

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034177.D
 Acq On : 09 Mar 2022 09:06
 Operator : CG/JU
 Sample : N1698-02
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGZR0

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034177.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.266	3	5	7	rVB2	21121	16877	0.87%	0.158%
2	3.307	7	12	16	rVB	1879845	1948048	100.00%	18.261%
3	3.619	63	65	67	rBV3	4238	5016	0.26%	0.047%
4	3.654	67	71	76	rVB	54972	68405	3.51%	0.641%
5	3.749	84	87	90	rVB3	3886	3882	0.20%	0.036%
6	4.137	149	153	158	rVB	14913	19913	1.02%	0.187%
7	4.596	226	231	237	rVB	10178	14915	0.77%	0.140%
8	5.601	396	402	403	rBV5	1317	2362	0.12%	0.022%
9	6.001	468	470	474	rVB4	1731	2038	0.10%	0.019%
10	7.142	661	664	667	rVB4	1826	2054	0.11%	0.019%
11	7.989	801	808	815	rBV	593681	939635	48.23%	8.808%
12	8.160	834	837	843	rVB5	1422	2274	0.12%	0.021%
13	9.154	1003	1006	1011	rVB5	1256	2602	0.13%	0.024%
14	10.801	1278	1286	1294	rBV	700923	1167337	59.92%	10.943%
15	10.942	1305	1310	1312	rBV3	2406	2382	0.12%	0.022%
16	14.624	1928	1936	1946	rBV	1063136	1516503	77.85%	14.216%
17	15.407	2066	2069	2071	rBV	1732	2134	0.11%	0.020%
18	15.871	2143	2148	2152	rBV3	9805	13869	0.71%	0.130%
19	16.795	2302	2305	2309	rVB2	1527	2053	0.11%	0.019%
20	16.901	2315	2323	2326	rBV3	14390	20188	1.04%	0.189%
21	17.024	2342	2344	2348	rBV3	1518	2511	0.13%	0.024%
22	17.148	2359	2365	2369	rVB5	4700	7074	0.36%	0.066%
23	17.248	2377	2382	2385	rBV4	2525	2949	0.15%	0.028%
24	17.283	2385	2388	2392	rVB4	6163	7433	0.38%	0.070%
25	17.365	2396	2402	2413	rVV	1202777	1682001	86.34%	15.767%
26	17.459	2416	2418	2422	rVB4	4254	5076	0.26%	0.048%
27	17.548	2429	2433	2436	rBV4	1538	2311	0.12%	0.022%
28	17.748	2464	2467	2473	rVB8	2846	4482	0.23%	0.042%
29	17.924	2493	2497	2500	rBV6	1929	2944	0.15%	0.028%
30	18.083	2520	2524	2529	rBV6	6074	8675	0.45%	0.081%
31	18.165	2535	2538	2542	rVB3	7756	10583	0.54%	0.099%
32	18.283	2556	2558	2562	rBV5	2157	3442	0.18%	0.032%
33	18.383	2573	2575	2576	rBV2	2610	2205	0.11%	0.021%
34	18.436	2582	2584	2587	rVB4	3768	3542	0.18%	0.033%
35	18.495	2591	2594	2597	rBV3	11329	11859	0.61%	0.111%
36	18.542	2597	2602	2605	rBV2	25753	32394	1.66%	0.304%

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37	18.689	2625	2627	2630	rBV3	2223	2198	0.11%	0.021%
38	19.048	2686	2688	2691	rVB3	6656	5547	0.28%	0.052%
39	19.183	2709	2711	2714	rBV4	3686	4972	0.26%	0.047%
40	19.242	2718	2721	2725	rVV6	4797	6754	0.35%	0.063%
41	19.559	2772	2775	2778	rVB4	7853	9010	0.46%	0.084%
42	19.624	2783	2786	2791	rVV7	13051	17071	0.88%	0.160%
43	19.683	2793	2796	2800	rVB3	22481	26460	1.36%	0.248%
44	20.271	2893	2896	2899	rBV4	12963	12842	0.66%	0.120%
45	20.542	2940	2942	2947	rVB5	22924	24614	1.26%	0.231%
46	21.012	3019	3022	3027	rVB6	27515	31780	1.63%	0.298%
47	21.536	3106	3111	3122	rBV	1156842	1509398	77.48%	14.149%
48	23.977	3520	3526	3536	rVB	715631	1475248	75.73%	13.829%

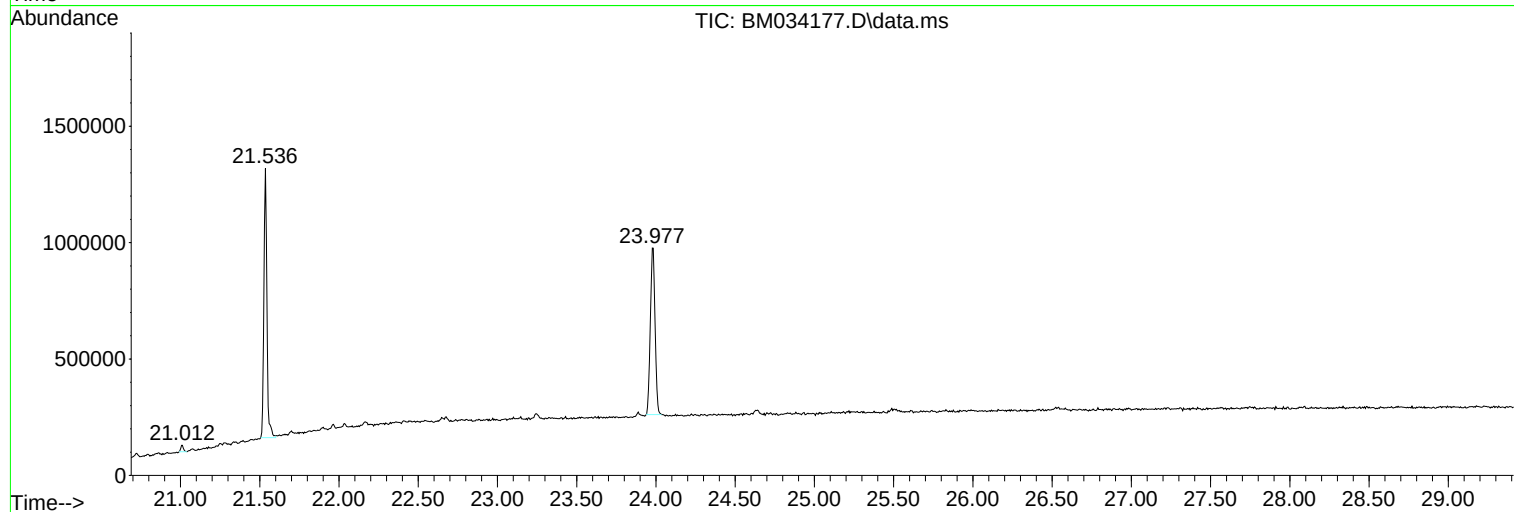
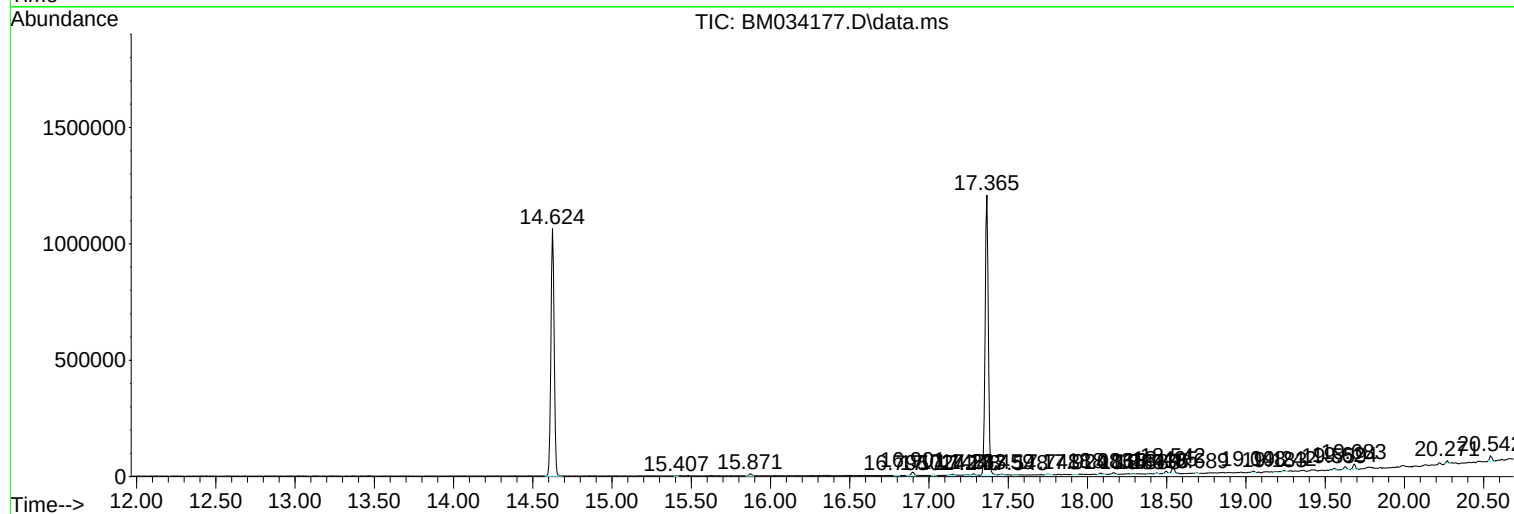
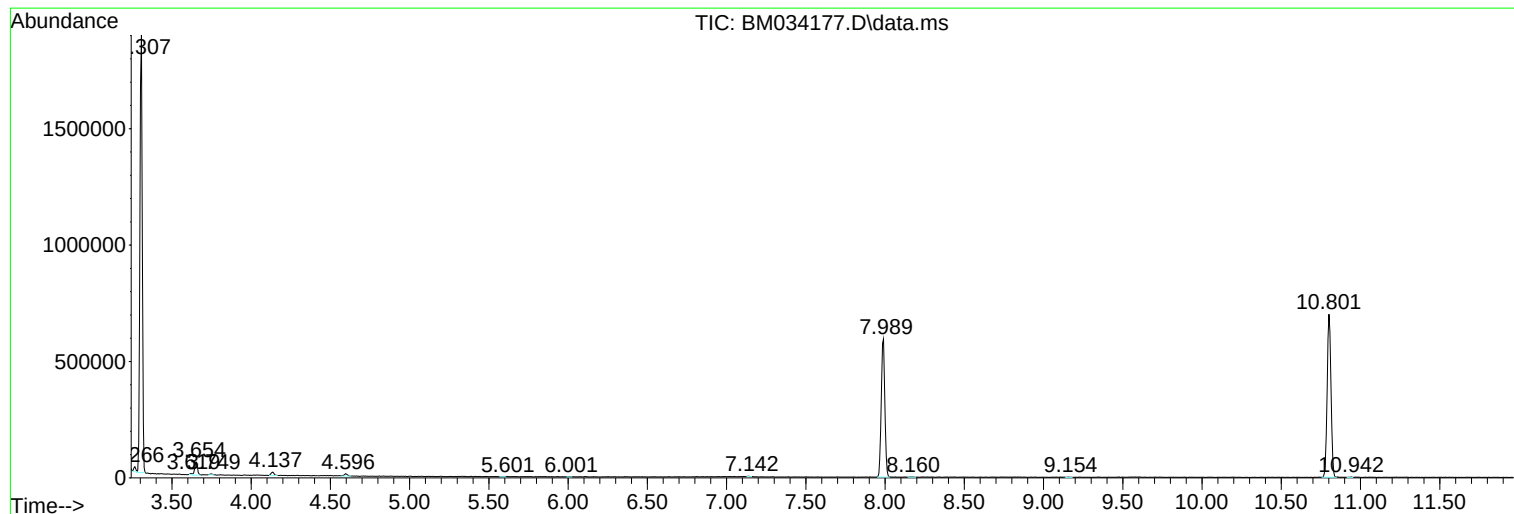
Sum of corrected areas: 10667862

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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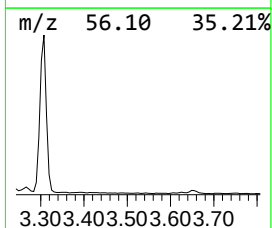
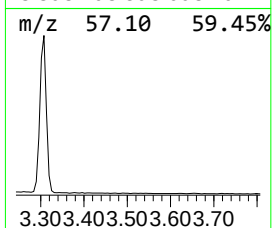
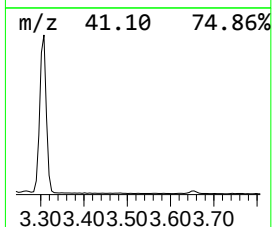
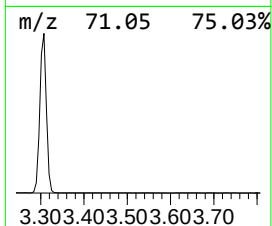
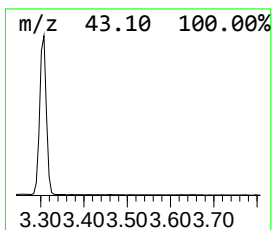
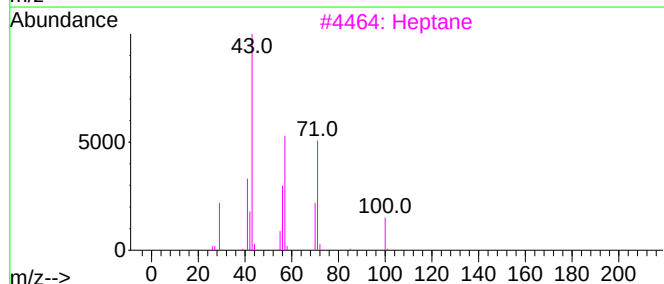
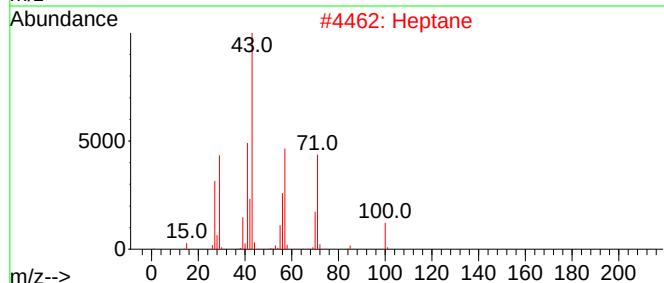
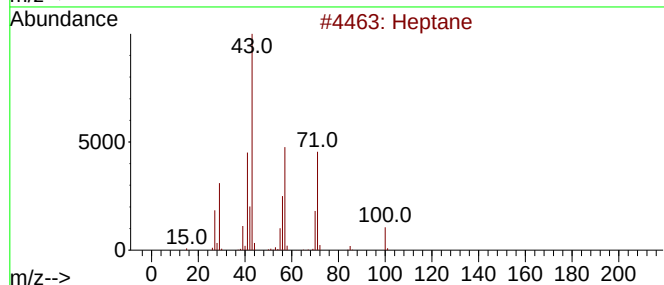
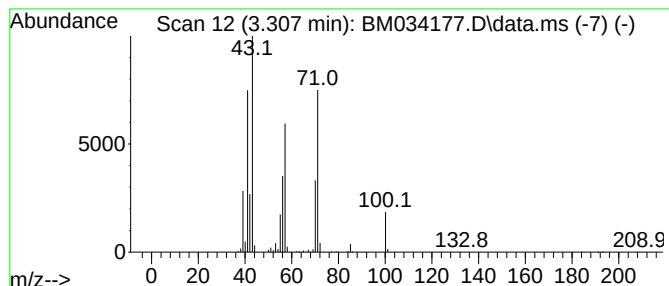
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.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.
3.307	41.46 ng/ul	1948050	1,4-Dichlorobenzene-d4	7.989

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	87
4	Heptane			100	C7H16	000142-82-5	64
5	Pentane, 2-methyl-			86	C6H14	000107-83-5	50



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	3.307	41.5	ng/u1	1948050	1	7.989	939635	20.0