

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081722\
 Data File : BP011468.D
 Acq On : 17 Aug 2022 19:20
 Operator : CG/JU
 Sample : PB147029TB 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB029

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-BP081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011468.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.070	11	14	20	rVB	43068	49206	2.09%	0.179%
2	3.475	80	83	102	rBV	492976	745851	31.67%	2.715%
3	3.781	131	135	144	rVB	52798	70378	2.99%	0.256%
4	4.411	239	242	247	rBV6	1460	2474	0.11%	0.009%
5	4.569	268	269	274	rBV5	1903	2663	0.11%	0.010%
6	6.746	632	639	652	rBV	562513	870789	36.98%	3.170%
7	6.911	660	667	681	rVB	427209	645502	27.41%	2.350%
8	7.093	691	698	707	rBV	572220	879390	37.34%	3.201%
9	7.558	770	777	788	rBV	396721	604673	25.68%	2.201%
10	8.281	892	900	911	rBV	647084	1028043	43.65%	3.742%
11	8.722	968	975	986	rBV	538010	838857	35.62%	3.054%
12	9.440	1087	1097	1107	rBV	453104	728368	30.93%	2.651%
13	9.975	1180	1188	1203	rBV	613088	997193	42.34%	3.630%
14	10.346	1242	1251	1261	rBV	530518	849357	36.07%	3.092%
15	10.499	1270	1277	1287	rBV	603277	1019859	43.31%	3.712%
16	11.951	1520	1524	1532	rVB2	9056	16775	0.71%	0.061%
17	13.204	1730	1737	1741	rBV9	2807	5961	0.25%	0.022%
18	13.657	1807	1814	1826	rBV	1074023	1423236	60.43%	5.181%
19	13.745	1826	1829	1834	rVV7	1988	3155	0.13%	0.011%
20	13.922	1850	1859	1867	rBV	1139039	1635994	69.47%	5.955%
21	14.234	1904	1912	1921	rBV	713814	1023709	43.47%	3.726%
22	14.298	1921	1923	1927	rVB5	2940	2747	0.12%	0.010%
23	14.463	1944	1951	1970	rBV	447594	752346	31.95%	2.739%
24	14.675	1985	1987	1994	rVB8	4622	6750	0.29%	0.025%
25	15.175	2067	2072	2074	rBV6	1777	2375	0.10%	0.009%
26	15.234	2076	2082	2090	rBV	1363712	1927129	81.83%	7.015%
27	15.375	2100	2106	2118	rBV	530542	715621	30.39%	2.605%
28	15.475	2120	2123	2128	rVB6	7347	10788	0.46%	0.039%
29	15.810	2176	2180	2184	rBV7	2038	3272	0.14%	0.012%
30	16.139	2231	2236	2239	rBV5	3961	5752	0.24%	0.021%
31	16.422	2280	2284	2285	rBV4	1950	2484	0.11%	0.009%
32	16.675	2325	2327	2330	rVB4	2277	2575	0.11%	0.009%
33	16.786	2341	2346	2352	rBV10	4584	9358	0.40%	0.034%
34	16.945	2371	2373	2377	rBV5	2744	3805	0.16%	0.014%
35	17.004	2377	2383	2389	rVV	884504	1168644	49.62%	4.254%
36	17.104	2393	2400	2410	rVV	1574874	2109236	89.56%	7.678%

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Peak Location: TOP

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Peak separation: 5

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 Title : SVOA CALIBRATION

37	17.698	2498	2501	2506	rVB7	5692	10023	0.43%	0.036%
38	17.922	2535	2539	2543	rBV3	16213	24857	1.06%	0.090%
39	18.939	2708	2712	2717	rBV6	20127	29389	1.25%	0.107%
40	19.010	2720	2724	2727	rBV	20944	30584	1.30%	0.111%
41	19.045	2727	2730	2732	rVB	11703	12793	0.54%	0.047%
42	19.369	2779	2785	2793	rBV	1827054	2355003	100.00%	8.573%
43	21.133	3080	3085	3089	rBV	1052445	1290036	54.78%	4.696%
44	23.280	3444	3450	3457	rBV2	1300637	2284777	97.02%	8.317%
45	23.427	3469	3475	3485	rVB2	659902	1269737	53.92%	4.622%

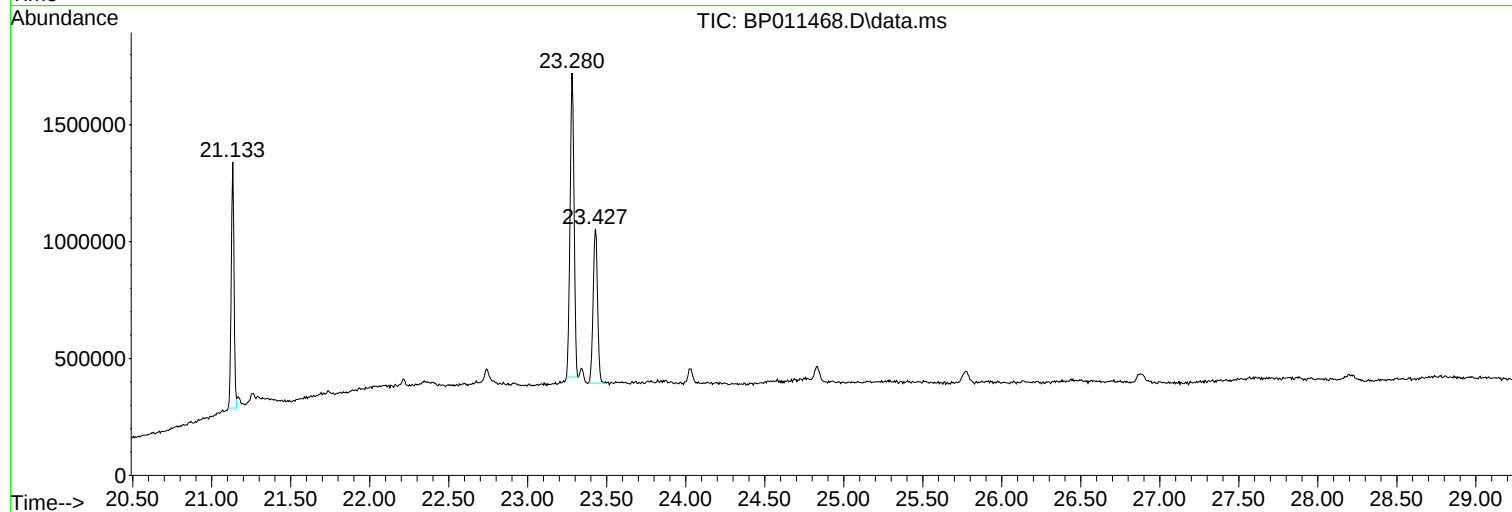
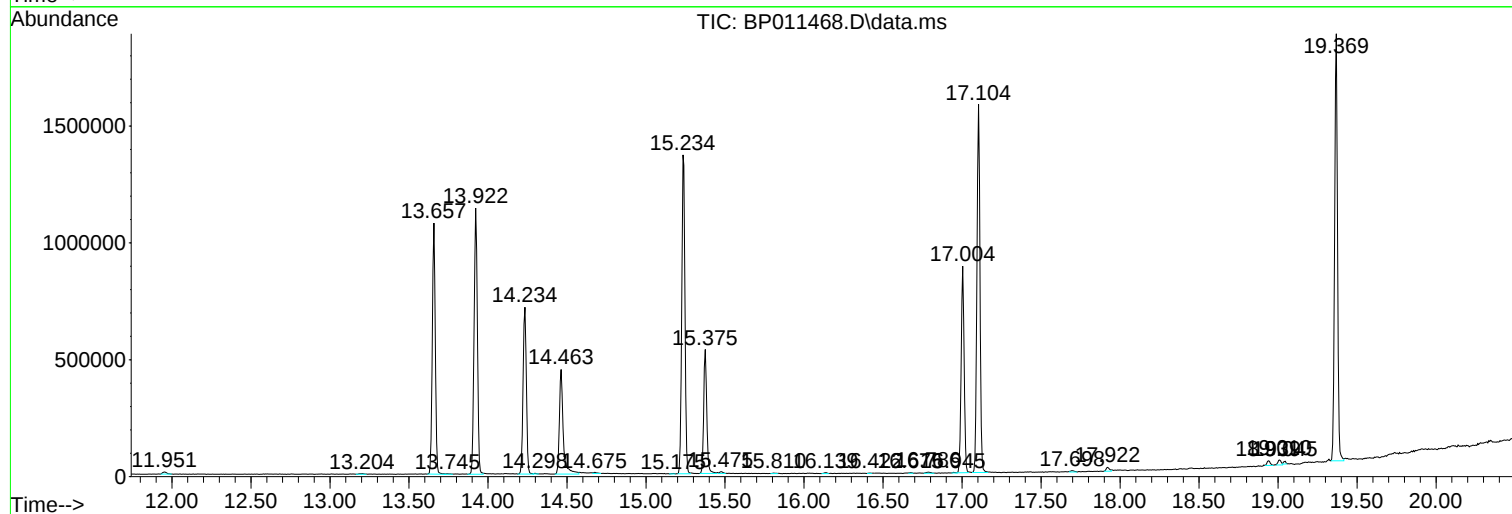
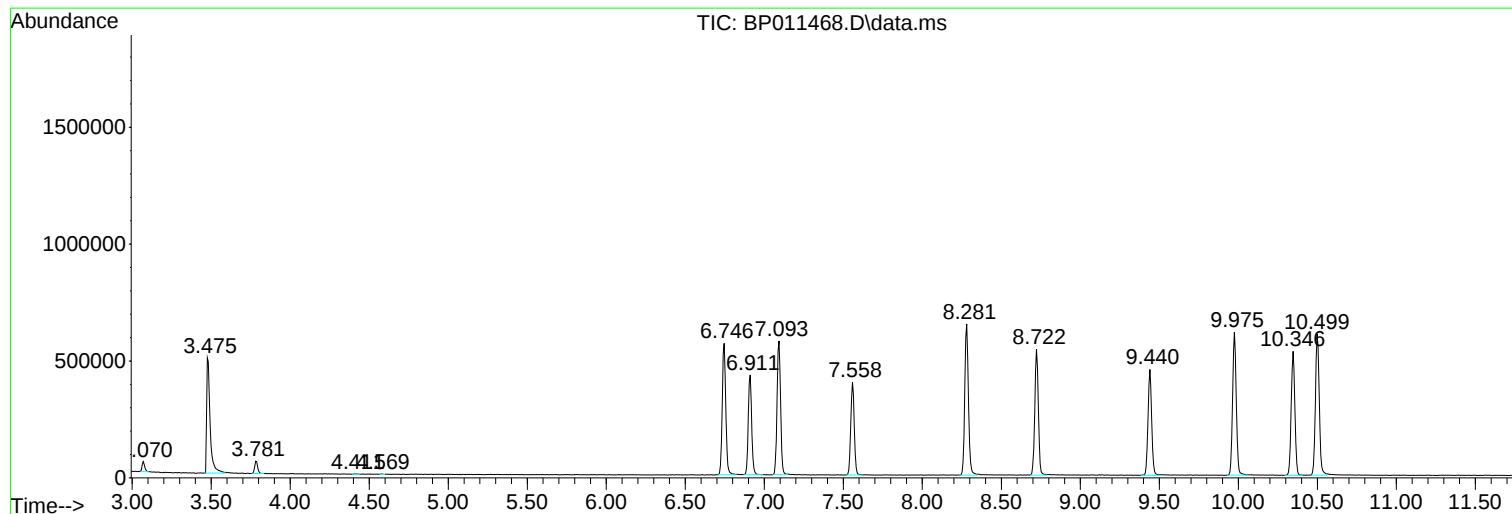
Sum of corrected areas: 27471514

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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



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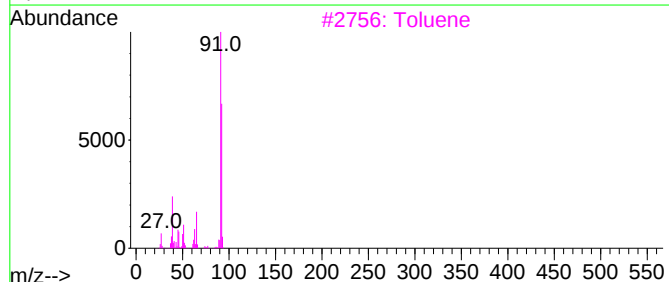
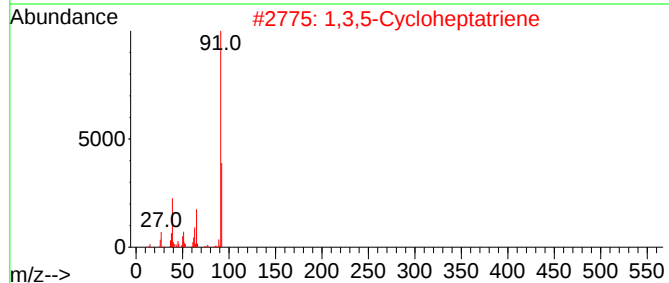
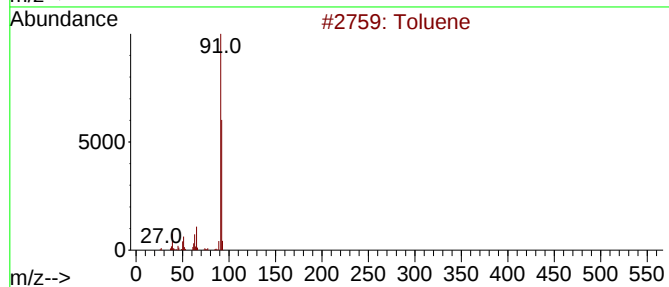
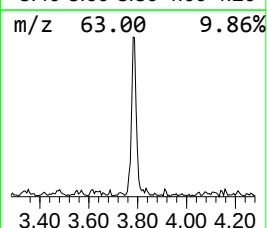
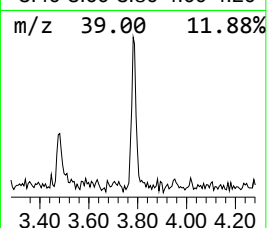
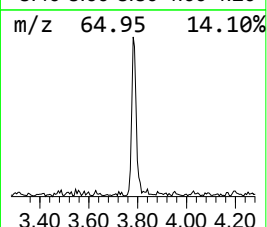
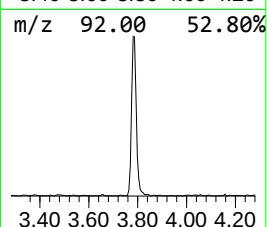
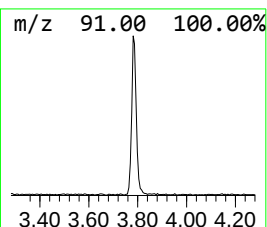
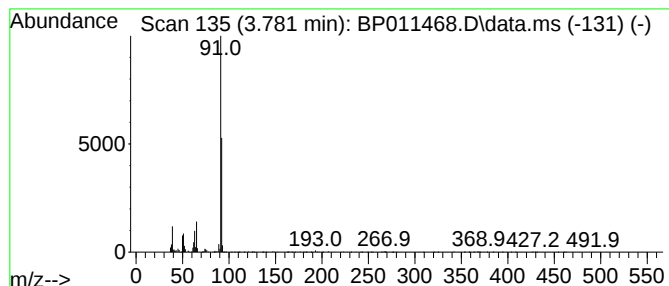
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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.
3.781	2.33 ng/ul	70378	1,4-Dichlorobenzene-d4	7.558

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	90
2			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	90
3			Toluene	92	C7H8	000108-88-3	90
4			Toluene	92	C7H8	000108-88-3	87
5			Spiro[2,4]hepta-4,6-diene	92	C7H8	000765-46-8	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	3.781	2.3	ng/ul	70378	1	7.558	604673	20.0