

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010162.D
 Acq On : 29 Apr 2022 21:45
 Operator : CG/JU
 Sample : PB144429BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK429

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_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010162.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.111	3	4	12	rVB	107847	110596	4.76%	0.866%
2	3.193	15	18	21	rVV	31487	33445	1.44%	0.262%
3	3.223	21	23	25	rVV3	13440	14195	0.61%	0.111%
4	3.252	25	28	30	rVV2	18942	20276	0.87%	0.159%
5	3.287	30	34	40	rVB	1503135	1519233	65.42%	11.898%
6	3.587	84	85	88	rBV3	3846	4083	0.18%	0.032%
7	3.628	88	92	97	rVB2	39192	47433	2.04%	0.371%
8	3.723	106	108	110	rVB4	3649	3261	0.14%	0.026%
9	3.952	144	147	149	rBV4	2194	2556	0.11%	0.020%
10	4.105	169	173	176	rVB	15730	20240	0.87%	0.159%
11	4.305	204	207	211	rBV6	2542	2552	0.11%	0.020%
12	4.393	220	222	227	rVB6	2128	2827	0.12%	0.022%
13	5.046	329	333	338	rVB6	3469	7084	0.31%	0.055%
14	6.058	503	505	510	rVB6	1764	2489	0.11%	0.019%
15	7.922	813	822	834	rBV	801937	1298474	55.91%	10.170%
16	8.216	869	872	879	rVB9	2038	3958	0.17%	0.031%
17	8.281	879	883	888	rBV8	1893	4086	0.18%	0.032%
18	8.875	982	984	989	rVB6	1741	2478	0.11%	0.019%
19	9.516	1090	1093	1097	rVB6	1590	2442	0.11%	0.019%
20	10.646	1282	1285	1289	rBV6	1237	2612	0.11%	0.020%
21	10.728	1291	1299	1315	rBV	993268	1708016	73.55%	13.377%
22	13.304	1732	1737	1740	rBV7	1887	3085	0.13%	0.024%
23	14.557	1941	1950	1961	rBV	1415238	2122837	91.41%	16.626%
24	15.334	2080	2082	2085	rBV4	1826	2353	0.10%	0.018%
25	15.398	2088	2093	2098	rVB8	2116	3934	0.17%	0.031%
26	15.563	2118	2121	2124	rBV5	2105	2626	0.11%	0.021%
27	15.945	2183	2186	2190	rBV6	2286	3502	0.15%	0.027%
28	16.081	2207	2209	2213	rBV5	2071	2637	0.11%	0.021%
29	16.175	2221	2225	2227	rBV5	1979	2335	0.10%	0.018%
30	17.098	2378	2382	2387	rVB6	2710	4631	0.20%	0.036%
31	17.310	2411	2418	2433	rBV	1497552	2322303	100.00%	18.188%
32	18.198	2567	2569	2574	rBV6	2580	2941	0.13%	0.023%
33	21.392	3107	3112	3120	rBV	1330982	1890807	81.42%	14.809%
34	23.845	3522	3529	3542	rVB2	724737	1591967	68.55%	12.468%

Sum of corrected areas: 12768294

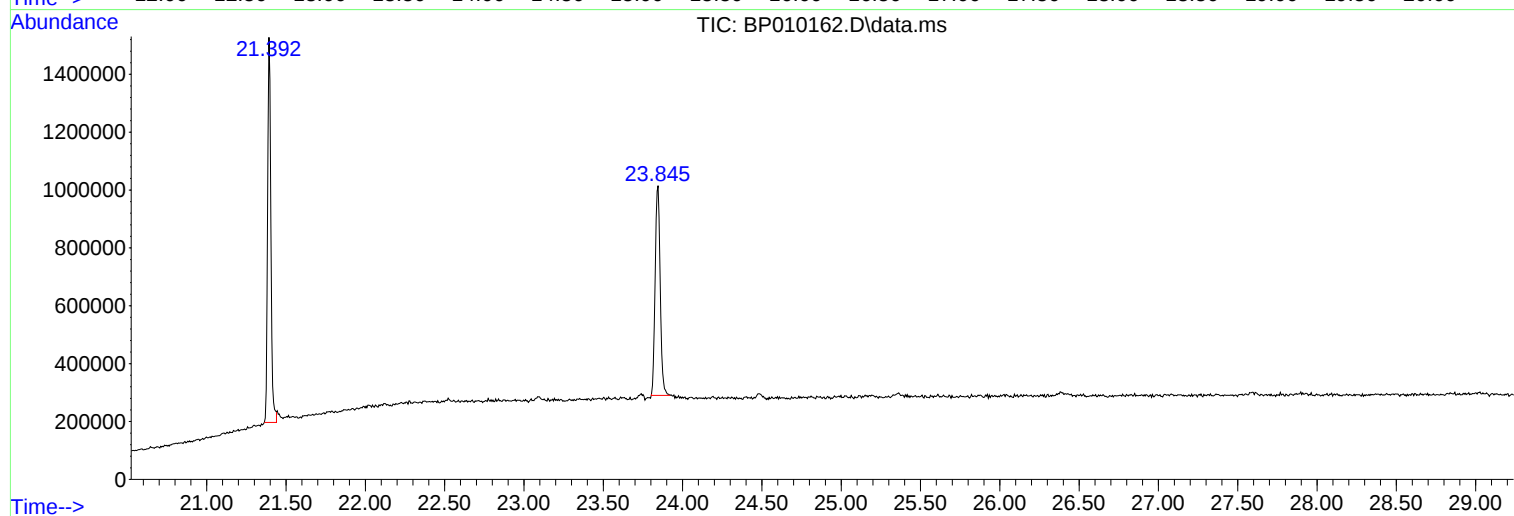
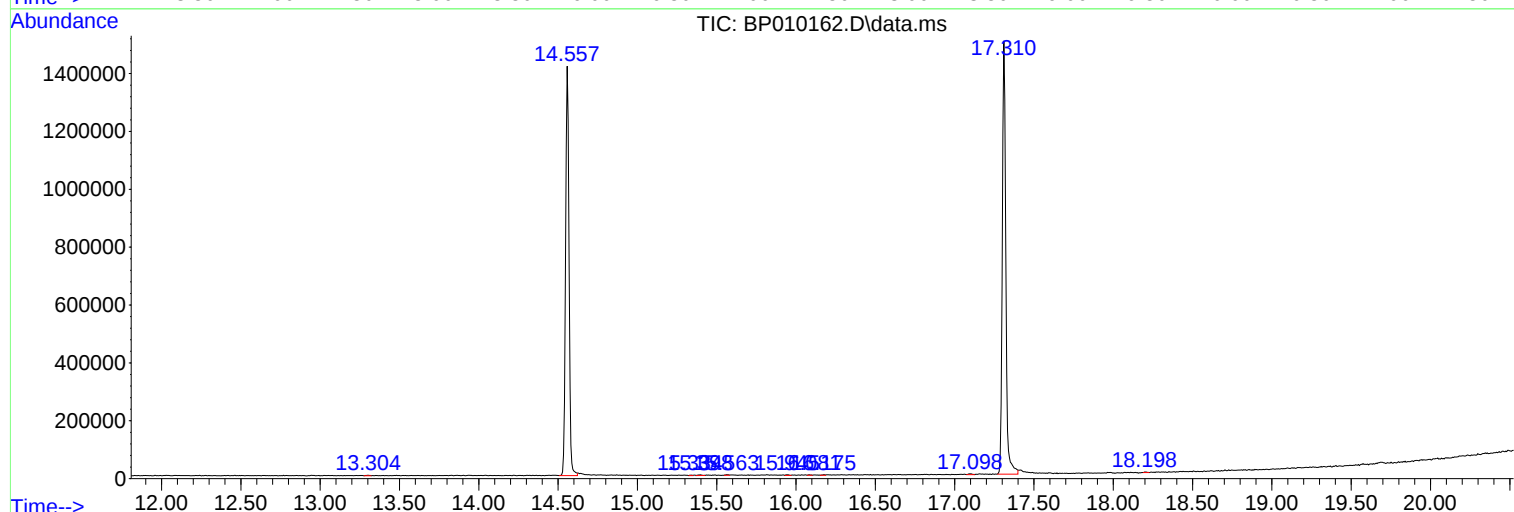
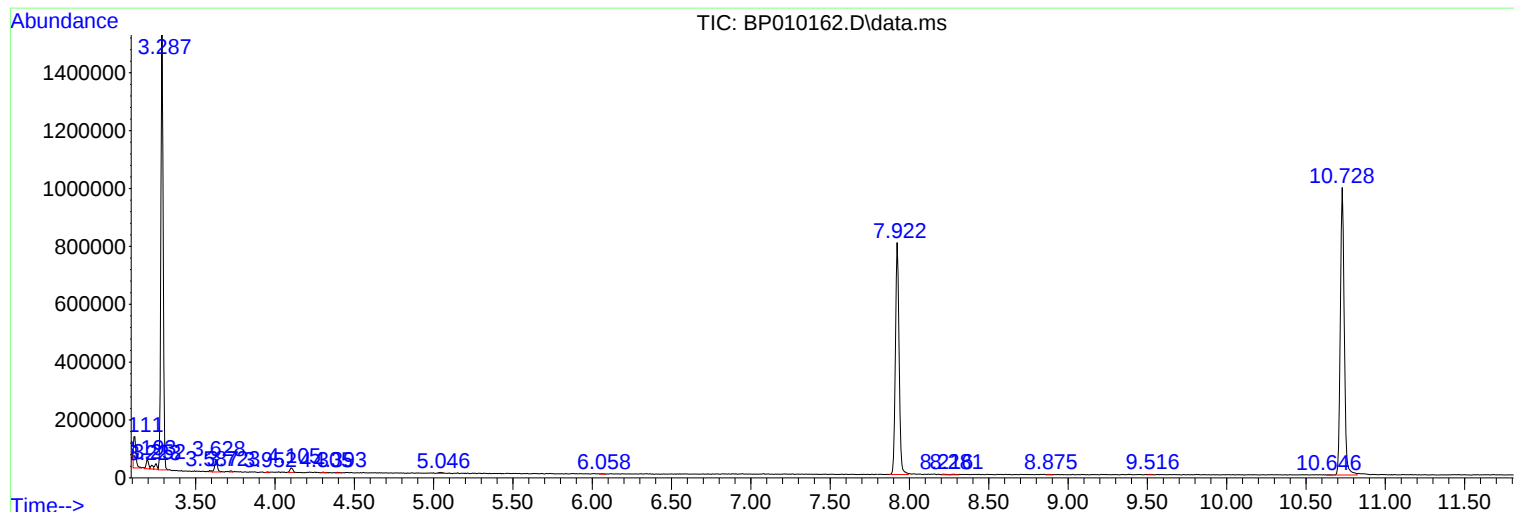
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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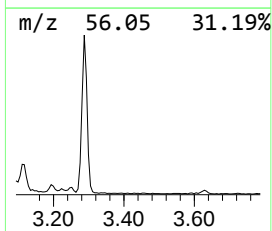
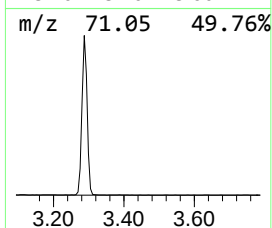
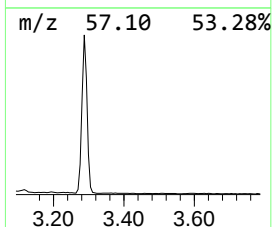
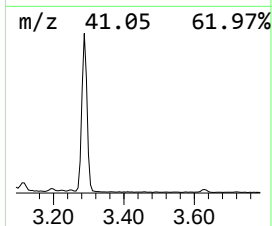
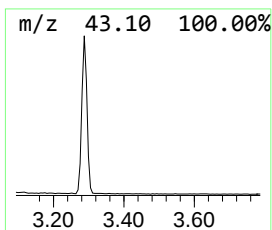
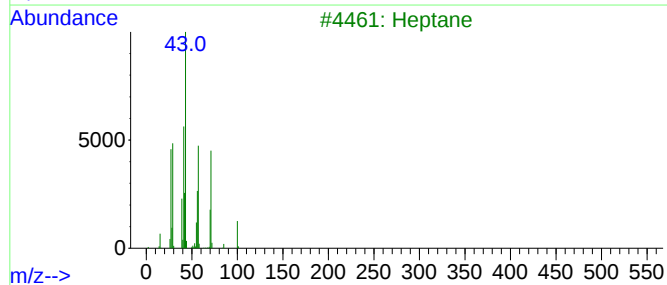
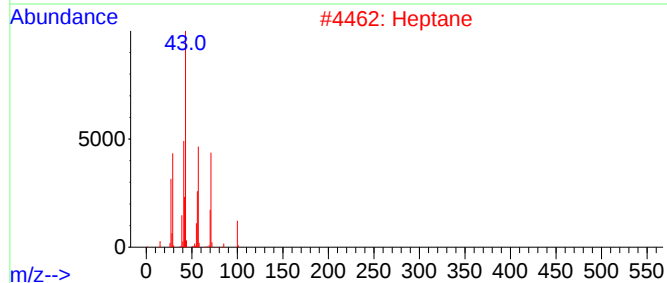
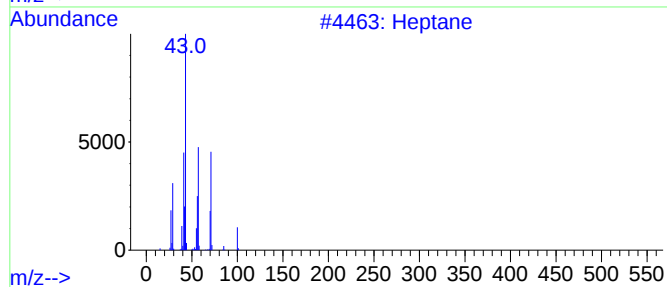
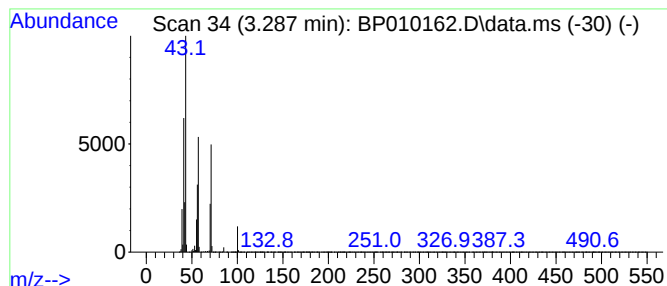
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Quant Title : SVOA CALIBRATION

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.
3.287	23.40 ng/ul	1519230	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	87
5	Hexane, 3-methyl-			100	C7H16	000589-34-4	42



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	3.287	23.4	ng/u1	1519230	1	7.922	1298470	20.0