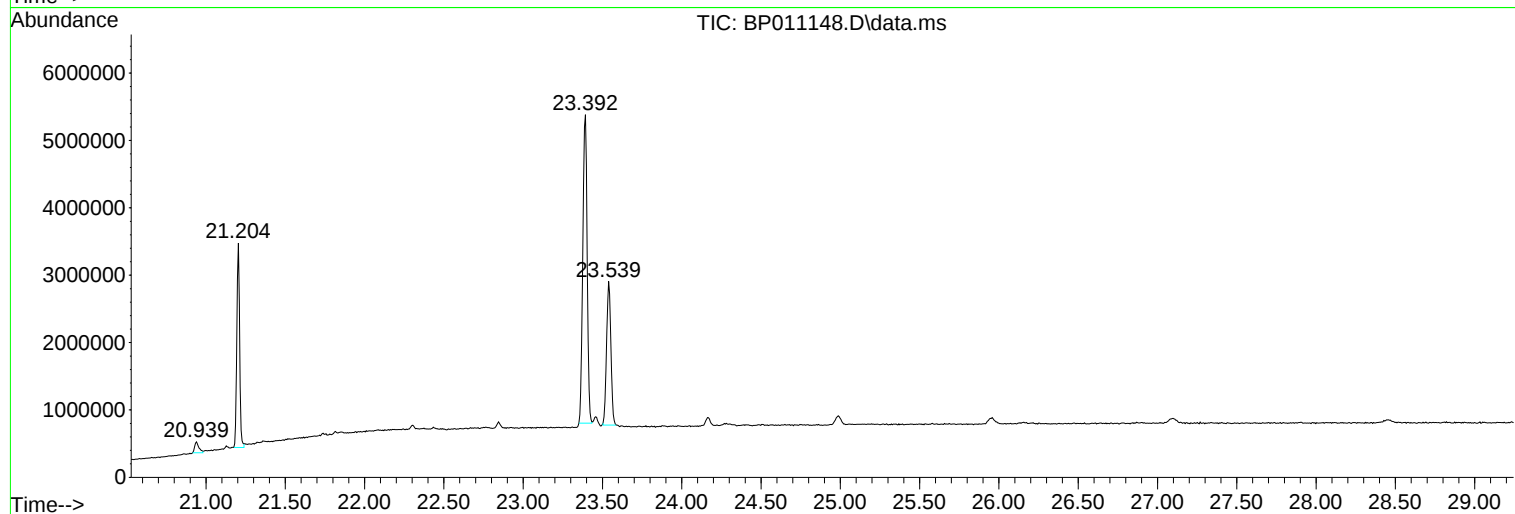
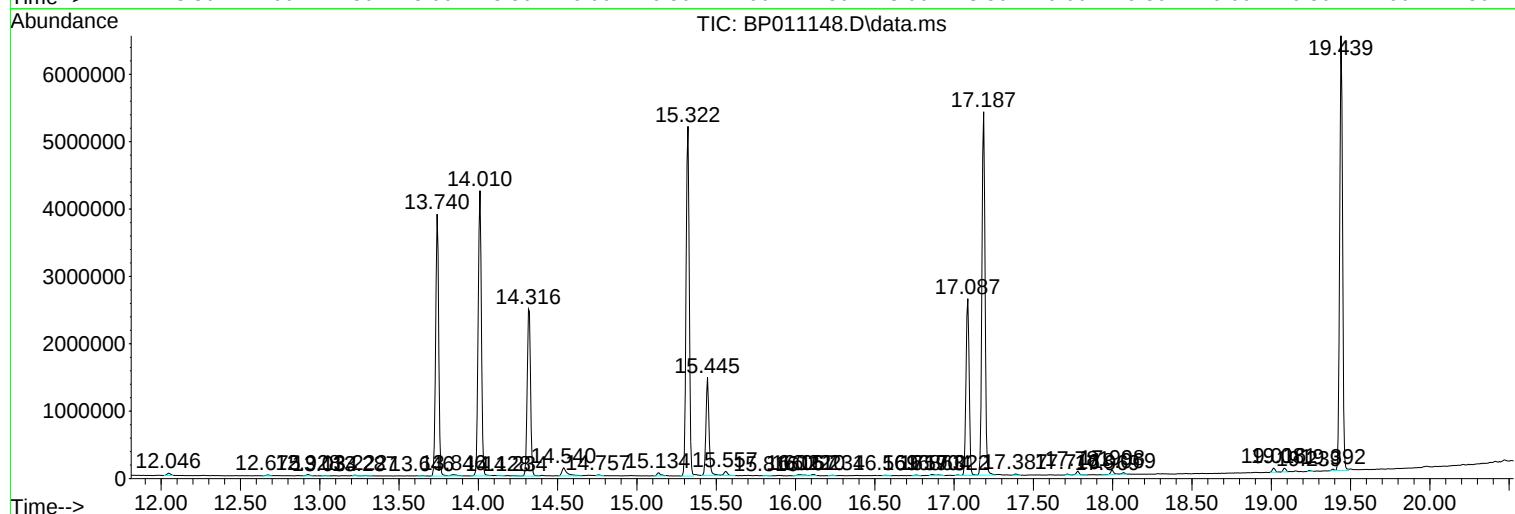
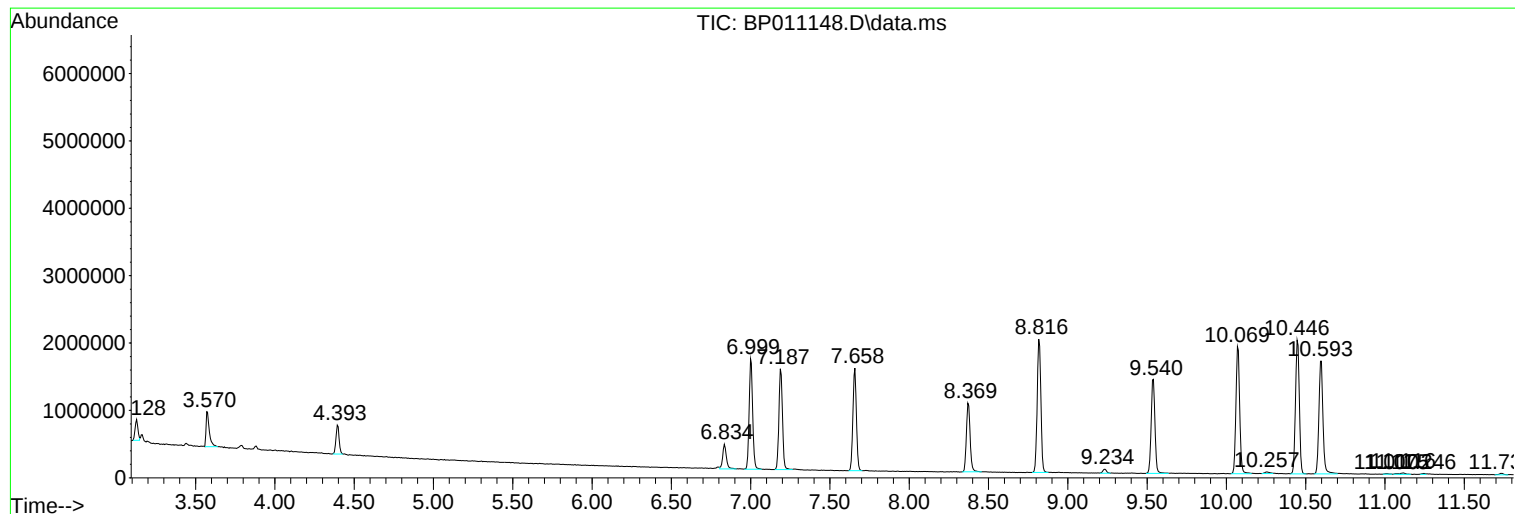
Sum of corrected areas: 87729834

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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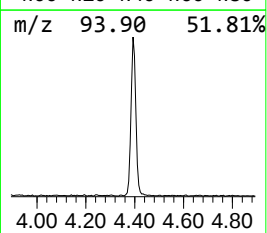
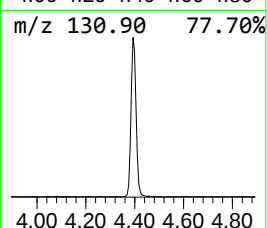
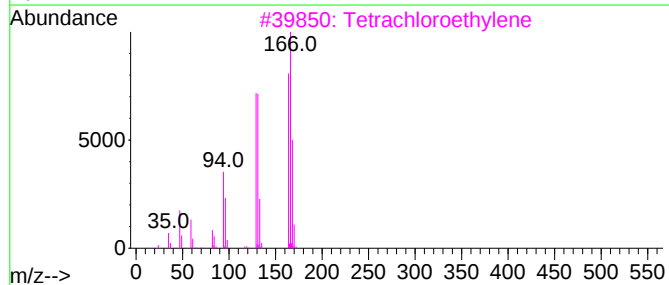
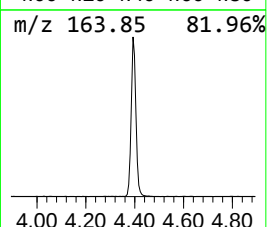
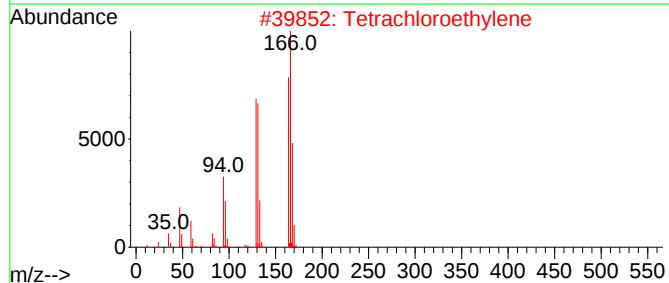
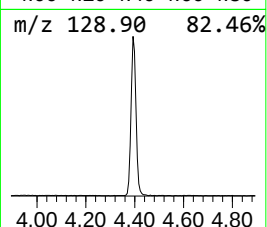
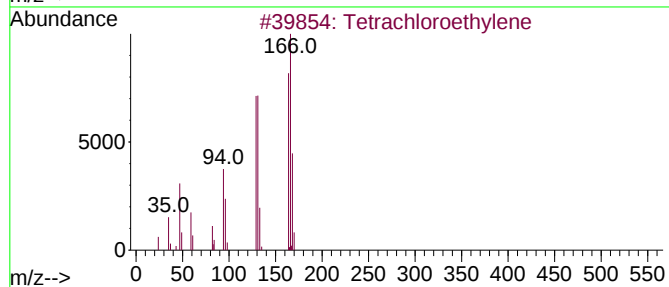
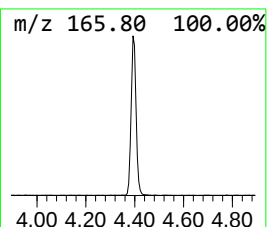
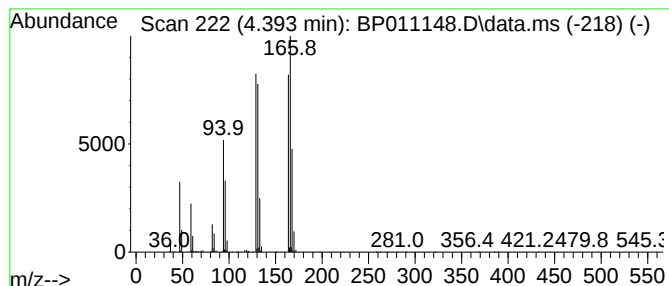
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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.
4.393	5.07 ng/ul	593144	1,4-Dichlorobenzene-d4	7.658

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HCl2FN2	002927-71-1	25



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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...	4.393	5.1	ng/u1	593144	1	7.658	2338230	20.0