

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092317\
 Data File : BG028902.D
 Acq On : 23 Sep 2017 17:38
 Operator : SJ/JU
 Sample : I5298-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 091817-SIDI-F-D-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-BG091217.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	5.380	298	305	312	rBV2	23659	38950	1.17%	0.368%
2	5.844	377	384	393	rBV	130548	232550	6.99%	2.194%
3	7.431	648	654	666	rBV	79375	162226	4.88%	1.531%
4	7.812	710	719	728	rBV	241652	449730	13.52%	4.243%
5	8.277	791	798	805	rBV2	80467	144842	4.35%	1.367%
6	8.600	846	853	862	rVB	352685	634781	19.08%	5.990%
7	9.446	990	997	1010	rBV	280112	540032	16.23%	5.096%
8	11.097	1270	1278	1289	rBV2	102373	202203	6.08%	1.908%
9	13.512	1682	1689	1699	rBV	919857	1458198	43.84%	13.759%
10	14.892	1918	1924	1932	rBV3	189052	318430	9.57%	3.005%
11	16.385	2171	2178	2190	rBV	665583	1063584	31.97%	10.036%
12	17.642	2386	2392	2405	rBV2	349131	570907	17.16%	5.387%
13	20.233	2825	2833	2844	rBV	2519294	3326360	100.00%	31.386%
14	21.961	3120	3127	3137	rBV2	406156	721821	21.70%	6.811%
15	25.415	3704	3715	3731	rVB2	235265	733503	22.05%	6.921%

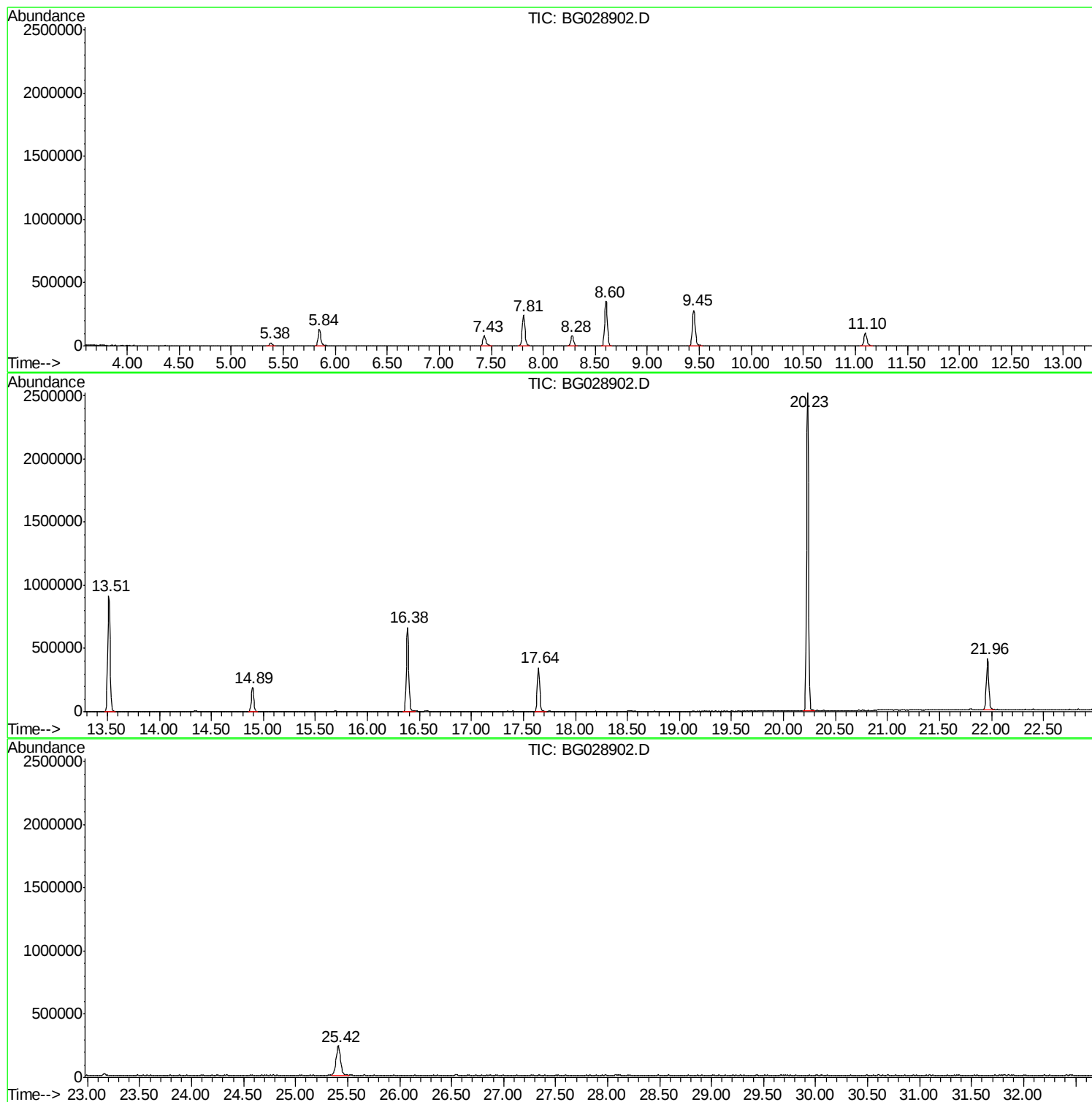
Sum of corrected areas: 10598117

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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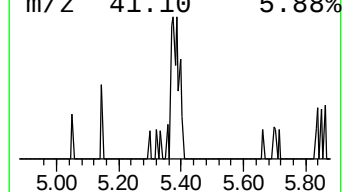
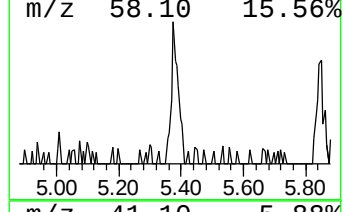
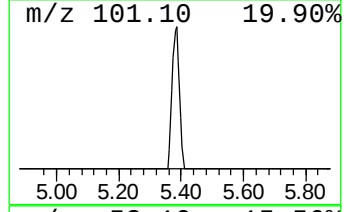
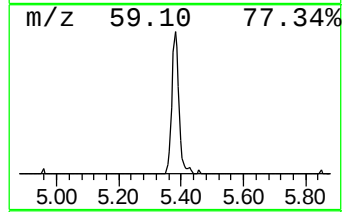
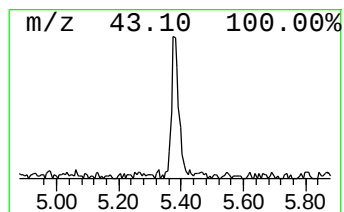
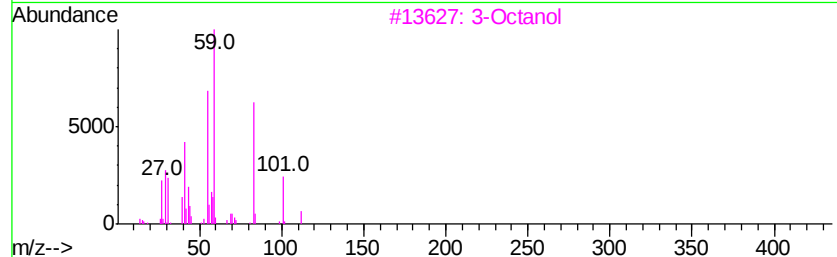
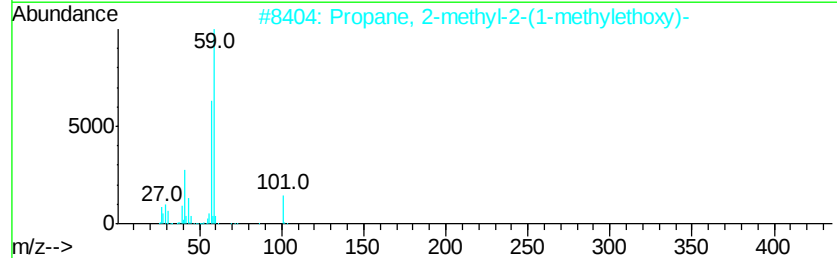
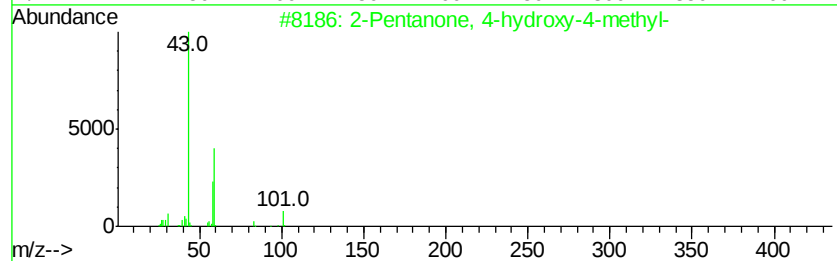
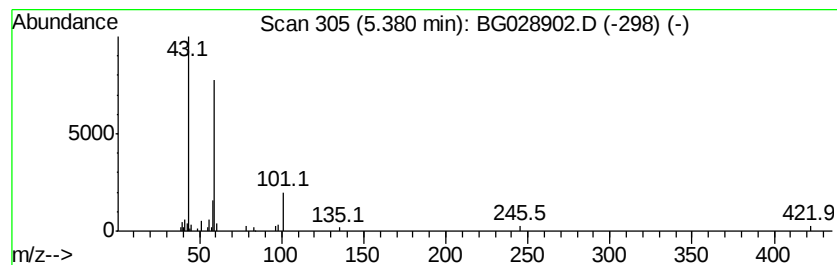
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.38	5.38 ng	38950	1,4-Dichlorobenzene-d4	8.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	45
3			3-Octanol	130	C8H18O	000589-98-0	38
4			2,3-Dimethyloxirane-2-carboxylic...	130	C6H10O3	1000306-64-4	36
5			3-Hexanol	102	C6H14O	000623-37-0	33



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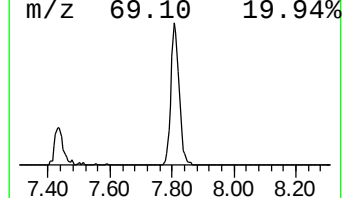
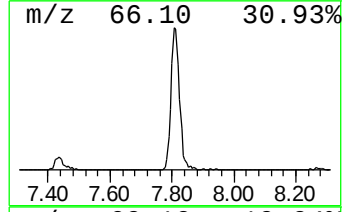
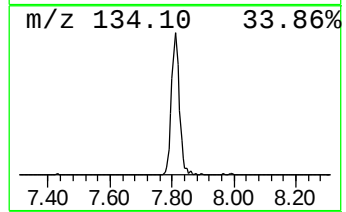
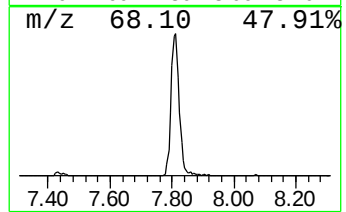
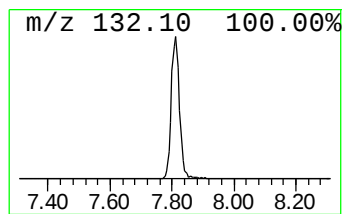
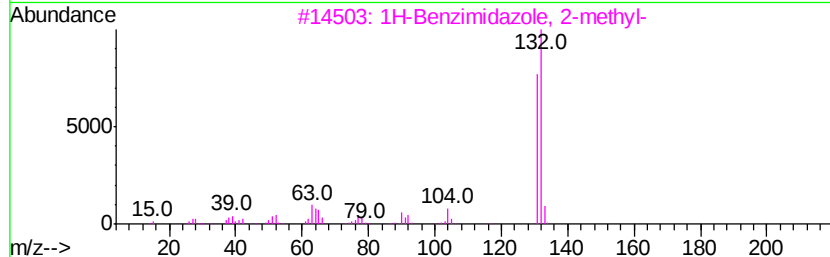
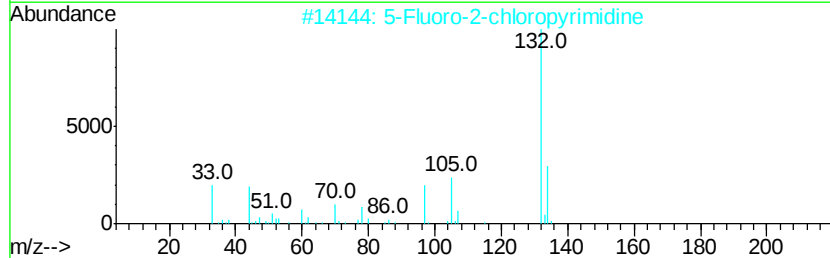
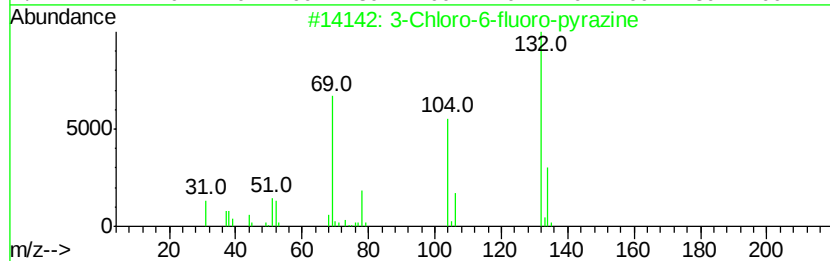
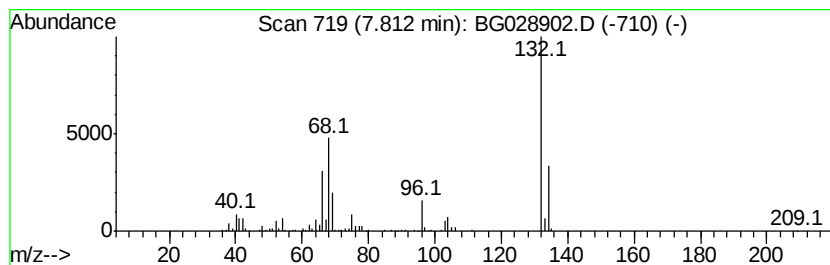
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Peak Number 2 unknown7.81 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	62.10 ng	449730	1,4-Dichlorobenzene-d4	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		3-Ethylphenyl isocyanate	147	C9H9NO	023138-58-1	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.38	5.4	ng	38950	1	8.28	144842	20.0
unknown7.81	7.81	62.1	ng	449730	1	8.28	144842	20.0