

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021722\
 Data File : BN018643.D
 Acq On : 18 Feb 2022 00:29
 Operator : CG/JU
 Sample : N1506-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGZD6

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_N\Methods\SFAM-EPA-BN021622.M
 Title : SVOA CALIBRATION

Signal : TIC: BN018643.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.040	6	9	16	rVB	210933	235509	6.48%	1.386%
2	3.123	20	23	26	rVV	60020	65951	1.81%	0.388%
3	3.152	26	28	30	rVV	26803	27910	0.77%	0.164%
4	3.181	30	33	35	rVV	41555	41475	1.14%	0.244%
5	3.223	35	40	45	rVB	3538567	3635072	100.00%	21.394%
6	3.570	95	99	104	rVB	103352	122651	3.37%	0.722%
7	3.664	112	115	120	rVB5	9142	12863	0.35%	0.076%
8	4.058	178	182	189	rVB	17073	26812	0.74%	0.158%
9	4.522	257	261	268	rVB	8140	11482	0.32%	0.068%
10	4.728	293	296	300	rVB4	3963	3715	0.10%	0.022%
11	5.352	397	402	406	rBV6	2410	4212	0.12%	0.025%
12	7.058	688	692	698	rBV7	2730	5724	0.16%	0.034%
13	7.916	831	838	852	rBV	959184	1513800	41.64%	8.909%
14	10.722	1307	1315	1328	rBV	1196352	2023107	55.66%	11.907%
15	14.551	1959	1966	1975	rBV	1549282	2295087	63.14%	13.507%
16	16.292	2258	2262	2265	rVB6	2916	3840	0.11%	0.023%
17	16.598	2311	2314	2322	rVB7	2572	4614	0.13%	0.027%
18	17.075	2387	2395	2399	rBV5	7172	14154	0.39%	0.083%
19	17.287	2424	2431	2443	rBV	1570202	2250815	61.92%	13.247%
20	17.381	2444	2447	2454	rVB5	5873	11512	0.32%	0.068%
21	17.957	2542	2545	2550	rVV7	4081	5607	0.15%	0.033%
22	18.010	2550	2554	2557	rVB6	6365	8170	0.22%	0.048%
23	18.086	2564	2567	2571	rVB5	8848	11041	0.30%	0.065%
24	18.363	2610	2614	2618	rBV7	9538	14552	0.40%	0.086%
25	18.422	2621	2624	2628	rVV5	8410	11286	0.31%	0.066%
26	18.463	2628	2631	2636	rVB5	14505	21014	0.58%	0.124%
27	18.910	2705	2707	2712	rVB6	5793	6717	0.18%	0.040%
28	19.075	2732	2735	2739	rBV5	8665	13739	0.38%	0.081%
29	19.210	2754	2758	2764	rVB6	29809	45119	1.24%	0.266%
30	19.339	2777	2780	2787	rVB3	9506	13781	0.38%	0.081%
31	19.404	2788	2791	2795	rVV4	17341	22988	0.63%	0.135%
32	19.486	2801	2805	2810	rVB3	43447	60194	1.66%	0.354%
33	19.551	2813	2816	2819	rVB4	15922	18223	0.50%	0.107%
34	19.916	2875	2878	2879	rBV3	19063	21349	0.59%	0.126%
35	20.057	2898	2902	2906	rBV4	46286	57445	1.58%	0.338%
36	20.198	2921	2926	2930	rBV	99636	133808	3.68%	0.788%

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37	20.475	2969	2973	2977	rVB3	104347	124190	3.42%	0.731%
38	20.639	2996	3001	3006	rBV4	40635	76076	2.09%	0.448%
39	20.939	3048	3052	3057	rBV3	101990	122486	3.37%	0.721%
40	20.998	3060	3062	3069	rVB7	31342	41503	1.14%	0.244%
41	21.204	3095	3097	3102	rVB2	59532	70028	1.93%	0.412%
42	21.463	3136	3141	3145	rBV	1466659	1914671	52.67%	11.269%
43	23.857	3542	3548	3562	rVB2	870370	1867017	51.36%	10.988%

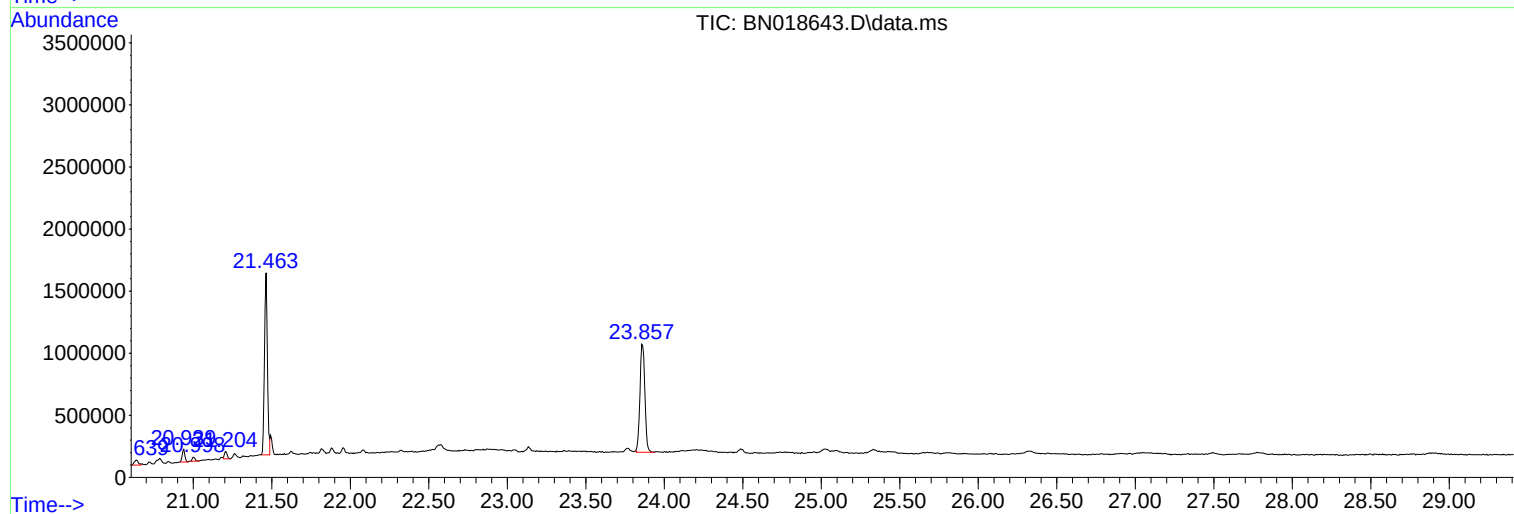
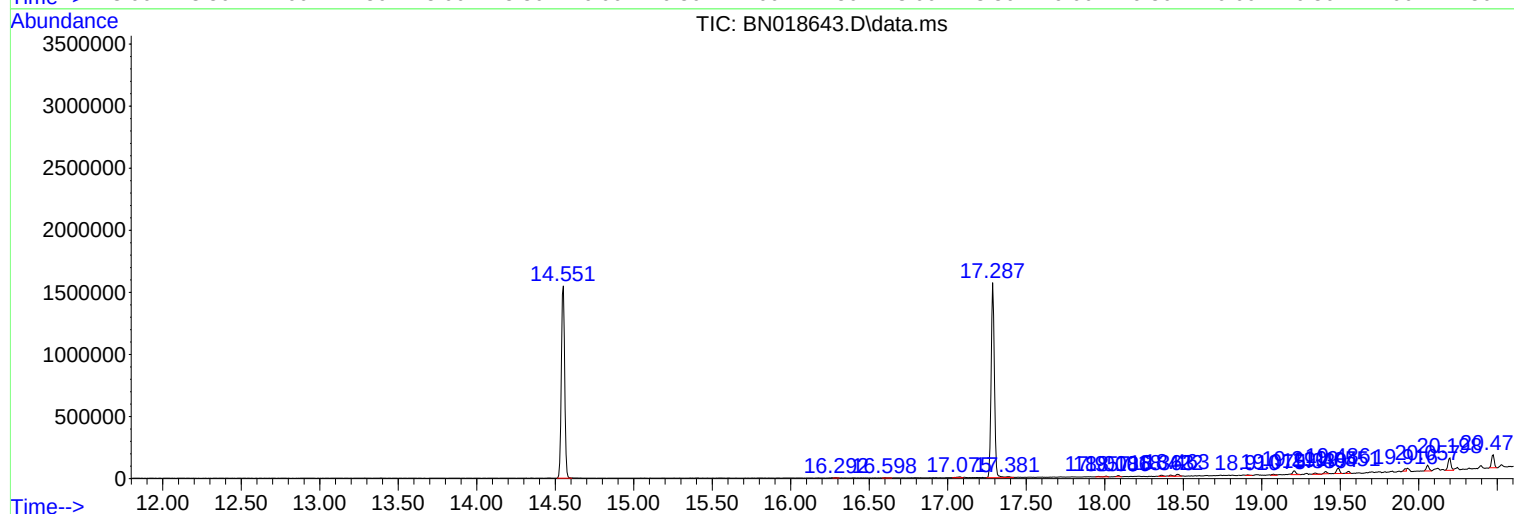
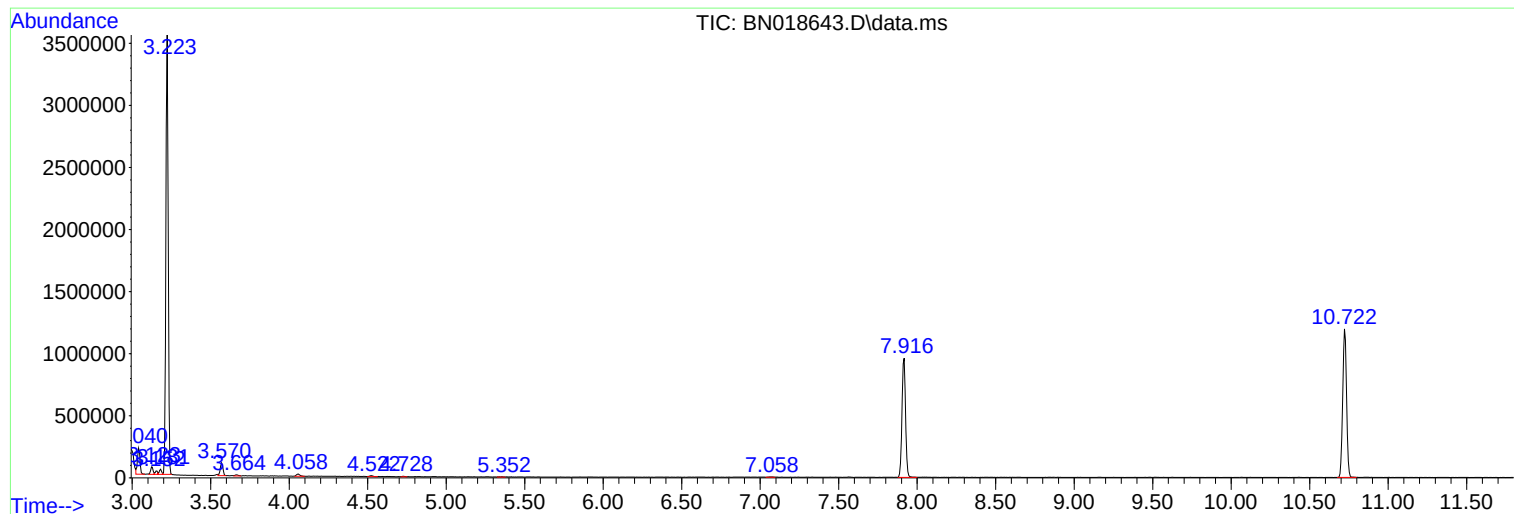
Sum of corrected areas: 16991309

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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



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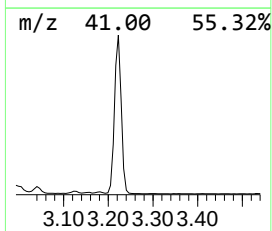
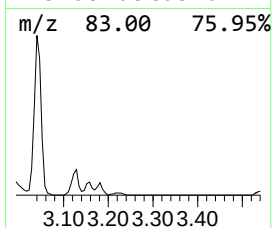
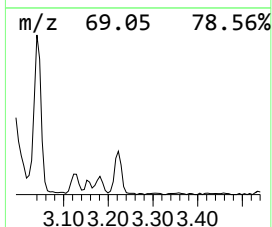
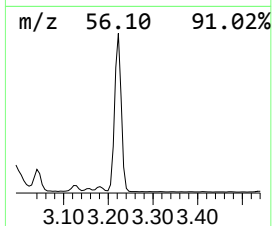
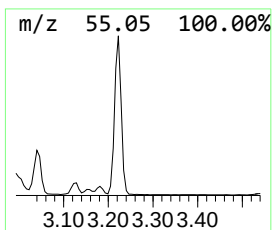
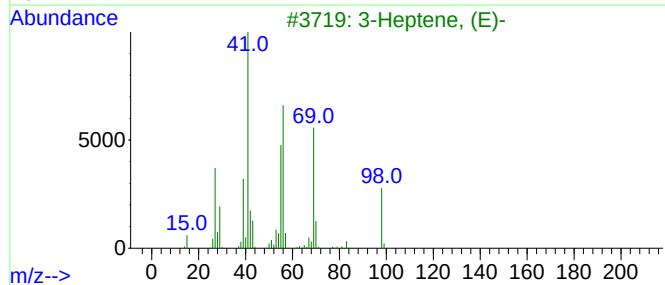
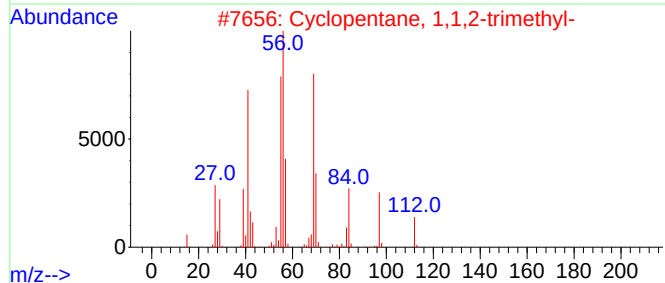
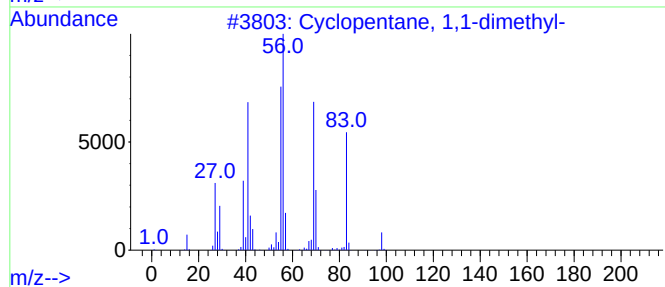
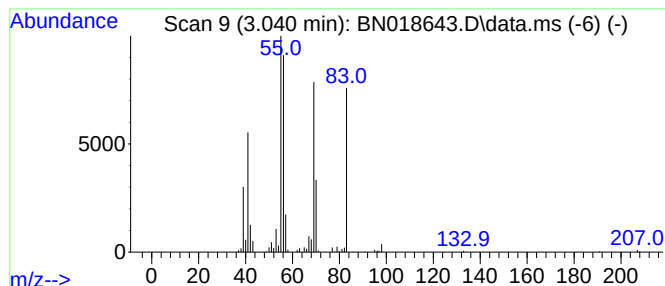
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.040 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	3.11 ng/ul	235509	1,4-Dichlorobenzene-d4	7.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	64
2			Cyclopentane, 1,1,2-trimethyl-	112	C8H16	004259-00-1	53
3			3-Heptene, (E)-	98	C7H14	014686-14-7	50
4			4-Nonene, 5-methyl-	140	C10H20	015918-07-7	50
5			3-Undecene, 8-methyl-	168	C12H24	1000061-84-4	50



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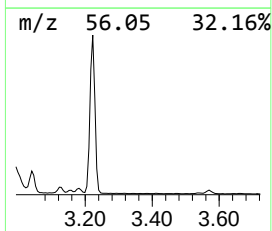
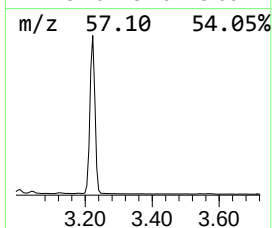
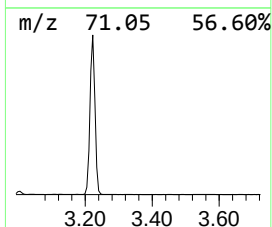
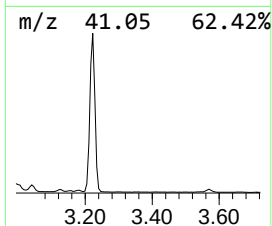
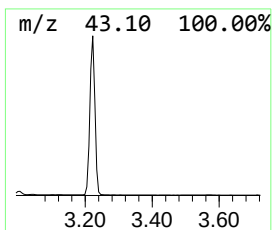
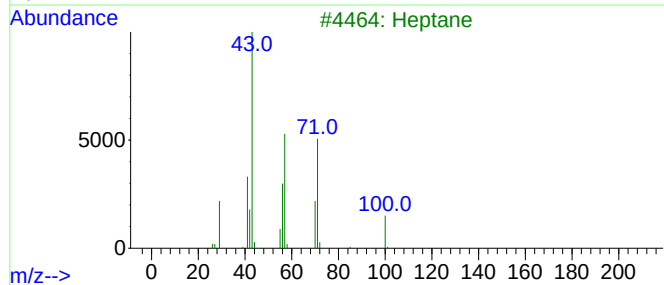
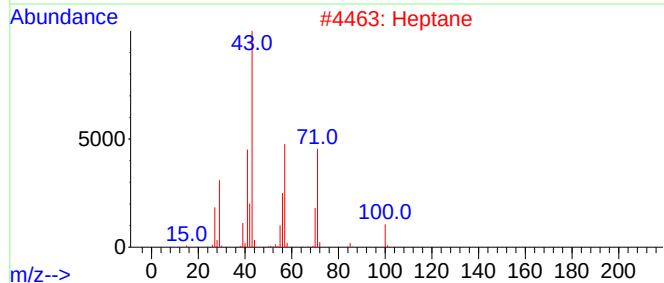
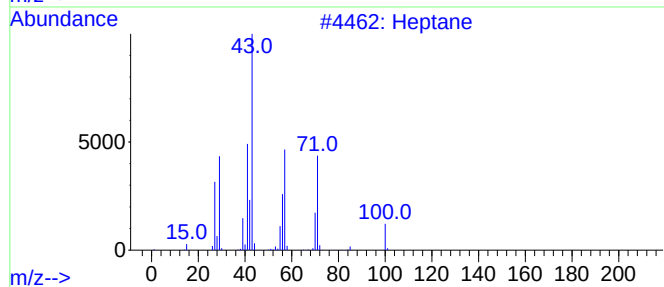
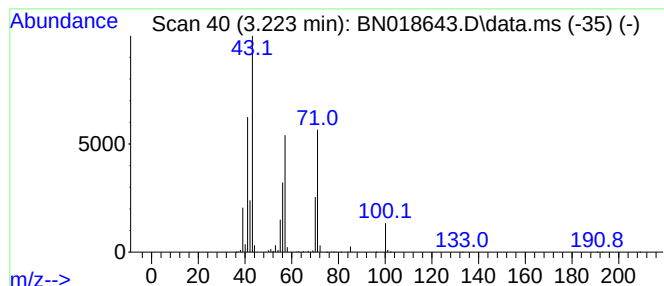
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.223	48.03 ng/ul	3635070	1,4-Dichlorobenzene-d4	7.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	94
2	Heptane			100	C7H16	000142-82-5	91
3	Heptane			100	C7H16	000142-82-5	91
4	Heptane			100	C7H16	000142-82-5	87
5	2-Butanone, 3,3-dimethyl-			100	C6H12O	000075-97-8	43



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.040	3.1	ng/u1	235509	1	7.916	1513800	20.0
(DEL) Alkane: S...	3.223	48.0	ng/u1	3635070	1	7.916	1513800	20.0