

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121818\
 Data File : BG038701.D
 Acq On : 18 Dec 2018 19:19
 Operator : JU/SJ
 Sample : PB115799BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115799BL

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-BG121218.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	5.142	342	349	356	rBV2	18053	31246	1.94%	0.418%
2	5.606	421	428	437	rBV	318851	534091	33.12%	7.140%
3	7.210	693	701	713	rBV	340969	596443	36.99%	7.973%
4	7.586	758	765	777	rBV	318642	562777	34.90%	7.523%
5	8.056	838	845	852	rBV	77807	132417	8.21%	1.770%
6	8.379	893	900	909	rVB	240253	431084	26.73%	5.763%
7	9.231	1038	1045	1057	rBV	187209	360527	22.36%	4.820%
8	10.876	1318	1325	1336	rBV	95198	180168	11.17%	2.409%
9	13.315	1733	1740	1751	rBV	550387	889203	55.14%	11.887%
10	14.695	1966	1975	1989	rVB2	165540	272342	16.89%	3.641%
11	16.182	2222	2228	2247	rBV2	401612	684096	42.42%	9.145%
12	17.439	2436	2442	2452	rBV2	235498	368676	22.86%	4.929%
13	20.042	2879	2885	2895	rBV	1247967	1612579	100.00%	21.557%
14	21.722	3163	3171	3183	rBV2	253883	434899	26.97%	5.814%
15	25.019	3721	3732	3747	rVB2	127754	389919	24.18%	5.212%

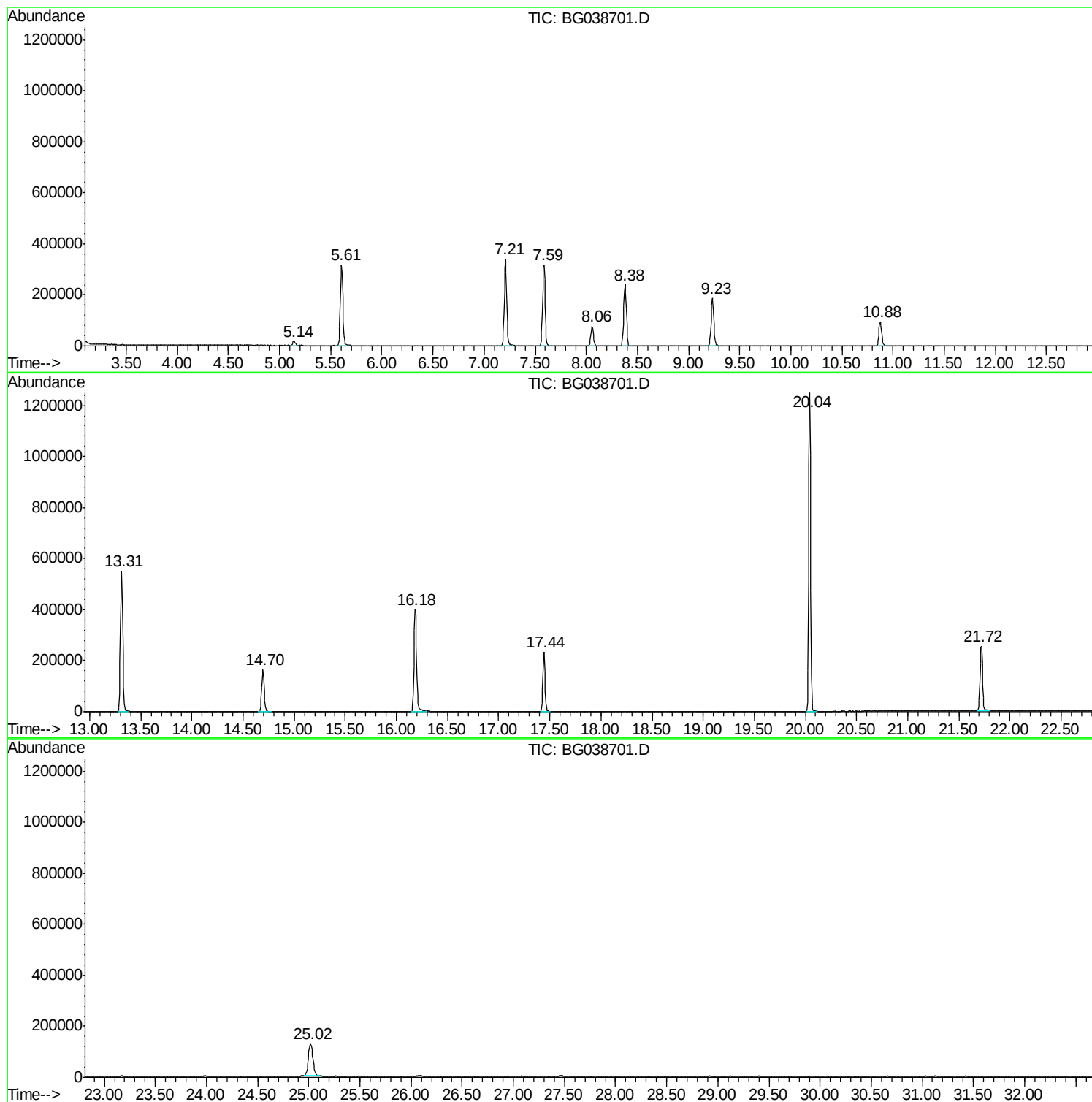
Sum of corrected areas: 7480467

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



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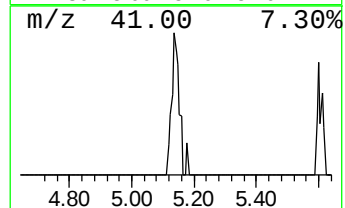
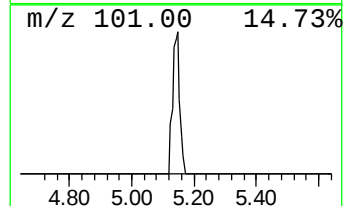
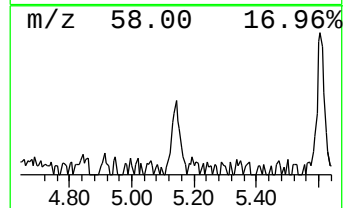
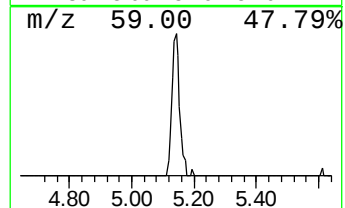
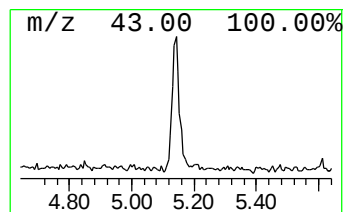
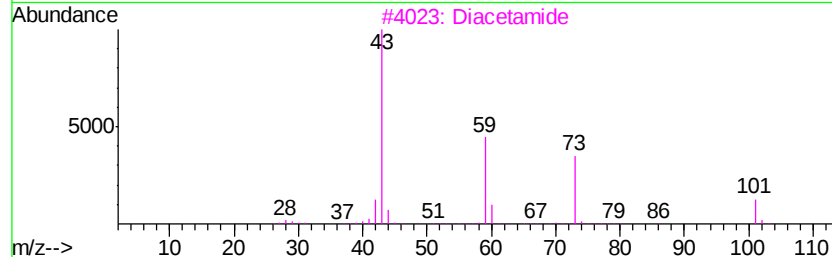
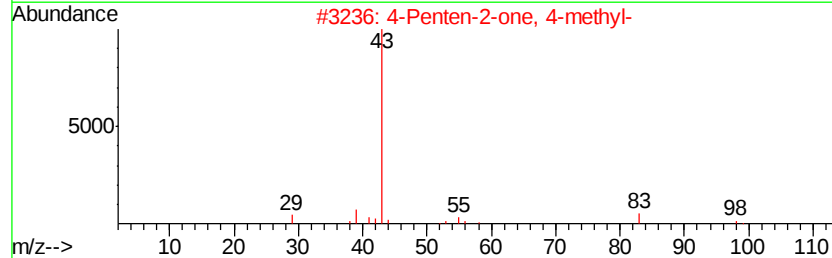
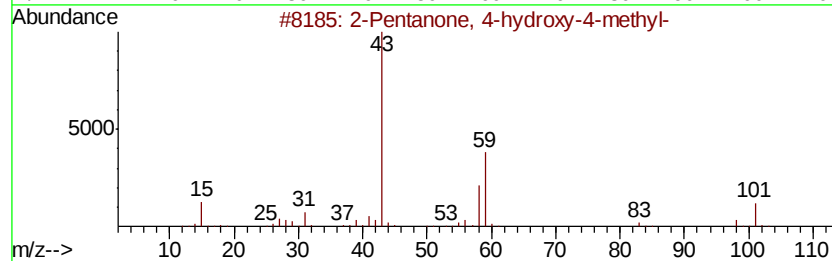
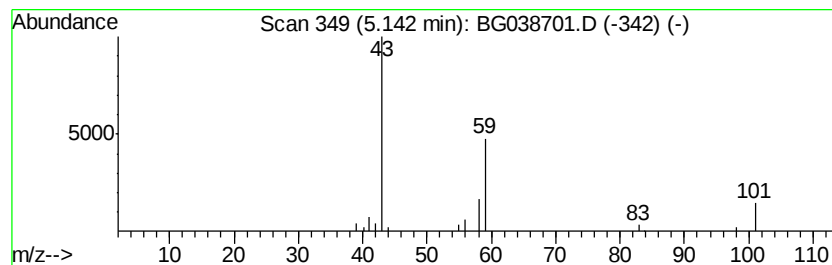
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.72 ng	31246	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
3			Diacetamide	101	C4H7NO2	000625-77-4	9
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5			5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9



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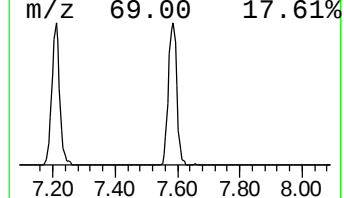
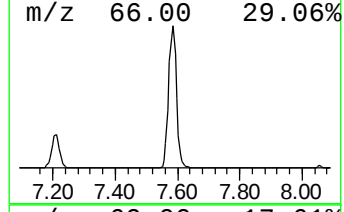
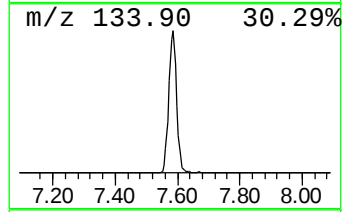
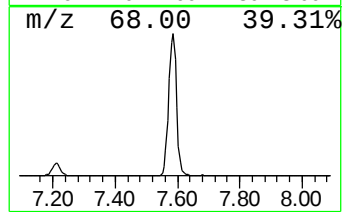
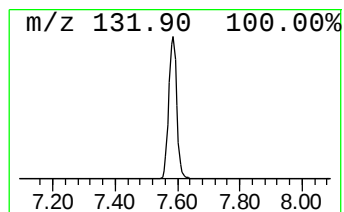
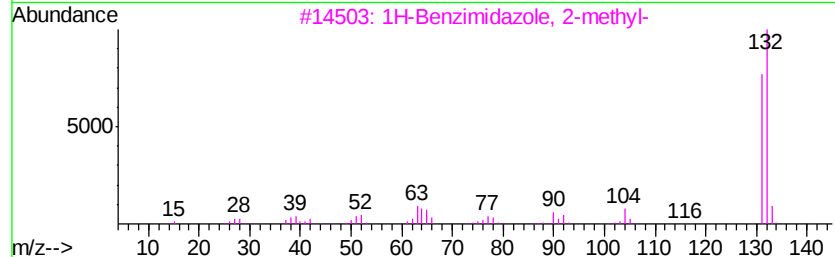
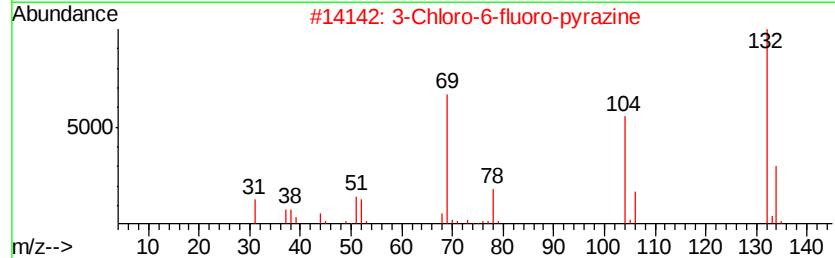
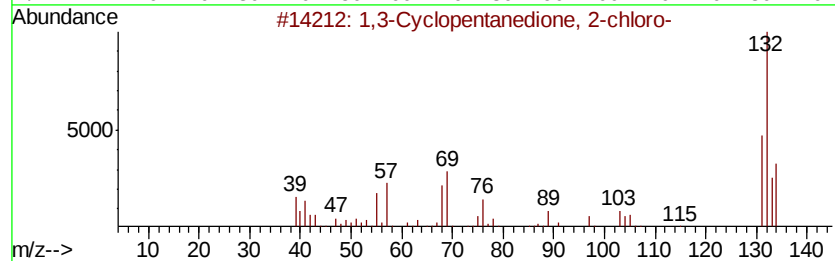
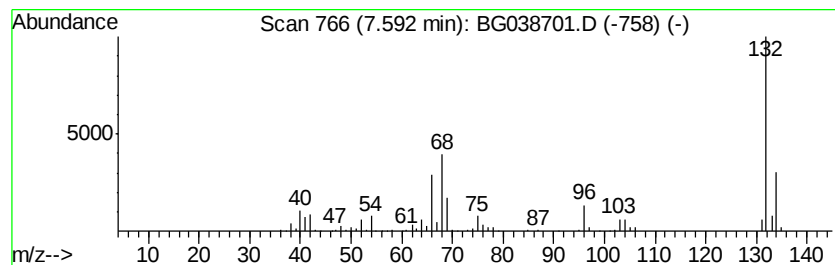
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	85.00 ng	562777	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	37
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4			5-Aminoindole	132	C8H8N2	005192-03-0	22
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	4.7	ng	31246	1	8.06	132417	20.0
unknown7.59	7.59	85.0	ng	562777	1	8.06	132417	20.0