

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
 Data File : BP019231.D
 Acq On : 03 Jan 2024 04:14
 Operator : MA/JU
 Sample : 05859-17
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGRY7

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-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019231.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	7	rVB	93309	73492	2.17%	0.396%
2	3.152	7	11	20	rVB	2080531	2081553	61.54%	11.210%
3	3.287	31	34	39	rVB5	8215	11186	0.33%	0.060%
4	3.458	60	63	65	rBV4	6766	7674	0.23%	0.041%
5	3.487	65	68	76	rVB2	38068	54291	1.61%	0.292%
6	3.969	146	150	156	rVB	19442	26903	0.80%	0.145%
7	4.117	170	175	180	rBV	80929	100111	2.96%	0.539%
8	4.169	180	184	194	rVB	105302	135799	4.01%	0.731%
9	4.275	197	202	205	rBV	77643	97562	2.88%	0.525%
10	4.317	205	209	217	rVB	74526	99658	2.95%	0.537%
11	4.911	304	310	317	rVB5	9861	16564	0.49%	0.089%
12	4.981	319	322	326	rVB6	3694	5680	0.17%	0.031%
13	5.105	340	343	345	rVB4	3301	3406	0.10%	0.018%
14	5.299	371	376	380	rBV8	3119	5799	0.17%	0.031%
15	5.499	407	410	415	rVB2	6785	10093	0.30%	0.054%
16	5.799	459	461	465	rBV5	3311	4086	0.12%	0.022%
17	5.899	474	478	484	rVB9	5463	10467	0.31%	0.056%
18	6.087	507	510	511	rBV3	3262	3864	0.11%	0.021%
19	6.446	564	571	577	rBV3	11691	20638	0.61%	0.111%
20	6.587	591	595	599	rBV	11538	16192	0.48%	0.087%
21	6.881	643	645	650	rVB6	4248	6744	0.20%	0.036%
22	7.434	737	739	745	rVB5	4774	7755	0.23%	0.042%
23	7.799	793	801	808	rBV	927377	1501730	44.40%	8.087%
24	9.552	1098	1099	1104	rVB5	3356	3634	0.11%	0.020%
25	10.593	1268	1276	1289	rBV	1150516	1924965	56.91%	10.367%
26	12.016	1515	1518	1522	rBV5	2962	4757	0.14%	0.026%
27	12.298	1562	1566	1571	rVB8	2682	4241	0.13%	0.023%
28	12.798	1648	1651	1655	rBV6	2454	4125	0.12%	0.022%
29	13.263	1728	1730	1735	rBV5	2640	3415	0.10%	0.018%
30	13.851	1827	1830	1833	rBV5	2902	4182	0.12%	0.023%
31	13.916	1837	1841	1847	rVB9	2260	5034	0.15%	0.027%
32	14.445	1924	1931	1942	rVV	1546399	2333386	68.99%	12.566%
33	15.239	2062	2066	2069	rBV6	2809	4067	0.12%	0.022%
34	15.416	2093	2096	2100	rBV6	2830	4529	0.13%	0.024%
35	16.857	2338	2341	2346	rBV5	12933	17891	0.53%	0.096%
36	16.986	2361	2363	2367	rVB5	6845	6702	0.20%	0.036%

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 Title : SVOA CALIBRATION

37	17.198	2393	2399	2409	rBV	1924074	2632825	77.84%	14.179%
38	17.892	2515	2517	2520	rBV4	10115	10244	0.30%	0.055%
39	18.092	2547	2551	2560	rBV2	51806	85548	2.53%	0.461%
40	18.575	2630	2633	2640	rVB9	30396	40505	1.20%	0.218%
41	18.804	2668	2672	2677	rBV8	24090	42745	1.26%	0.230%
42	18.869	2679	2683	2689	rVV7	43134	73136	2.16%	0.394%
43	19.522	2791	2794	2798	rBV5	60422	81403	2.41%	0.438%
44	19.704	2822	2825	2829	rBV5	62319	71695	2.12%	0.386%
45	19.863	2849	2852	2856	rBV2	100329	133927	3.96%	0.721%
46	19.904	2857	2859	2863	rVB4	47915	51043	1.51%	0.275%
47	20.045	2880	2883	2891	rBV9	70847	139019	4.11%	0.749%
48	20.180	2904	2906	2909	rVB3	47730	44760	1.32%	0.241%
49	20.651	2983	2986	2993	rBV8	81688	116940	3.46%	0.630%
50	21.304	3092	3097	3104	rBV	2393573	3040393	89.89%	16.374%
51	23.668	3493	3499	3506	rVB	1771637	3382415	100.00%	18.216%

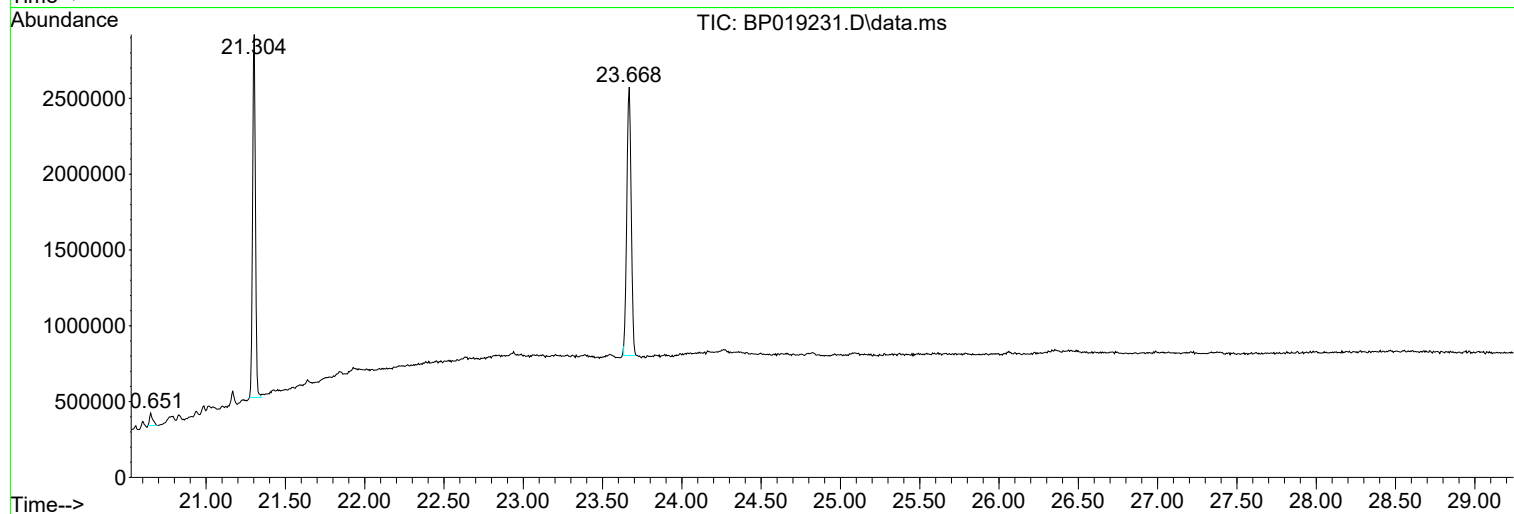
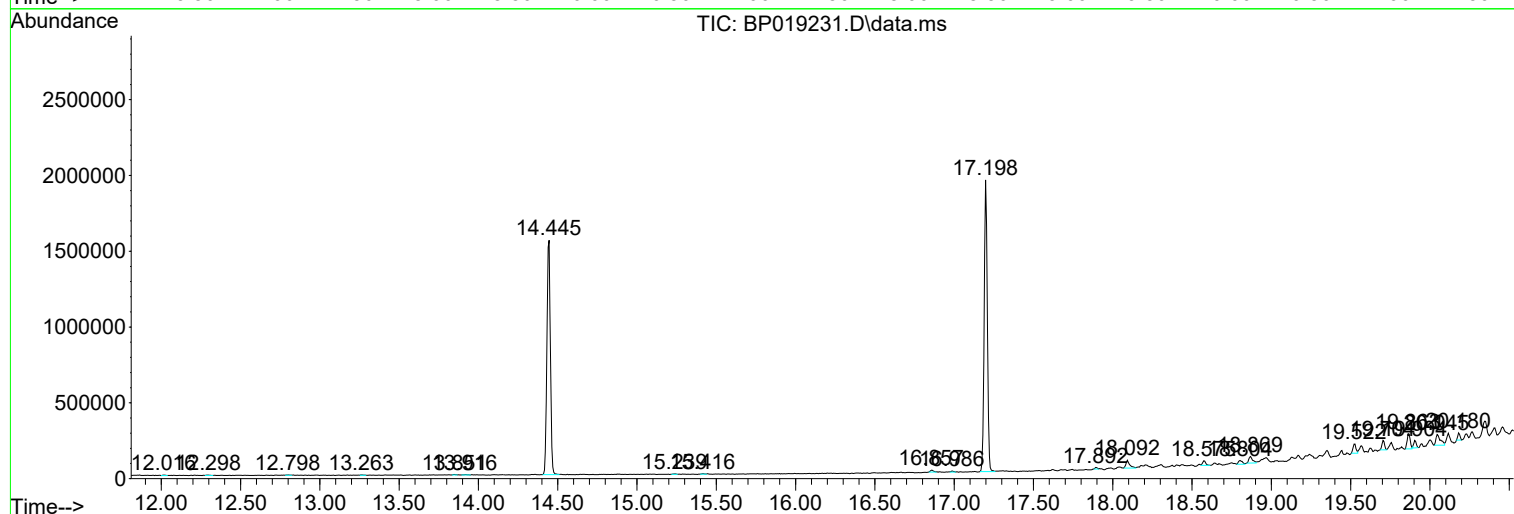
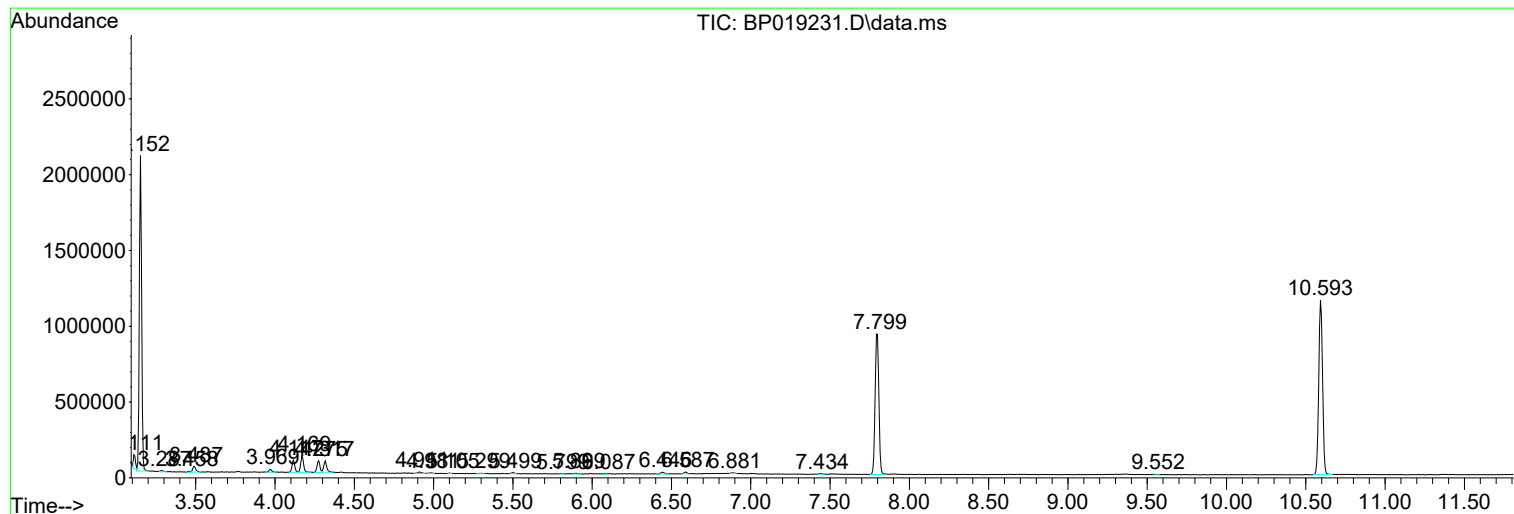
Sum of corrected areas: 18568773

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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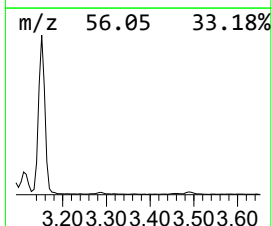
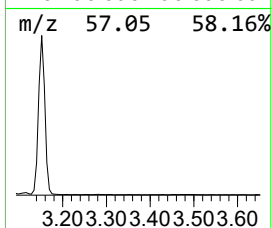
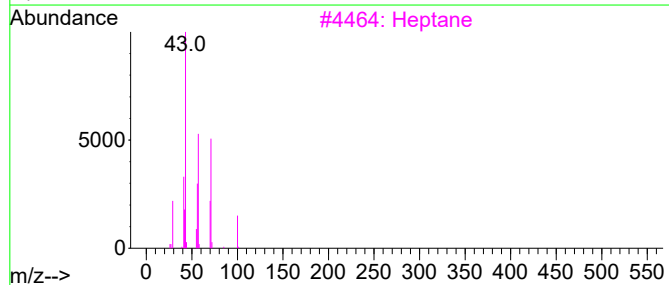
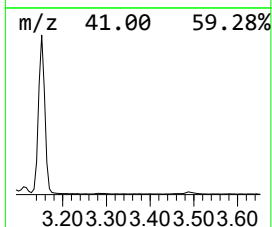
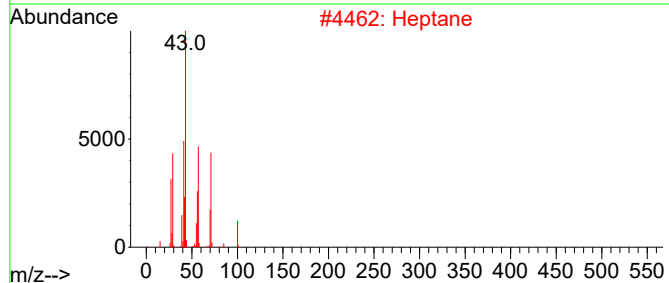
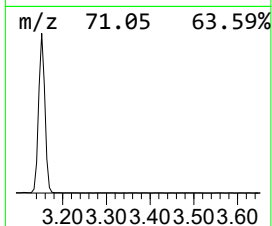
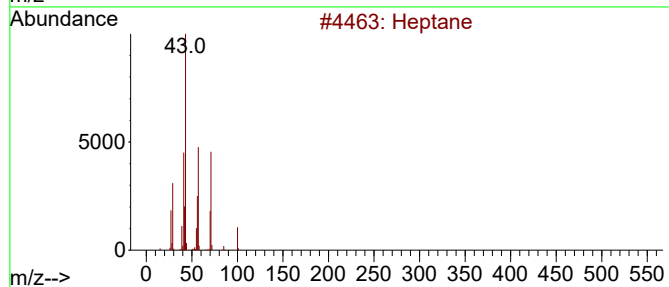
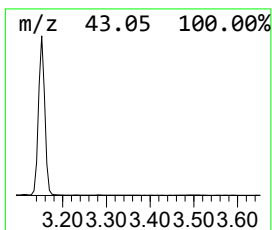
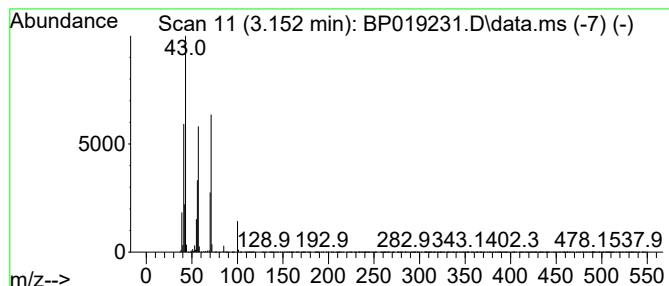
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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.
3.152	27.72 ng/ul	2081550	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Heptane	100	C7H16	000142-82-5	91
3			Heptane	100	C7H16	000142-82-5	91
4			Heptane	100	C7H16	000142-82-5	90
5			Hexane, 3-methyl-	100	C7H16	000589-34-4	43



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	3.152	27.7	ng/ul	2081550	1	7.799	1501730	20.0