

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG040519\
 Data File : BG040381.D
 Acq On : 5 Apr 2019 17:45
 Operator : JU/SJ
 Sample : PB118645BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118645BL

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	5.136	343	348	359	rVB	81781	128033	3.93%	0.758%
2	5.600	421	427	441	rBV	631827	1047077	32.15%	6.199%
3	7.198	692	699	716	rBV	611207	1082197	33.23%	6.406%
4	7.574	753	763	771	rBV	623575	1079656	33.15%	6.391%
5	8.044	836	843	850	rBV	216450	369188	11.34%	2.186%
6	8.367	890	898	910	rBV	715660	1228741	37.73%	7.274%
7	9.219	1035	1043	1054	rBV	543779	996430	30.60%	5.899%
8	10.858	1315	1322	1338	rBV	261226	506508	15.55%	2.998%
9	13.297	1730	1737	1749	rBV	1442159	2287240	70.23%	13.540%
10	14.678	1964	1972	1984	rBV2	413923	698157	21.44%	4.133%
11	16.164	2219	2225	2241	rBV	567646	1013313	31.11%	5.999%
12	17.421	2433	2439	2452	rBV2	564675	888033	27.27%	5.257%
13	19.783	2829	2841	2842	rBV3	18009	45675	1.40%	0.270%
14	19.824	2842	2848	2855	rVB6	20351	62747	1.93%	0.371%
15	20.018	2875	2881	2891	rVB	2323109	3256764	100.00%	19.280%
16	21.693	3159	3166	3177	rBV	658719	1121152	34.43%	6.637%
17	24.971	3714	3724	3745	rVB3	357914	1081322	33.20%	6.401%

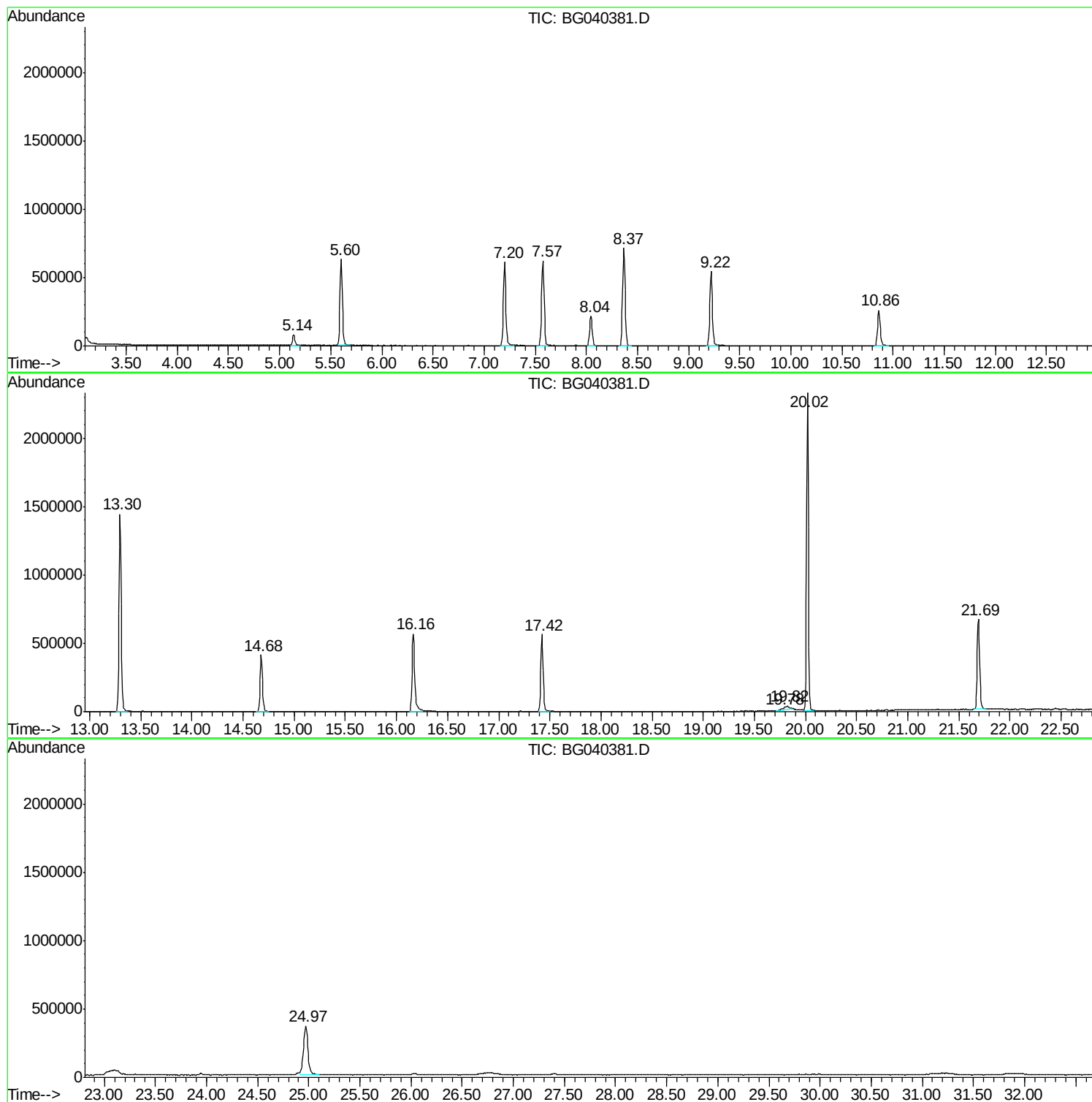
Sum of corrected areas: 16892233

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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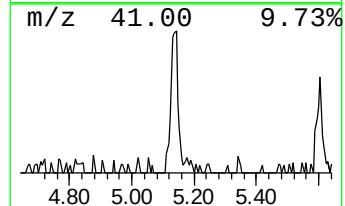
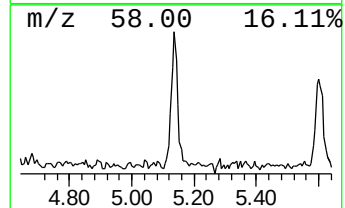
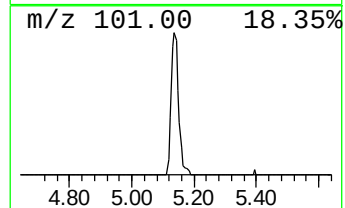
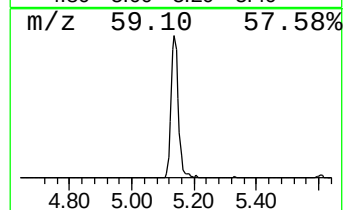
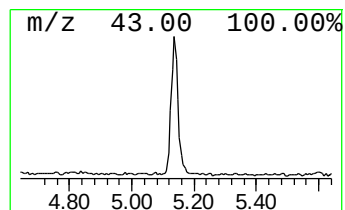
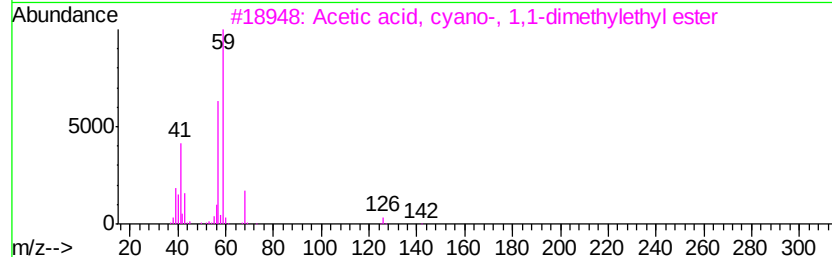
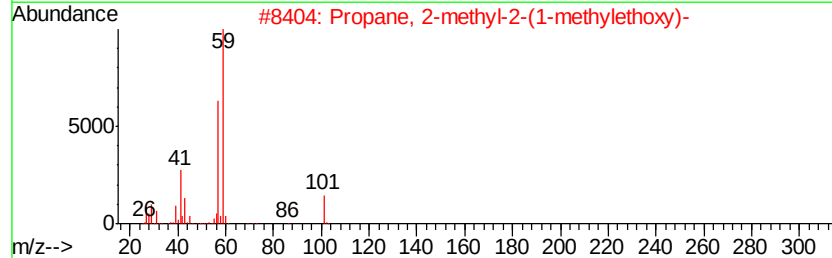
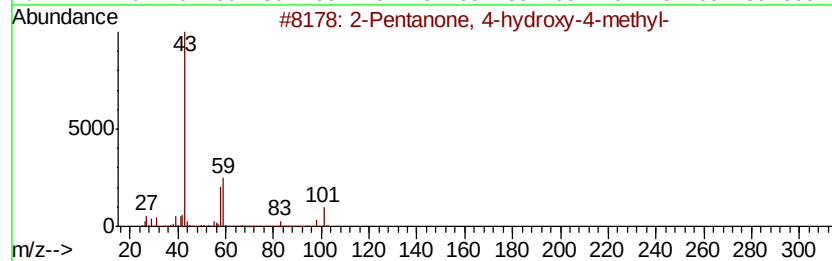
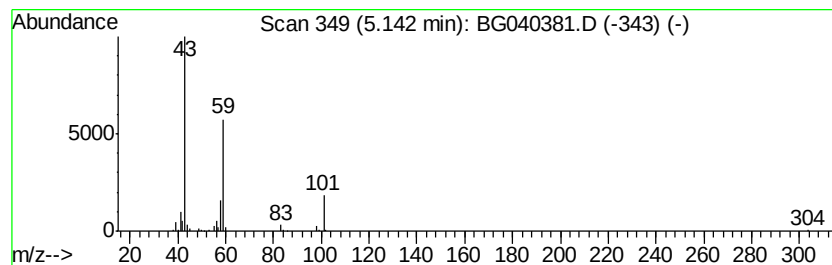
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	6.94 ng	128033	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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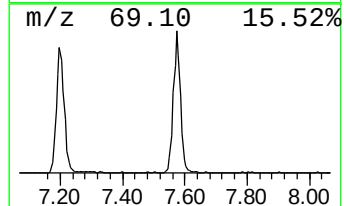
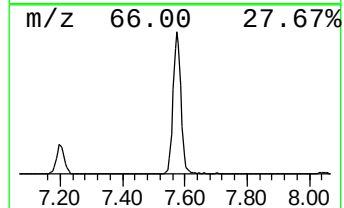
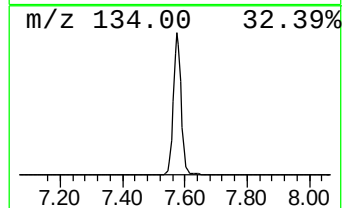
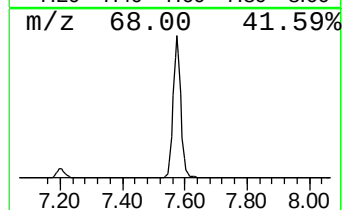
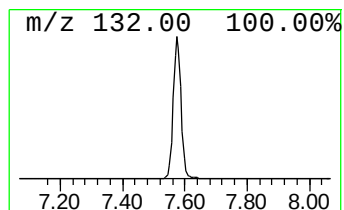
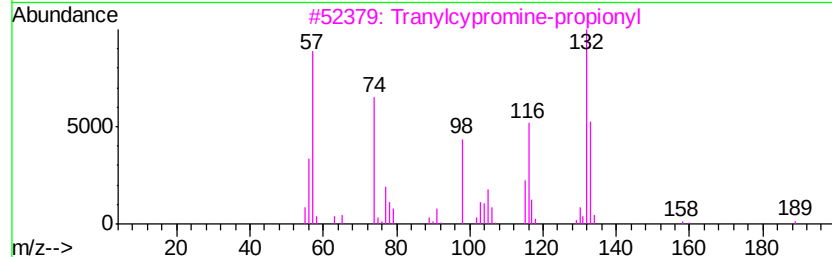
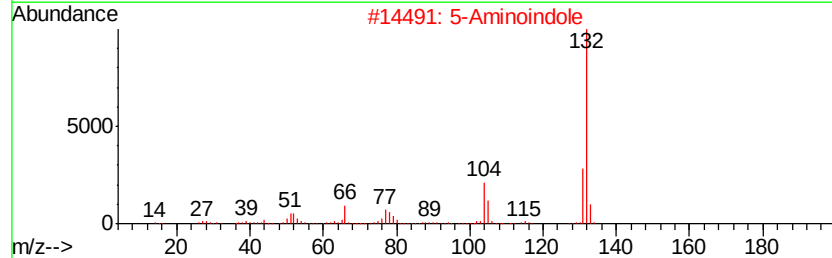
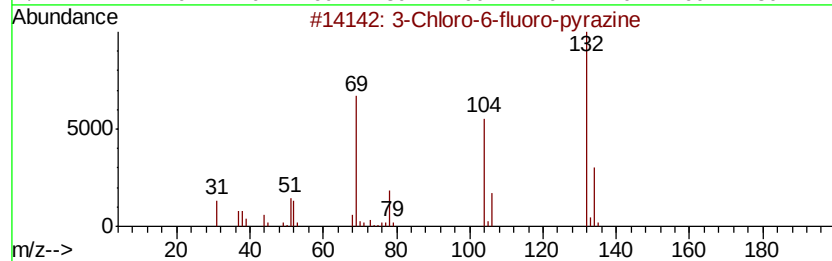
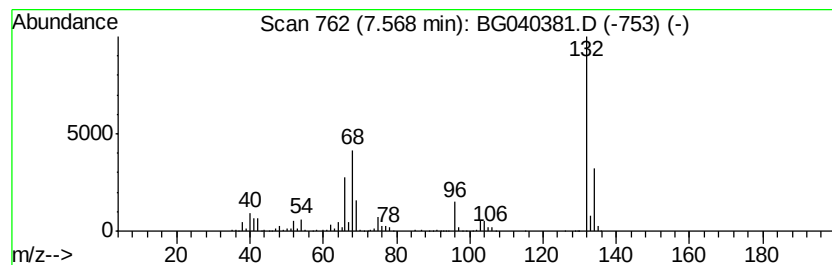
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Peak Number 2 unknown7.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	58.49 ng	1079660	1,4-Dichlorobenzene-d4	8.04

Hit#	of	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	12
3	Tranvlcy	promine-propionyl	189	C12H15NO	1000123-86-3	10
4	5-Fluoro-2-chloro	pyrimidine	132	C4H2ClFN2	062802-42-0	9
5	1-Methylpyrrolo[1,2-a]	pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	6.9	ng	128033	1	8.04	369188	20.0
unknown7.57	7.57	58.5	ng	1079660	1	8.04	369188	20.0