

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062117\
 Data File : BG027579.D
 Acq On : 22 Jun 2017 3:49
 Operator : SJ/JU
 Sample : PB99928BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99928BL

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	5.645	292	299	309	rVB	83941	134566	5.66%	1.356%
2	6.097	369	376	384	rBV	429757	731373	30.78%	7.370%
3	7.654	632	641	654	rBV	419410	767918	32.32%	7.738%
4	8.048	699	708	717	rBV	401594	733754	30.88%	7.394%
5	8.512	781	787	797	rBV	74607	133385	5.61%	1.344%
6	8.841	835	843	851	rVB	287131	508760	21.41%	5.126%
7	9.675	976	985	995	rBV	230284	432466	18.20%	4.358%
8	11.326	1259	1266	1275	rBV	100973	194399	8.18%	1.959%
9	13.718	1666	1673	1681	rBV	645215	1028402	43.28%	10.363%
10	15.092	1900	1907	1916	rVB2	166207	275537	11.60%	2.776%
11	16.579	2153	2160	2171	rBV	666620	1041900	43.85%	10.499%
12	17.842	2368	2375	2383	rBV	291085	448312	18.87%	4.517%
13	20.422	2807	2814	2822	rBV	1731014	2376228	100.00%	23.944%
14	22.214	3110	3119	3128	rBV	305451	572092	24.08%	5.765%
15	25.850	3724	3738	3755	rBV3	159219	545076	22.94%	5.492%

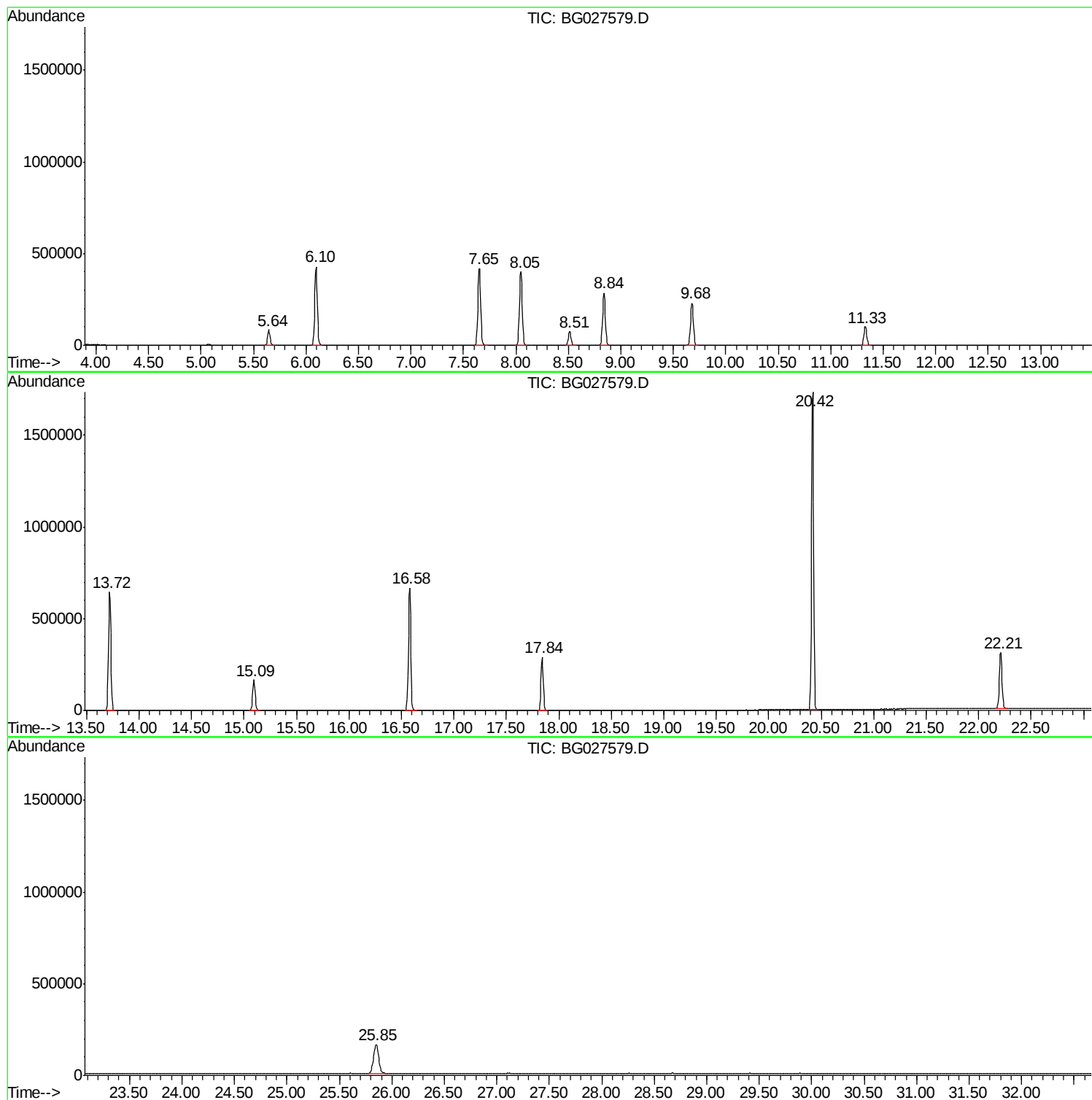
Sum of corrected areas: 9924168

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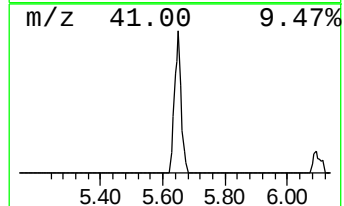
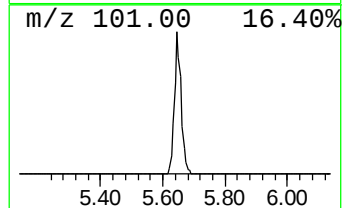
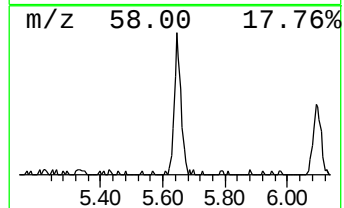
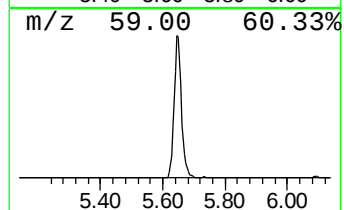
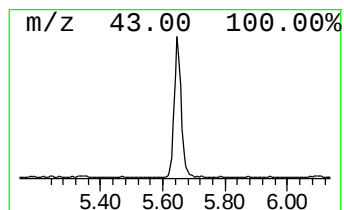
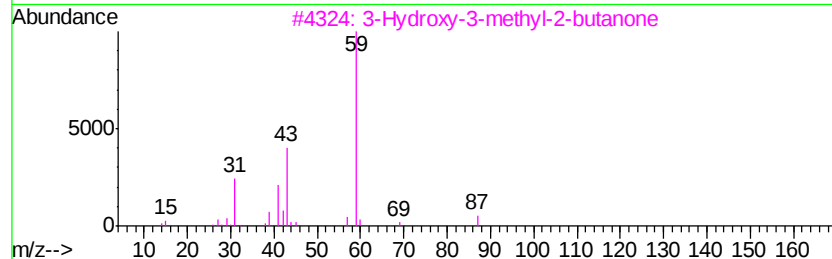
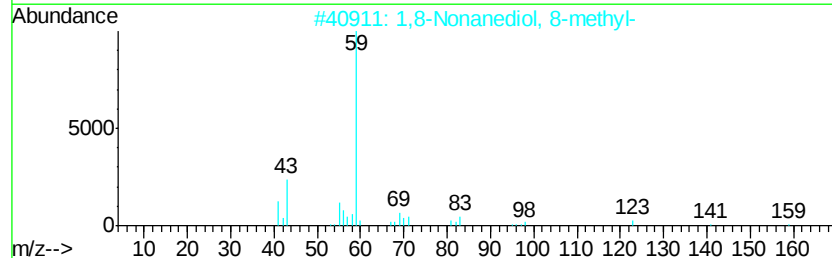
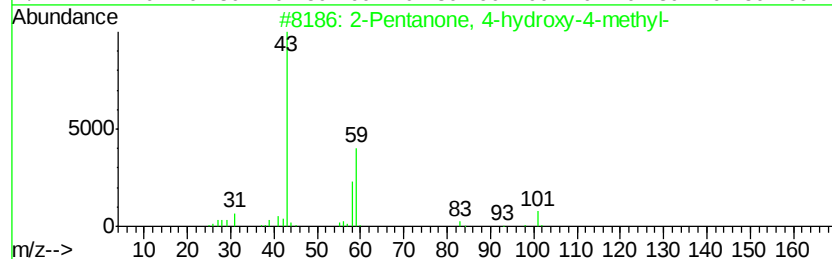
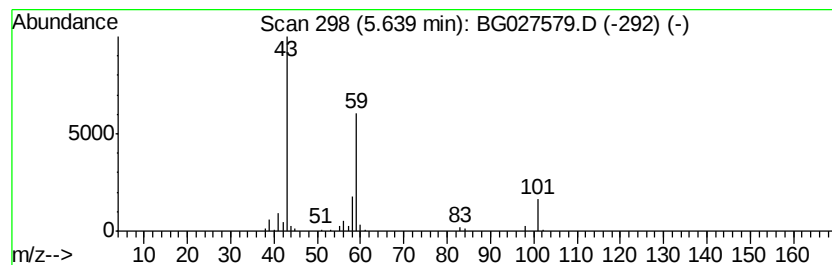
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	20.18 ng	134566	1,4-Dichlorobenzene-d4	8.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25
3		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5		Acetone	58	C3H6O	000067-64-1	12



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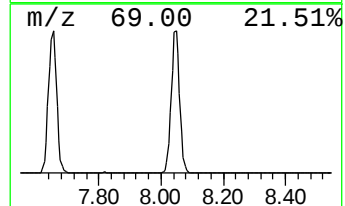
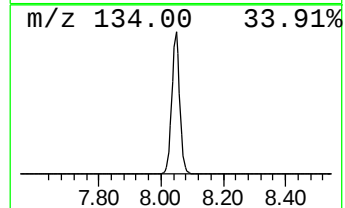
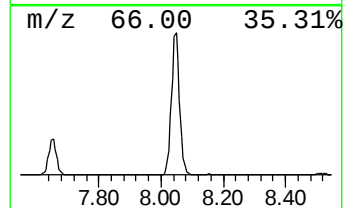
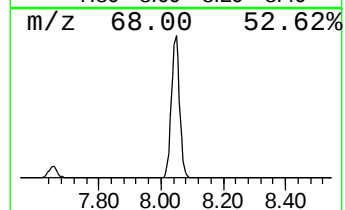
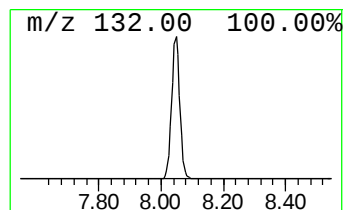
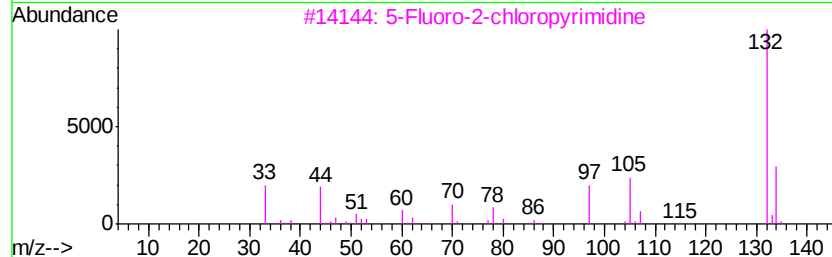
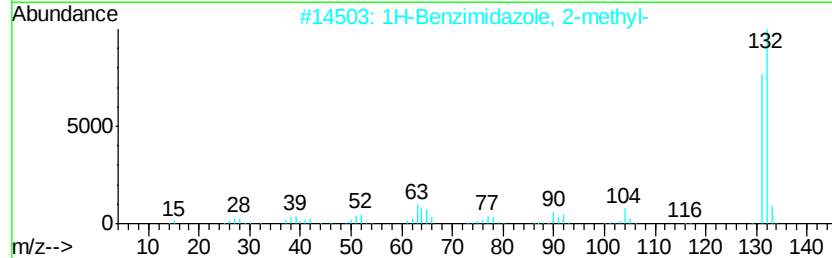
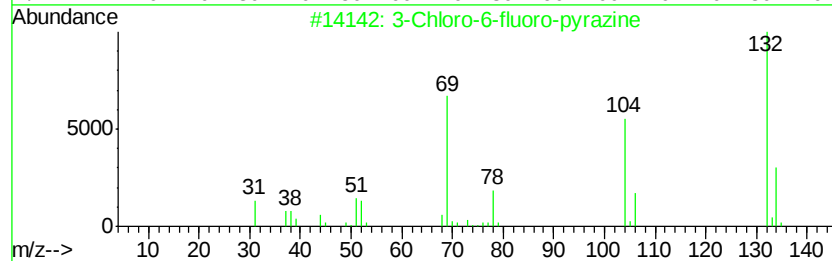
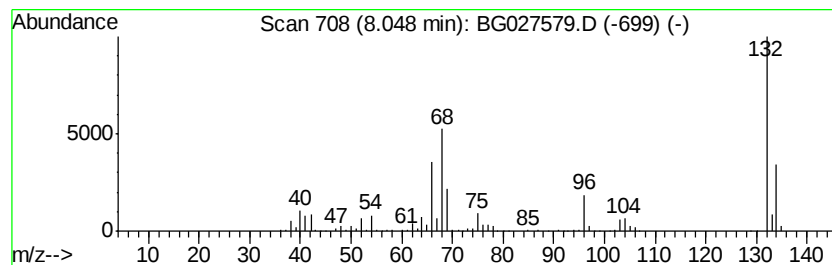
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Peak Number 2 unknown8.05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.05	110.02 ng	733754	1,4-Dichlorobenzene-d4	8.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
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	20.2	ng	134566	1	8.52	133385	20.0
unknown8.05	8.05	110.0	ng	733754	1	8.52	133385	20.0