

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
 Data File : BP016683.D
 Acq On : 21 Aug 2023 20:13
 Operator : MA/JU
 Sample : PB154937BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK937

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-BP081423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016683.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.387	13	17	31	rVB	193611	301318	2.22%	0.224%
2	3.823	87	91	113	rBV	2142938	3437392	25.35%	2.557%
3	4.146	140	146	158	rVB	254723	391958	2.89%	0.292%
4	7.187	654	663	677	rBV	2597701	4014382	29.61%	2.986%
5	7.381	689	696	708	rVB	2049568	3123022	23.04%	2.323%
6	7.563	720	727	741	rBV	2409181	3744459	27.62%	2.786%
7	8.046	800	809	818	rBV	1386867	2139747	15.78%	1.592%
8	8.746	921	928	941	rBV	3036702	4966949	36.64%	3.695%
9	9.246	1003	1013	1024	rBV	2410339	3922844	28.93%	2.918%
10	9.957	1126	1134	1150	rBV	2005421	3238744	23.89%	2.409%
11	10.481	1214	1223	1243	rBV	2897860	4919518	36.29%	3.660%
12	10.875	1281	1290	1298	rBV	2051032	3351368	24.72%	2.493%
13	11.034	1308	1317	1331	rBV	3188669	5349056	39.45%	3.979%
14	12.451	1553	1558	1565	rVB	47016	81086	0.60%	0.060%
15	13.546	1738	1744	1750	rBV	39490	65275	0.48%	0.049%
16	13.657	1760	1763	1768	rVB5	9659	15635	0.12%	0.012%
17	14.110	1832	1840	1852	rBV	5297707	7460774	55.03%	5.550%
18	14.398	1881	1889	1902	rBV	5970228	8949075	66.01%	6.657%
19	14.698	1932	1940	1953	rBV	3029873	4562857	33.66%	3.394%
20	14.898	1967	1974	1992	rBV	2642880	4147229	30.59%	3.085%
21	15.692	2102	2109	2121	rBV	7787591	11050868	81.51%	8.221%
22	15.810	2122	2129	2143	rVV	2340281	3140377	23.16%	2.336%
23	15.928	2145	2149	2157	rVB2	37518	62369	0.46%	0.046%
24	17.128	2349	2353	2358	rBV6	11365	18647	0.14%	0.014%
25	17.463	2403	2410	2420	rBV	3986374	5597150	41.28%	4.164%
26	17.563	2420	2427	2439	rBV2	8765165	12091709	89.19%	8.995%
27	18.304	2549	2553	2556	rBV6	20723	33462	0.25%	0.025%
28	19.369	2729	2734	2738	rBV2	105611	141670	1.04%	0.105%
29	19.433	2741	2745	2751	rVB	94750	127962	0.94%	0.095%
30	19.792	2801	2806	2818	rBV	10089193	13557486	100.00%	10.086%
31	21.557	3101	3106	3113	rBV	4028361	5182224	38.22%	3.855%
32	23.969	3507	3516	3530	rVB2	4953130	10708766	78.99%	7.966%
33	24.133	3537	3544	3557	rVB	2073804	4528895	33.41%	3.369%

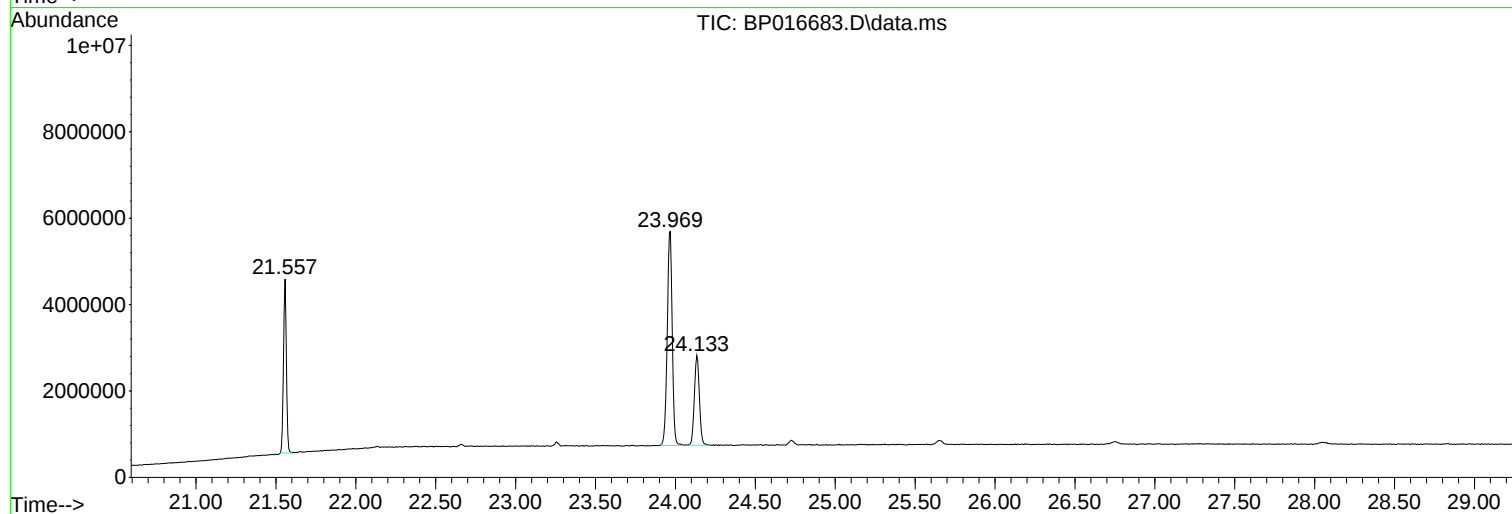
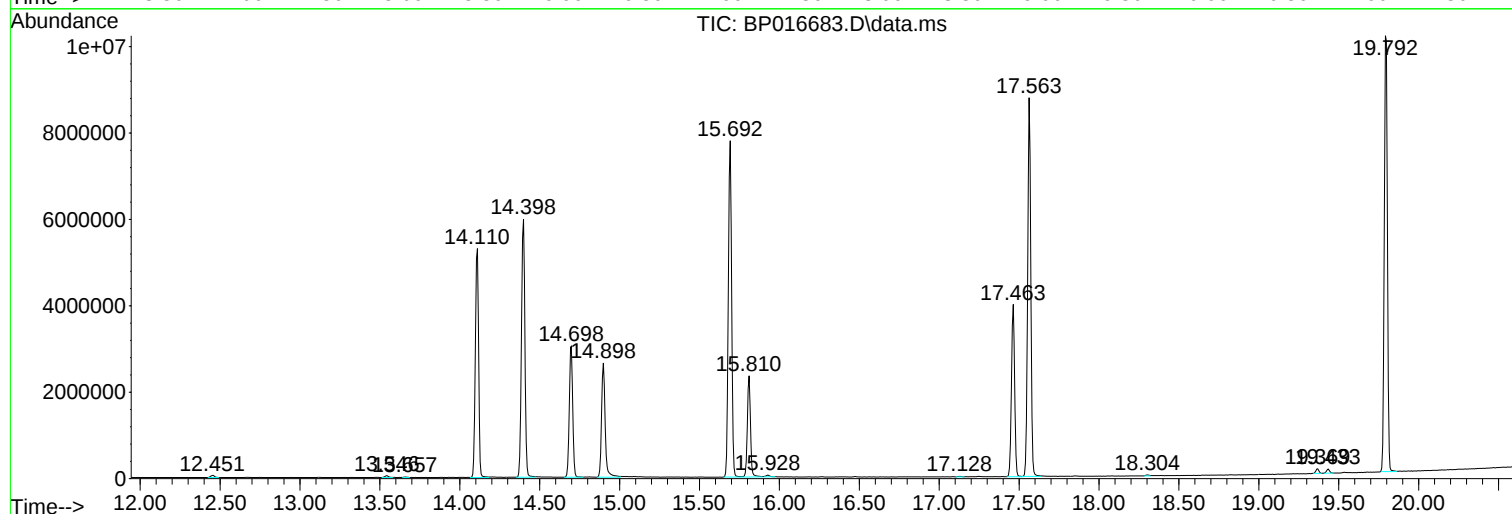
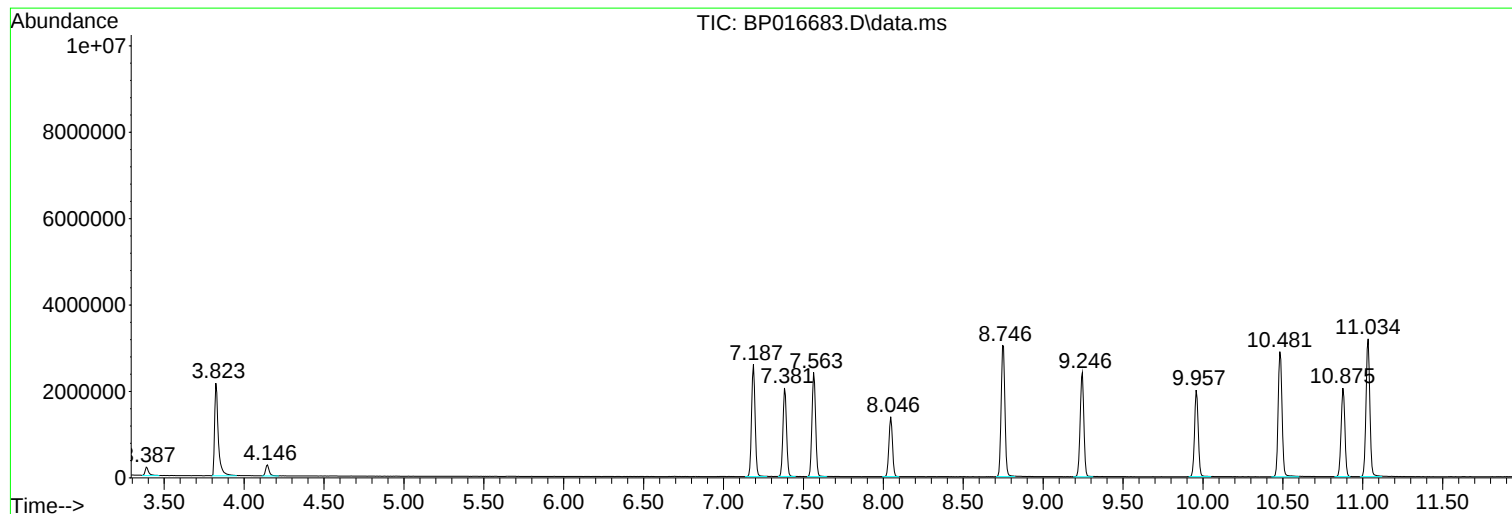
Sum of corrected areas: 134424273

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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



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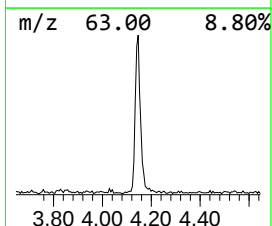
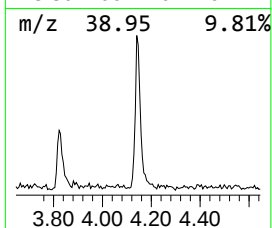
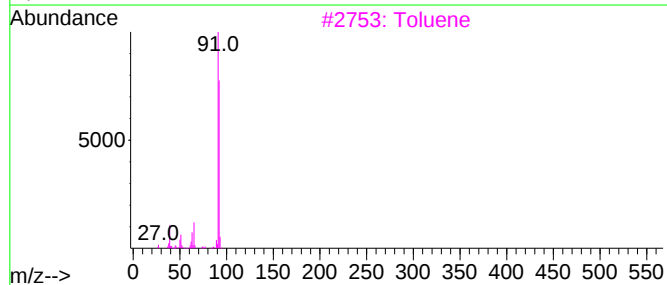
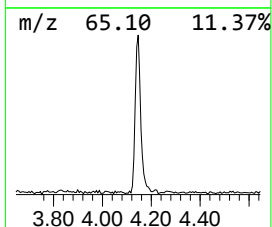
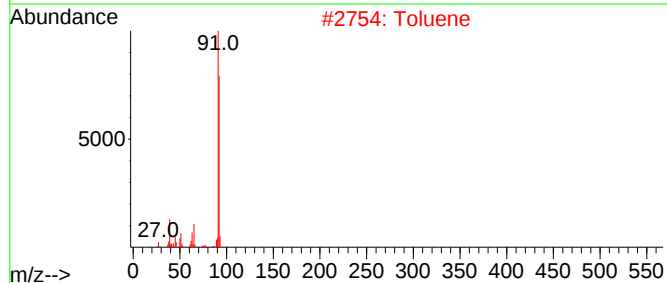
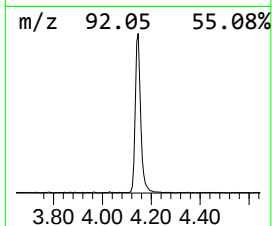
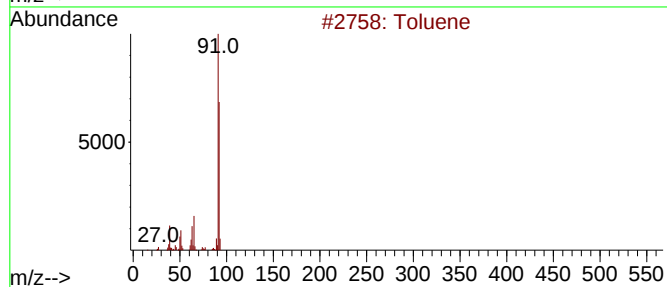
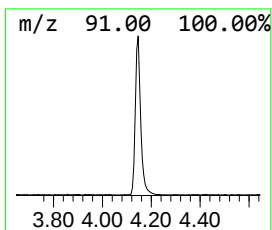
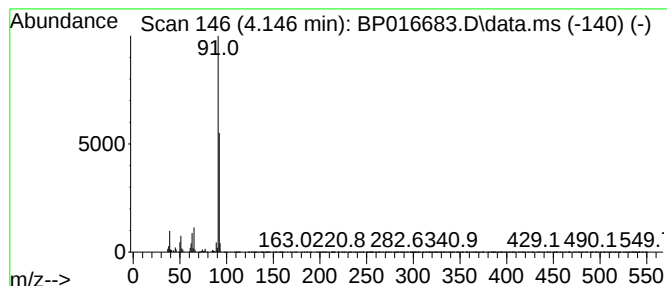
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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.
4.146	3.66 ng/ul	391958	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	93
2	Toluene			92	C7H8	000108-88-3	91
3	Toluene			92	C7H8	000108-88-3	90
4	Toluene			92	C7H8	000108-88-3	87
5	Toluene			92	C7H8	000108-88-3	87



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Toluene	4.146	3.7	ng/u1	391958	1	8.046	2139750	20.0