

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041521\
 Data File : BM029487.D
 Acq On : 15 Apr 2021 05:13
 Operator : CG/JU
 Sample : PB135575BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK75

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-BM040721MA.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	2.934	6	8	15	rVB	80212	81996	5.39%	0.912%
2	3.010	19	21	24	rVV3	21380	21420	1.41%	0.238%
3	3.040	24	26	28	rVV2	11048	10267	0.67%	0.114%
4	3.063	28	30	33	rVV3	13547	15369	1.01%	0.171%
5	3.104	33	37	46	rVB	1010482	1058372	69.57%	11.770%
6	3.434	89	93	98	rVB	28163	31716	2.08%	0.353%
7	3.899	168	172	180	rVB2	10605	15322	1.01%	0.170%
8	4.428	260	262	268	rVB5	1516	1780	0.12%	0.020%
9	4.846	329	333	338	rBV7	2504	4285	0.28%	0.048%
10	6.557	623	624	629	rVB5	1269	1682	0.11%	0.019%
11	7.575	793	797	800	rBV5	886	1769	0.12%	0.020%
12	7.675	807	814	828	rBV	525549	833888	54.81%	9.274%
13	8.651	977	980	985	rBV3	1601	2192	0.14%	0.024%
14	9.439	1111	1114	1117	rBV4	1301	1925	0.13%	0.021%
15	9.969	1200	1204	1209	rVB4	876	1597	0.10%	0.018%
16	10.457	1280	1287	1300	rBV	654599	1106440	72.73%	12.305%
17	14.316	1936	1943	1953	rBV	944516	1387330	91.19%	15.429%
18	16.845	2369	2373	2378	rVB4	2942	4690	0.31%	0.052%
19	17.057	2402	2409	2419	rBV	1074432	1521350	100.00%	16.919%
20	17.157	2423	2426	2428	rVB4	2501	2732	0.18%	0.030%
21	17.421	2468	2471	2474	rVB2	1557	1918	0.13%	0.021%
22	17.721	2518	2522	2523	rBV4	1245	1526	0.10%	0.017%
23	17.980	2561	2566	2573	rVB8	1836	4883	0.32%	0.054%
24	18.039	2573	2576	2579	rBV4	1210	1849	0.12%	0.021%
25	18.221	2604	2607	2608	rVB3	1802	1603	0.11%	0.018%
26	18.256	2611	2613	2614	rBV2	2292	1682	0.11%	0.019%
27	18.362	2630	2631	2633	rBV2	1848	1592	0.10%	0.018%
28	18.539	2660	2661	2664	rBV3	1892	1687	0.11%	0.019%
29	18.568	2664	2666	2668	rBV3	1987	2188	0.14%	0.024%
30	18.927	2725	2727	2729	rBV3	2472	1553	0.10%	0.017%
31	18.962	2731	2733	2736	rBV4	2153	3127	0.21%	0.035%
32	19.392	2802	2806	2811	rBV	33421	42846	2.82%	0.476%
33	19.509	2823	2826	2829	rBV5	6740	7708	0.51%	0.086%
34	20.439	2981	2984	2988	rVB2	31075	31363	2.06%	0.349%

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35	20.503	2992	2995	2997	rBV2	16148	15468	1.02%	0.172%
36	21.233	3114	3119	3124	rBV	1161224	1449814	95.30%	16.123%
37	23.438	3488	3494	3505	rVB2	707620	1315015	86.44%	14.624%

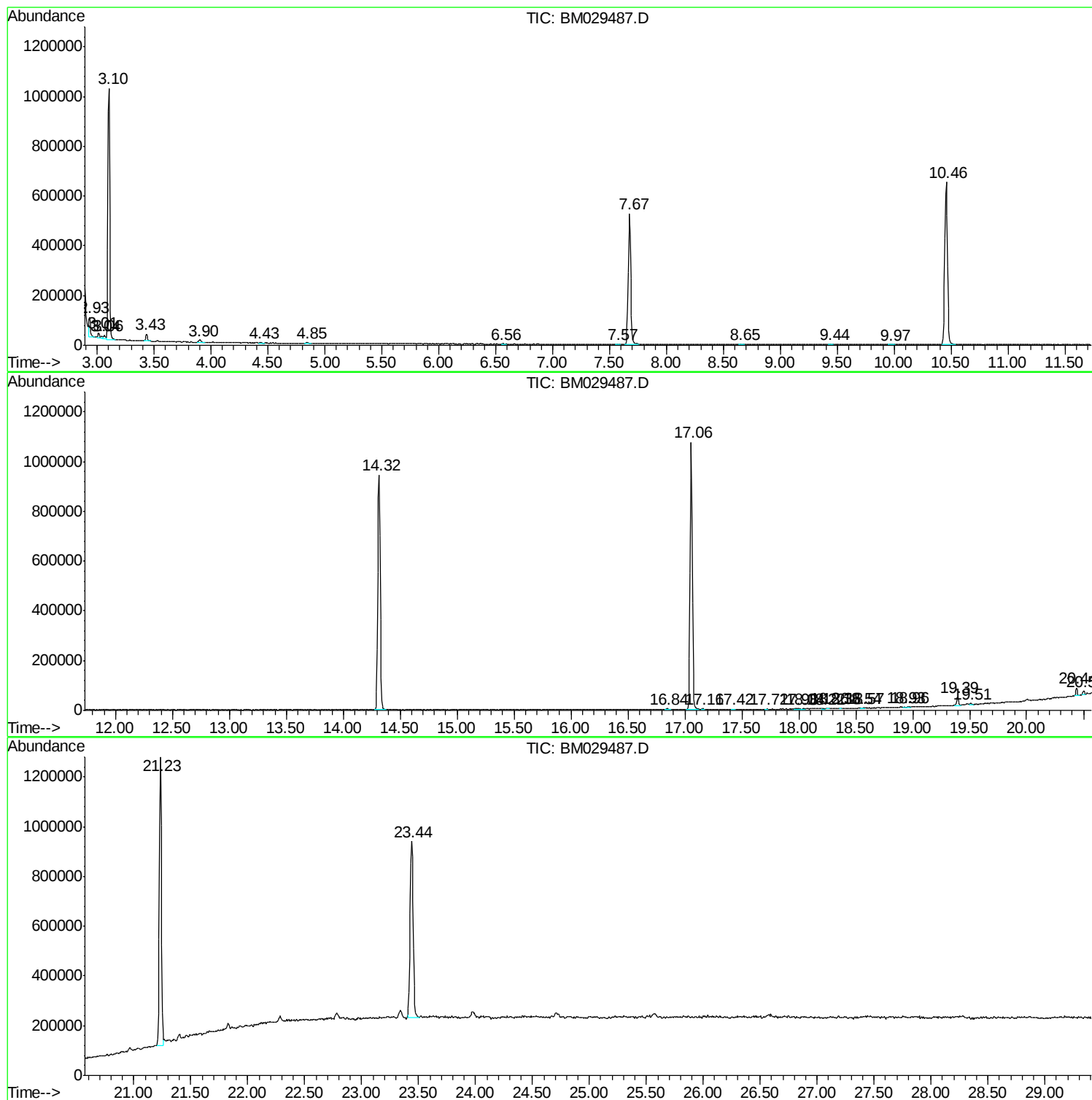
Sum of corrected areas: 8991944

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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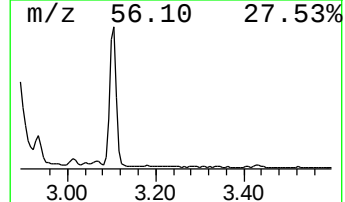
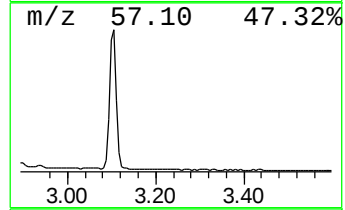
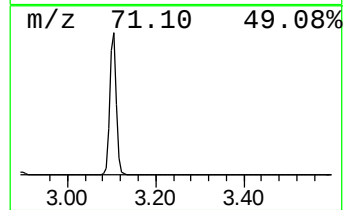
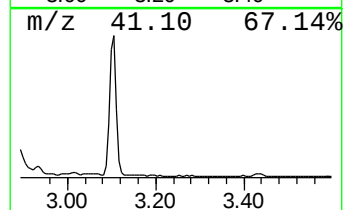
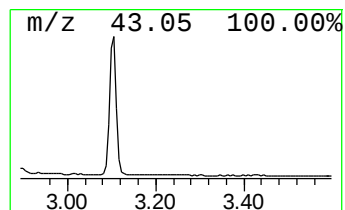
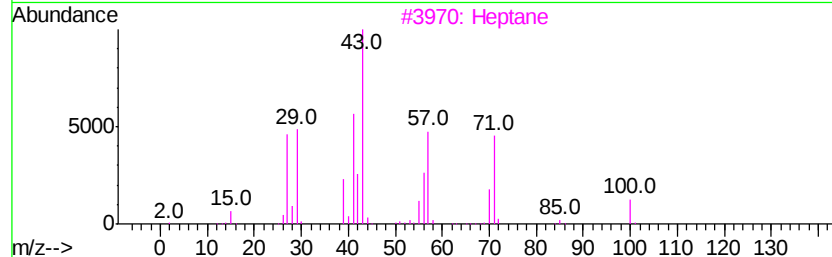
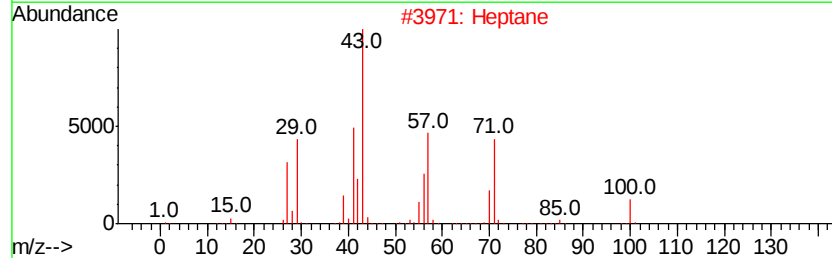
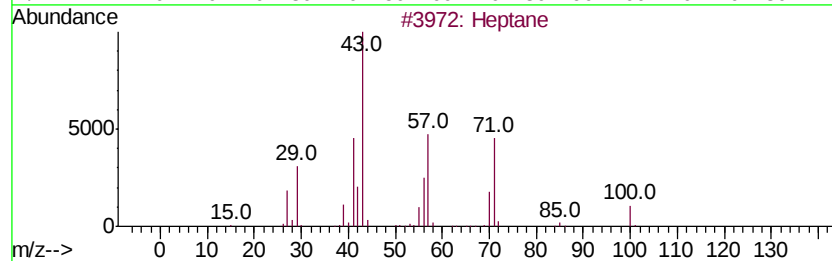
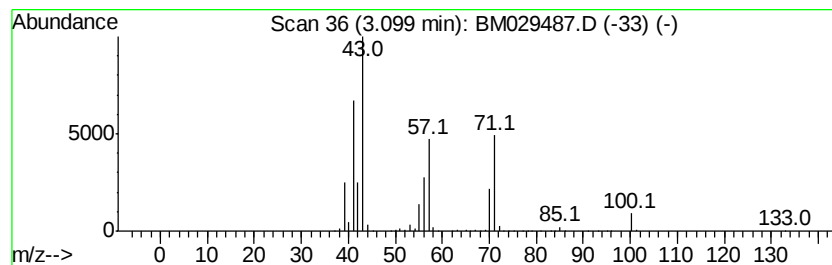
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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.10	25.38 ng/ul	1058370	1,4-Dichlorobenzene-d4	7.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Heptane	100	C7H16	000142-82-5	90
3		Heptane	100	C7H16	000142-82-5	87
4		Heptane	100	C7H16	000142-82-5	80
5		Hexane, 3-methyl-	100	C7H16	000589-34-4	45



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.10	25.4	ng/ul	1058370	1	7.67	833888	20.0