

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062117\
Data File : BG027571.D
Acq On : 21 Jun 2017 22:05
Operator : SJ/JU
Sample : PB99899BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB99899BL

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-BG062117.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	5.638	291	298	310	rVB	101326	167076	5.30%	1.330%
2	6.084	367	374	387	rBV	544559	950017	30.12%	7.563%
3	7.641	631	639	649	rBV	552597	1000596	31.72%	7.966%
4	8.035	697	706	720	rBV	522738	962107	30.50%	7.659%
5	8.505	779	786	793	rBV	84334	150022	4.76%	1.194%
6	8.828	833	841	850	rVB	356538	654434	20.75%	5.210%
7	9.668	975	984	993	rBV	303345	574933	18.23%	4.577%
8	11.325	1256	1266	1274	rBV	117423	226712	7.19%	1.805%
9	13.716	1665	1673	1682	rBV	840922	1350280	42.81%	10.750%
10	15.091	1900	1907	1915	rVB3	187525	316756	10.04%	2.522%
11	16.572	2152	2159	2167	rBV	849939	1346522	42.69%	10.720%
12	17.841	2368	2375	2381	rBV2	304471	497530	15.77%	3.961%
13	20.414	2807	2813	2821	rBV	2269833	3154262	100.00%	25.111%
14	22.206	3110	3118	3128	rBV	334749	627435	19.89%	4.995%
15	25.843	3725	3737	3753	rVB2	176426	582499	18.47%	4.637%

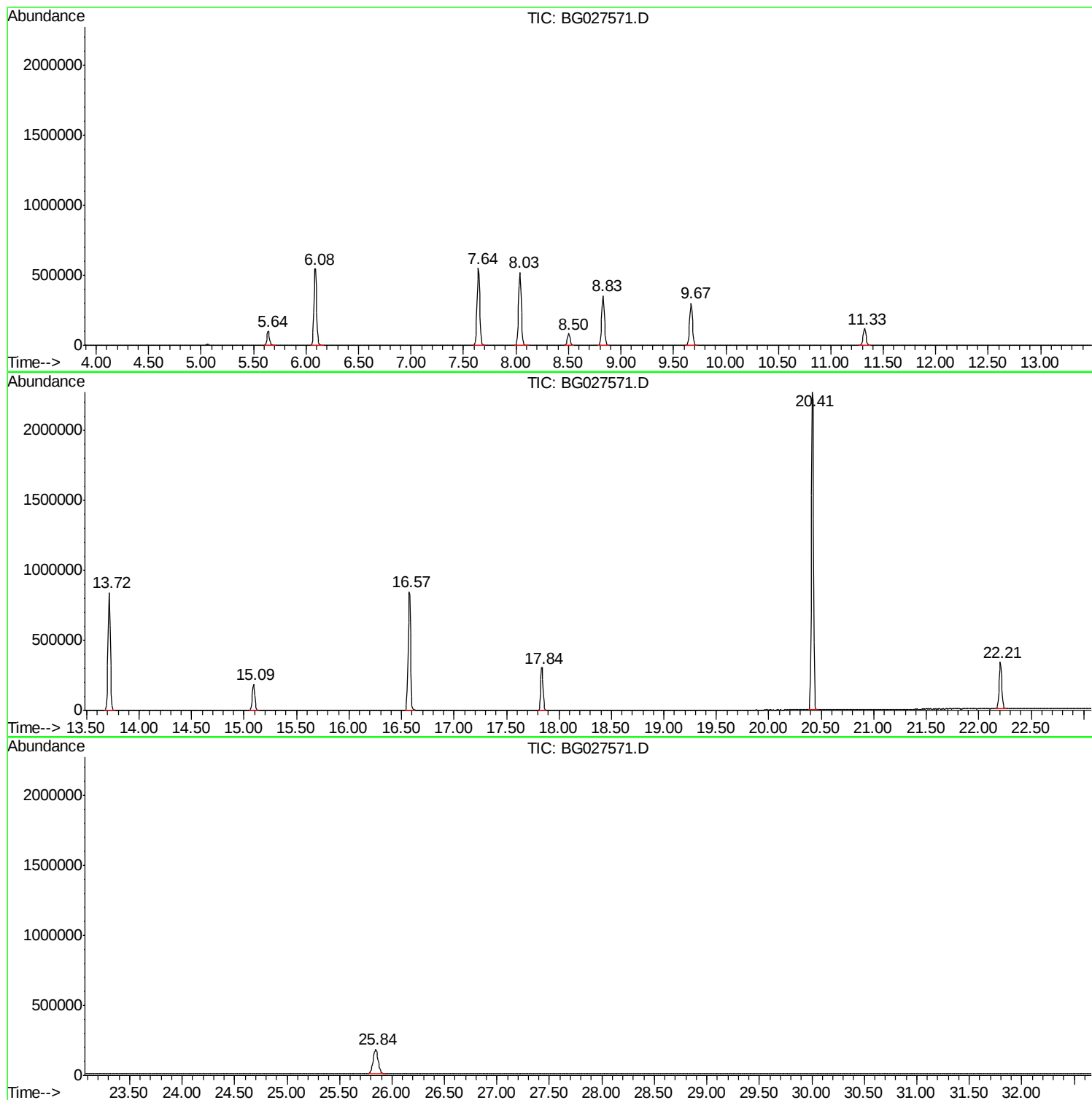
Sum of corrected areas: 12561181

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062117.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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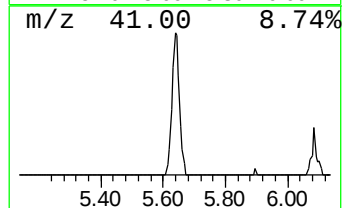
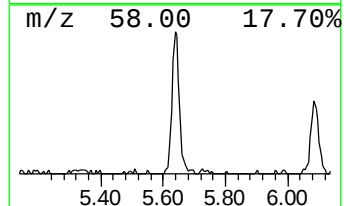
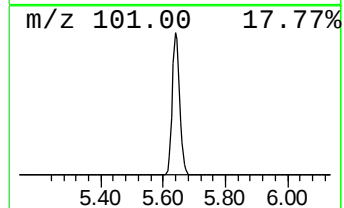
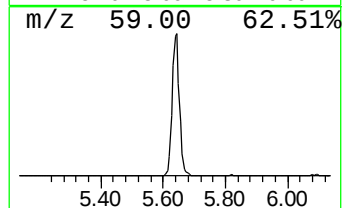
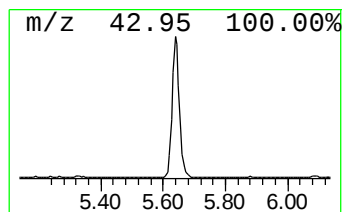
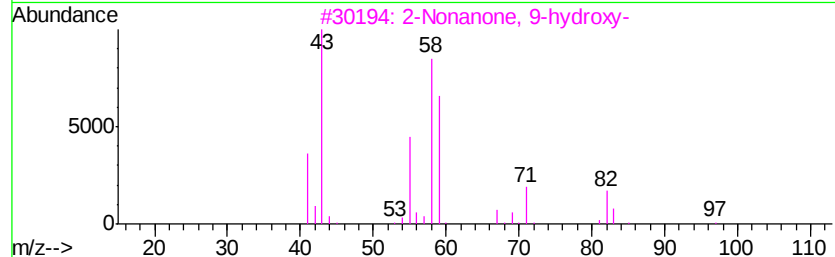
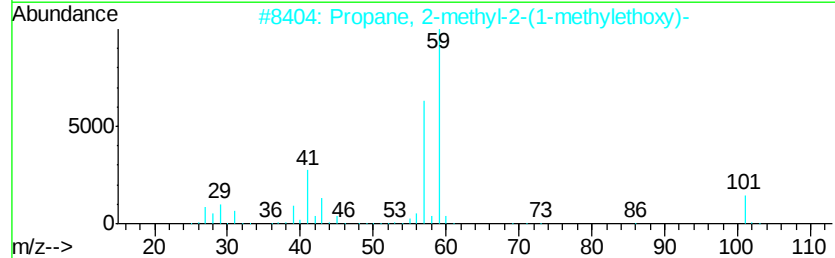
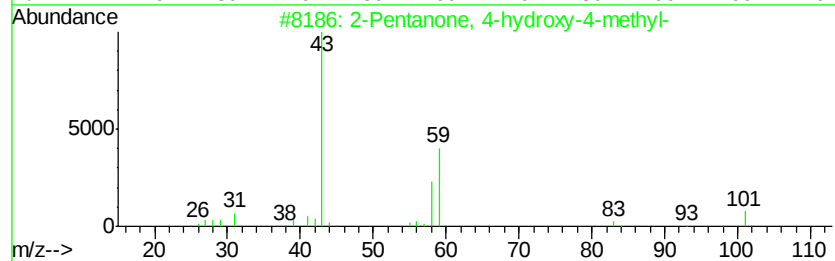
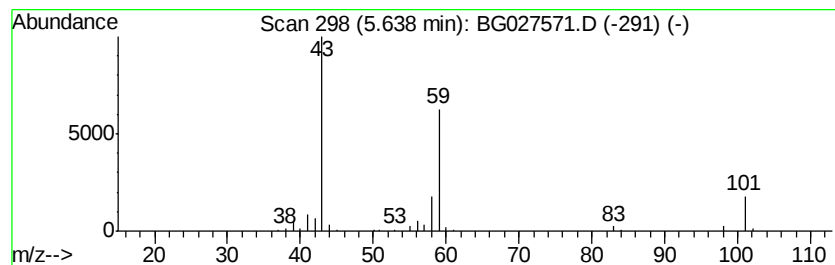
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	22.27 ng	167076	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	36
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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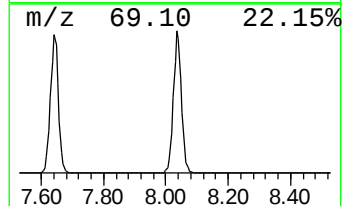
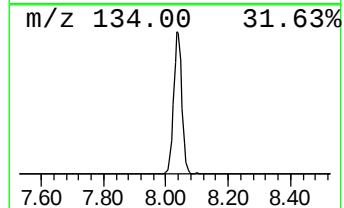
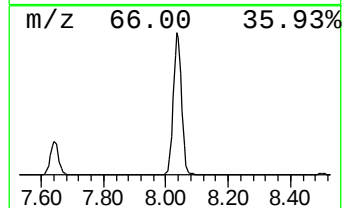
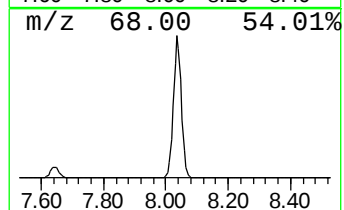
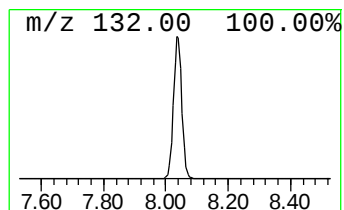
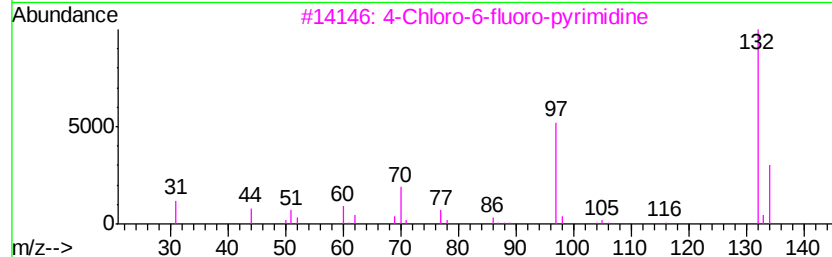
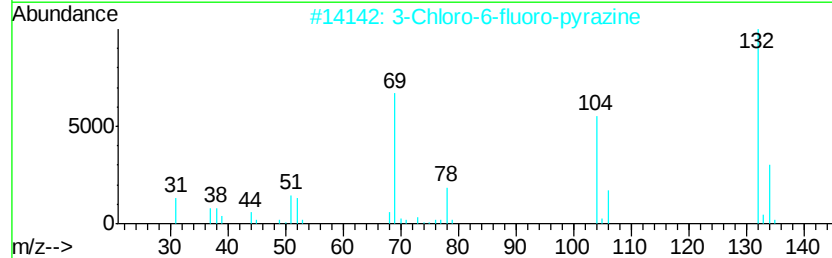
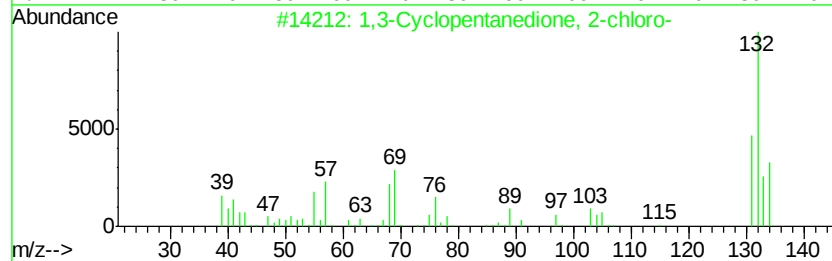
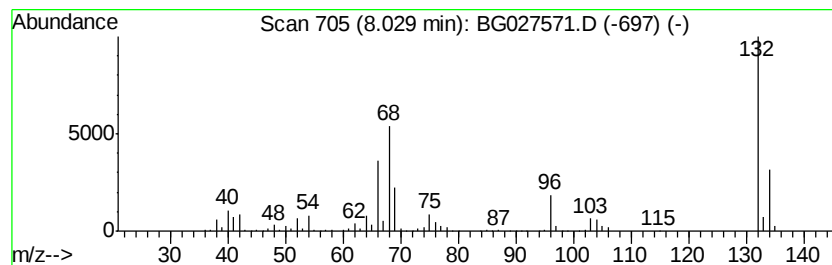
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Peak Number 2 unknown8.03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.03	128.26 ng	962107	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.64	22.3	ng	167076	1	8.50	150022	20.0
unknown8.03	8.03	128.3	ng	962107	1	8.50	150022	20.0