

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG091117\
 Data File : BG028661.D
 Acq On : 12 Sep 2017 5:01
 Operator : SJ/JU
 Sample : I5181-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 090817-TIDO-F-U-M

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:\HPCHEM1\BNA_G\METHODS\8270-BG090717.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG028662.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.445	312	316	322	rBV	36940	55220	1.45%	0.391%
2	5.909	388	395	405	rBV	241448	411345	10.82%	2.915%
3	7.489	657	664	678	rVB2	140396	255146	6.71%	1.808%
4	7.871	721	729	743	rBV	402851	733717	19.30%	5.200%
5	8.335	799	808	815	rBV2	140984	249479	6.56%	1.768%
6	8.659	854	863	874	rVB	563644	1012706	26.63%	7.177%
7	9.505	998	1007	1018	rBV	462280	859566	22.60%	6.092%
8	11.150	1280	1287	1296	rBV	180516	339276	8.92%	2.404%
9	13.565	1690	1698	1710	rBV	1395400	2164699	56.93%	15.341%
10	14.945	1926	1933	1943	rVB2	300806	525664	13.82%	3.725%
11	16.432	2179	2186	2200	rBV	795740	1255974	33.03%	8.901%
12	17.695	2393	2401	2411	rBV2	452103	719272	18.92%	5.098%
13	20.280	2834	2841	2850	rBV	2925669	3802580	100.00%	26.949%
14	22.025	3131	3138	3146	rBV	496951	877789	23.08%	6.221%
15	25.533	3724	3735	3748	rVB2	267563	847682	22.29%	6.008%

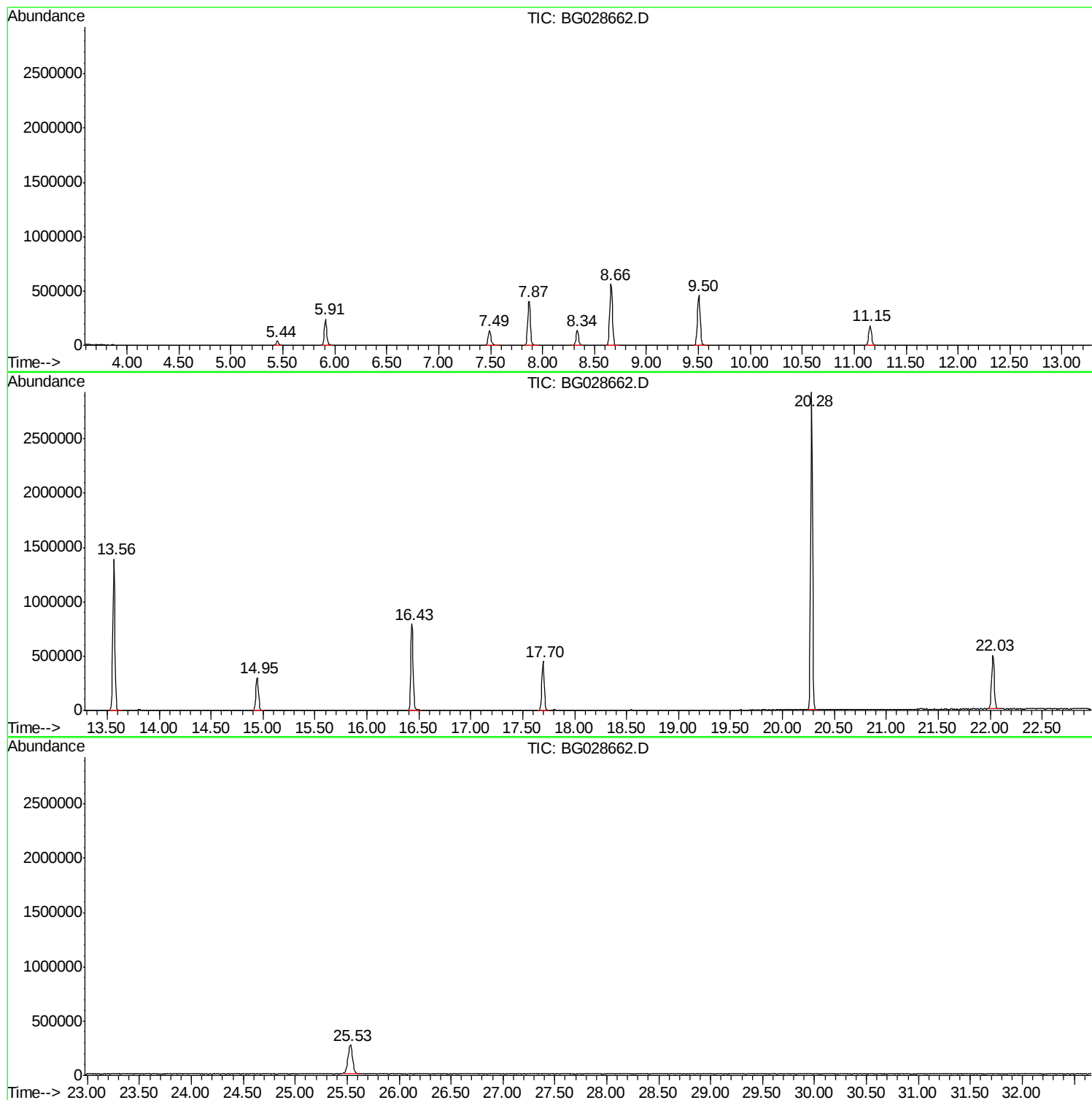
Sum of corrected areas: 14110115

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



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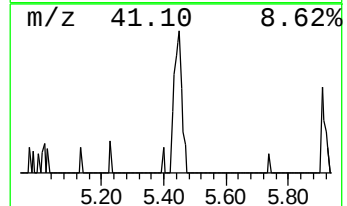
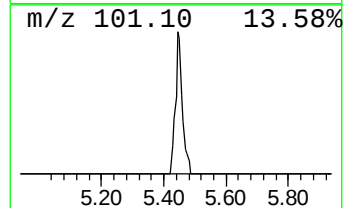
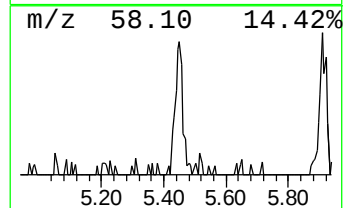
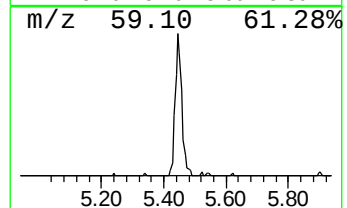
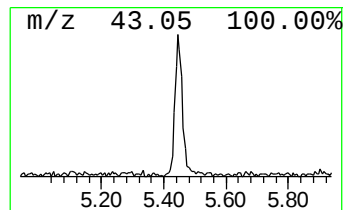
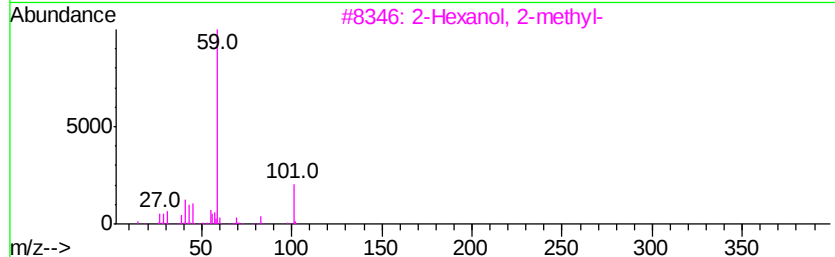
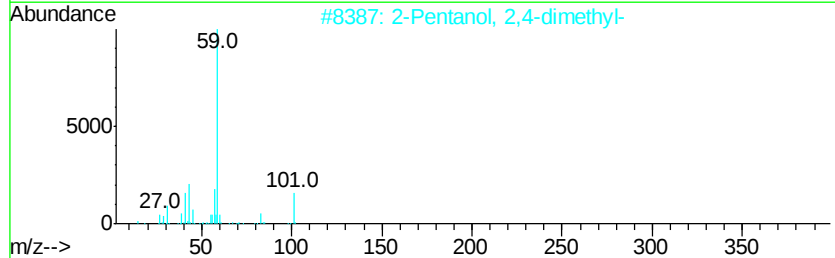
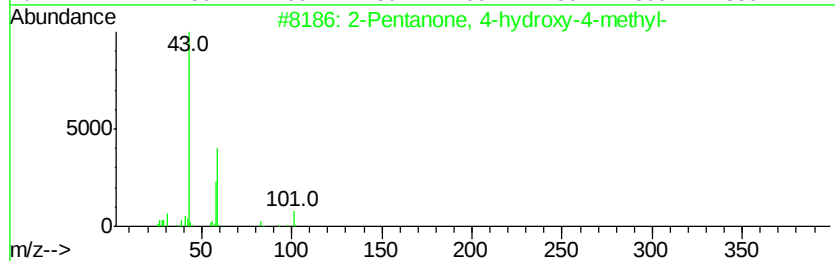
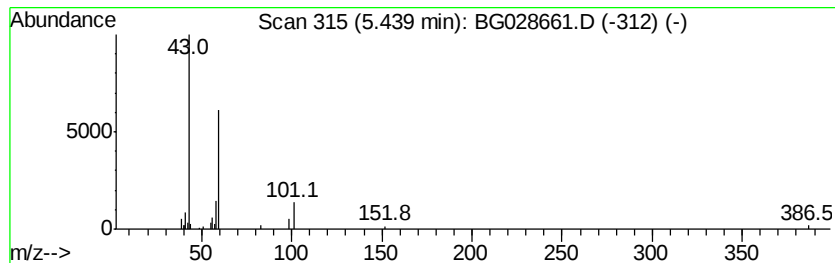
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.44	4.43 ng	55220	1,4-Dichlorobenzene-d4	8.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	23
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	12



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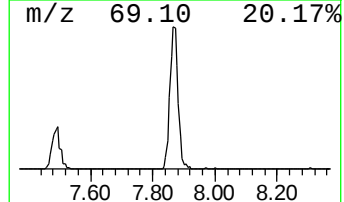
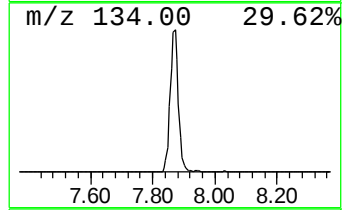
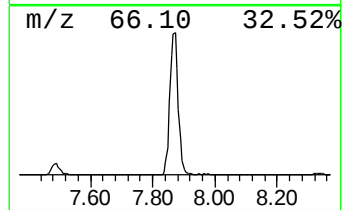
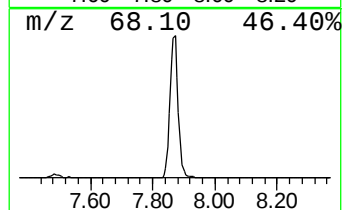
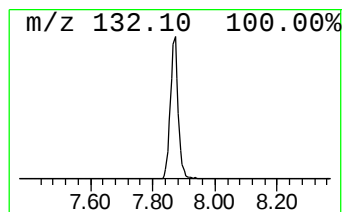
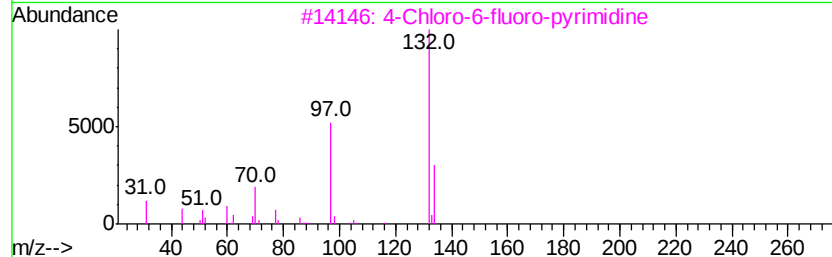
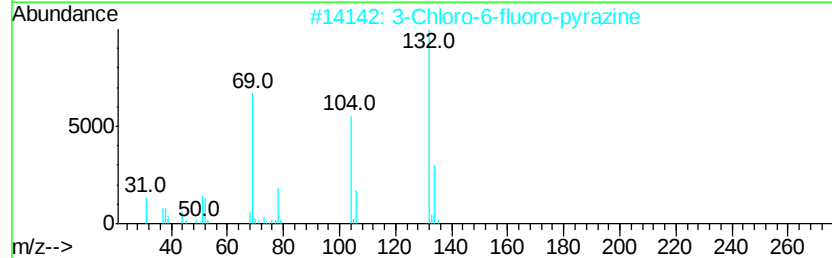
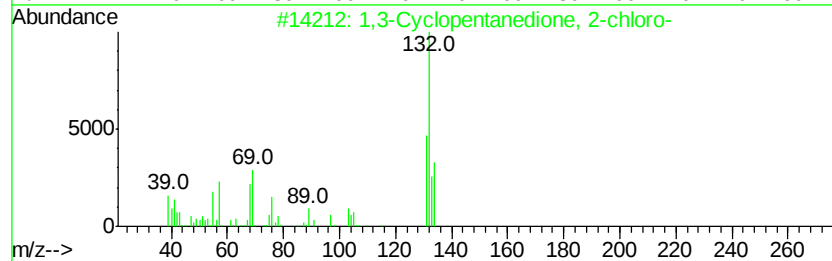
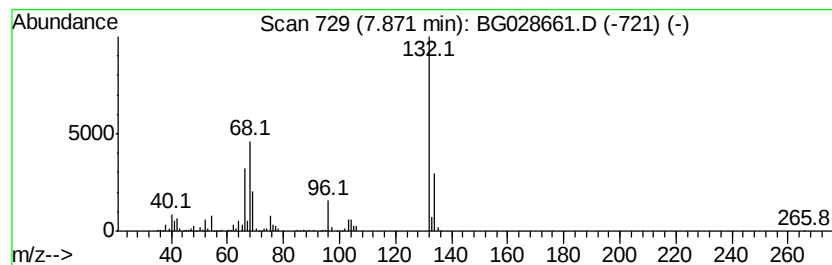
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Peak Number 2 unknown7.87 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	58.82 ng	733717	1,4-Dichlorobenzene-d4	8.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3			4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
4			5-Aminoindole	132	C8H8N2	005192-03-0	22
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.44	4.4	ng	55220	1	8.34	249479	20.0
unknown7.87	7.87	58.8	ng	733717	1	8.34	249479	20.0