

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
 Data File : BP016669.D
 Acq On : 21 Aug 2023 11:39
 Operator : MA/JU
 Sample : PB154934BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK934

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: BP016669.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	14	18	30	rVB	156268	218483	2.59%	0.220%
2	3.828	87	92	107	rBV	1655710	2519590	29.85%	2.537%
3	4.146	141	146	156	rVB	229364	321047	3.80%	0.323%
4	7.187	656	663	674	rBV	1896853	2930217	34.71%	2.950%
5	7.381	689	696	705	rBV	1488521	2303672	27.29%	2.320%
6	7.563	720	727	737	rBV	1746905	2775999	32.89%	2.795%
7	8.046	800	809	817	rBV	1480768	2291302	27.15%	2.307%
8	8.752	920	929	945	rBV	2212741	3633583	43.05%	3.659%
9	9.246	1002	1013	1029	rBV	1742661	2889959	34.24%	2.910%
10	9.957	1125	1134	1148	rBV	1443402	2349196	27.83%	2.365%
11	10.487	1216	1224	1244	rBV	2111356	3614885	42.83%	3.640%
12	10.875	1282	1290	1300	rBV	2114445	3568782	42.28%	3.593%
13	11.034	1309	1317	1336	rBV	2297802	3899784	46.20%	3.927%
14	12.451	1552	1558	1566	rBV	37014	64481	0.76%	0.065%
15	13.545	1739	1744	1747	rBV	25255	35045	0.42%	0.035%
16	13.657	1756	1763	1769	rVB6	8532	15852	0.19%	0.016%
17	14.110	1833	1840	1853	rBV	3820856	5497789	65.13%	5.536%
18	14.198	1853	1855	1863	rVB9	5129	11397	0.14%	0.011%
19	14.398	1881	1889	1909	rBV	4496217	6619929	78.43%	6.665%
20	14.698	1933	1940	1954	rBV	3423290	4883904	57.86%	4.917%
21	14.898	1967	1974	1996	rBV	1507208	2501779	29.64%	2.519%
22	15.128	2011	2013	2021	rVB8	6443	11569	0.14%	0.012%
23	15.692	2101	2109	2123	rBV	5791784	8117673	96.17%	8.173%
24	15.810	2123	2129	2145	rVV	1376403	1876781	22.23%	1.890%
25	15.928	2145	2149	2159	rVB2	29974	53139	0.63%	0.054%
26	16.263	2203	2206	2211	rBV7	5901	10533	0.12%	0.011%
27	16.475	2238	2242	2247	rBV8	7918	15337	0.18%	0.015%
28	17.128	2350	2353	2359	rBV7	8833	15503	0.18%	0.016%
29	17.251	2370	2374	2378	rBV4	8685	13619	0.16%	0.014%
30	17.463	2404	2410	2420	rBV	4066397	5669198	67.16%	5.708%
31	17.563	2421	2427	2444	rVV2	5842159	8440820	100.00%	8.499%
32	18.310	2551	2554	2562	rVB9	13021	24439	0.29%	0.025%
33	19.369	2730	2734	2739	rBV3	70168	93351	1.11%	0.094%
34	19.439	2742	2746	2749	rVV	49268	68498	0.81%	0.069%
35	19.798	2801	2807	2815	rBV	6373765	8352215	98.95%	8.410%
36	21.563	3102	3107	3114	rBV	2893692	3848349	45.59%	3.875%

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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

37	23.968	3509	3516	3527	rVB2	2810545	5961611	70.63%	6.003%
38	24.139	3539	3545	3563	rVB	1726838	3798103	45.00%	3.824%

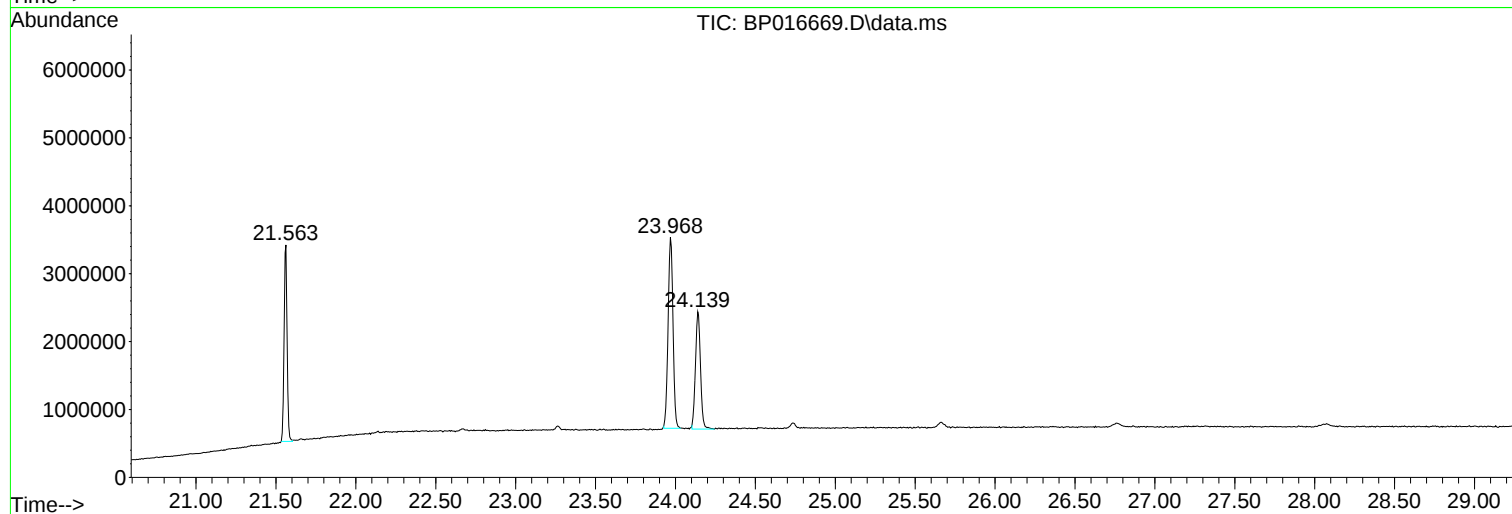
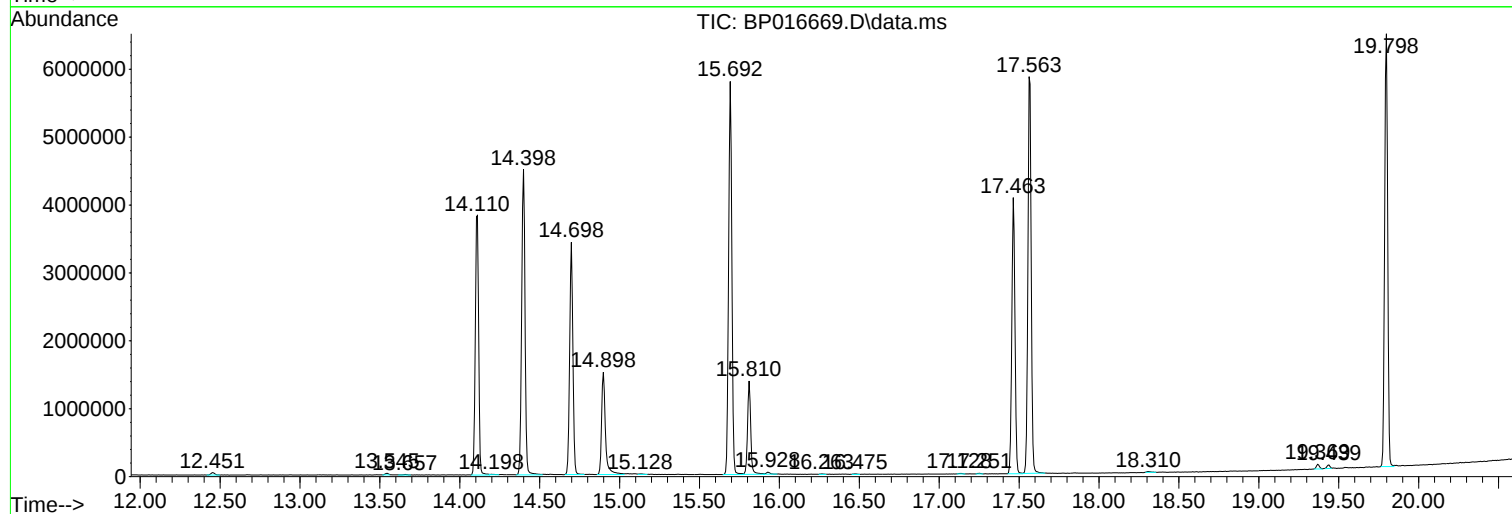
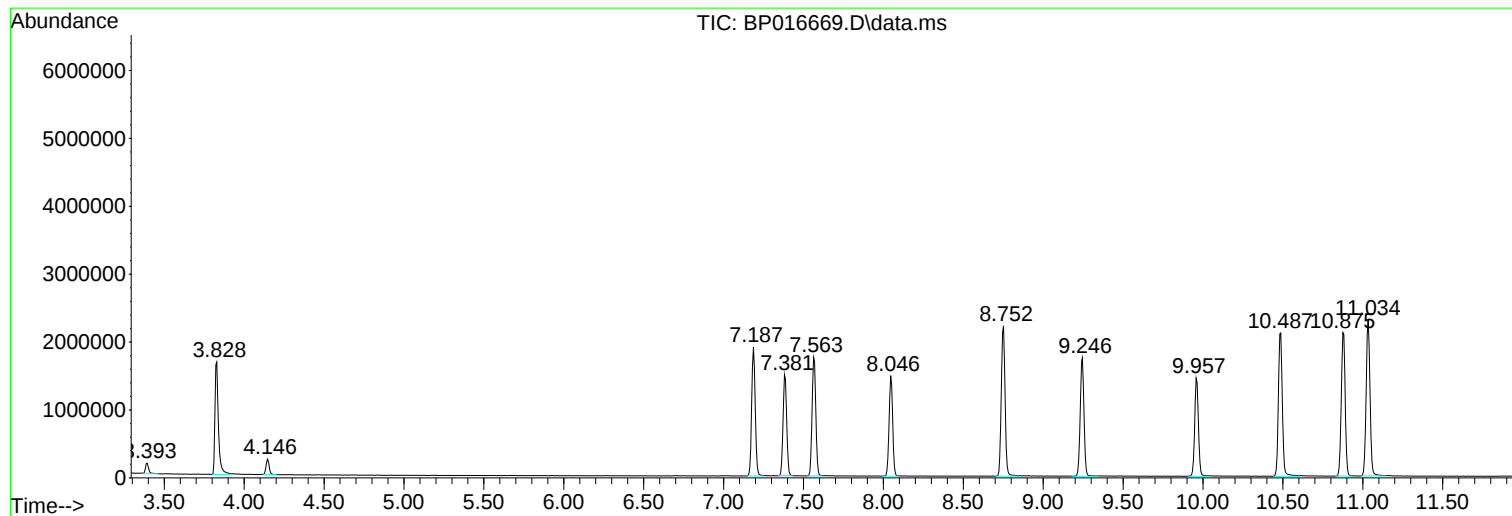
Sum of corrected areas: 99317413

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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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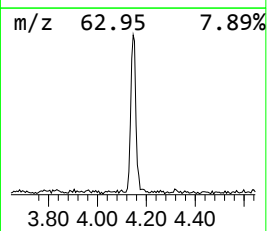
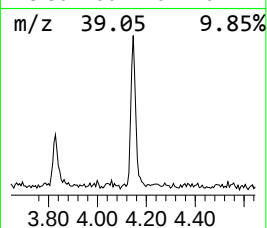
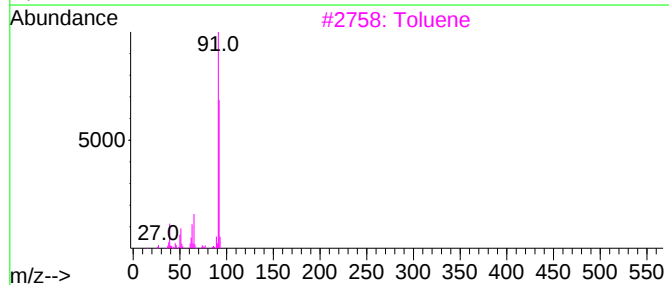
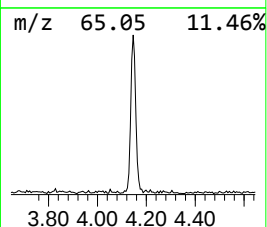
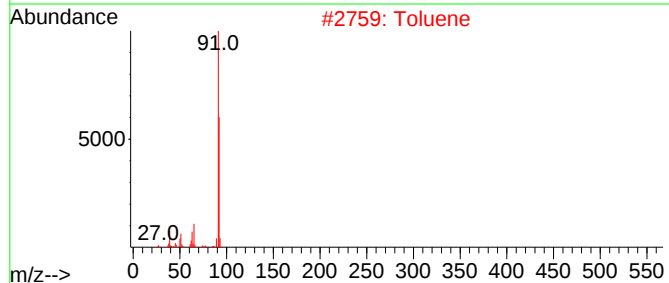
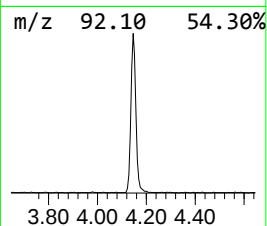
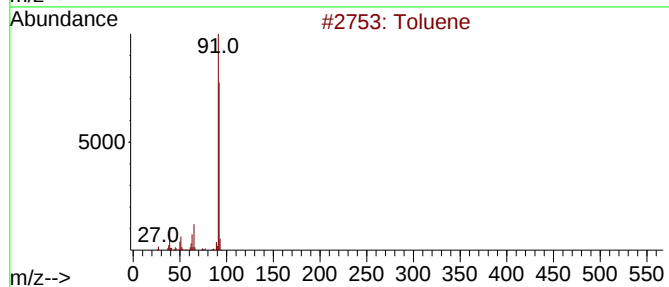
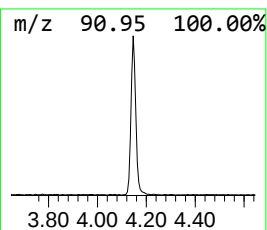
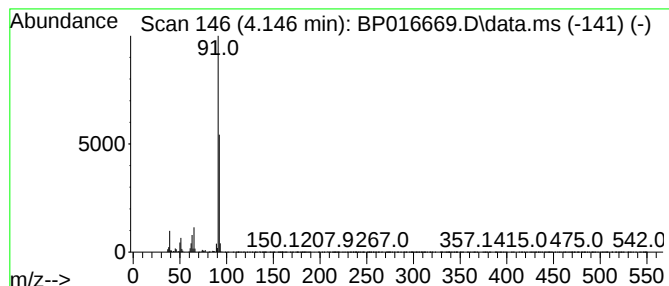
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Toluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.146	2.80 ng/ul	321047	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	95
2	Toluene			92	C7H8	000108-88-3	95
3	Toluene			92	C7H8	000108-88-3	94
4	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	90
5	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	90



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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	4.146	2.8	ng/ul	321047	1	8.046	2291300	20.0