

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
 Data File : BM026337.D
 Acq On : 10 Jun 2020 19:28
 Operator : JU/CG
 Sample : PB129356BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK56

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-BM051320MA.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	2.922	3	6	13	rVB	216286	200915	11.64%	2.296%
2	3.240	56	60	64	rVB3	12539	14170	0.82%	0.162%
3	3.552	112	113	118	rBV5	2415	2484	0.14%	0.028%
4	3.699	134	138	143	rVB3	11478	15381	0.89%	0.176%
5	4.046	195	197	201	rVB4	2984	2033	0.12%	0.023%
6	4.228	224	228	229	rBV3	1828	2548	0.15%	0.029%
7	4.328	241	245	248	rBV5	3829	4568	0.26%	0.052%
8	4.628	293	296	301	rVB6	2540	4852	0.28%	0.055%
9	4.722	311	312	315	rVB3	1905	1809	0.10%	0.021%
10	4.822	326	329	333	rVB5	1893	2846	0.16%	0.033%
11	4.928	344	347	349	rBV4	1728	1972	0.11%	0.023%
12	5.075	370	372	375	rBV5	2049	2423	0.14%	0.028%
13	6.040	533	536	537	rBV3	1759	1813	0.11%	0.021%
14	6.910	682	684	689	rVB4	1362	2078	0.12%	0.024%
15	7.416	763	770	781	rBV	628916	982288	56.92%	11.224%
16	7.593	798	800	803	rBV4	1915	2408	0.14%	0.028%
17	8.004	866	870	872	rBV3	1430	1814	0.11%	0.021%
18	8.698	985	988	991	rVB4	1332	1873	0.11%	0.021%
19	8.940	1026	1029	1031	rVV3	1667	2115	0.12%	0.024%
20	9.034	1040	1045	1048	rBV4	1427	1968	0.11%	0.022%
21	9.498	1121	1124	1126	rBV2	1270	1766	0.10%	0.020%
22	9.845	1178	1183	1185	rBV4	1098	1835	0.11%	0.021%
23	10.181	1232	1240	1252	rBV	771991	1268560	73.50%	14.496%
24	10.516	1293	1297	1300	rBV4	1726	2705	0.16%	0.031%
25	10.669	1317	1323	1328	rBV5	1337	3250	0.19%	0.037%
26	11.951	1539	1541	1544	rBV4	1573	1842	0.11%	0.021%
27	12.928	1704	1707	1708	rBV3	1669	1900	0.11%	0.022%
28	14.075	1895	1902	1922	rBV	1055956	1543200	89.42%	17.634%
29	15.774	2187	2191	2193	rBV3	1470	2194	0.13%	0.025%
30	16.221	2266	2267	2272	rBV4	1306	1844	0.11%	0.021%
31	16.616	2327	2334	2340	rBV3	3658	7508	0.44%	0.086%
32	16.827	2363	2370	2384	rBV	1249618	1725840	100.00%	19.721%
33	16.921	2384	2386	2395	rVB7	5177	12612	0.73%	0.144%
34	17.174	2425	2429	2431	rVB4	1891	2416	0.14%	0.028%

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35	17.292	2446	2449	2451	rBV4	1941	2183	0.13%	0.025%
36	17.345	2453	2458	2459	rBV3	1682	1899	0.11%	0.022%
37	17.363	2459	2461	2465	rBV5	1969	2196	0.13%	0.025%
38	17.586	2497	2499	2502	rVB4	1944	2001	0.12%	0.023%
39	17.933	2555	2558	2559	rBV3	2151	2523	0.15%	0.029%
40	18.262	2611	2614	2617	rBV5	2124	2671	0.15%	0.031%
41	18.427	2641	2642	2644	rBV2	1829	1739	0.10%	0.020%
42	21.033	3080	3085	3095	rBV	1197735	1530140	88.66%	17.485%
43	23.127	3436	3441	3452	rVB	792024	1376188	79.74%	15.725%

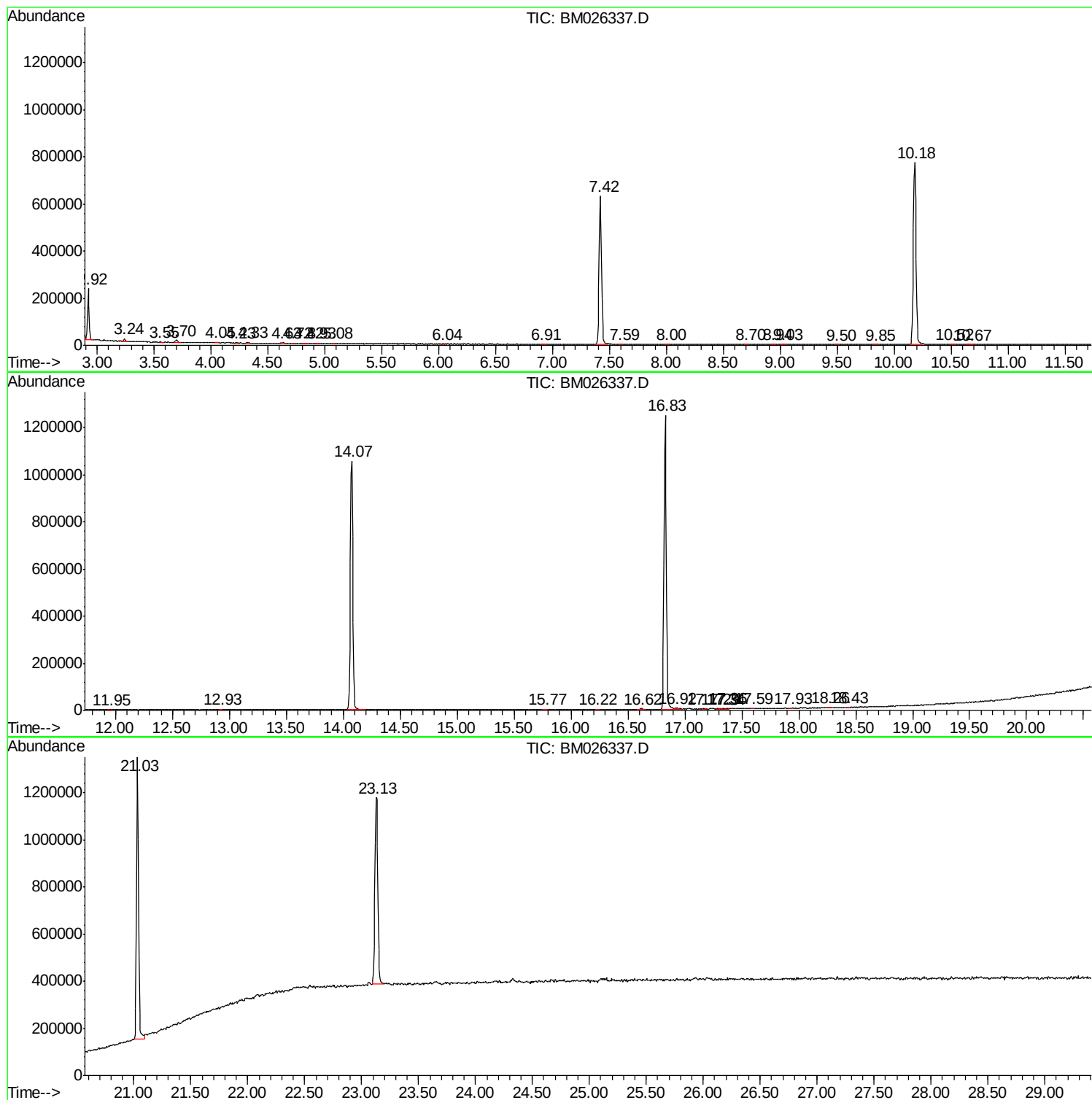
Sum of corrected areas: 8751370

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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



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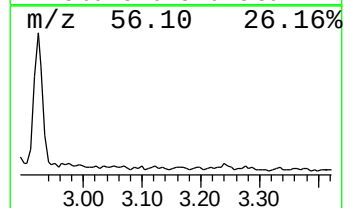
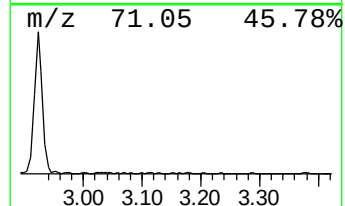
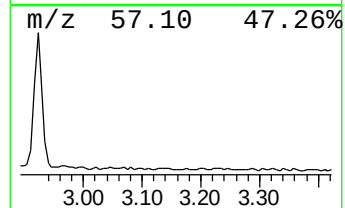
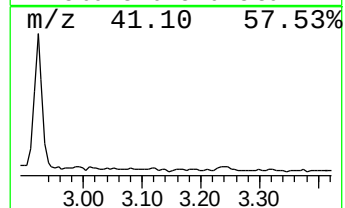
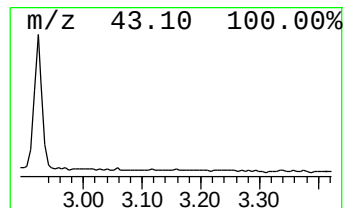
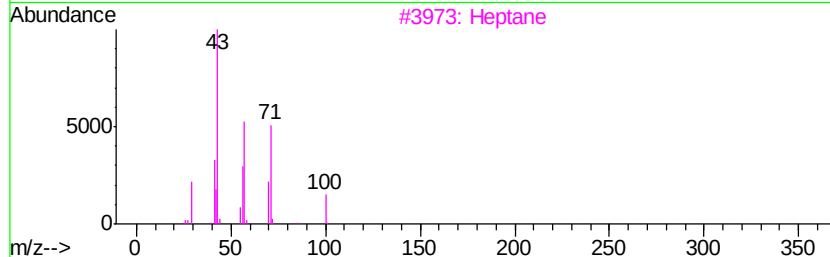
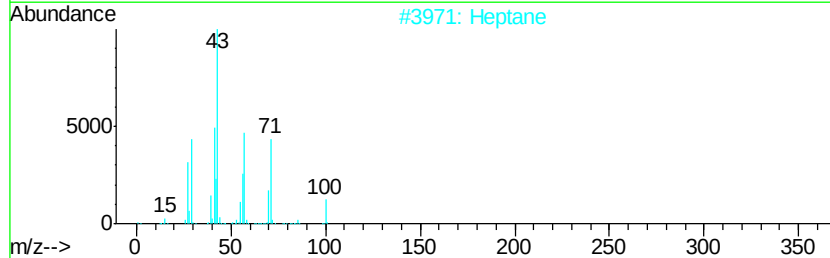
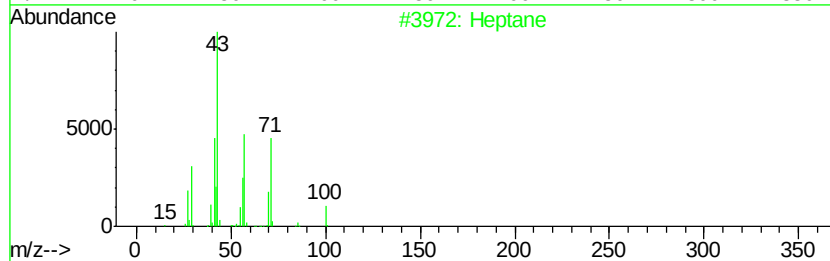
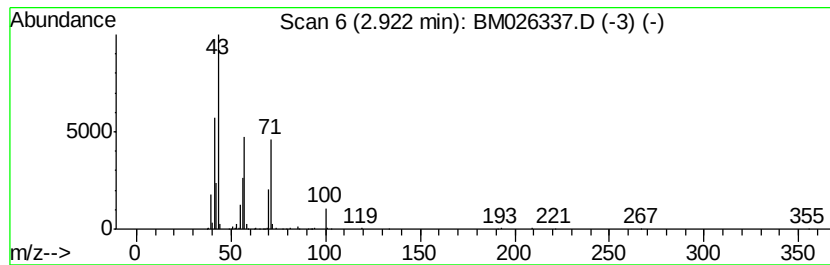
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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	4.09 ng/ul	200915	1,4-Dichlorobenzene-d4	7.42

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	78
5		Morpholine, 2,6-dimethyl-	115	C6H13NO	000141-91-3	42



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	2.92	4.1	ng/ul	200915	1	7.42	982288	20.0