

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082418\
 Data File : BG036382.D
 Acq On : 23 Aug 2018 18:12
 Operator : JU/SJ
 Sample : PB112153BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112153BL

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-BG082318.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	5.174	314	321	329	rBV	36669	61959	1.76%	0.371%
2	5.632	393	399	409	rBV	762611	1204262	34.22%	7.209%
3	7.218	662	669	677	rBV	775728	1305199	37.09%	7.813%
4	7.600	726	734	742	rBV	708714	1217106	34.58%	7.286%
5	8.070	807	814	821	rBV	177916	298350	8.48%	1.786%
6	8.394	861	869	877	rBV	535828	916954	26.05%	5.489%
7	9.228	1003	1011	1019	rBV	436120	755564	21.47%	4.523%
8	10.879	1283	1292	1300	rVB	238274	427560	12.15%	2.559%
9	13.323	1700	1708	1717	rVB	1219833	1892272	53.77%	11.327%
10	14.692	1934	1941	1950	rBV2	394698	618874	17.58%	3.705%
11	16.179	2187	2194	2211	rBV	921678	1440321	40.92%	8.622%
12	17.436	2401	2408	2417	rVB2	579483	854077	24.27%	5.113%
13	20.045	2846	2852	2863	rBV	2610802	3519460	100.00%	21.068%
14	21.701	3127	3134	3142	rBV	671110	1130144	32.11%	6.765%
15	24.927	3672	3683	3693	rBV2	381655	1063264	30.21%	6.365%

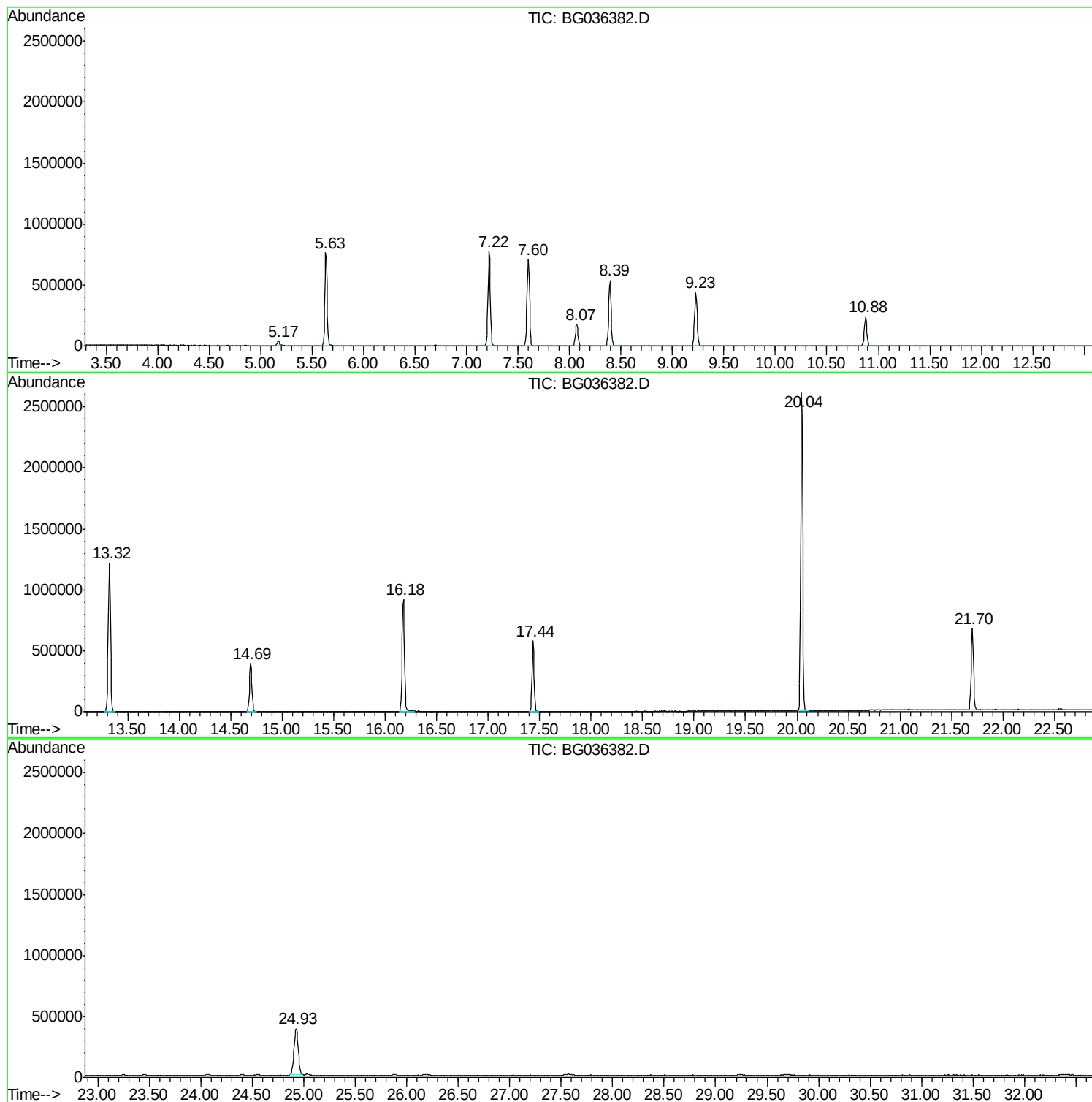
Sum of corrected areas: 16705366

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



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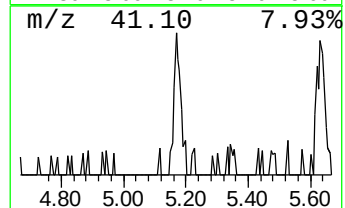
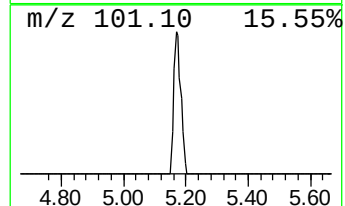
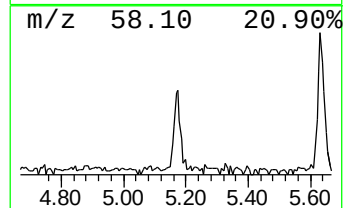
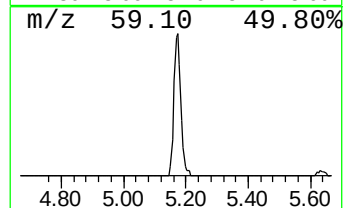
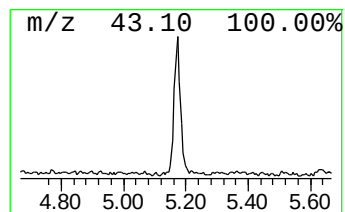
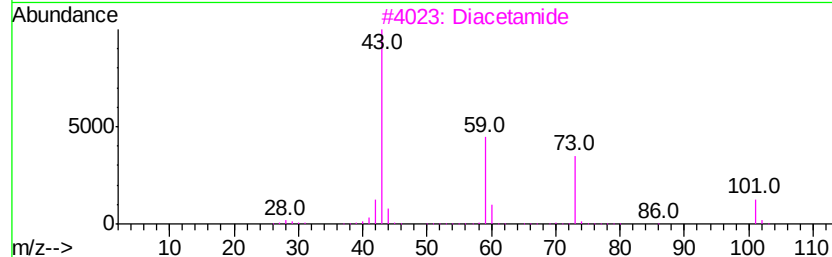
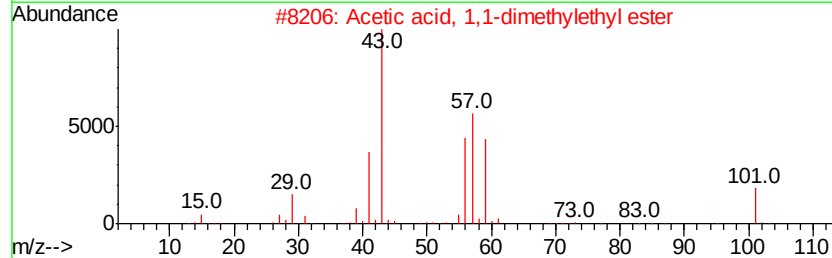
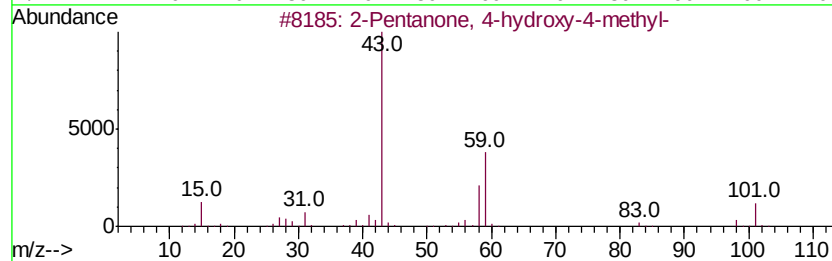
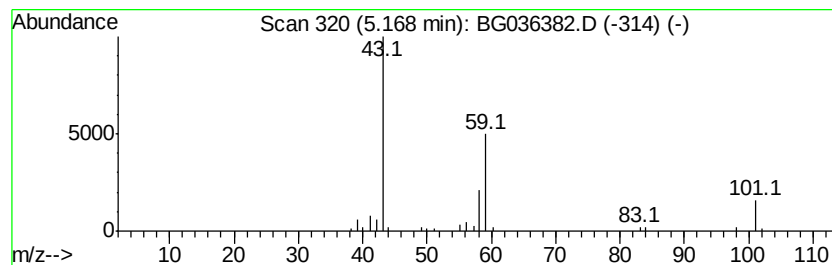
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.15 ng	61959	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Diacetamide	101	C4H7NO2	000625-77-4	36
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7



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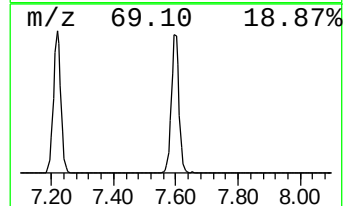
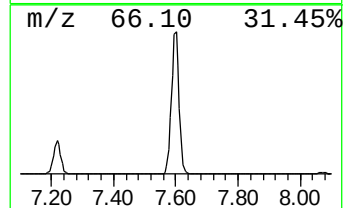
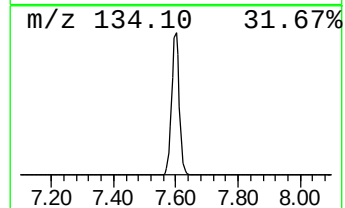
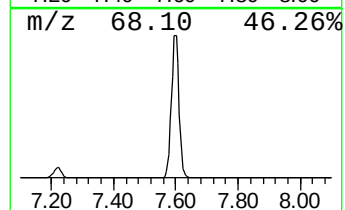
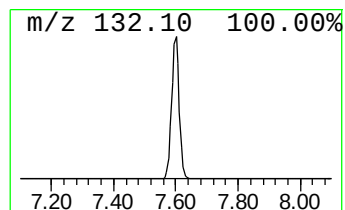
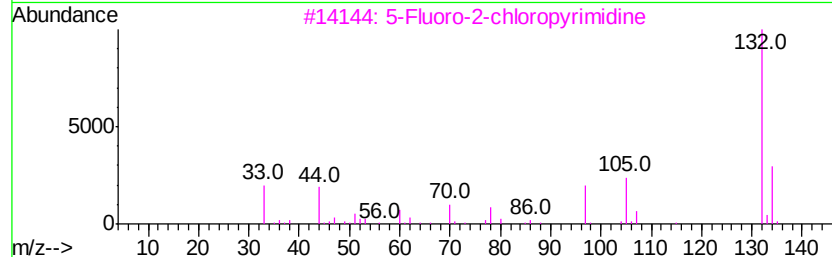
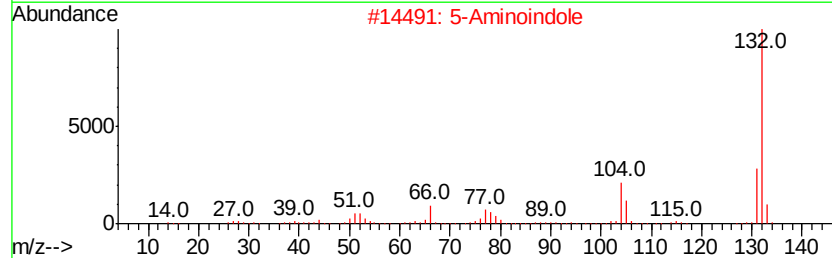
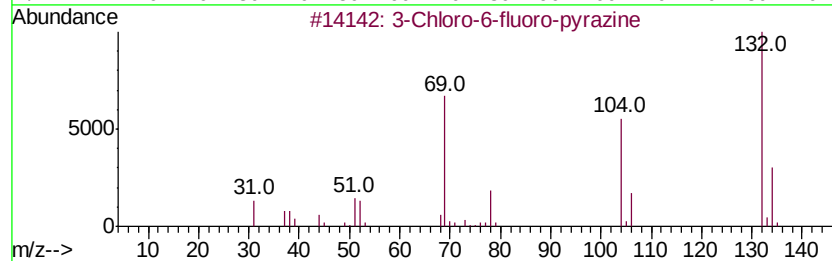
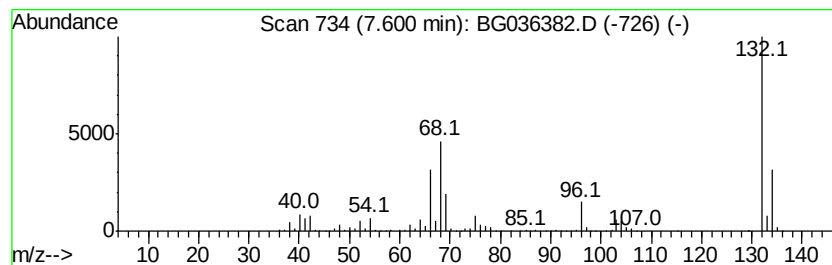
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	81.59 ng	1217110	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	22
3			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	4.2	ng	61959	1	8.07	298350	20.0
unknown7.60	7.60	81.6	ng	1217110	1	8.07	298350	20.0