

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\
 Data File : BM016952.D
 Acq On : 02 Oct 2018 09:40
 Operator : MJ/SJ
 Sample : PB113341BL
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK41

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	30	34	39	rVB	689963	705238	20.74%	3.831%
2	4.864	316	319	323	rVB2	35963	49178	1.45%	0.267%
3	7.710	796	803	811	rVB	1225179	1930072	56.76%	10.484%
4	8.034	853	858	863	rVB5	4596	9117	0.27%	0.050%
5	10.487	1267	1275	1285	rBV	1604161	2639098	77.62%	14.335%
6	14.351	1925	1932	1940	rBV2	2052956	3119443	91.74%	16.944%
7	15.345	2097	2101	2106	rBV3	6656	9540	0.28%	0.052%
8	16.080	2223	2226	2229	rBV2	2841	3704	0.11%	0.020%
9	16.298	2259	2263	2266	rBV3	4331	5612	0.17%	0.030%
10	16.880	2358	2362	2367	rBV4	6131	9389	0.28%	0.051%
11	17.098	2392	2399	2410	rBV	2249881	3210098	94.41%	17.436%
12	21.286	3106	3111	3119	rBV	2615593	3319687	97.63%	18.032%
13	23.515	3484	3490	3500	rVB2	1833268	3400238	100.00%	18.469%

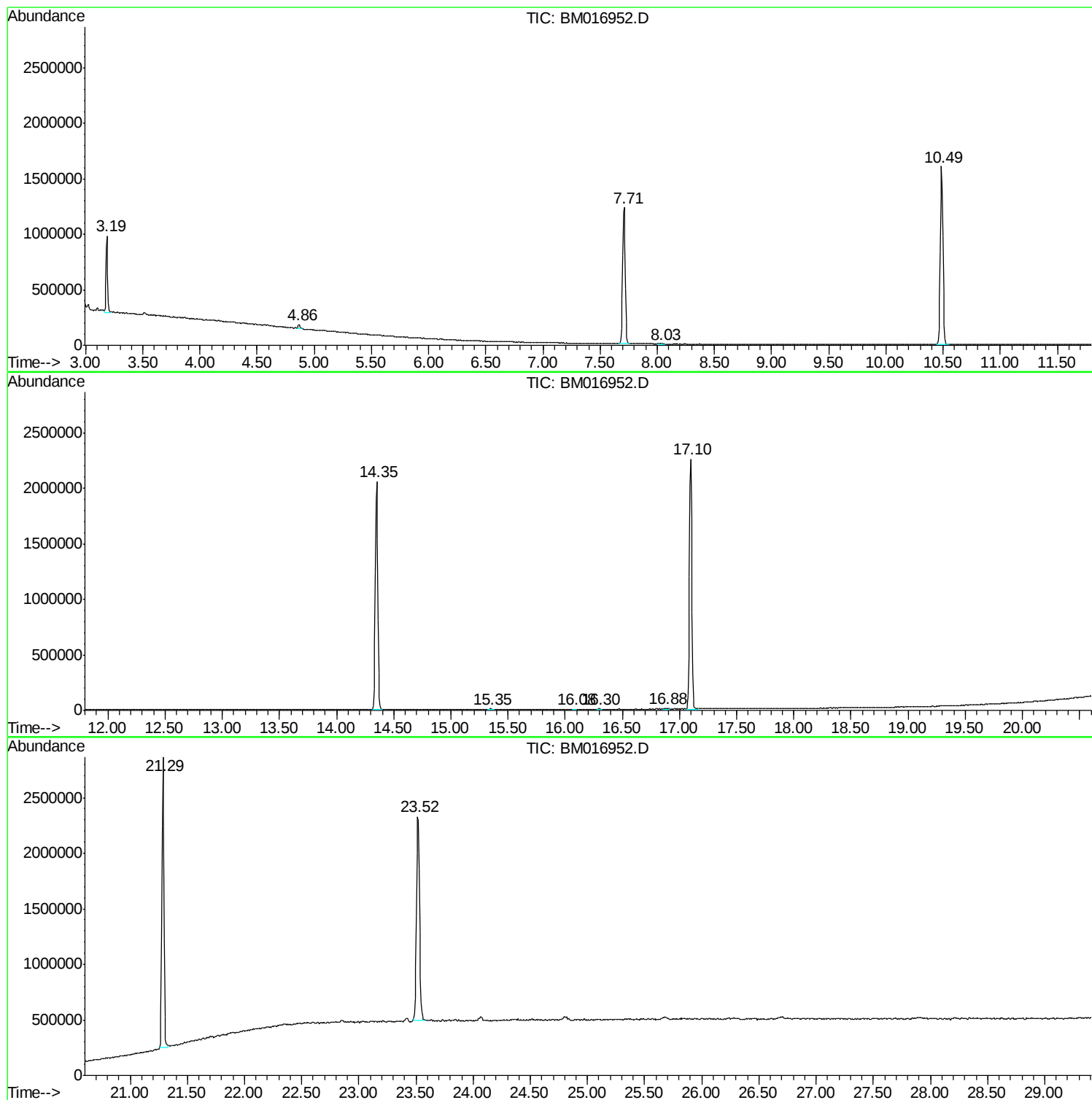
Sum of corrected areas: 18410414

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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ClientSampled :
SBLK41

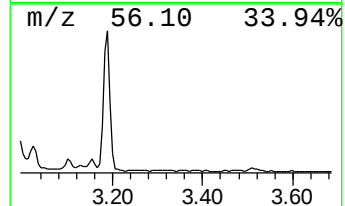
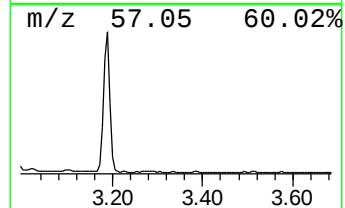
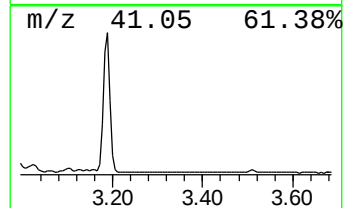
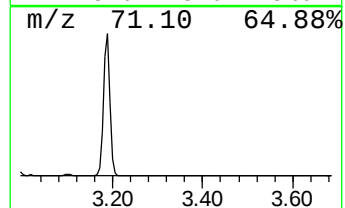
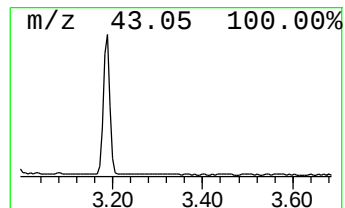
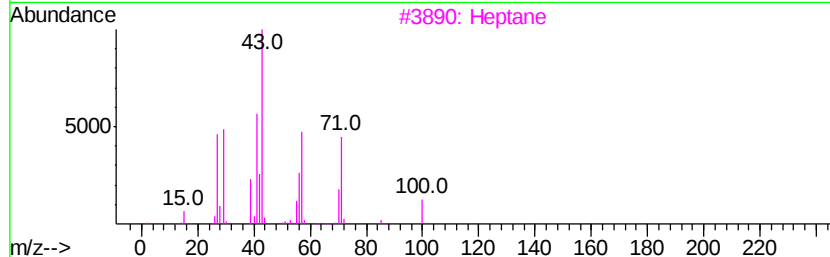
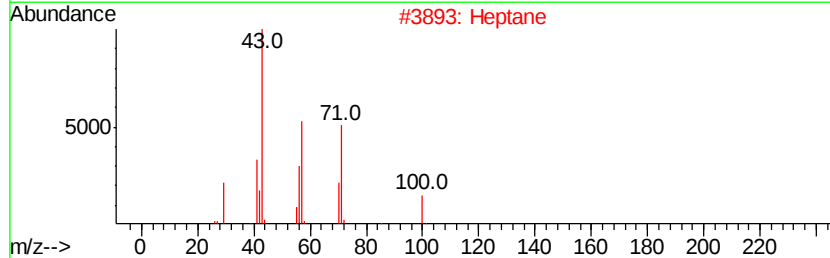
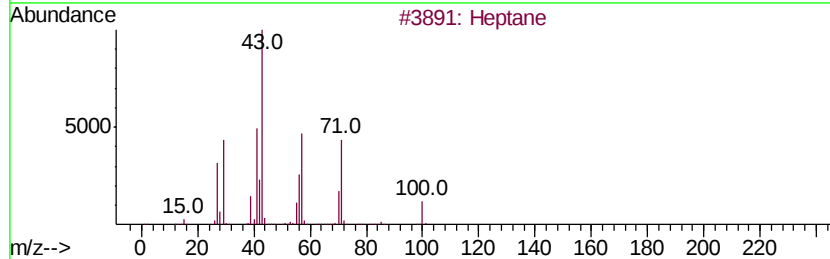
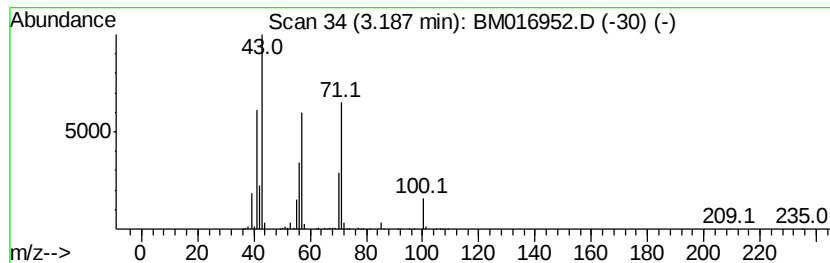
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.19	7.31 ng/ul	705238	1,4-Dichlorobenzene-d4	7.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane		100	C7H16	000142-82-5	91
2	Heptane		100	C7H16	000142-82-5	91
3	Heptane		100	C7H16	000142-82-5	91
4	Heptane		100	C7H16	000142-82-5	87
5	Hexane, 3-methyl-		100	C7H16	000589-34-4	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.19	7.3	ng/ul	705238	1	7.71	1930070	20.0