

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG032819\
 Data File : BG040198.D
 Acq On : 28 Mar 2019 13:42
 Operator : JU/SJ
 Sample : PB118345BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118345BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % 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\8270-BG032719.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.141	3	9	19	rBV2	50803	100130	3.28%	0.650%
2	5.620	424	431	448	rBV	706176	1146148	37.52%	7.445%
3	7.218	694	703	714	rBV	704183	1216492	39.82%	7.902%
4	7.600	758	768	778	rBV	670110	1178909	38.59%	7.658%
5	8.070	840	848	857	rBV	182120	312109	10.22%	2.027%
6	8.393	891	903	911	rBV	501235	894081	29.27%	5.808%
7	9.245	1040	1048	1064	rBV	390834	716144	23.44%	4.652%
8	10.885	1319	1327	1335	rBV	235606	416178	13.62%	2.703%
9	13.323	1734	1742	1749	rBV	1069439	1661426	54.39%	10.792%
10	14.698	1970	1976	1987	rVB2	337043	567559	18.58%	3.687%
11	16.184	2223	2229	2243	rBV	748585	1181397	38.67%	7.674%
12	17.447	2436	2444	2454	rVB2	504445	770457	25.22%	5.005%
13	20.044	2879	2886	2892	rBV	2357204	3054894	100.00%	19.843%
14	21.725	3164	3172	3180	rBV	669909	1075284	35.20%	6.985%
15	23.999	3550	3559	3564	rBV6	20203	42940	1.41%	0.279%
16	25.027	3717	3734	3746	rBV2	353887	1060871	34.73%	6.891%

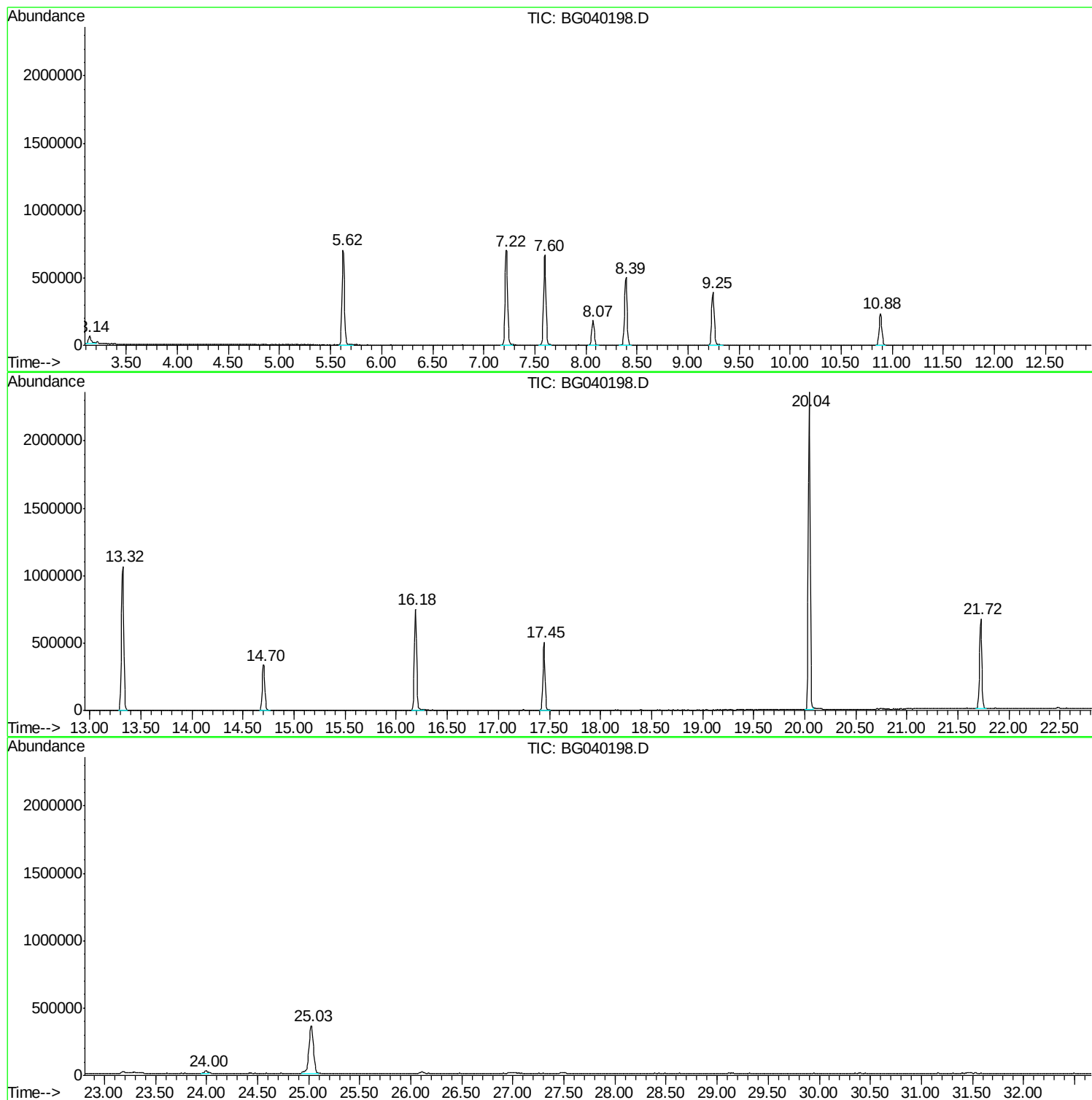
Sum of corrected areas: 15395019

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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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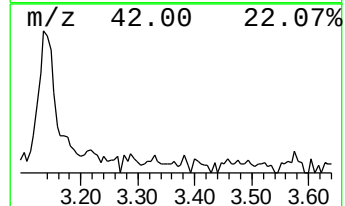
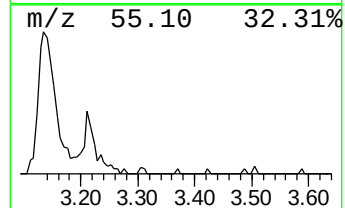
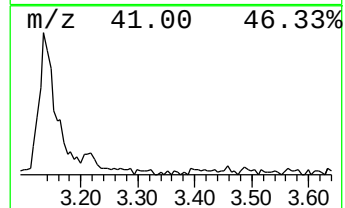
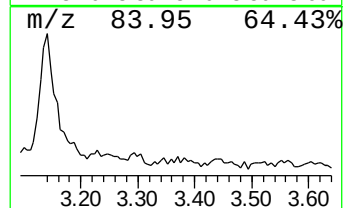
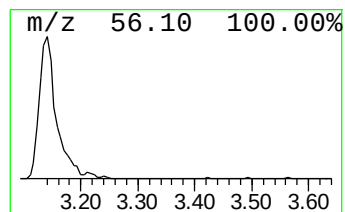
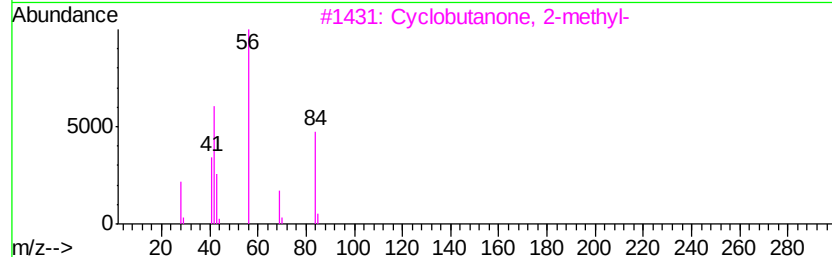
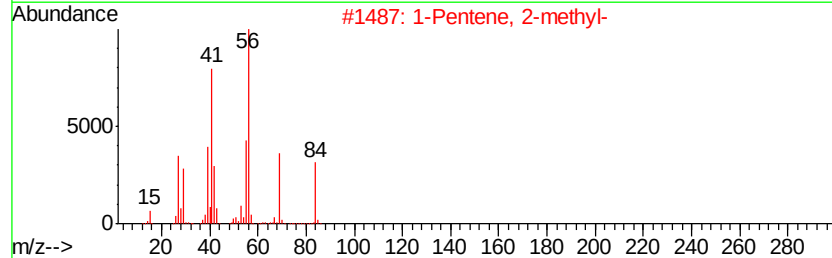
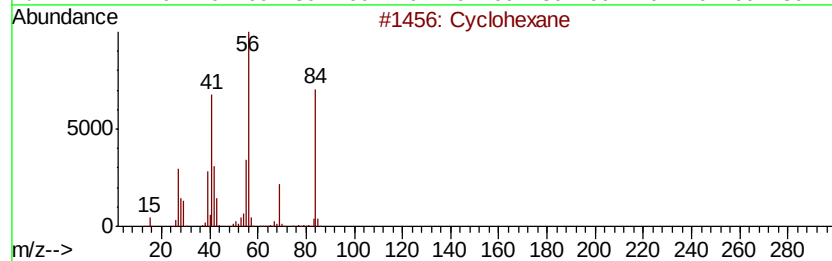
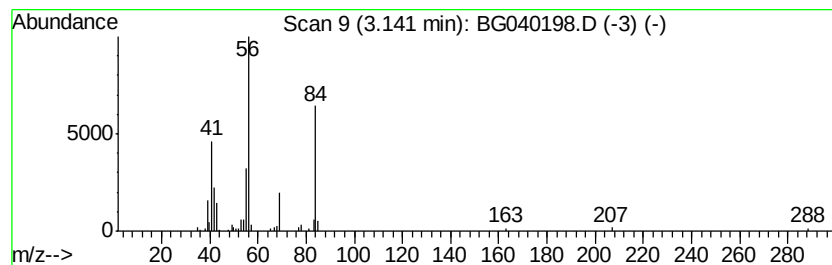
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Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.14	6.42 ng	100130	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	90
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
3		Cyclobutanone, 2-methyl-	84	C5H8O	001517-15-3	59
4		Cyclopropane, propyl-	84	C6H12	002415-72-7	50
5		Cyclopentane, methyl-	84	C6H12	000096-37-7	43



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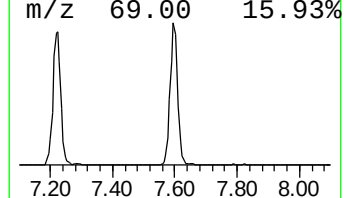
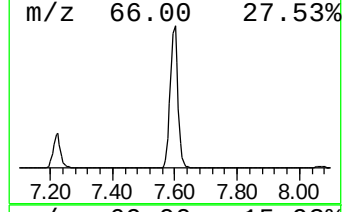
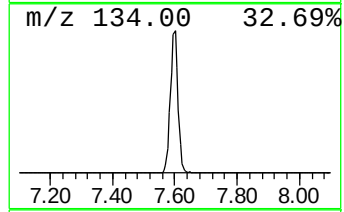
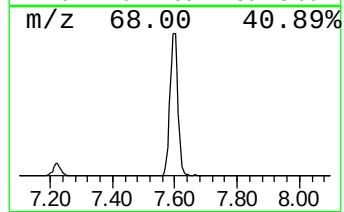
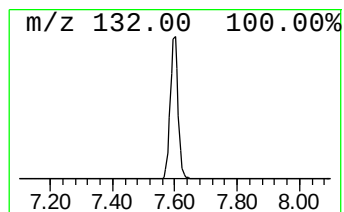
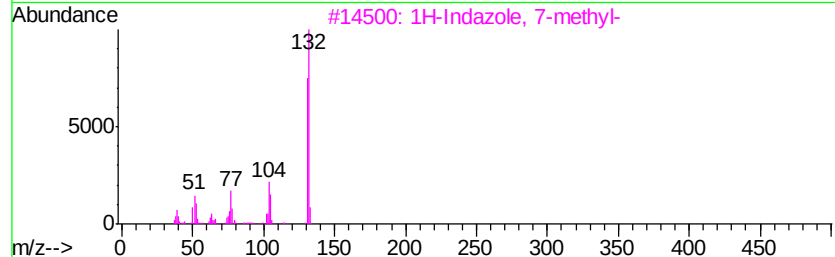
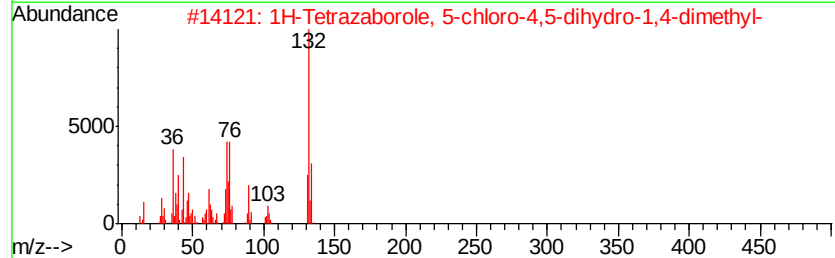
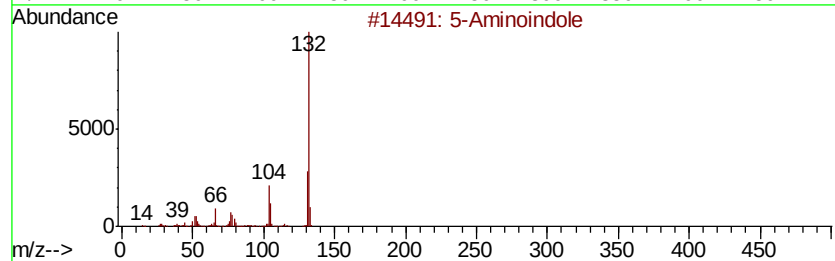
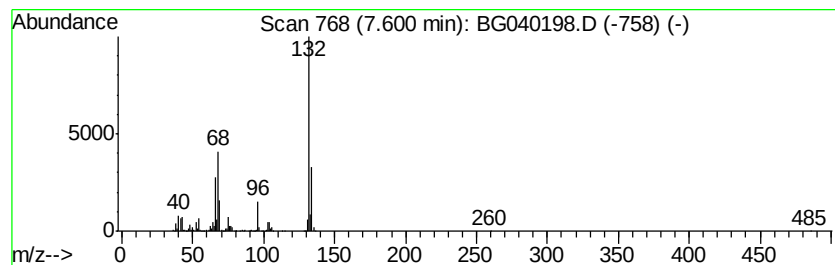
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	75.54 ng	1178910	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		1H-Tetrazaborole, 5-chloro-4,5-d...	132	C2H6BClN4	021960-49-6	12
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	12
4		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Pentene, 2-methyl-	3.14	6.4	ng	100130	1	8.07	312109	20.0
unknown7.60	7.60	75.5	ng	1178910	1	8.07	312109	20.0