

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009993.D
 Acq On : 21 Apr 2022 01:08
 Operator : CG/JU
 Sample : PB144233BL
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK233

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009993.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	44	48	53	rBV	33207	45981	0.94%	0.104%
2	3.787	115	119	135	rBV	410350	736798	15.00%	1.659%
3	4.122	171	176	185	rVB	57535	89337	1.82%	0.201%
4	7.099	674	682	698	rBV	585112	981693	19.99%	2.210%
5	7.269	704	711	724	rBV	489702	790256	16.09%	1.779%
6	7.475	738	746	758	rBV	611700	1007990	20.53%	2.269%
7	7.940	819	825	834	rBV	441557	739689	15.06%	1.665%
8	8.645	936	945	959	rBV	793618	1334745	27.18%	3.005%
9	9.098	1015	1022	1034	rBV	652691	1104308	22.49%	2.486%
10	9.822	1137	1145	1158	rBV	530228	937677	19.09%	2.111%
11	10.363	1229	1237	1252	rBV	803563	1379551	28.09%	3.106%
12	10.745	1294	1302	1312	rVV	683273	1179847	24.03%	2.656%
13	10.881	1316	1325	1339	rBV	807132	1471761	29.97%	3.314%
14	12.334	1564	1572	1579	rBV3	14445	28691	0.58%	0.065%
15	13.981	1845	1852	1864	rBV	1760846	2492776	50.76%	5.612%
16	14.263	1893	1900	1916	rBV	1790055	2730320	55.60%	6.147%
17	14.575	1946	1953	1965	rBV	1220864	1812591	36.91%	4.081%
18	14.757	1976	1984	2003	rBV	786331	1265333	25.77%	2.849%
19	15.216	2057	2062	2065	rBV7	4104	5594	0.11%	0.013%
20	15.386	2087	2091	2097	rVB8	3142	6442	0.13%	0.015%
21	15.563	2109	2121	2134	rBV	2458687	3587325	73.05%	8.077%
22	15.675	2134	2140	2152	rBV	830803	1175271	23.93%	2.646%
23	15.804	2158	2162	2170	rVB	30110	48310	0.98%	0.109%
24	16.251	2235	2238	2242	rBV6	3371	5255	0.11%	0.012%
25	16.351	2249	2255	2260	rBV8	10350	19441	0.40%	0.044%
26	17.004	2364	2366	2371	rVB5	5678	6869	0.14%	0.015%
27	17.104	2379	2383	2389	rVB5	5794	10646	0.22%	0.024%
28	17.322	2413	2420	2429	rBV	1612007	2309024	47.02%	5.199%
29	17.422	2431	2437	2457	rVB2	2819937	4105469	83.60%	9.243%
30	17.857	2508	2511	2514	rVB5	3698	5398	0.11%	0.012%
31	17.904	2514	2519	2521	rBV5	2976	5659	0.12%	0.013%
32	19.227	2740	2744	2751	rBV2	36713	50599	1.03%	0.114%
33	19.298	2752	2756	2762	rVB	36641	49623	1.01%	0.112%
34	19.651	2810	2816	2832	rBV	3712159	4910890	100.00%	11.056%
35	21.404	3108	3114	3125	rVB	1687964	2373809	48.34%	5.344%
36	23.698	3496	3504	3511	rBV2	1726486	3589200	73.09%	8.081%

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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-BP040722.M
Title : SVOA CALIBRATION

37	23.856	3523	3531	3545	rVB2	958054	2022365	41.18%	4.553%
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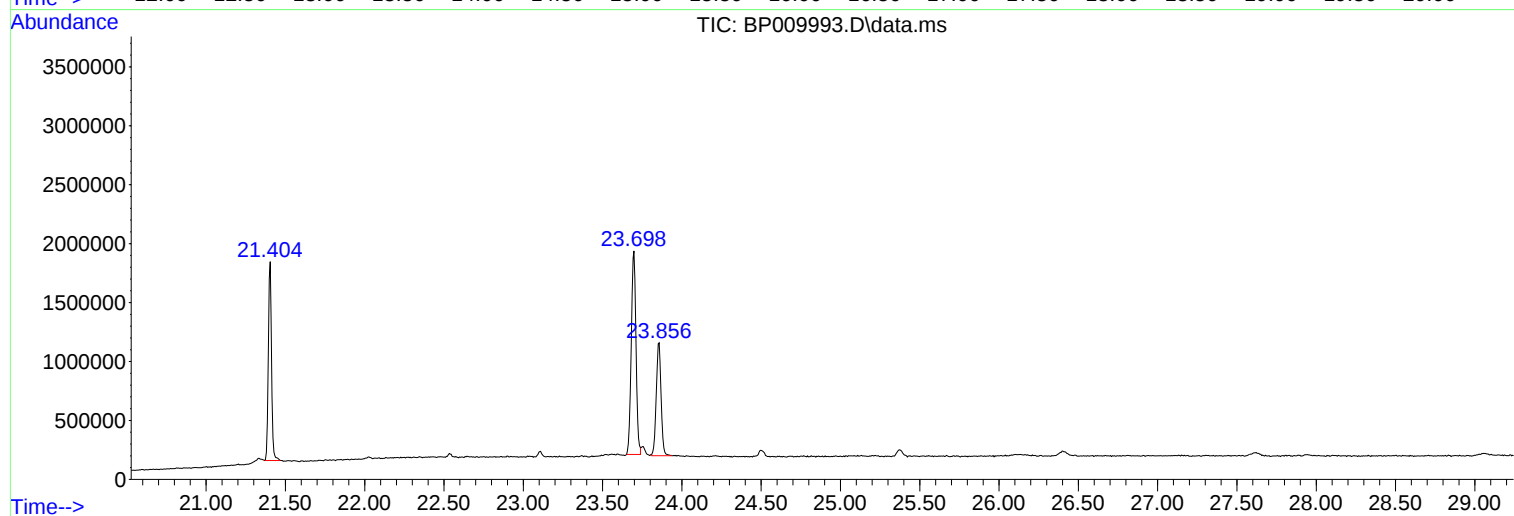
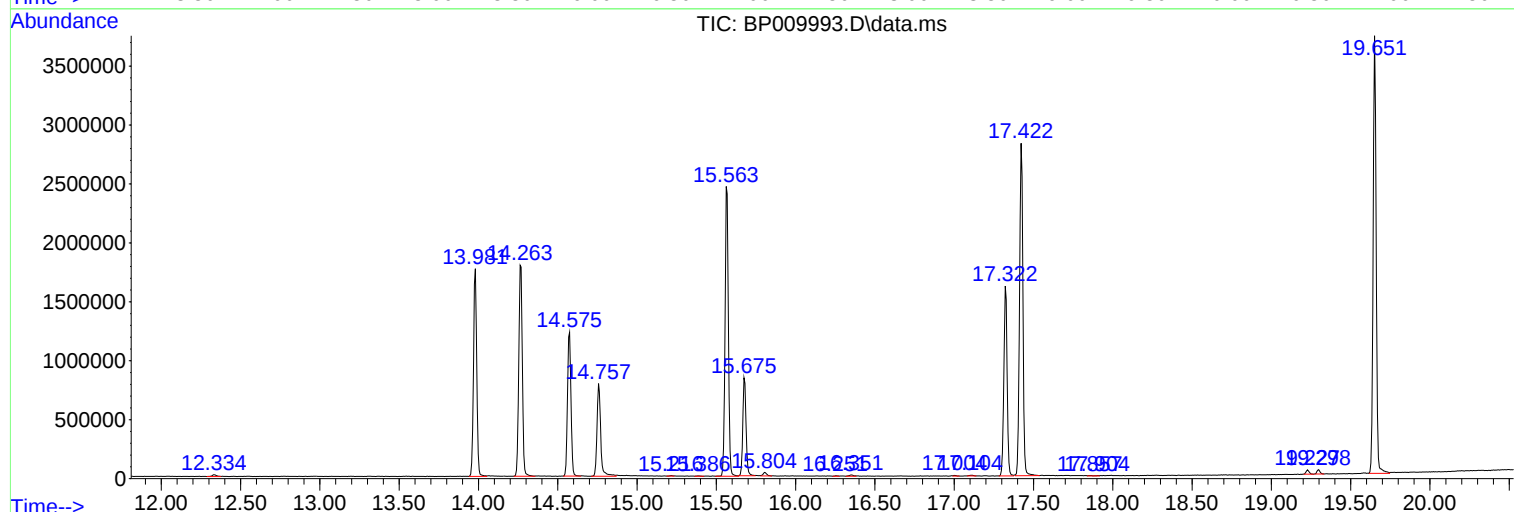
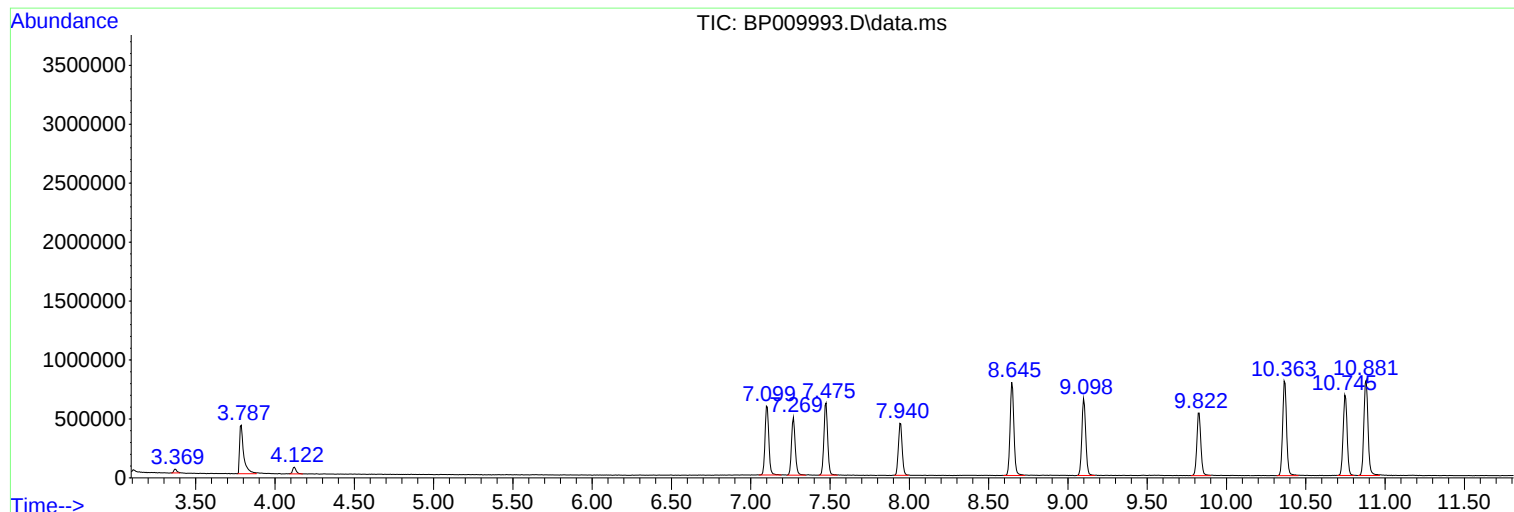
Sum of corrected areas: 44416533

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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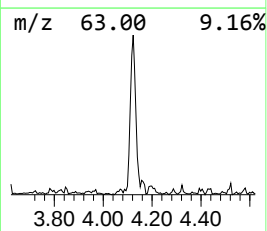
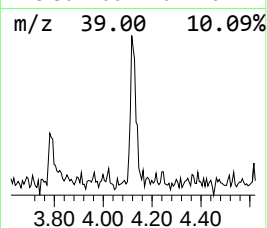
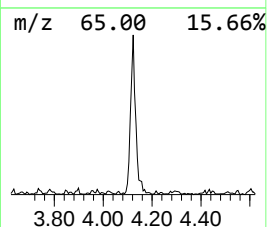
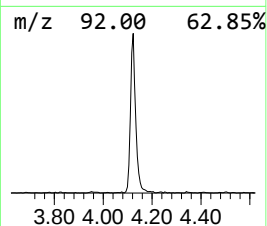
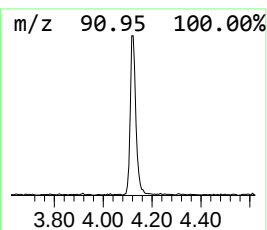
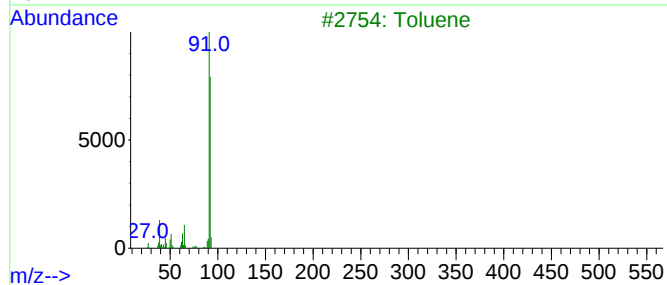
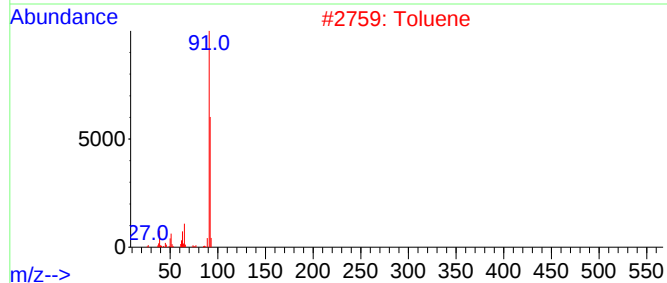
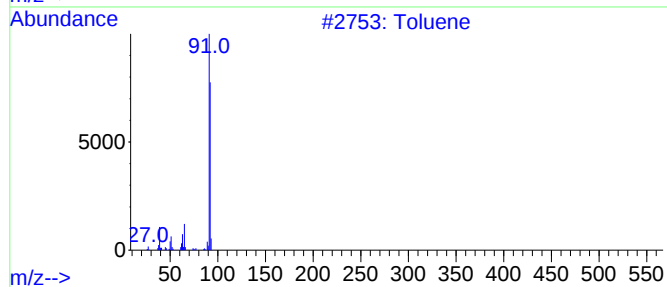
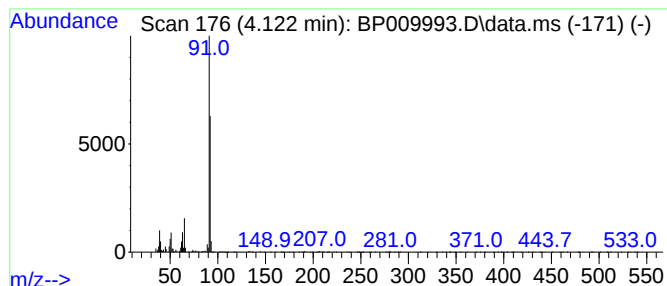
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Toluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.122	2.42 ng/ul	89337	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	91
2	Toluene			92	C7H8	000108-88-3	91
3	Toluene			92	C7H8	000108-88-3	91
4	Toluene			92	C7H8	000108-88-3	90
5	Toluene			92	C7H8	000108-88-3	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	4.122	2.4	ng/u1	89337	1	7.946	739689	20.0