

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
 Data File : BP019342.D
 Acq On : 12 Jan 2024 22:04
 Operator : MA/JU
 Sample : P1091-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AZ7

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-BP010824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019342.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	18	22	30	rVB	24530	32530	0.66%	0.084%
2	3.640	90	94	106	rBV	143453	232763	4.76%	0.600%
3	3.864	128	132	137	rVB3	5526	10291	0.21%	0.027%
4	3.958	143	148	153	rBV	51007	69815	1.43%	0.180%
5	4.875	300	304	309	rBV7	3473	6821	0.14%	0.018%
6	6.969	654	660	668	rBV	114727	187342	3.83%	0.483%
7	7.122	679	686	697	rVB	481790	747284	15.27%	1.928%
8	7.311	711	718	729	rBV	444032	712769	14.56%	1.839%
9	7.405	729	734	744	rVV	19404	37873	0.77%	0.098%
10	7.775	791	797	806	rBV	415354	666643	13.62%	1.720%
11	8.505	913	921	934	rBV	362272	593318	12.12%	1.531%
12	8.616	936	940	947	rVB4	7648	14865	0.30%	0.038%
13	8.946	989	996	1007	rBV	630446	1014792	20.74%	2.618%
14	9.287	1048	1054	1058	rBV9	4839	10023	0.20%	0.026%
15	9.663	1108	1118	1130	rBV	480121	805159	16.45%	2.077%
16	10.205	1201	1210	1227	rBV	702979	1215317	24.83%	3.135%
17	10.575	1265	1273	1281	rVB	570531	962897	19.68%	2.484%
18	10.728	1291	1299	1313	rBV	662917	1214374	24.81%	3.133%
19	12.175	1539	1545	1553	rBV2	10022	19471	0.40%	0.050%
20	13.322	1735	1740	1748	rVB	17426	27896	0.57%	0.072%
21	13.846	1822	1829	1840	rBV	1486215	2089560	42.70%	5.390%
22	14.116	1868	1875	1887	rBV	1620708	2340855	47.83%	6.038%
23	14.422	1920	1927	1938	rBV	986361	1417451	28.96%	3.656%
24	14.669	1964	1969	1984	rBV	69309	166516	3.40%	0.430%
25	14.875	2000	2004	2009	rVB8	6268	9079	0.19%	0.023%
26	15.422	2087	2097	2107	rBV	2255600	3178369	64.95%	8.199%
27	15.557	2114	2120	2134	rBV	560250	843564	17.24%	2.176%
28	15.663	2134	2138	2144	rVB5	10566	16460	0.34%	0.042%
29	16.210	2226	2231	2236	rVB6	7985	13696	0.28%	0.035%
30	16.857	2337	2341	2346	rVB7	4880	7352	0.15%	0.019%
31	16.916	2346	2351	2355	rBV6	4736	8564	0.17%	0.022%
32	17.181	2389	2396	2406	rBV	1402813	1942696	39.70%	5.011%
33	17.281	2406	2413	2428	rVV	2833766	3921471	80.13%	10.116%
34	18.075	2544	2548	2554	rBV8	14861	27837	0.57%	0.072%
35	18.486	2615	2618	2622	rVB5	6720	8682	0.18%	0.022%
36	19.098	2718	2722	2728	rBV	39668	53204	1.09%	0.137%

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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-BP010824.MA.M
Title : SVOA CALIBRATION

37	19.169	2731	2734	2739	rBV	33835	47068	0.96%	0.121%
38	19.316	2756	2759	2764	rBV5	14590	19710	0.40%	0.051%
39	19.457	2780	2783	2787	rBV4	12466	15685	0.32%	0.040%
40	19.528	2789	2795	2804	rVV	3723392	4893860	100.00%	12.624%
41	21.286	3089	3094	3101	rBV	1882333	2221925	45.40%	5.732%
42	23.492	3461	3469	3479	rVB2	2401309	4744712	96.95%	12.240%
43	23.639	3487	3494	3508	rVB	1087519	2194991	44.85%	5.662%

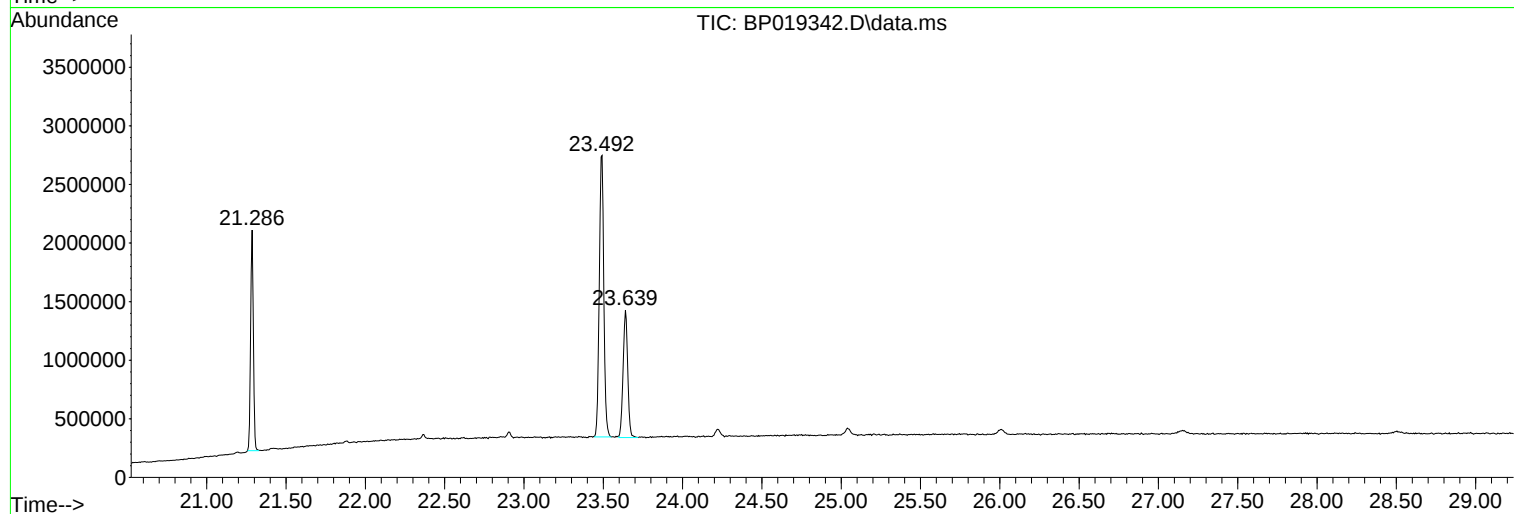
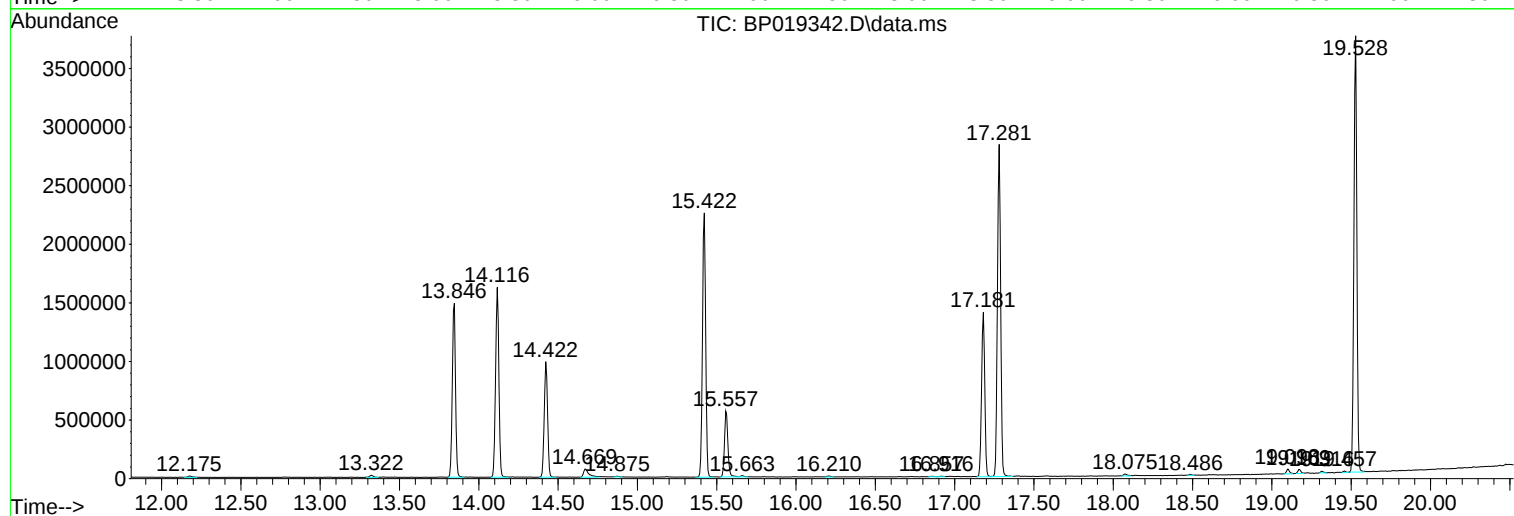
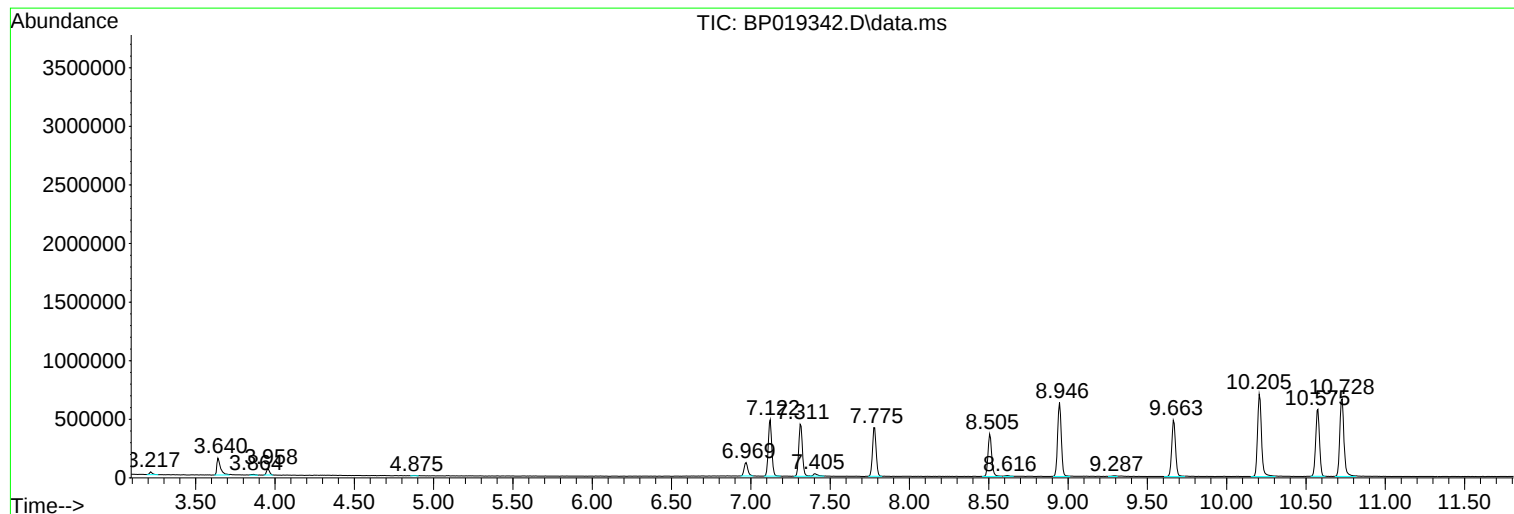
Sum of corrected areas: 38765550

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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



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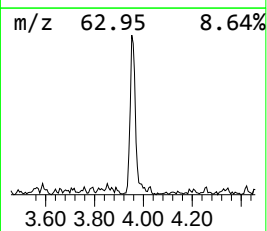
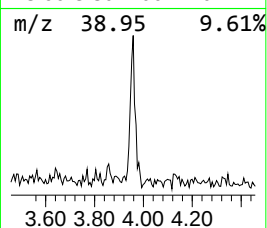
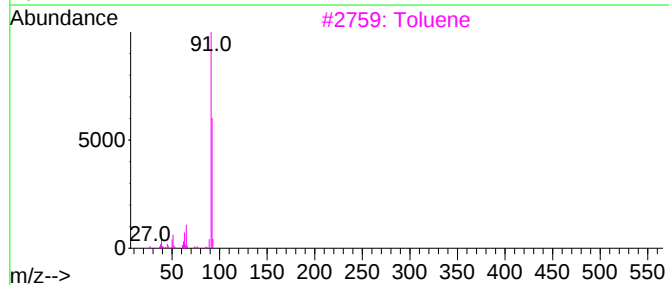
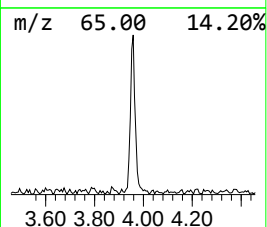
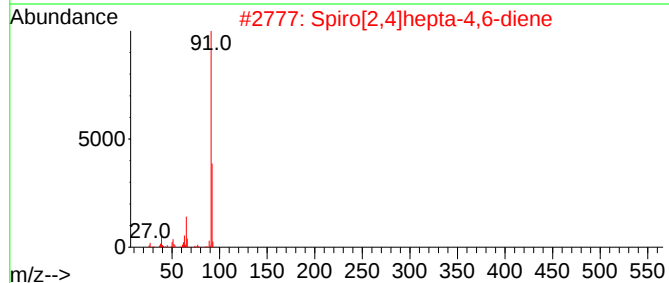
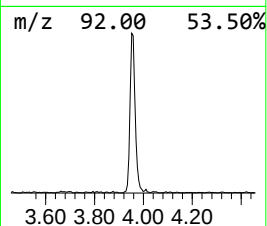
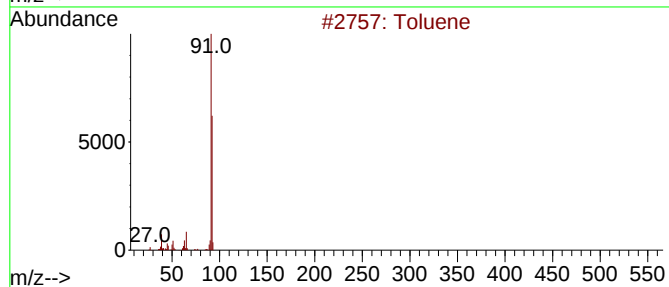
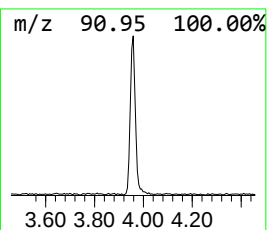
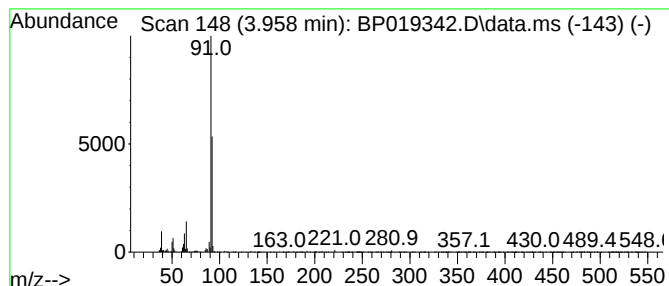
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Toluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.958	2.09 ng/ul	69815	1,4-Dichlorobenzene-d4	7.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	87
2			Spiro[2,4]hepta-4,6-diene	92	C7H8	000765-46-8	87
3			Toluene	92	C7H8	000108-88-3	83
4			Toluene	92	C7H8	000108-88-3	83
5			Toluene	92	C7H8	000108-88-3	76



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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	3.958	2.1	ng/u1	69815	1	7.781	666643	20.0