

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072421\
 Data File : BG049443.D
 Acq On : 24 Jul 2021 3:24
 Operator : CG/JU
 Sample : M3036-08
 Misc : GCMS CONFIRMATION
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0CW7

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_G\METHODS\SFAM-EPA-BG072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049443.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.091	3	9	22	rBV	14046619	18387017	100.00%	90.255%
2	3.349	48	53	62	rVB	410500	479243	2.61%	2.352%
3	8.043	845	852	859	rBV2	87187	147340	0.80%	0.723%
4	10.869	1324	1333	1341	rBV	103810	189573	1.03%	0.931%
5	14.693	1978	1984	1993	rBV	150754	246221	1.34%	1.209%
6	17.442	2446	2452	2459	rBV	179517	284475	1.55%	1.396%
7	21.725	3172	3181	3189	rVB	192641	317742	1.73%	1.560%
8	24.997	3727	3738	3750	rVB2	106331	320680	1.74%	1.574%

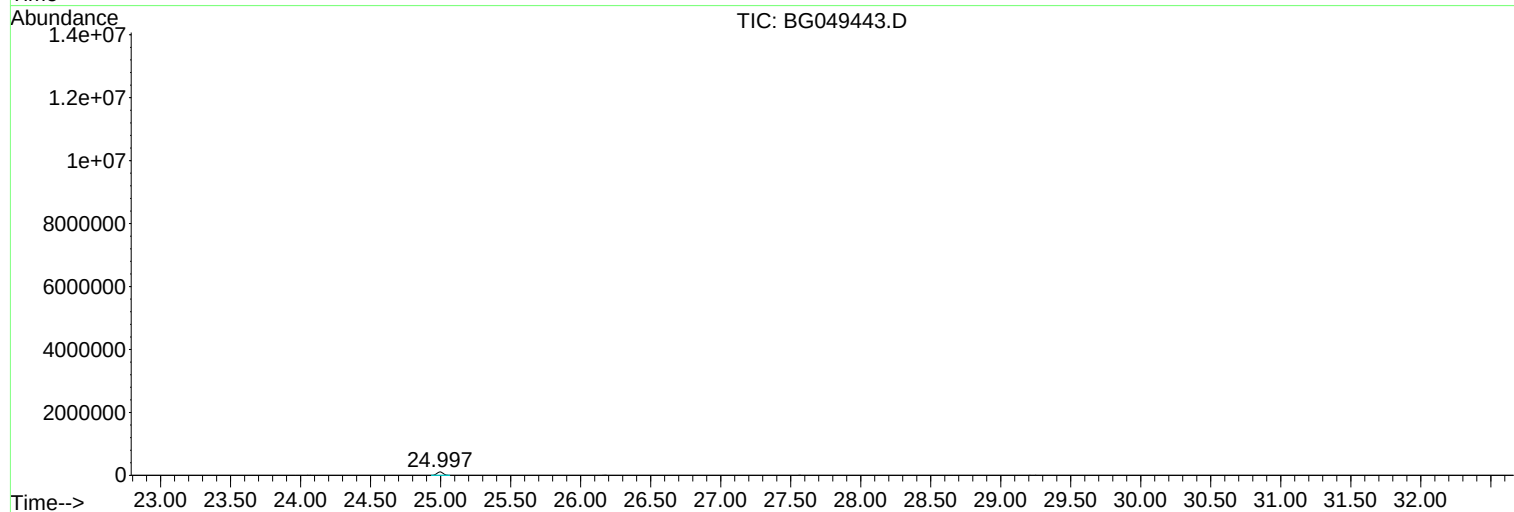
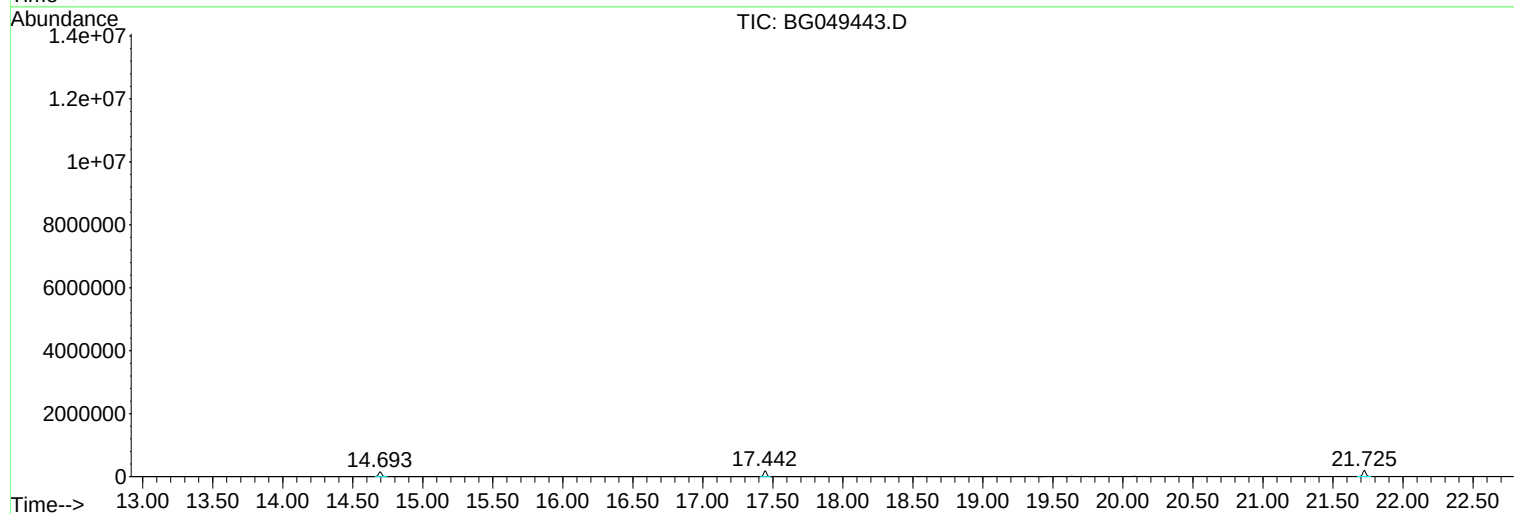
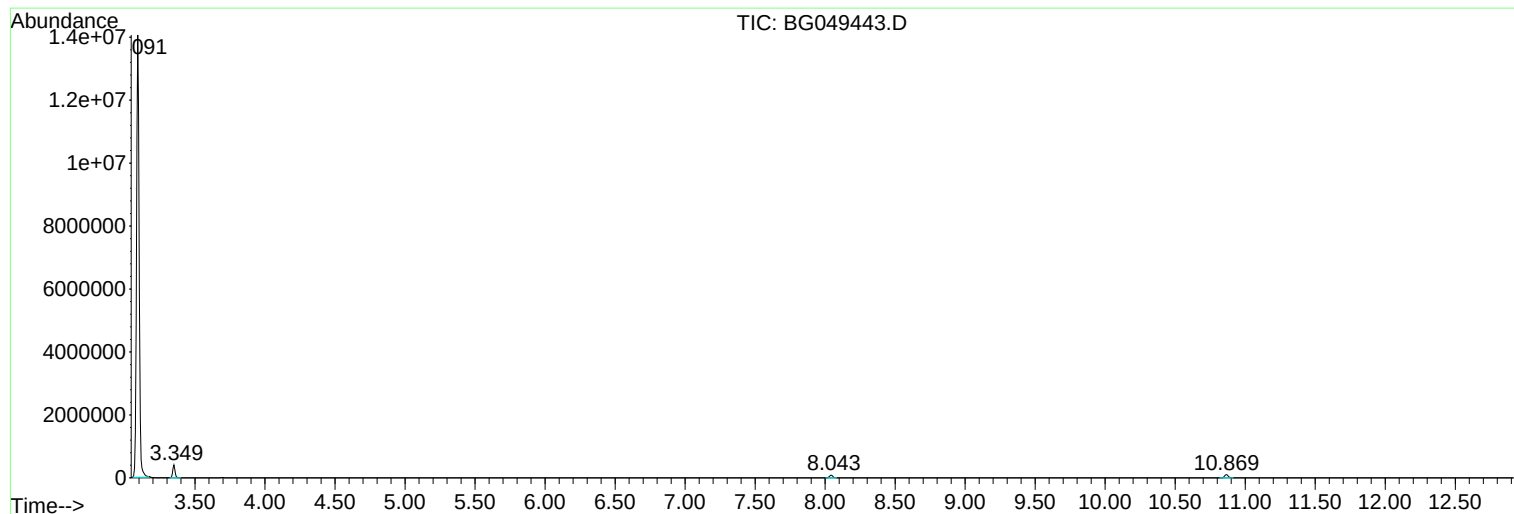
Sum of corrected areas: 20372291

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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



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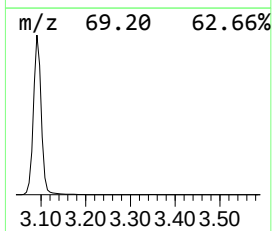
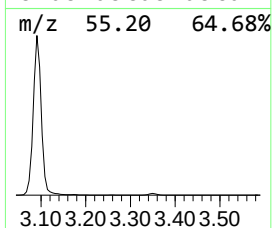
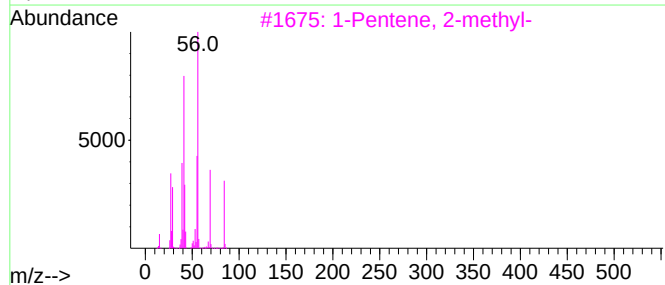
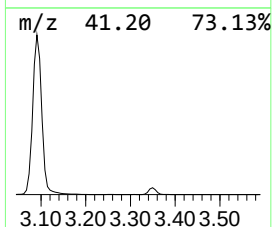
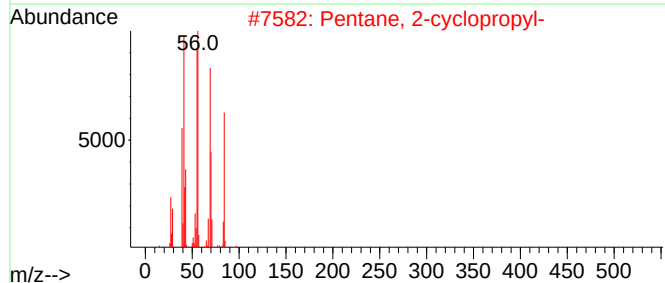
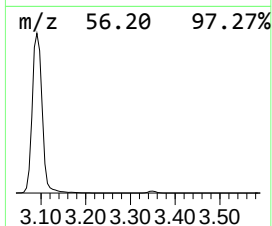
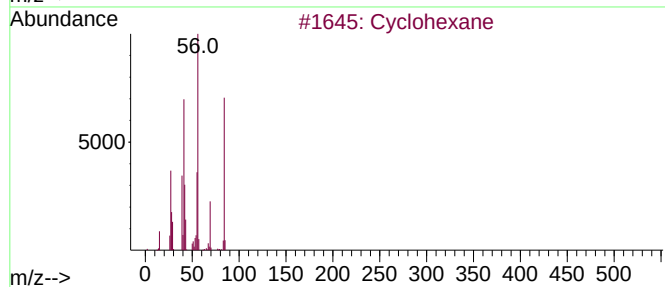
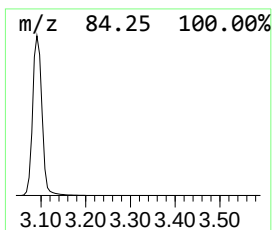
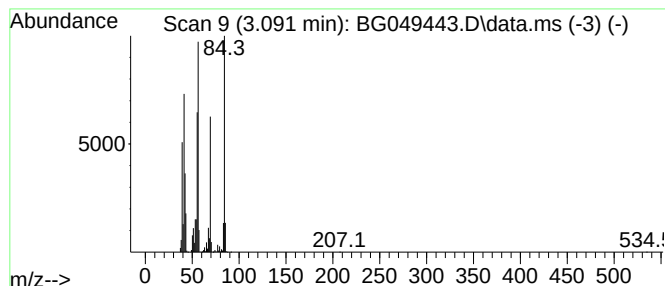
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Cyclic3.09 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.090	2495.86 ng/ul	18387000	1,4-Dichlorobenzene-d4	8.043

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane	84	C6H12	000110-82-7	59
2		Pentane, 2-cyclopropyl-	112	C8H16	005458-16-2	59
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
4		1-Silacyclo-3-pentene	84	C4H8Si	007049-25-4	58
5		2-Amino-oxazole	84	C3H4N2O	004570-45-0	53



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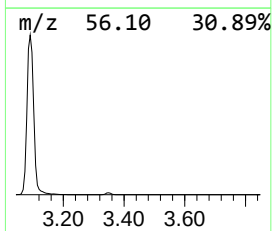
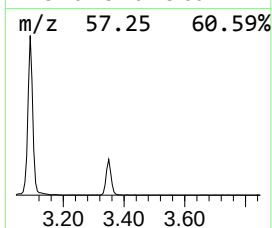
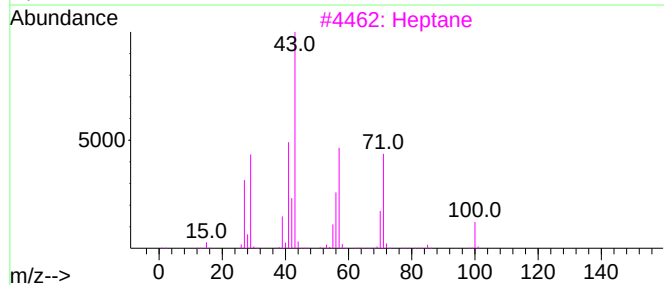
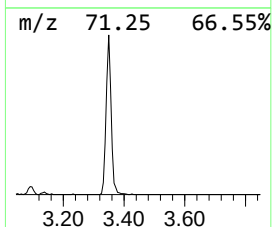
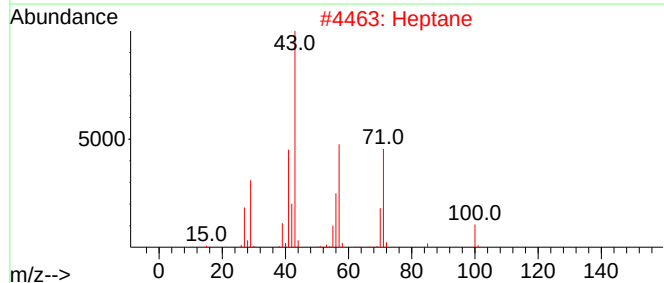
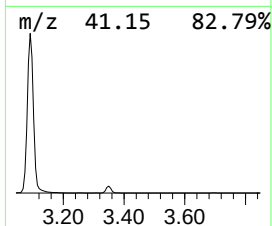
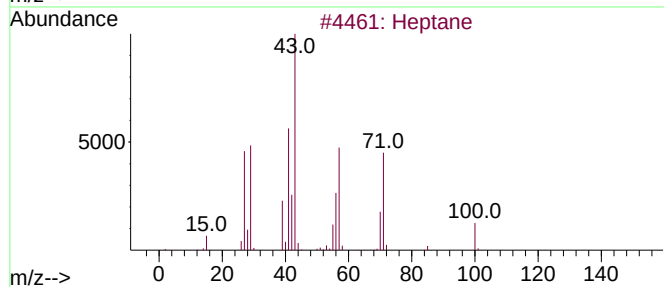
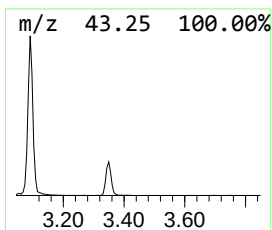
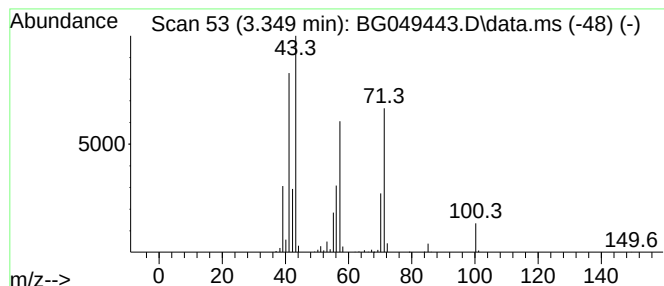
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.350	65.05 ng/ul	479243	1,4-Dichlorobenzene-d4	8.043

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane	100	C7H16	000142-82-5	91
2		Heptane	100	C7H16	000142-82-5	87
3		Heptane	100	C7H16	000142-82-5	80
4		Heptane	100	C7H16	000142-82-5	72
5		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.090	2495.9	ng/ul	18387000	1	8.043	147340	20.0
(DEL) Alkane: S...	3.350	65.0	ng/ul	479243	1	8.043	147340	20.0