

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012149.D
 Acq On : 14 Oct 2022 22:10
 Operator : CG/JU
 Sample : N4984-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGNC3

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-BP101122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012149.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.011	170	174	187	rVB	366873	527447	14.28%	1.370%
2	6.775	641	644	649	rVB5	9250	12844	0.35%	0.033%
3	7.010	678	684	701	rVB	474595	836532	22.65%	2.172%
4	7.175	706	712	723	rVB	421682	677107	18.34%	1.758%
5	7.375	739	746	757	rBV	486469	811024	21.96%	2.106%
6	7.840	818	825	833	rBV	833397	1337924	36.23%	3.474%
7	8.557	940	947	959	rBV	578757	979658	26.53%	2.544%
8	8.681	963	968	974	rVB	20751	37992	1.03%	0.099%
9	9.010	1016	1024	1039	rBV	457532	774374	20.97%	2.011%
10	9.399	1088	1090	1095	rVB5	4937	7161	0.19%	0.019%
11	9.734	1139	1147	1158	rBV	328458	566367	15.34%	1.471%
12	10.269	1231	1238	1259	rBV	510339	934021	25.29%	2.425%
13	10.440	1261	1267	1282	rBV	51808	149503	4.05%	0.388%
14	10.645	1294	1302	1317	rVB	1122199	1921658	52.04%	4.990%
15	10.793	1320	1327	1347	rBV	463023	861480	23.33%	2.237%
16	11.493	1441	1446	1454	rVB	2312	6928	0.19%	0.018%
17	12.240	1569	1573	1578	rBV2	8699	14252	0.39%	0.037%
18	12.845	1672	1676	1680	rVB6	2833	4564	0.12%	0.012%
19	13.098	1713	1719	1723	rBV8	3024	5402	0.15%	0.014%
20	13.216	1736	1739	1742	rBV5	3100	4306	0.12%	0.011%
21	13.387	1763	1768	1775	rVB2	17412	30335	0.82%	0.079%
22	13.457	1778	1780	1787	rVB7	3513	5285	0.14%	0.014%
23	13.822	1837	1842	1846	rVB6	4526	7537	0.20%	0.020%
24	13.904	1848	1856	1861	rBV	956427	1296791	35.12%	3.367%
25	13.951	1861	1864	1873	rVB	45522	73085	1.98%	0.190%
26	14.186	1897	1904	1922	rBV	1116057	1686652	45.67%	4.380%
27	14.492	1949	1956	1969	rBV	1698706	2454805	66.47%	6.374%
28	14.710	1986	1993	2019	rBV	180376	433496	11.74%	1.126%
29	14.922	2024	2029	2040	rVB5	33433	62633	1.70%	0.163%
30	15.339	2098	2100	2105	rVB6	3321	3784	0.10%	0.010%
31	15.486	2118	2125	2138	rBV	1587247	2266761	61.38%	5.886%
32	15.616	2141	2147	2161	rVV	192949	315322	8.54%	0.819%
33	15.733	2163	2167	2175	rVB10	9374	20347	0.55%	0.053%
34	15.945	2200	2203	2208	rBV7	3572	5350	0.14%	0.014%
35	16.210	2244	2248	2249	rBV4	6620	7767	0.21%	0.020%
36	16.998	2380	2382	2385	rBV3	5720	6648	0.18%	0.017%

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Filtering: 5

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Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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Title : SVOA CALIBRATION

37	17.251	2418	2425	2435	rBV	2083716	2979834	80.69%	7.737%
38	17.351	2436	2442	2454	rVV	1753508	2540936	68.80%	6.598%
39	18.133	2571	2575	2585	rBV2	78604	129335	3.50%	0.336%
40	19.369	2782	2785	2789	rBV	51765	70969	1.92%	0.184%
41	19.586	2817	2822	2834	rBV	2328332	3136128	84.92%	8.143%
42	21.345	3116	3121	3132	rBV	2713815	3524424	95.44%	9.152%
43	23.598	3499	3504	3517	rVB2	1590071	3290030	89.09%	8.543%
44	23.757	3525	3531	3544	rVB2	1783250	3692963	100.00%	9.589%

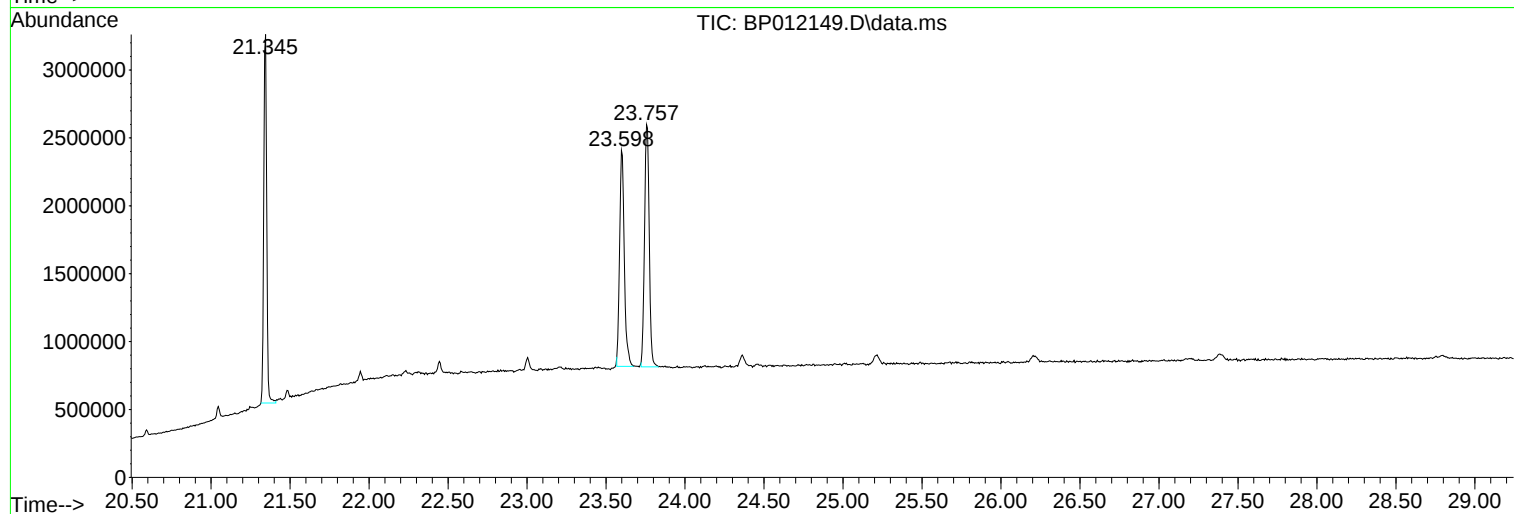
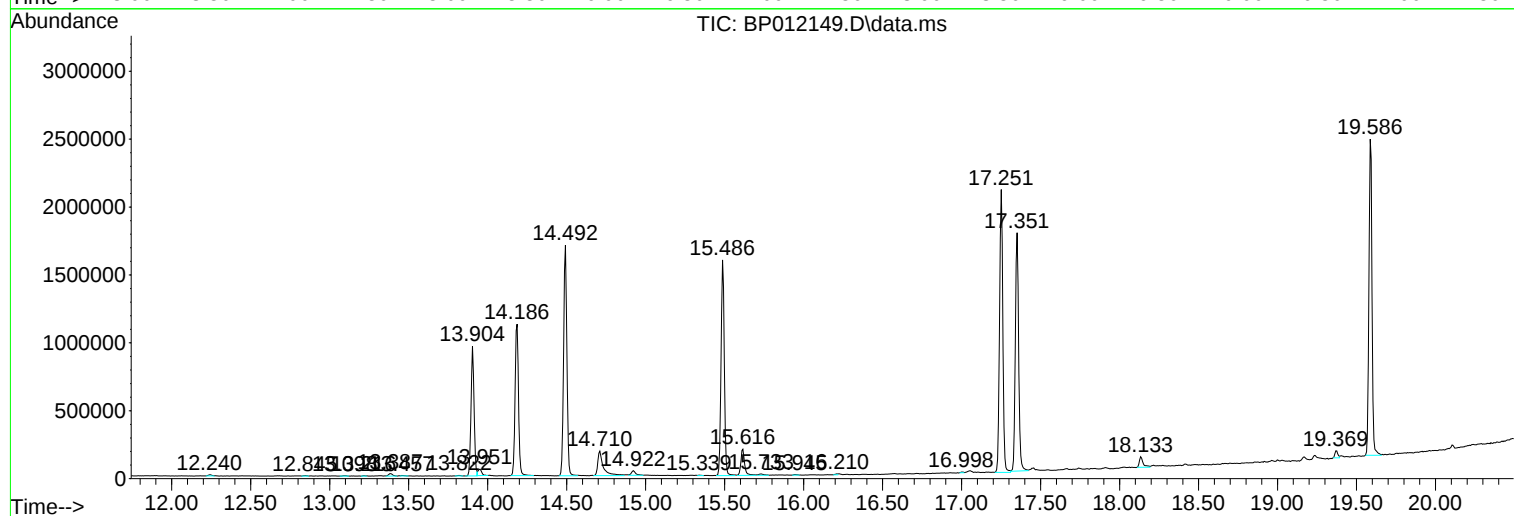
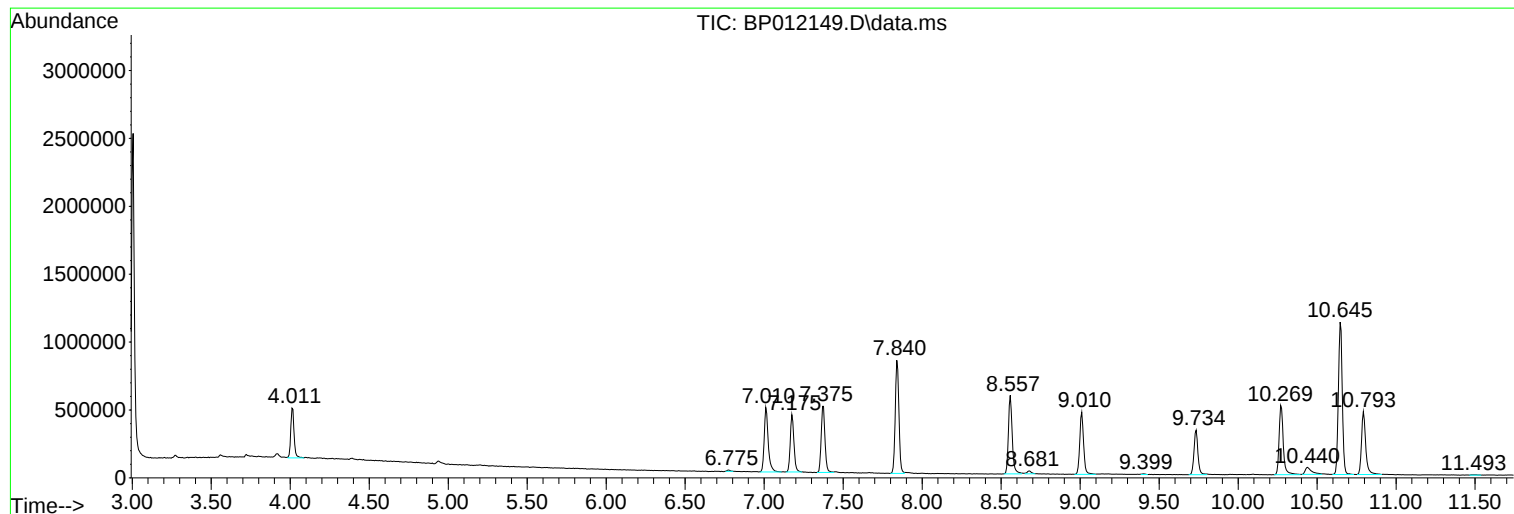
Sum of corrected areas: 38511761

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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



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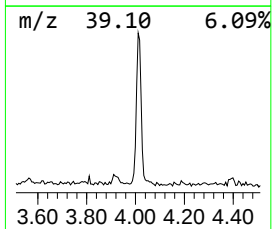
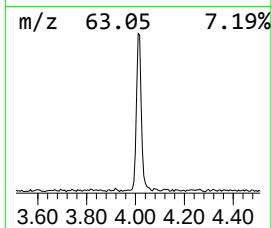
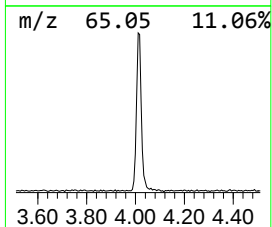
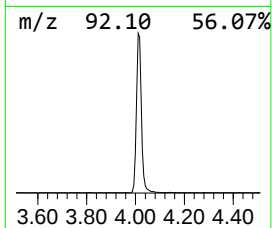
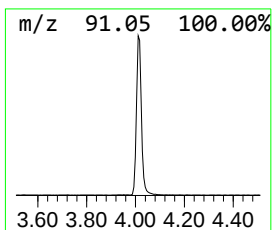
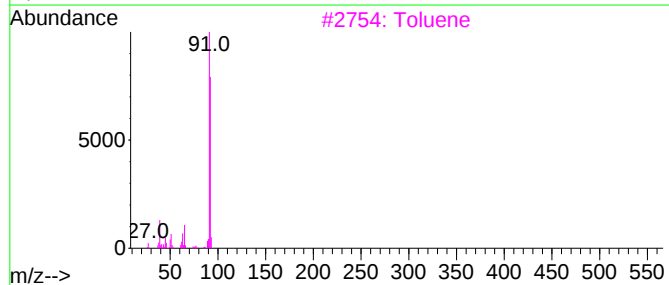
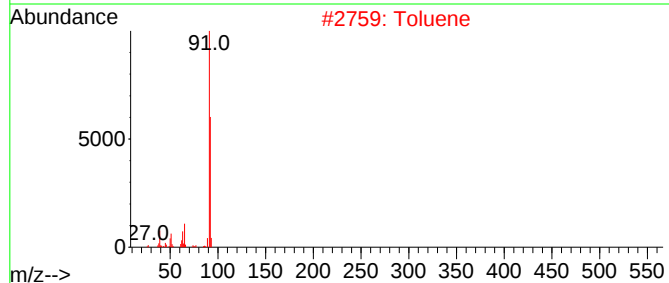
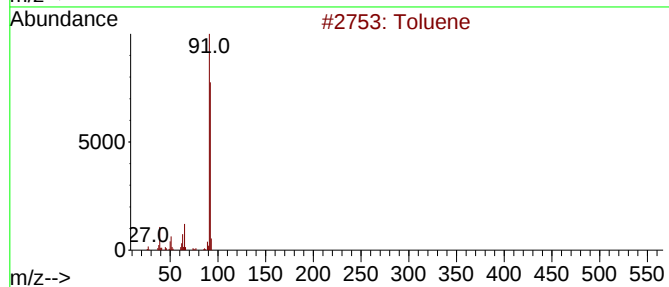
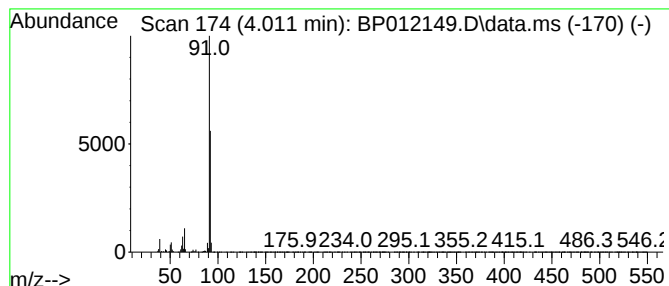
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Toluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.011	7.88 ng/ul	527447	1,4-Dichlorobenzene-d4	7.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	95
2	Toluene			92	C7H8	000108-88-3	94
3	Toluene			92	C7H8	000108-88-3	91
4	Toluene			92	C7H8	000108-88-3	91
5	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	87



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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
Toluene	4.011	7.9	ng/ul	527447	1	7.840	1337920	20.0