

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112918\
 Data File : BM017809.D
 Acq On : 29 Nov 2018 23:10
 Operator : JU/SJ
 Sample : J6041-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB4H4

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-BM111918MA.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	3.016	3	5	8	rBV	28373	28398	0.93%	0.156%
2	3.046	8	10	12	rVV2	17069	15856	0.52%	0.087%
3	3.069	12	14	16	rVV	26768	23360	0.76%	0.129%
4	3.104	16	20	30	rVB	618638	631756	20.62%	3.479%
5	3.228	38	41	43	rBV3	3276	3350	0.11%	0.018%
6	3.422	70	74	78	rVB2	16214	19557	0.64%	0.108%
7	3.869	148	150	154	rVB3	6774	8320	0.27%	0.046%
8	4.081	185	186	191	rVB3	3169	3521	0.11%	0.019%
9	5.269	383	388	389	rBV3	2267	3093	0.10%	0.017%
10	5.628	445	449	453	rVB5	2930	4084	0.13%	0.022%
11	5.981	506	509	515	rBV3	1707	3130	0.10%	0.017%
12	6.269	556	558	563	rVB3	2542	3492	0.11%	0.019%
13	7.204	712	717	721	rBV3	5202	7481	0.24%	0.041%
14	7.239	721	723	726	rVB2	3981	3745	0.12%	0.021%
15	7.587	774	782	797	rBV	941549	1617520	52.80%	8.907%
16	7.934	835	841	845	rVB4	2269	4804	0.16%	0.026%
17	8.634	956	960	962	rBV2	2073	3091	0.10%	0.017%
18	8.810	984	990	997	rVB6	2573	6371	0.21%	0.035%
19	10.357	1242	1253	1275	rBV	1313016	2374894	77.52%	13.078%
20	10.498	1275	1277	1281	rVV3	6331	9025	0.29%	0.050%
21	10.528	1281	1282	1287	rVB2	3472	3711	0.12%	0.020%
22	12.069	1540	1544	1545	rBV3	2407	3284	0.11%	0.018%
23	12.269	1568	1578	1583	rBV7	4941	16664	0.54%	0.092%
24	13.051	1710	1711	1716	rBV5	2940	4768	0.16%	0.026%
25	13.239	1740	1743	1746	rBV3	1961	3507	0.11%	0.019%
26	13.404	1768	1771	1777	rVB5	3341	6580	0.21%	0.036%
27	13.551	1793	1796	1802	rBV3	5458	12793	0.42%	0.070%
28	13.621	1802	1808	1809	rVV4	4122	8460	0.28%	0.047%
29	13.639	1809	1811	1815	rVB3	4748	5384	0.18%	0.030%
30	13.786	1830	1836	1837	rBV3	2627	3957	0.13%	0.022%
31	13.963	1862	1866	1869	rVV4	3627	4741	0.15%	0.026%
32	14.057	1877	1882	1886	rBV4	4774	6181	0.20%	0.034%
33	14.233	1906	1912	1922	rBV2	2007582	3008186	98.19%	16.565%
34	14.468	1946	1952	1954	rBV3	3444	5096	0.17%	0.028%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
 Title : SVOA CALIBRATION

35	14.533	1960	1963	1965	rBV3	4086	4443	0.15%	0.024%
36	14.657	1979	1984	1986	rBV3	6774	10040	0.33%	0.055%
37	14.721	1993	1995	2000	rVB4	5518	6200	0.20%	0.034%
38	14.857	2016	2018	2021	rBV4	5042	6534	0.21%	0.036%
39	14.939	2027	2032	2034	rBV5	5227	7267	0.24%	0.040%
40	15.204	2075	2077	2079	rVB3	4052	3144	0.10%	0.017%
41	15.251	2079	2085	2087	rBV5	6593	12646	0.41%	0.070%
42	15.298	2087	2093	2099	rVV5	18808	44611	1.46%	0.246%
43	15.462	2116	2121	2125	rBV6	6305	11560	0.38%	0.064%
44	15.504	2126	2128	2131	rVB4	3850	4247	0.14%	0.023%
45	15.610	2142	2146	2151	rVB5	9375	14730	0.48%	0.081%
46	15.674	2153	2157	2158	rBV4	3332	3670	0.12%	0.020%
47	15.751	2168	2170	2175	rBV4	6383	8657	0.28%	0.048%
48	15.839	2183	2185	2190	rVB6	5948	8049	0.26%	0.044%
49	15.915	2195	2198	2201	rBV5	3109	4491	0.15%	0.025%
50	15.974	2203	2208	2210	rBV5	11791	19812	0.65%	0.109%
51	16.110	2227	2231	2234	rBV6	5239	9029	0.29%	0.050%
52	16.292	2256	2262	2268	rBV9	18993	40647	1.33%	0.224%
53	16.533	2301	2303	2304	rBV2	5187	4460	0.15%	0.025%
54	16.815	2347	2351	2354	rVB4	14562	20388	0.67%	0.112%
55	16.880	2359	2362	2363	rBV3	6387	5216	0.17%	0.029%
56	16.898	2363	2365	2367	rBV3	5183	5022	0.16%	0.028%
57	16.992	2374	2381	2386	rBV	2006600	3063633	100.00%	16.871%
58	17.033	2386	2388	2397	rVB	275426	369198	12.05%	2.033%
59	17.239	2421	2423	2425	rBV3	7383	6632	0.22%	0.037%
60	17.374	2443	2446	2449	rVB5	23408	24707	0.81%	0.136%
61	17.492	2464	2466	2469	rBV4	13928	16675	0.54%	0.092%
62	17.709	2499	2503	2509	rBV4	59701	83300	2.72%	0.459%
63	17.786	2511	2516	2521	rBV	126731	176886	5.77%	0.974%
64	17.874	2527	2531	2535	rBV2	37373	64862	2.12%	0.357%
65	17.921	2536	2539	2545	rVB	40530	66826	2.18%	0.368%
66	18.062	2559	2563	2567	rBV4	49548	84719	2.77%	0.467%
67	18.098	2567	2569	2570	rBV2	19424	13799	0.45%	0.076%
68	18.168	2577	2581	2592	rVB3	121516	184930	6.04%	1.018%
69	19.062	2728	2733	2737	rBV	246393	387944	12.66%	2.136%
70	19.921	2876	2879	2882	rVB4	144007	167738	5.48%	0.924%
71	21.192	3091	3095	3105	rVB	1829620	2870631	93.70%	15.808%

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Stop Thrs : 0 Peak Location: TOP

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72	23.380	3462	3467	3476	rVB	1421100	2454036	80.10%	13.514%
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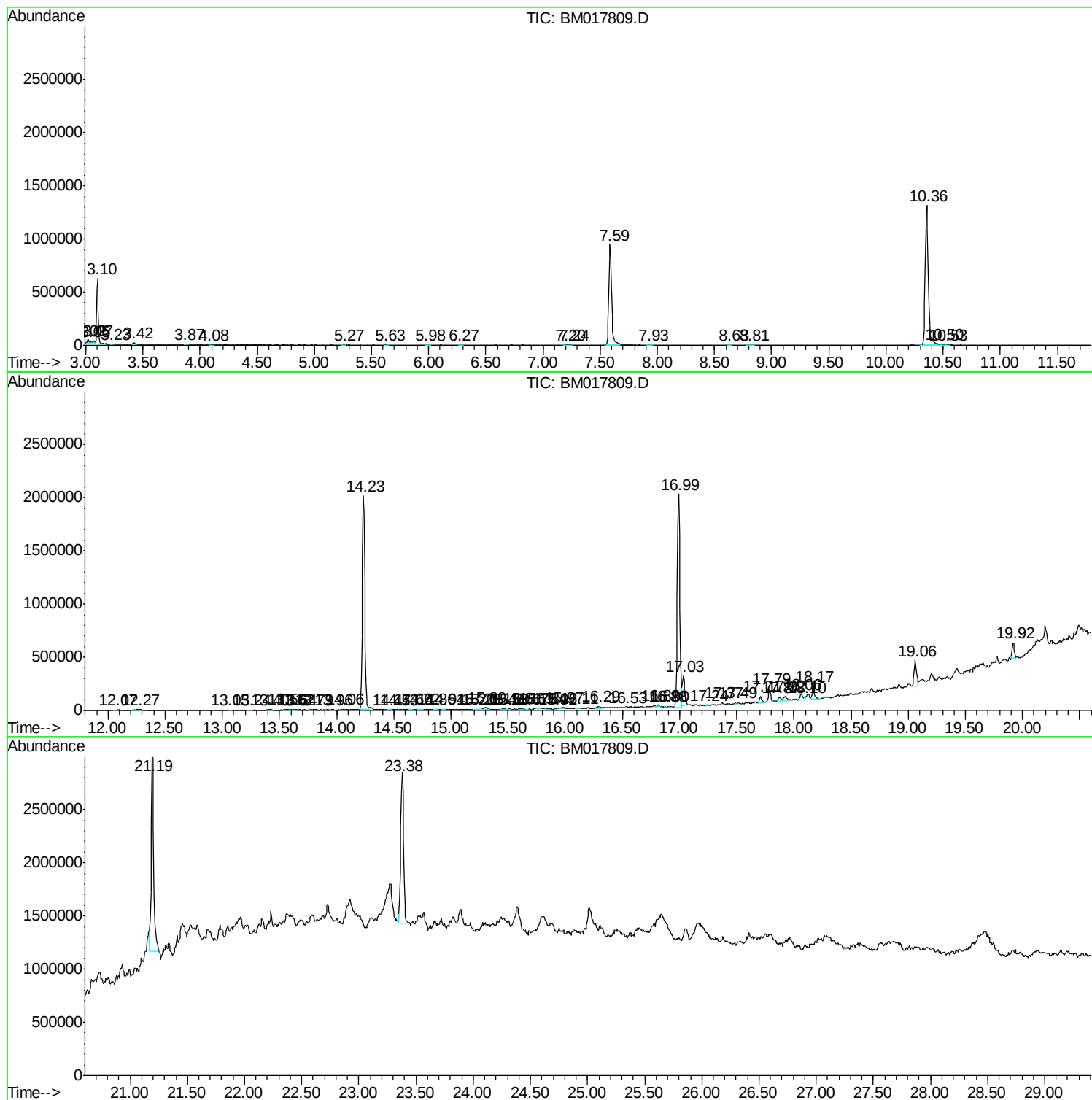
Sum of corrected areas: 18159539

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_M
ClientSampleId :
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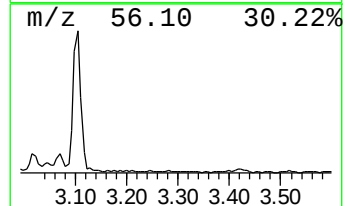
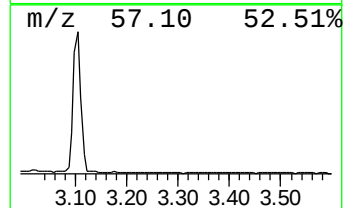
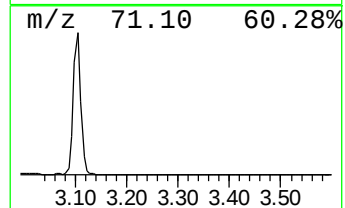
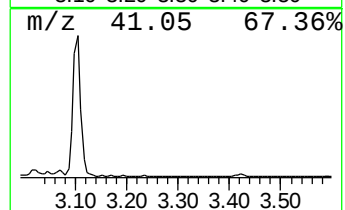
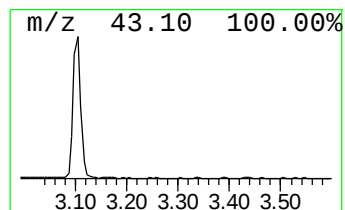
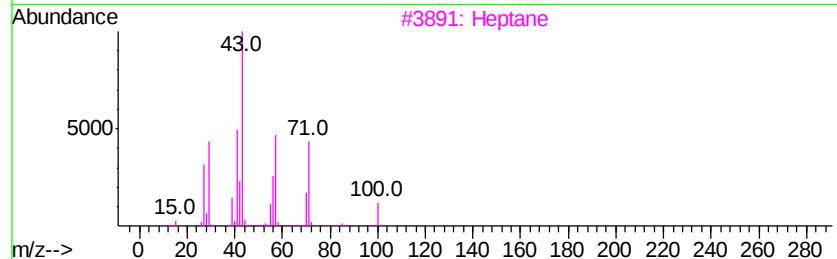
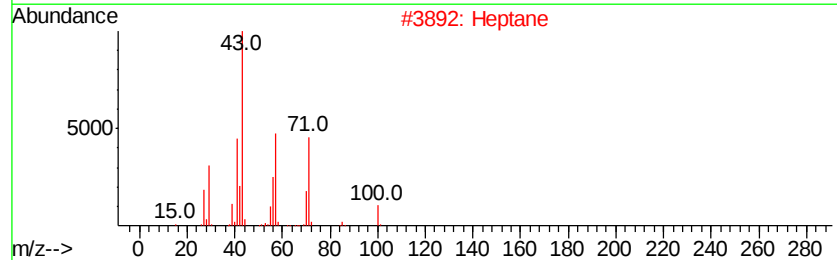
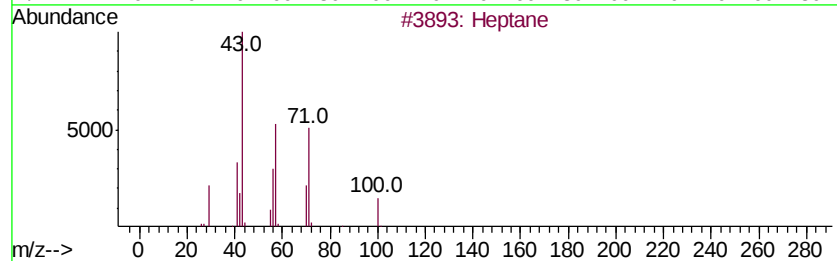
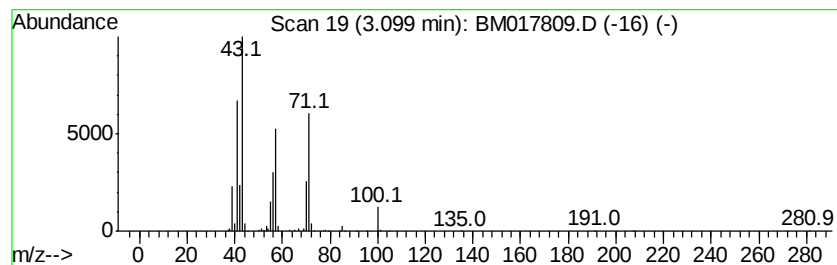
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.10	7.81 ng/ul	631756	1,4-Dichlorobenzene-d4	7.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	91
2	Heptane			100	C7H16	000142-82-5	90
3	Heptane			100	C7H16	000142-82-5	90
4	Heptane			100	C7H16	000142-82-5	86
5	Hexane, 3-methyl-			100	C7H16	000589-34-4	42



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.10	7.8	ng/ul	631756	1	7.59	1617520	20.0