

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013065.D
 Acq On : 11 Dec 2022 06:05
 Operator : CG/JU
 Sample : N5927-08
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXZ96

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: BP013065.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.317	17	22	26	rBV	645382	952907	4.14%	0.479%
2	3.781	98	101	122	rBV	591139	1036570	4.50%	0.521%
3	4.011	136	140	146	rVB2	53690	88091	0.38%	0.044%
4	4.111	153	157	163	rVB	33215	53190	0.23%	0.027%
5	4.217	170	175	182	rBV2	78379	114660	0.50%	0.058%
6	4.646	242	248	262	rVB	2447512	3963428	17.22%	1.994%
7	7.122	662	669	681	rBV	616179	986932	4.29%	0.496%
8	7.293	691	698	707	rBV	2439105	3808758	16.55%	1.916%
9	7.363	707	710	719	rVB	13787	25778	0.11%	0.013%
10	7.499	726	733	741	rBV	2196316	3551158	15.43%	1.786%
11	7.969	804	813	820	rBV	2058828	3230263	14.04%	1.625%
12	8.028	820	823	828	rVB3	29300	44532	0.19%	0.022%
13	8.675	924	933	942	rBV	1826261	2908814	12.64%	1.463%
14	9.134	1002	1011	1023	rBV	2962297	4770661	20.73%	2.400%
15	9.299	1035	1039	1044	rVB5	16715	25142	0.11%	0.013%
16	9.540	1075	1080	1086	rVB	29105	50823	0.22%	0.026%
17	9.863	1126	1135	1149	rBV	2276076	3811712	16.56%	1.917%
18	10.399	1218	1226	1247	rBV	3470254	5719158	24.85%	2.877%
19	10.557	1247	1253	1260	rBV2	49094	88908	0.39%	0.045%
20	10.781	1284	1291	1299	rBV	3038334	5019305	21.81%	2.525%
21	10.916	1307	1314	1332	rBV	2924724	5124287	22.27%	2.578%
22	11.316	1376	1382	1387	rBV	162870	265909	1.16%	0.134%
23	11.375	1387	1392	1397	rVV	119441	184625	0.80%	0.093%
24	11.557	1420	1423	1429	rBV6	16853	32878	0.14%	0.017%
25	12.363	1554	1560	1565	rVB2	68994	108895	0.47%	0.055%
26	13.210	1700	1704	1713	rVB6	17401	33522	0.15%	0.017%
27	13.428	1737	1741	1748	rVB3	13969	26308	0.11%	0.013%
28	13.498	1748	1753	1757	rVV2	40216	63136	0.27%	0.032%
29	13.575	1761	1766	1771	rVB6	16492	33437	0.15%	0.017%
30	14.016	1834	1841	1855	rBV	7873749	10672265	46.37%	5.369%
31	14.128	1855	1860	1869	rVB	17868	35529	0.15%	0.018%
32	14.245	1875	1880	1883	rBV5	66405	107501	0.47%	0.054%
33	14.304	1883	1890	1900	rVV	8114706	12223155	53.11%	6.149%
34	14.381	1901	1903	1908	rVB6	17614	26102	0.11%	0.013%
35	14.492	1918	1922	1928	rVB3	21743	34170	0.15%	0.017%
36	14.610	1935	1942	1951	rVB	4904700	7308260	31.76%	3.676%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	14.704	1953	1958	1965	rVB9	16129	32473	0.14%	0.016%
38	14.798	1967	1974	1993	rBV	501898	868488	3.77%	0.437%
39	15.016	2006	2011	2016	rVB8	32348	58237	0.25%	0.029%
40	15.239	2044	2049	2057	rVB	44161	68300	0.30%	0.034%
41	15.398	2072	2076	2078	rVV	52600	70617	0.31%	0.036%
42	15.445	2082	2084	2089	rVB2	33152	38088	0.17%	0.019%
43	15.604	2103	2111	2123	rBV	11180289	16349455	71.04%	8.225%
44	15.716	2124	2130	2142	rVV	3966419	5269487	22.90%	2.651%
45	15.839	2147	2151	2158	rVB2	61407	105473	0.46%	0.053%
46	15.957	2167	2171	2175	rBV3	29675	38138	0.17%	0.019%
47	16.004	2175	2179	2183	rVB4	29101	41451	0.18%	0.021%
48	16.057	2183	2188	2194	rBV7	19864	33490	0.15%	0.017%
49	16.145	2198	2203	2205	rBV6	22518	34716	0.15%	0.017%
50	16.239	2215	2219	2222	rBV4	28047	35767	0.16%	0.018%
51	16.275	2222	2225	2228	rVV2	43661	61035	0.27%	0.031%
52	16.310	2228	2231	2239	rVV4	48002	95166	0.41%	0.048%
53	16.381	2240	2243	2250	rVB7	33813	52858	0.23%	0.027%
54	16.492	2258	2262	2266	rBV4	40148	54401	0.24%	0.027%
55	16.581	2275	2277	2284	rVB5	13499	23959	0.10%	0.012%
56	16.657	2286	2290	2293	rBV6	19037	30615	0.13%	0.015%
57	16.834	2313	2320	2325	rBV9	32033	62563	0.27%	0.031%
58	17.104	2361	2366	2370	rBV5	41414	68142	0.30%	0.034%
59	17.157	2370	2375	2381	rVB5	41481	77458	0.34%	0.039%
60	17.269	2390	2394	2400	rVB	104869	129828	0.56%	0.065%
61	17.363	2403	2410	2417	rBV	6604638	9278017	40.32%	4.667%
62	17.463	2421	2427	2439	rVB2	13013211	18828287	81.81%	9.471%
63	17.645	2454	2458	2469	rBV	111889	210494	0.91%	0.106%
64	17.739	2471	2474	2477	rBV4	33046	42721	0.19%	0.021%
65	18.016	2517	2521	2527	rVB2	136424	190696	0.83%	0.096%
66	18.228	2553	2557	2564	rBV	249436	321845	1.40%	0.162%
67	18.304	2567	2570	2575	rVB	189338	224515	0.98%	0.113%
68	18.516	2603	2606	2610	rBV4	29803	43732	0.19%	0.022%
69	18.645	2624	2628	2632	rBV2	72776	86480	0.38%	0.044%
70	18.810	2654	2656	2659	rVB3	39886	38213	0.17%	0.019%
71	18.969	2680	2683	2687	rVB	73247	91987	0.40%	0.046%
72	19.057	2694	2698	2702	rBV	111562	134770	0.59%	0.068%
73	19.269	2730	2734	2738	rBV	211276	258043	1.12%	0.130%
74	19.339	2742	2746	2749	rBV	181606	257216	1.12%	0.129%
75	19.463	2763	2767	2770	rBV3	111919	145364	0.63%	0.073%
76	19.692	2800	2806	2818	rVB	16782246	22952288	99.73%	11.546%

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Title : SVOA CALIBRATION

77	21.139	3048	3052	3063	rBV2	565586	910890	3.96%	0.458%
78	21.451	3099	3105	3111	rBV	7259583	10192237	44.29%	5.127%
79	23.151	3390	3394	3401	rVB	316474	499702	2.17%	0.251%
80	23.774	3492	3500	3519	rVB2	10865434	23013798	100.00%	11.577%
81	23.933	3520	3527	3540	rVB2	4651864	9897797	43.01%	4.979%
82	24.557	3629	3633	3641	rVB	483482	939404	4.08%	0.473%

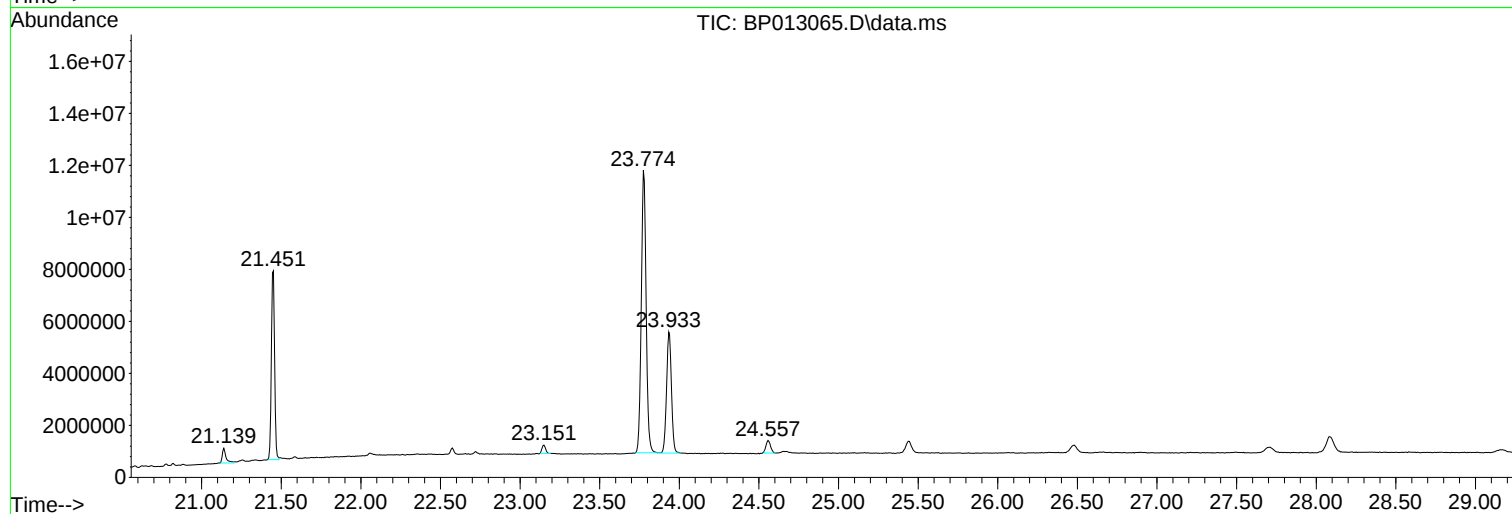
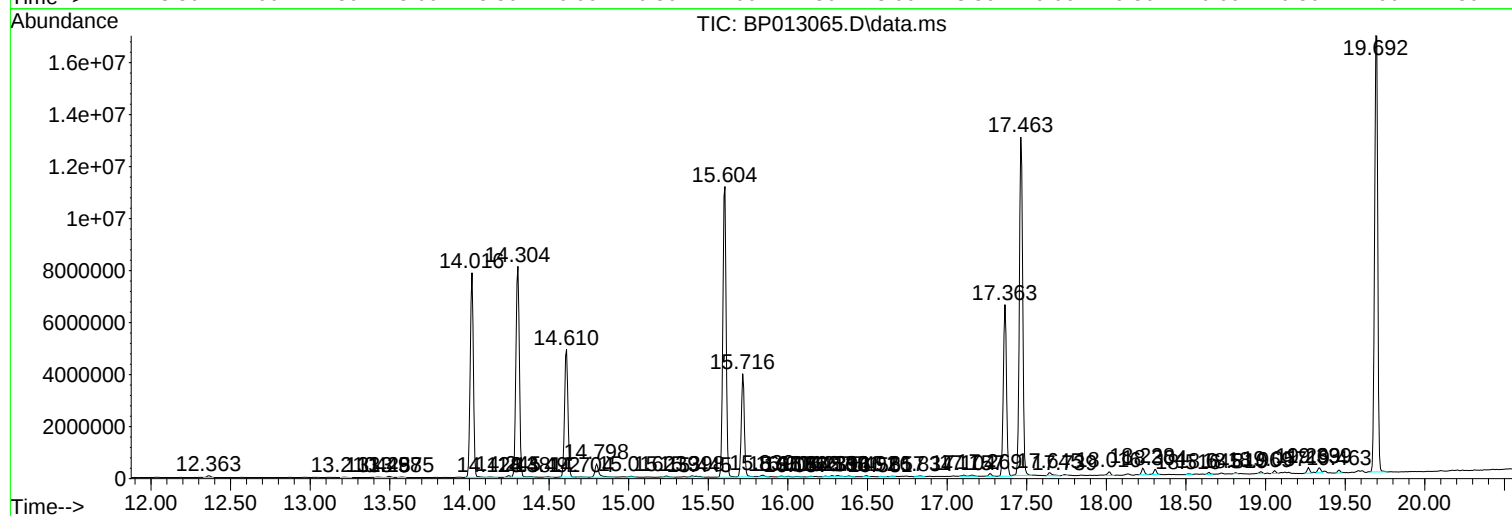
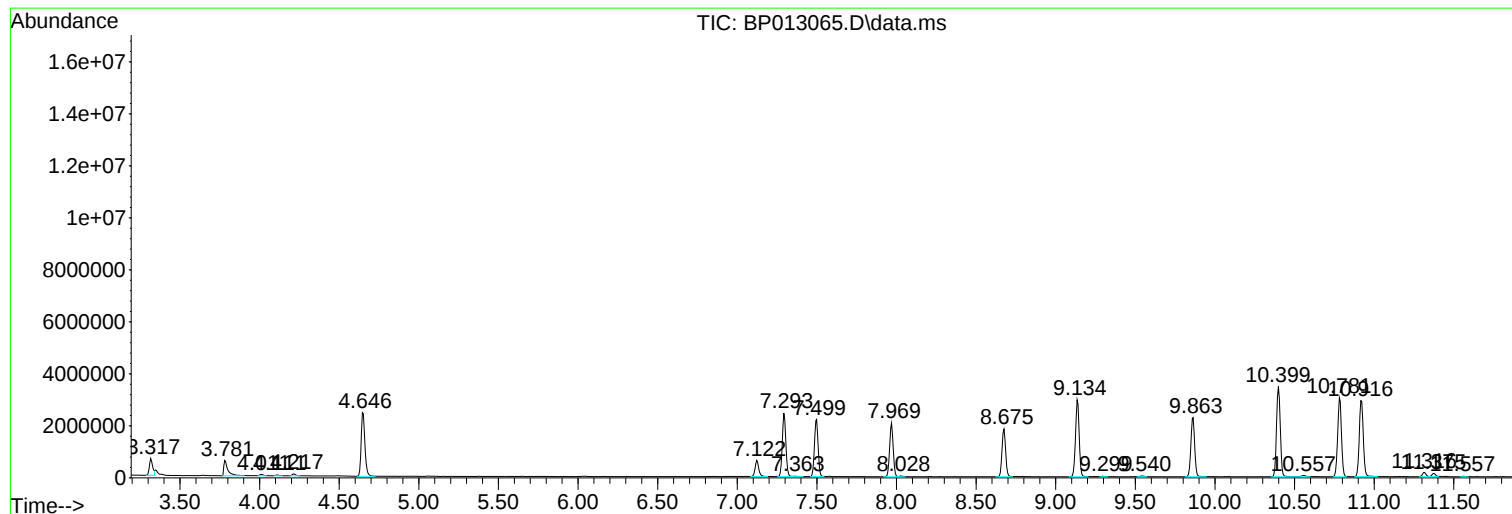
Sum of corrected areas: 198789470

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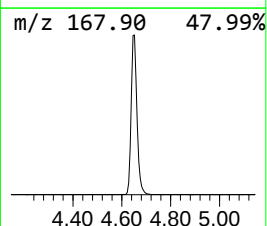
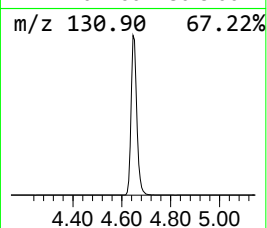
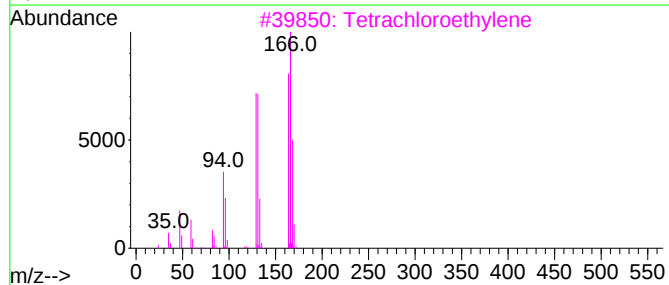
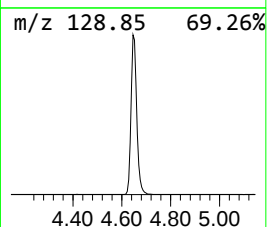
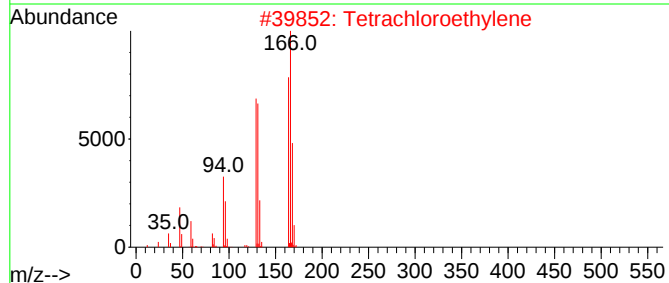
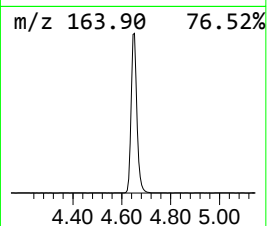
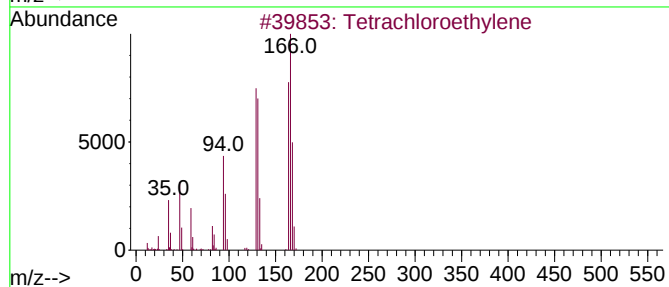
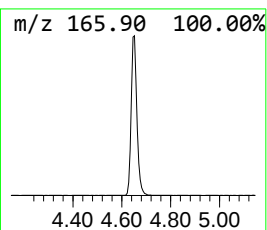
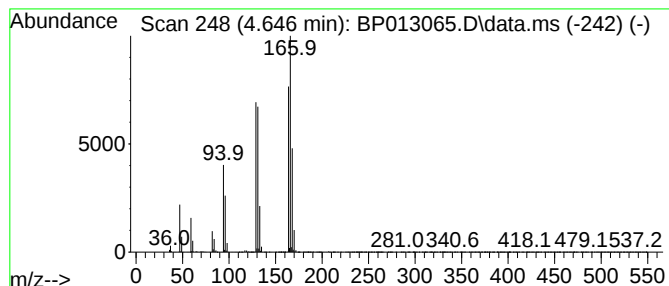
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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.
4.646	24.54 ng/ul	3963430	1,4-Dichlorobenzene-d4	7.969

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	97
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HCl2FN2	002927-71-1	38



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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.646	24.5	ng/u1	3963430	1	7.969	3230260	20.0