

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG062817\
 Data File : BG027719.D
 Acq On : 28 Jun 2017 21:28
 Operator : SJ/JU
 Sample : PB100124BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB100124BL

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-BG062817.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	5.604	285	292	303	rBV	119910	192567	4.70%	1.198%
2	6.056	362	369	385	rVB	687381	1167977	28.49%	7.263%
3	7.613	627	634	645	rBV	675583	1249220	30.47%	7.769%
4	8.007	693	701	712	rBV	652985	1194785	29.14%	7.430%
5	8.477	773	781	788	rBV	100868	183264	4.47%	1.140%
6	8.800	829	836	846	rBV	453067	840477	20.50%	5.227%
7	9.640	970	979	988	rBV	381436	724224	17.66%	4.504%
8	11.291	1253	1260	1270	rVB	154308	291992	7.12%	1.816%
9	13.688	1660	1668	1676	rBV	1142856	1814204	44.25%	11.282%
10	15.063	1893	1902	1911	rVB2	253361	417305	10.18%	2.595%
11	16.550	2147	2155	2170	rBV	1054211	1707903	41.66%	10.621%
12	17.813	2362	2370	2379	rBV2	412647	655028	15.98%	4.073%
13	20.392	2802	2809	2818	rBV	2859791	4099972	100.00%	25.496%
14	22.172	3104	3112	3121	rVB	425270	782163	19.08%	4.864%
15	25.780	3710	3726	3740	rBV2	237926	759469	18.52%	4.723%

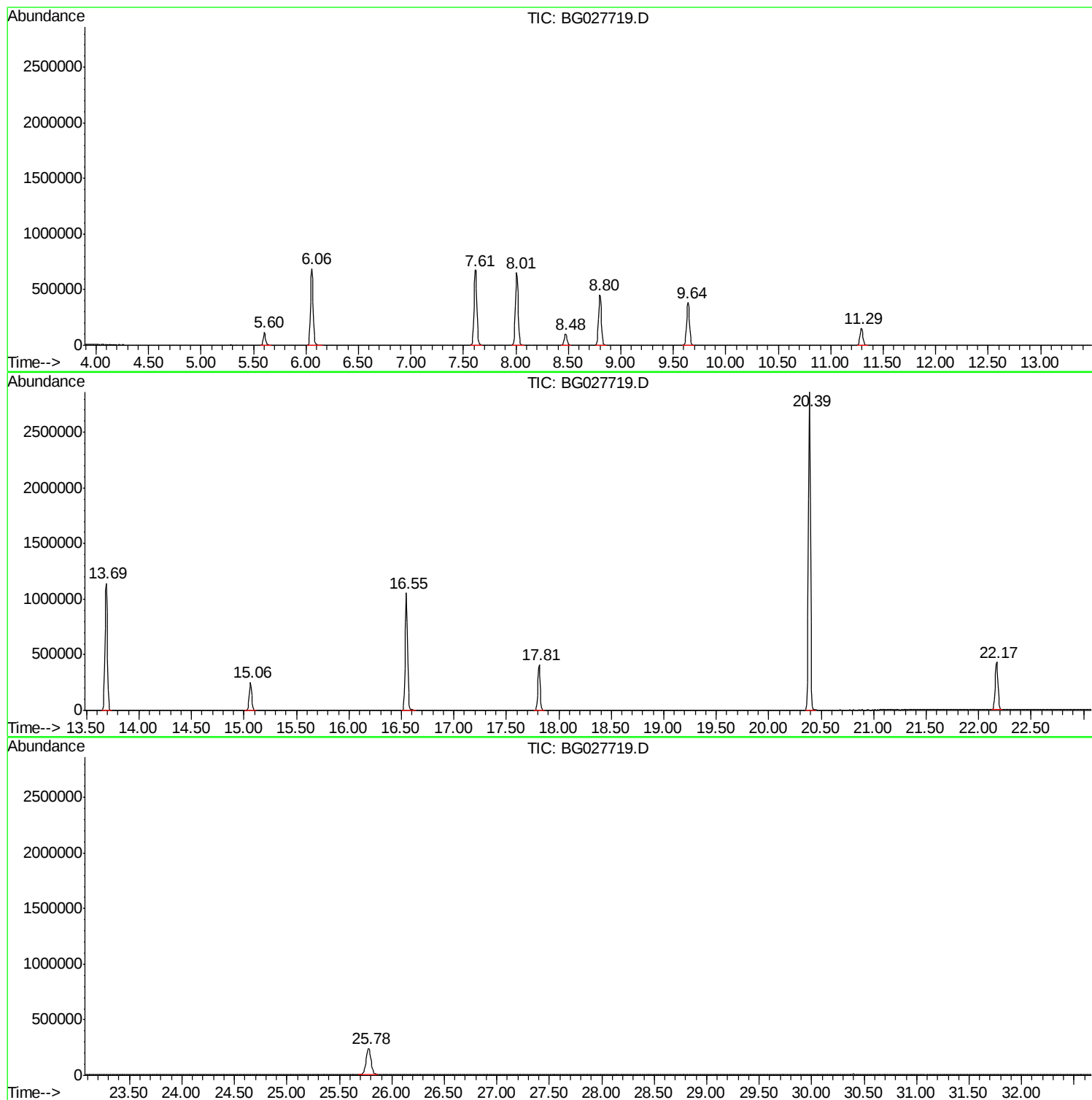
Sum of corrected areas: 16080550

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062817.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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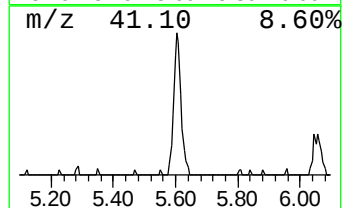
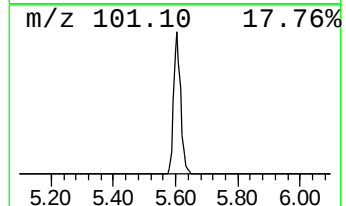
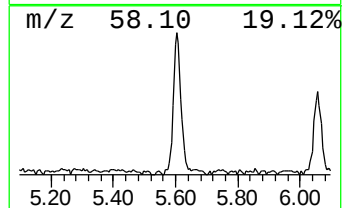
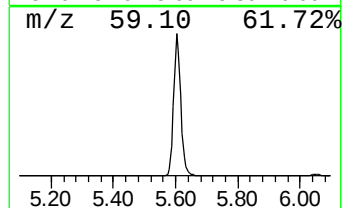
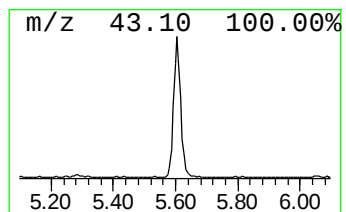
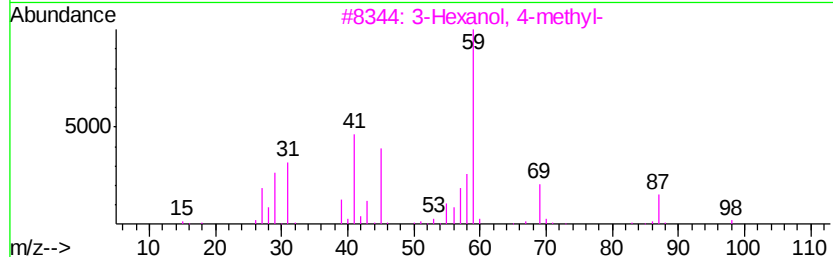
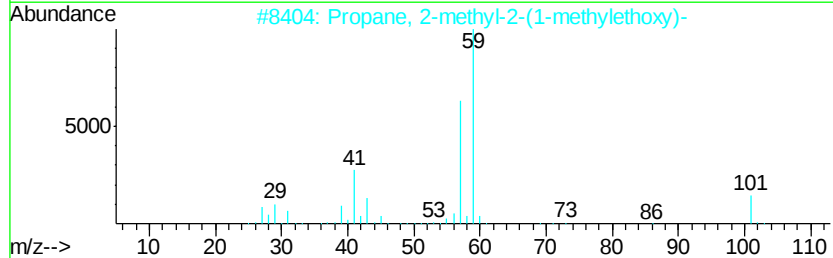
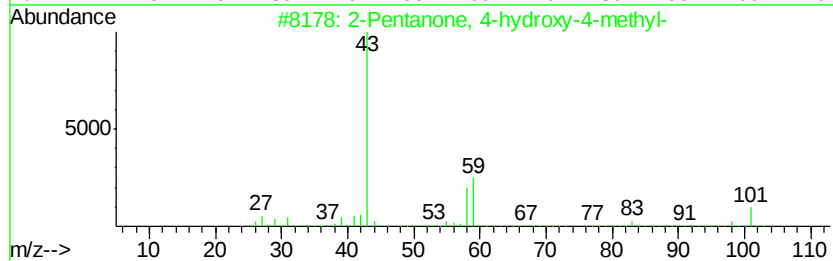
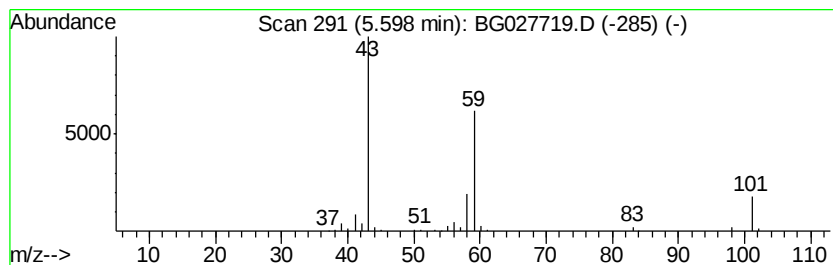
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.60	21.02 ng	192567	1,4-Dichlorobenzene-d4	8.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10



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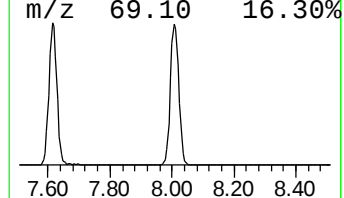
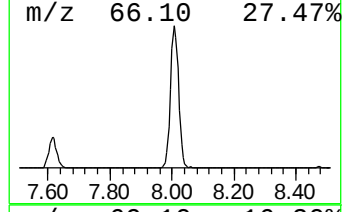
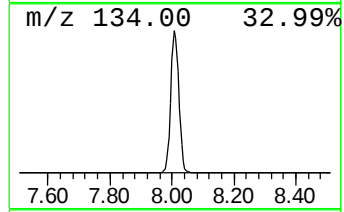
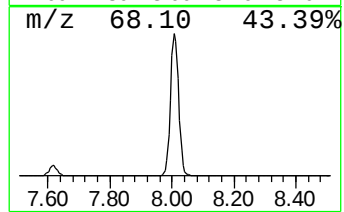
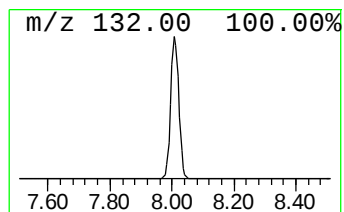
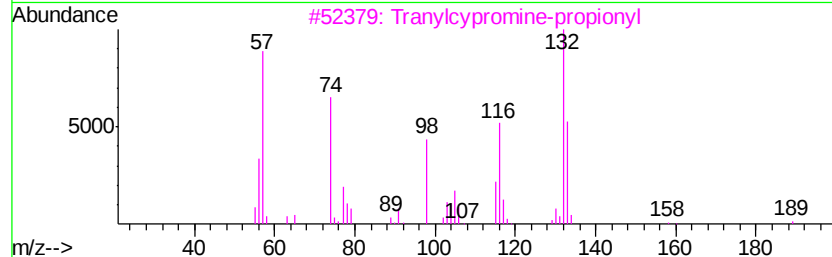
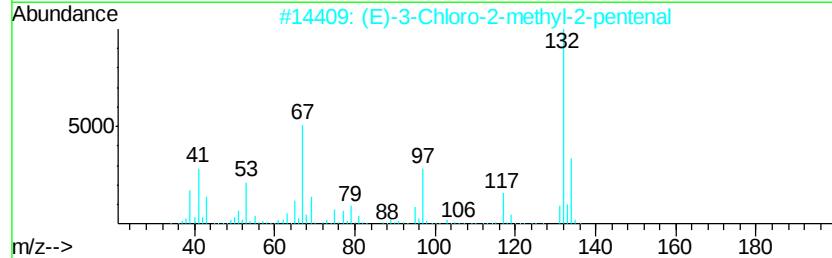
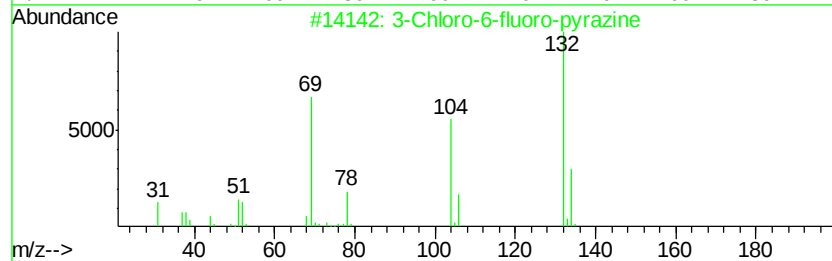
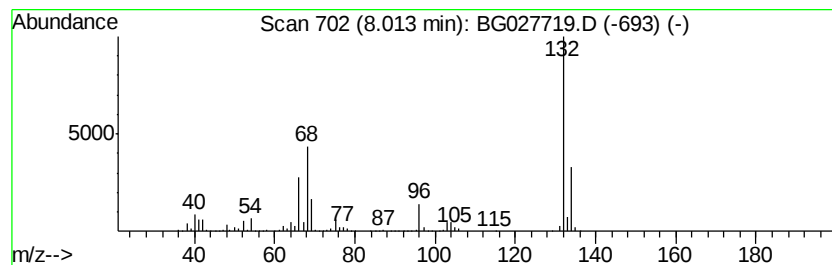
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Peak Number 2 unknown8.01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.01	130.39 ng	1194790	1,4-Dichlorobenzene-d4	8.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.60	21.0	ng	192567	1	8.48	183264	20.0
unknown8.01	8.01	130.4	ng	1194790	1	8.48	183264	20.0