

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017575.D
 Acq On : 05 Oct 2023 20:30
 Operator : MA/JU
 Sample : 04588-22
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJZ9

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-BP100423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017575.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.240	23	26	29	rVB5	9811	10601	0.25%	0.033%
2	3.287	31	34	39	rVV	20544	25923	0.61%	0.081%
3	3.352	43	45	49	rVB2	9149	9426	0.22%	0.029%
4	3.517	70	73	77	rVV3	22963	26153	0.62%	0.081%
5	3.552	77	79	85	rVB6	5497	6748	0.16%	0.021%
6	3.728	106	109	117	rBV	98402	144251	3.40%	0.448%
7	4.046	161	163	170	rVB7	15466	23533	0.55%	0.073%
8	4.481	233	237	239	rBV5	8819	12815	0.30%	0.040%
9	4.505	239	241	244	rVV4	8534	9931	0.23%	0.031%
10	4.552	244	249	262	rVB	149555	223532	5.26%	0.695%
11	4.781	284	288	292	rBV7	9686	14343	0.34%	0.045%
12	4.864	300	302	312	rVB7	3821	9047	0.21%	0.028%
13	5.069	334	337	347	rBV7	10998	25845	0.61%	0.080%
14	5.163	350	353	358	rVB7	5102	6688	0.16%	0.021%
15	5.216	358	362	369	rVB7	13917	21798	0.51%	0.068%
16	5.334	377	382	383	rBV5	2724	4701	0.11%	0.015%
17	5.740	447	451	454	rBV6	4082	6363	0.15%	0.020%
18	5.793	457	460	469	rVB5	9679	18334	0.43%	0.057%
19	5.946	481	486	489	rBV6	5058	7313	0.17%	0.023%
20	5.981	489	492	495	rVB4	4238	5704	0.13%	0.018%
21	6.052	499	504	506	rBV6	4613	7751	0.18%	0.024%
22	6.363	553	557	562	rBV4	13939	21966	0.52%	0.068%
23	6.822	631	635	641	rBV9	5671	11528	0.27%	0.036%
24	6.869	641	643	648	rVB6	3735	4495	0.11%	0.014%
25	7.099	676	682	695	rVB2	103954	216888	5.10%	0.674%
26	7.281	705	713	722	rBV	492601	765267	18.01%	2.379%
27	7.457	735	743	753	rBV	417222	667045	15.70%	2.074%
28	7.552	755	759	763	rVB6	4226	6033	0.14%	0.019%
29	7.934	815	824	835	rBV	339623	548151	12.90%	1.704%
30	8.199	866	869	874	rBV5	3404	5514	0.13%	0.017%
31	8.293	879	885	889	rBV8	4965	9287	0.22%	0.029%
32	8.593	933	936	941	rVB7	8881	14731	0.35%	0.046%
33	8.657	941	947	957	rVV	287567	478283	11.26%	1.487%
34	9.022	1003	1009	1011	rBV7	2794	4271	0.10%	0.013%
35	9.104	1017	1023	1025	rBV	60958	93365	2.20%	0.290%
36	9.146	1025	1030	1042	rVB	521381	854264	20.11%	2.655%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
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37	9.334	1060	1062	1065	rBV4	4450	5480	0.13%	0.017%
38	9.722	1120	1128	1131	rBV10	4530	9805	0.23%	0.030%
39	9.863	1145	1152	1166	rBV	418741	719261	16.93%	2.236%
40	10.004	1168	1176	1181	rVB9	3387	7065	0.17%	0.022%
41	10.134	1193	1198	1203	rBV3	12011	18436	0.43%	0.057%
42	10.387	1231	1241	1256	rBV	563986	1039338	24.46%	3.231%
43	10.775	1298	1307	1315	rBV	454237	773592	18.21%	2.405%
44	10.846	1315	1319	1325	rVV6	4759	9379	0.22%	0.029%
45	10.946	1328	1336	1349	rBV	490524	890997	20.97%	2.770%
46	11.845	1485	1489	1493	rVB7	3827	6911	0.16%	0.021%
47	12.363	1574	1577	1582	rVB3	9610	15382	0.36%	0.048%
48	12.687	1629	1632	1634	rBV4	3189	4418	0.10%	0.014%
49	13.369	1745	1748	1753	rVB7	3380	4944	0.12%	0.015%
50	13.463	1760	1764	1765	rBV3	4306	4633	0.11%	0.014%
51	13.845	1825	1829	1835	rVB7	5233	8204	0.19%	0.026%
52	13.957	1842	1848	1851	rBV7	2483	4513	0.11%	0.014%
53	14.045	1856	1863	1873	rBV	1273223	1747671	41.14%	5.433%
54	14.239	1890	1896	1903	rBV	3698	9172	0.22%	0.029%
55	14.328	1904	1911	1922	rBV	1456858	2006687	47.23%	6.238%
56	14.628	1955	1962	1969	rBV	734084	1032760	24.31%	3.210%
57	14.887	1998	2006	2021	rBV4	49490	133558	3.14%	0.415%
58	15.016	2027	2028	2035	rVB7	4636	6707	0.16%	0.021%
59	15.245	2063	2067	2074	rVB10	2631	5932	0.14%	0.018%
60	15.422	2093	2097	2099	rBV5	3973	4784	0.11%	0.015%
61	15.634	2126	2133	2147	rVV	1881456	2654423	62.48%	8.251%
62	15.775	2151	2157	2170	rBV	545601	754912	17.77%	2.347%
63	15.875	2170	2174	2178	rVB6	10039	15145	0.36%	0.047%
64	16.428	2264	2268	2271	rVB6	3921	4749	0.11%	0.015%
65	17.416	2430	2436	2443	rBV	970570	1326919	31.23%	4.125%
66	17.516	2447	2453	2468	rBV	2191842	3085512	72.62%	9.591%
67	17.828	2505	2506	2512	rVB6	4245	4739	0.11%	0.015%
68	18.263	2576	2580	2590	rBV	60171	89591	2.11%	0.278%
69	19.339	2759	2763	2767	rBV2	32279	37627	0.89%	0.117%
70	19.410	2772	2775	2779	rVB3	26687	36312	0.85%	0.113%
71	19.510	2788	2792	2797	rBV5	37346	56765	1.34%	0.176%
72	19.645	2813	2815	2820	rBV5	10028	9683	0.23%	0.030%
73	19.774	2831	2837	2844	rBV	2962719	3841722	90.42%	11.942%
74	21.551	3135	3139	3145	rBV	1166574	1554255	36.58%	4.831%
75	23.962	3541	3549	3563	rVB2	2054831	4248576	100.00%	13.207%
76	24.133	3571	3578	3588	rVB	803699	1657394	39.01%	5.152%

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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-BP100423.MA.M
Title : SVOA CALIBRATION

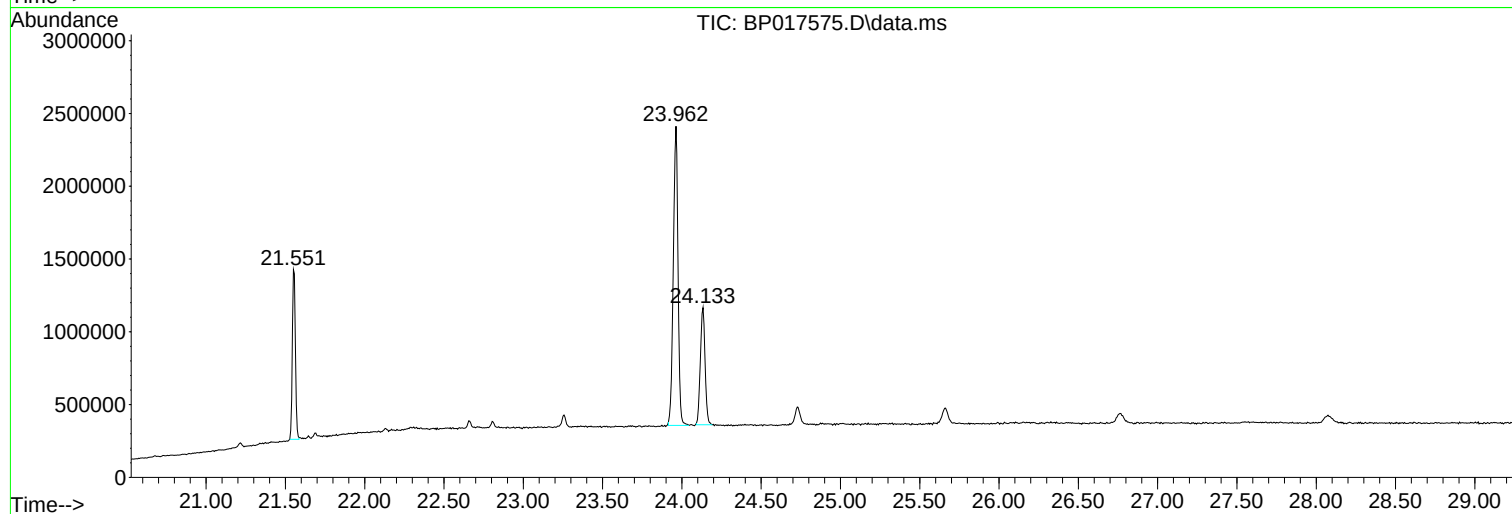
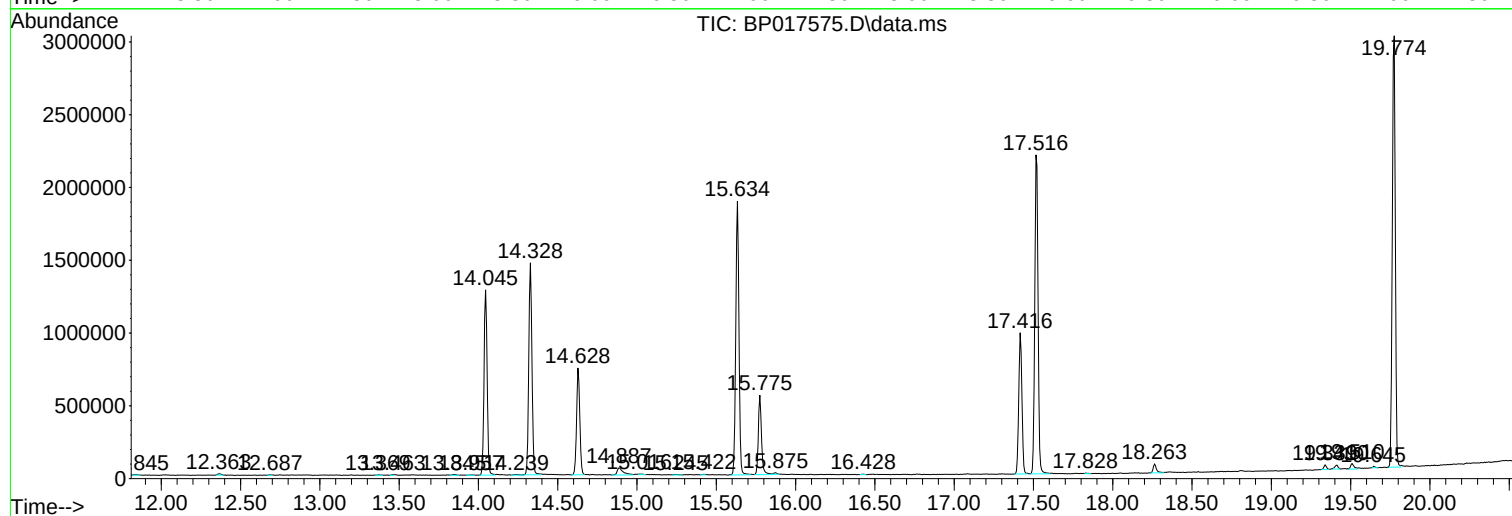
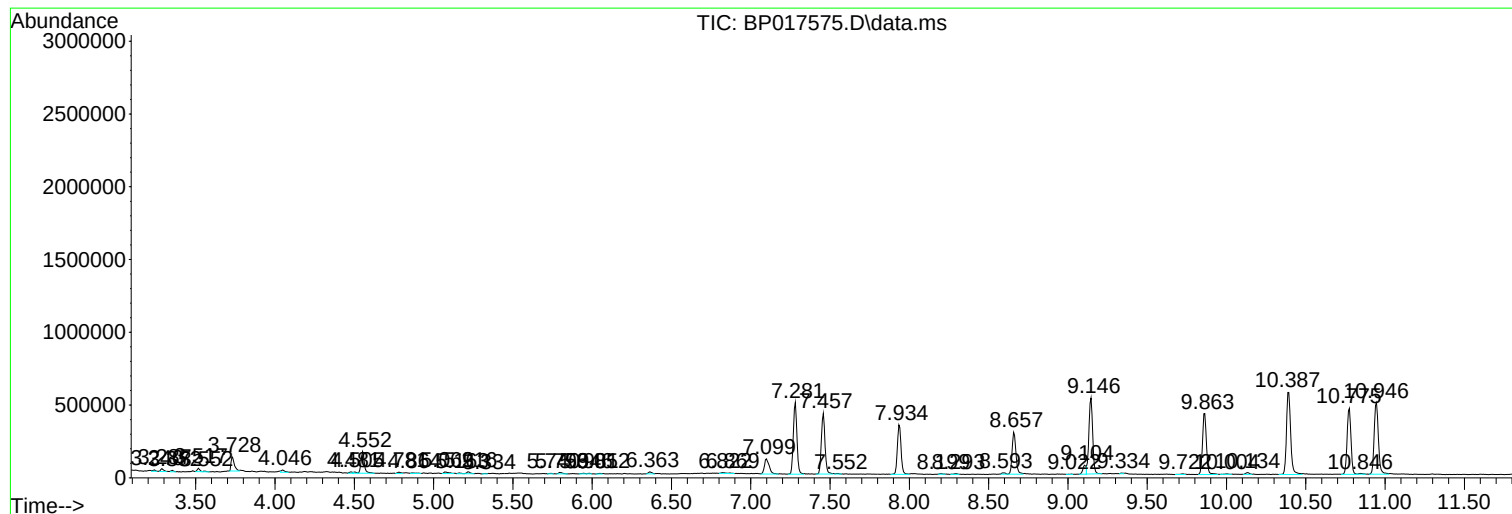
Sum of corrected areas: 32169841

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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

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



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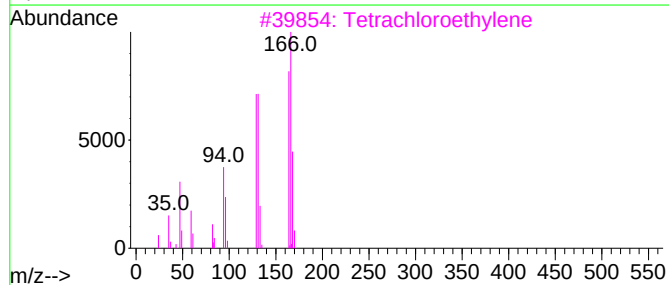
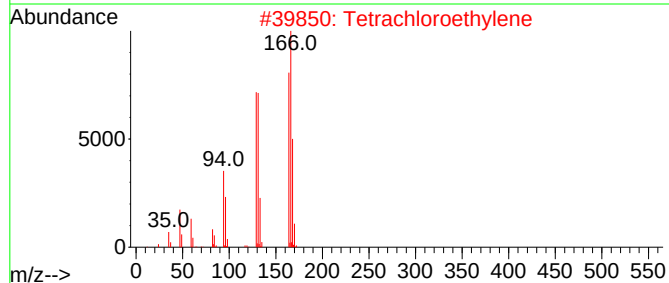
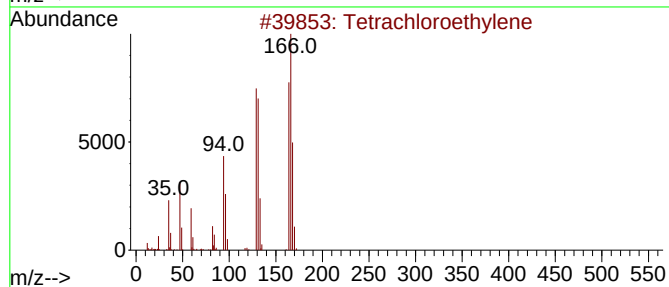
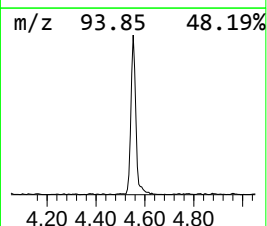
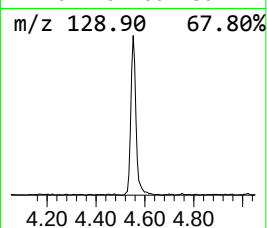
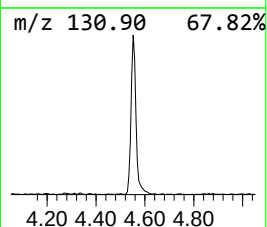
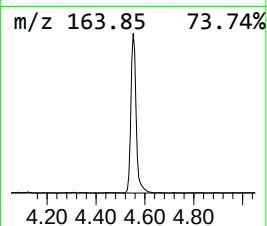
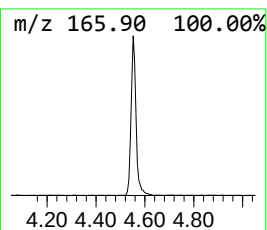
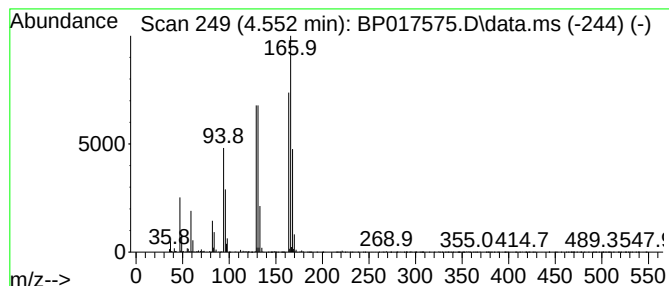
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

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.
4.552	8.16 ng/ul	223532	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	35



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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...	4.552	8.2	ng/u1	223532	1	7.934	548151	20.0