

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101720\
 Data File : BN012336.D
 Acq On : 18 Oct 2020 18:59
 Operator : CG/JU
 Sample : PB132369BL
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK69

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_N\METHODS\SOM-EPA-BN101720.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	3.023	4	6	8	rVV2	17508	14630	0.55%	0.100%
2	3.058	8	12	22	rVB	1050480	1053938	39.74%	7.197%
3	3.128	22	24	26	rVB3	4915	3499	0.13%	0.024%
4	3.193	33	35	39	rVB4	4588	4187	0.16%	0.029%
5	3.305	52	54	55	rBV2	3342	2699	0.10%	0.018%
6	3.387	65	68	71	rVB2	26068	30312	1.14%	0.207%
7	3.846	142	146	150	rVV	15234	20356	0.77%	0.139%
8	4.046	179	180	183	rVB3	3818	3551	0.13%	0.024%
9	4.105	187	190	195	rVB5	2339	3016	0.11%	0.021%
10	4.205	205	207	211	rVB5	3248	3047	0.11%	0.021%
11	4.287	218	221	226	rBV6	4612	8321	0.31%	0.057%
12	4.381	234	237	239	rBV3	2236	2688	0.10%	0.018%
13	4.422	242	244	247	rVB3	2995	2732	0.10%	0.019%
14	4.664	284	285	290	rVB4	4506	5932	0.22%	0.041%
15	4.711	290	293	296	rBV4	3854	5168	0.19%	0.035%
16	4.764	299	302	310	rVB6	11263	15858	0.60%	0.108%
17	5.211	375	378	382	rVB2	3694	3961	0.15%	0.027%
18	5.275	387	389	392	rBV2	2816	2920	0.11%	0.020%
19	5.405	408	411	414	rVB4	4410	4394	0.17%	0.030%
20	5.952	502	504	505	rVB2	4588	2678	0.10%	0.018%
21	5.987	509	510	513	rVV2	3041	2778	0.10%	0.019%
22	6.158	537	539	542	rVB4	3240	3417	0.13%	0.023%
23	6.246	551	554	556	rBV4	3332	3708	0.14%	0.025%
24	6.287	559	561	565	rVV4	2714	3898	0.15%	0.027%
25	6.399	575	580	585	rBV4	3366	6549	0.25%	0.045%
26	6.669	622	626	627	rBV3	2759	3391	0.13%	0.023%
27	6.699	630	631	636	rVV2	3136	5049	0.19%	0.034%
28	6.752	636	640	643	rVV3	4303	6652	0.25%	0.045%
29	6.781	643	645	650	rVB3	2644	3507	0.13%	0.024%
30	7.540	769	774	775	rBV3	2395	3075	0.12%	0.021%
31	7.593	777	783	791	rBV	970461	1489865	56.18%	10.174%
32	8.052	858	861	864	rBV4	2428	3387	0.13%	0.023%
33	8.152	876	878	882	rBV2	3061	3335	0.13%	0.023%
34	8.275	896	899	901	rBV2	2953	2834	0.11%	0.019%

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35	8.481	922	934	936	rBV3	3221	7853	0.30%	0.054%
36	9.328	1074	1078	1080	rVB4	2034	2685	0.10%	0.018%
37	10.363	1247	1254	1265	rBV	1152177	1900741	71.67%	12.980%
38	10.881	1338	1342	1344	rVB4	2118	2841	0.11%	0.019%
39	12.228	1566	1571	1577	rVB2	2317	4016	0.15%	0.027%
40	13.498	1784	1787	1792	rBV3	1670	2784	0.10%	0.019%
41	14.093	1885	1888	1891	rBV3	2000	3095	0.12%	0.021%
42	14.240	1906	1913	1924	rBV	1594997	2364250	89.15%	16.145%
43	15.945	2201	2203	2206	rVB2	2908	3516	0.13%	0.024%
44	15.981	2206	2209	2211	rBV3	2600	3184	0.12%	0.022%
45	16.187	2243	2244	2249	rBV3	3063	3862	0.15%	0.026%
46	16.781	2339	2345	2349	rBV4	4909	9908	0.37%	0.068%
47	16.986	2373	2380	2393	rBV	1876271	2651914	100.00%	18.110%
48	17.086	2393	2397	2401	rVB5	5931	8781	0.33%	0.060%
49	17.339	2437	2440	2441	rBV2	3681	3132	0.12%	0.021%
50	17.398	2449	2450	2453	rBV2	2764	3038	0.11%	0.021%
51	17.451	2457	2459	2462	rBV4	3191	3543	0.13%	0.024%
52	19.016	2723	2725	2727	rBV2	3998	2893	0.11%	0.020%
53	19.133	2743	2745	2747	rBV3	5332	4887	0.18%	0.033%
54	21.186	3089	3094	3099	rBV	2032064	2496504	94.14%	17.048%
55	23.363	3458	3464	3472	rVB2	1371036	2424969	91.44%	16.560%

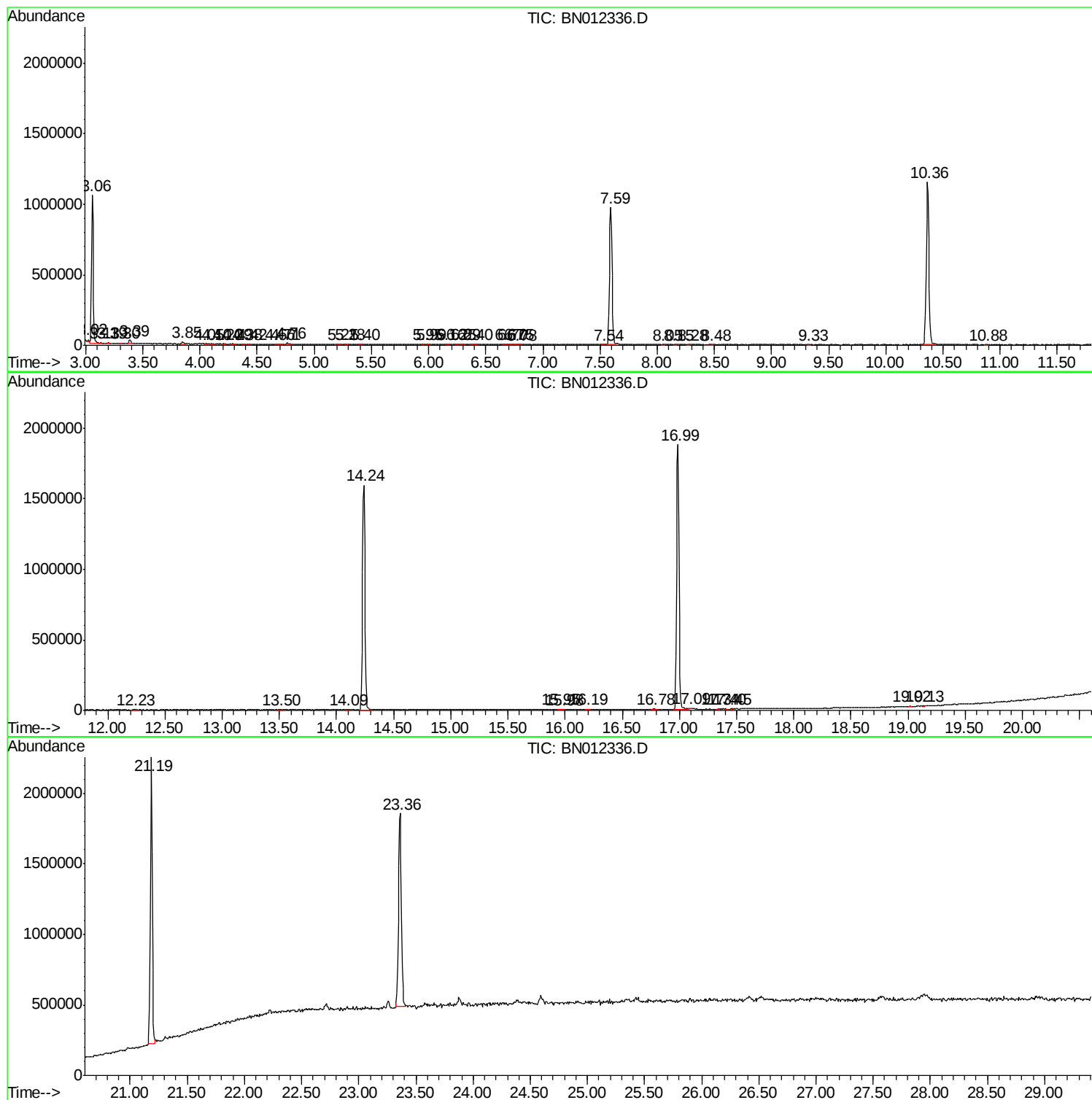
Sum of corrected areas: 14643728

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101720.M
Quant Title : SVOA CALIBRATION

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



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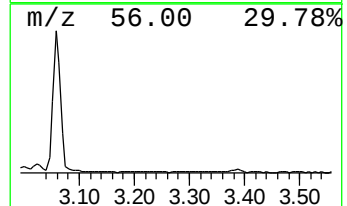
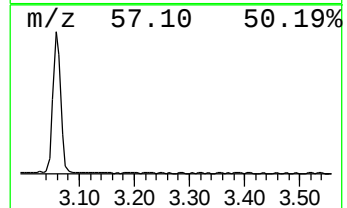
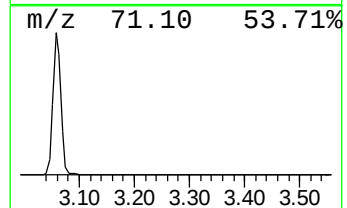
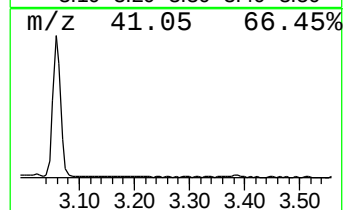
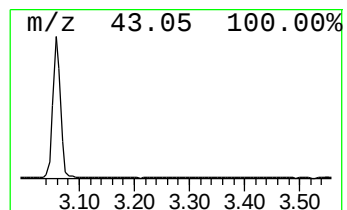
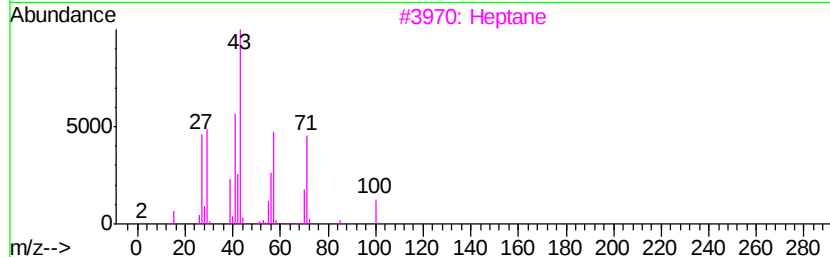
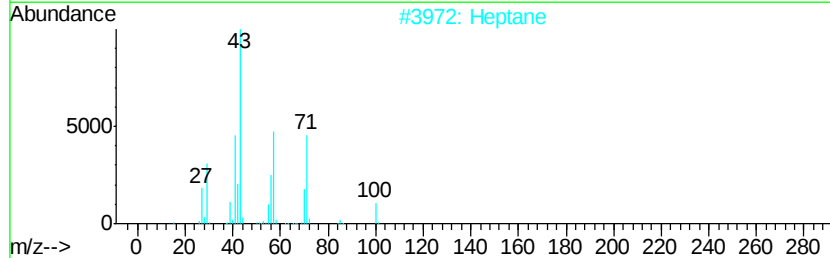
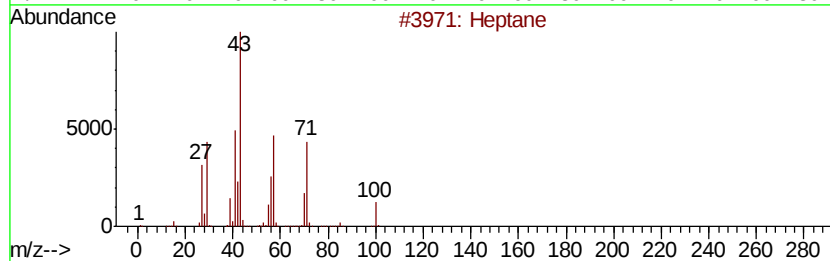
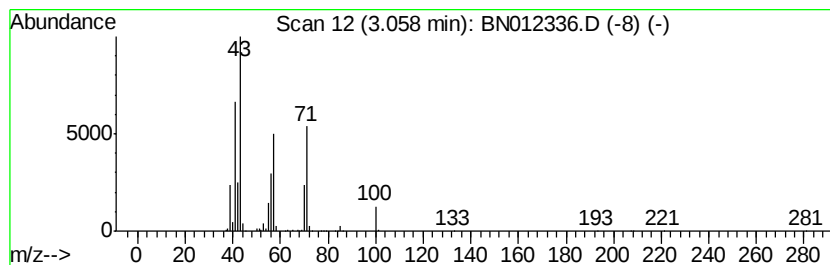
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	14.15 ng/ul	1053940	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	94
2		Heptane	100	C7H16	000142-82-5	94
3		Heptane	100	C7H16	000142-82-5	91
4		Heptane	100	C7H16	000142-82-5	83
5		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	45



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.06	14.2	ng/ul	1053940	1	7.59	1489870	20.0