

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG062217\
 Data File : BG027608.D
 Acq On : 22 Jun 2017 22:52
 Operator : SJ/JU
 Sample : I3657-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-FIELDBLANK

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	5.639	293	298	308	rVB	25072	38293	1.41%	0.379%
2	6.086	367	374	385	rVB	240330	399921	14.77%	3.959%
3	7.637	630	638	646	rBV	146308	271610	10.03%	2.689%
4	8.036	698	706	715	rBV	418858	750304	27.71%	7.428%
5	8.506	779	786	797	rBB	76347	135739	5.01%	1.344%
6	8.830	832	841	850	rBB	329879	597231	22.06%	5.913%
7	9.670	975	984	994	rBV	296615	563755	20.82%	5.581%
8	11.321	1257	1265	1275	rBV	106131	207145	7.65%	2.051%
9	13.712	1664	1672	1681	rBV	865219	1350702	49.89%	13.372%
10	15.087	1899	1906	1916	rBB2	174524	294766	10.89%	2.918%
11	16.573	2151	2159	2171	rBV	859212	1381751	51.04%	13.679%
12	17.837	2367	2374	2382	rVB2	278116	433386	16.01%	4.291%
13	20.416	2806	2813	2819	rBV	2094297	2707317	100.00%	26.803%
14	22.202	3110	3117	3125	rVB	273814	500297	18.48%	4.953%
15	25.839	3724	3736	3751	rVB3	144094	468752	17.31%	4.641%

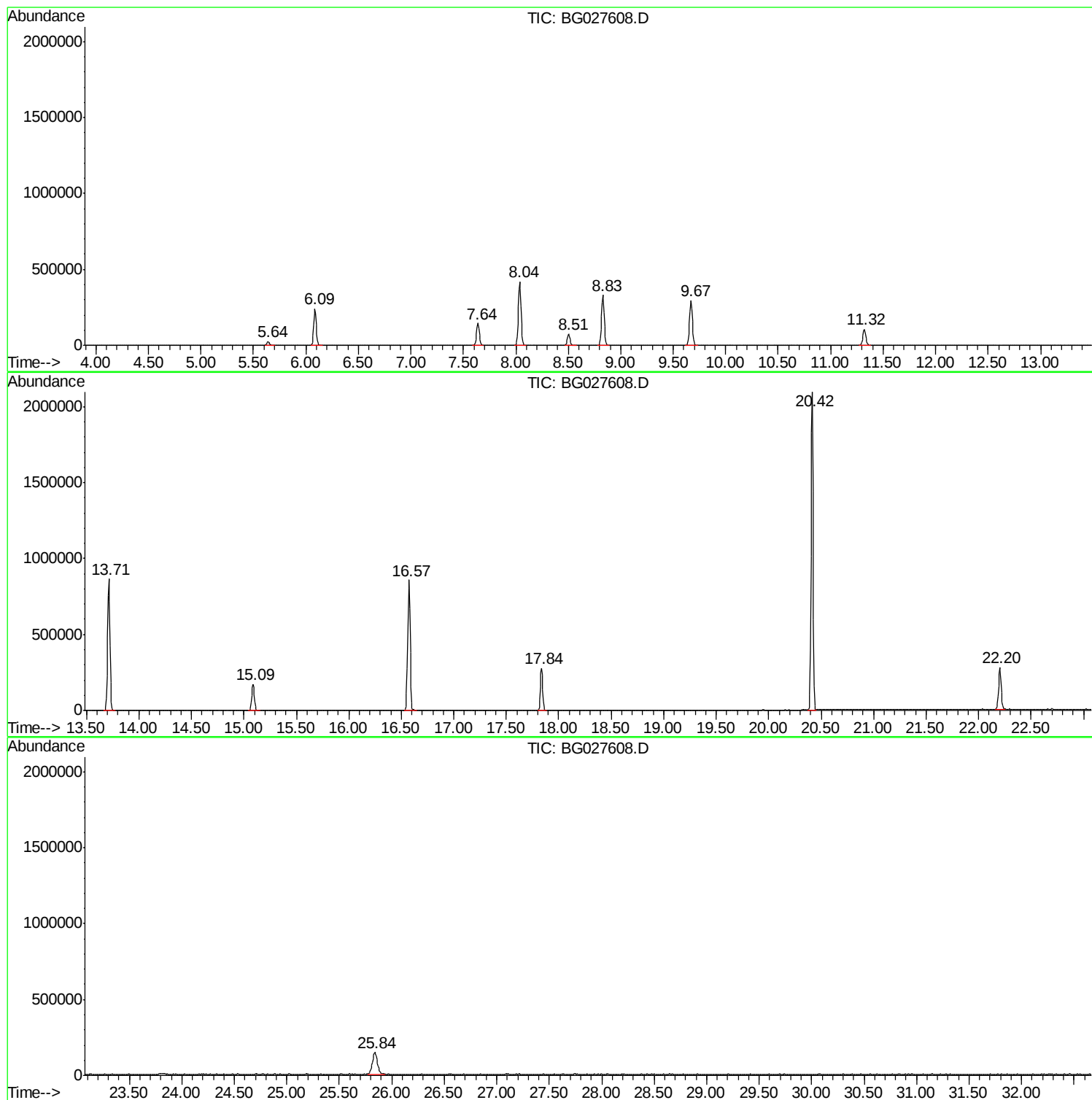
Sum of corrected areas: 10100969

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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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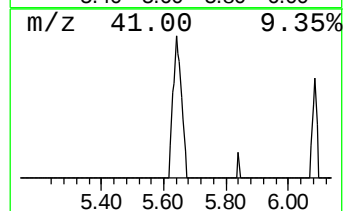
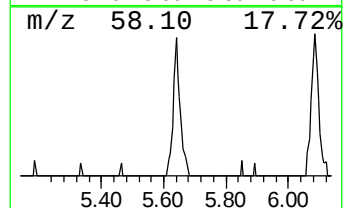
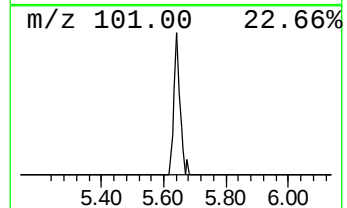
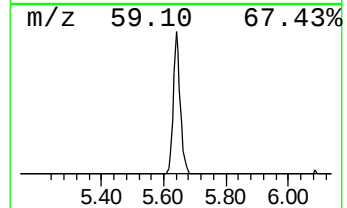
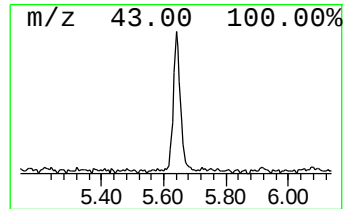
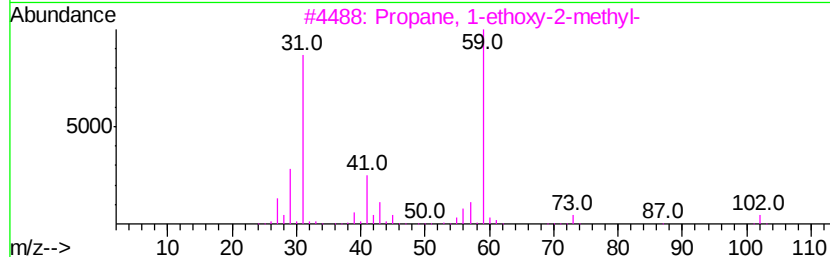
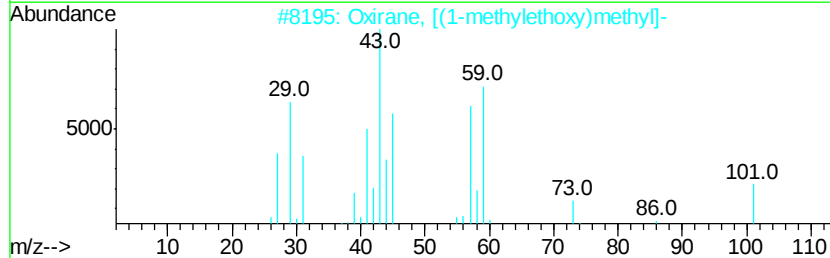
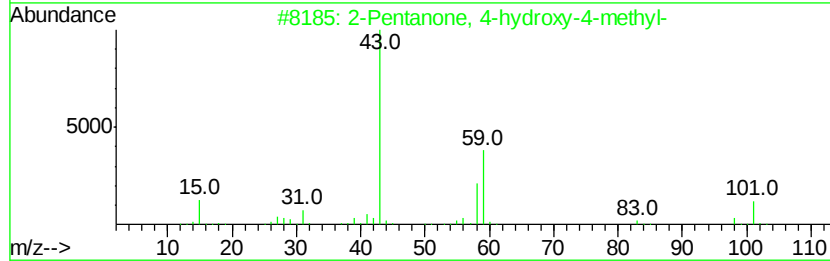
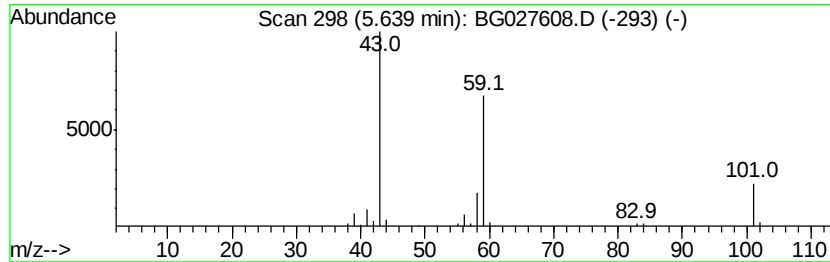
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	5.64 ng	38293	1,4-Dichlorobenzene-d4	8.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40	
2	Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	40	
3	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	35	
4	3-Octanol	130	C8H18O	000589-98-0	28	
5	3-Hexanol	102	C6H14O	000623-37-0	28	



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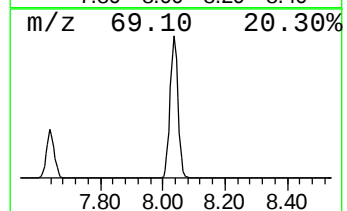
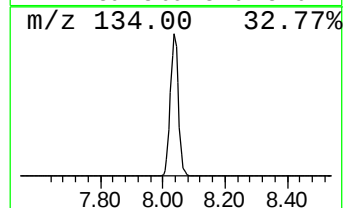
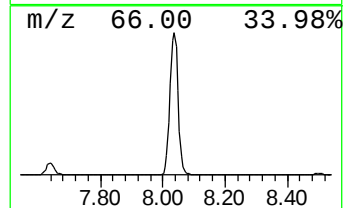
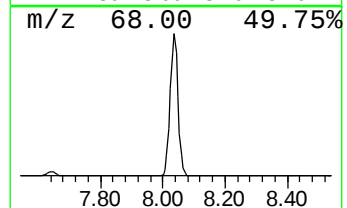
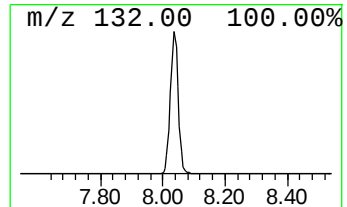
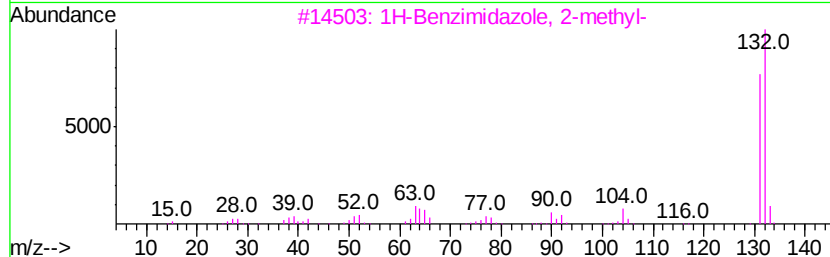
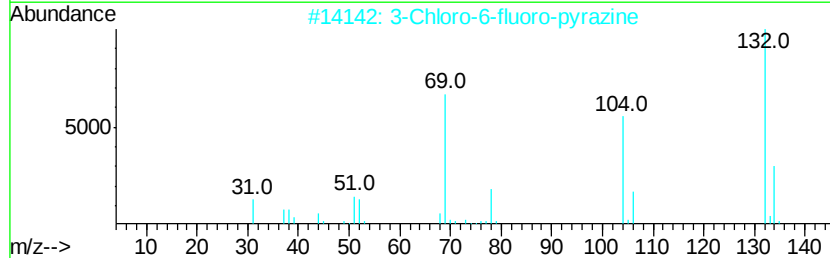
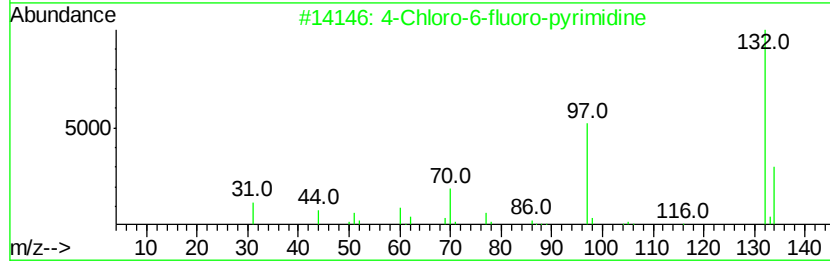
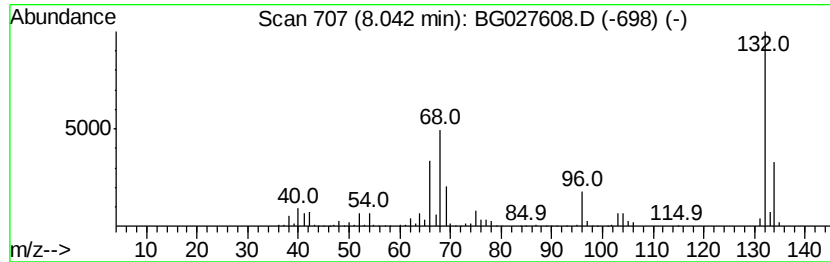
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Peak Number 2 unknown8.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.04	110.55 ng	750304	1,4-Dichlorobenzene-d4	8.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	25
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	16
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.64	5.6	ng	38293	1	8.51	135739	20.0
unknown8.04	8.04	110.6	ng	750304	1	8.51	135739	20.0