

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110417\
 Data File : BG030704.D
 Acq On : 3 Nov 2017 17:43
 Operator : SJ/JU
 Sample : PB103961BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103961BL

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-BG101617.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	5.228	325	330	341	rVB	116939	181541	4.83%	0.961%
2	5.709	405	412	425	rBV	702587	1136029	30.20%	6.015%
3	7.308	676	684	694	rBV	730988	1246573	33.14%	6.600%
4	7.684	740	748	758	rBV	717198	1236130	32.86%	6.545%
5	8.154	821	828	837	rVB	199538	337834	8.98%	1.789%
6	8.477	876	883	895	rVB	542724	952051	25.31%	5.041%
7	9.317	1019	1026	1036	rBV	415783	753607	20.03%	3.990%
8	10.968	1299	1307	1322	rBV	279424	517534	13.76%	2.740%
9	13.395	1713	1720	1736	rBV	1426701	2230373	59.29%	11.809%
10	14.769	1947	1954	1968	rVB2	519483	859321	22.84%	4.550%
11	16.250	2200	2206	2222	rBV	1118521	1775948	47.21%	9.403%
12	17.507	2412	2420	2428	rBV2	742457	1146918	30.49%	6.073%
13	20.099	2855	2861	2876	rBV	2858991	3761870	100.00%	19.918%
14	21.779	3140	3147	3159	rBV	837084	1371896	36.47%	7.264%
15	25.081	3698	3709	3722	rBV2	482776	1379000	36.66%	7.301%

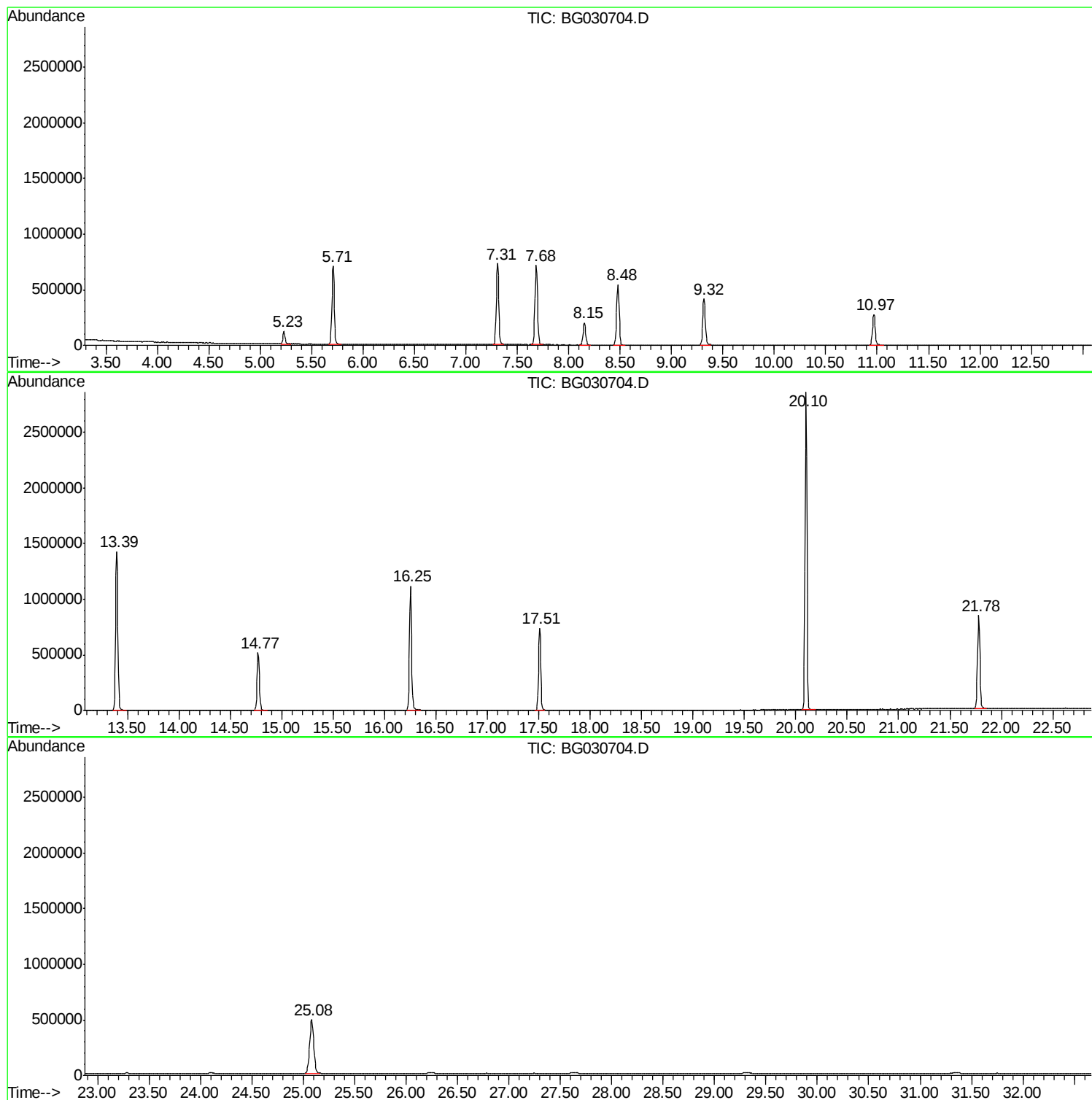
Sum of corrected areas: 18886625

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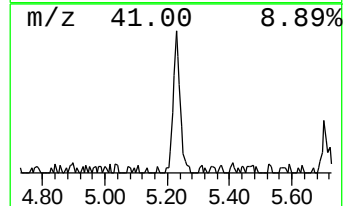
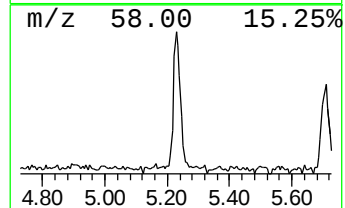
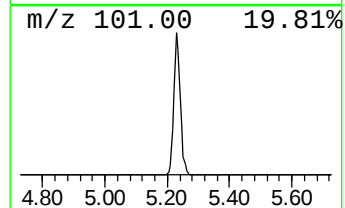
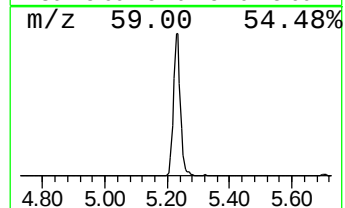
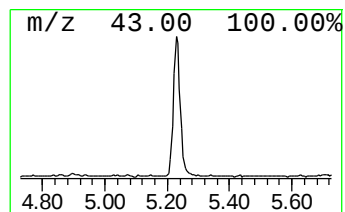
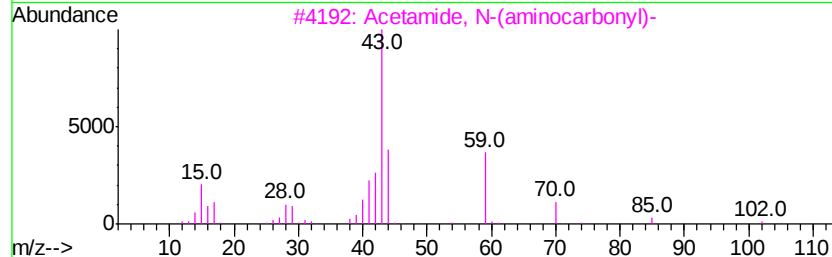
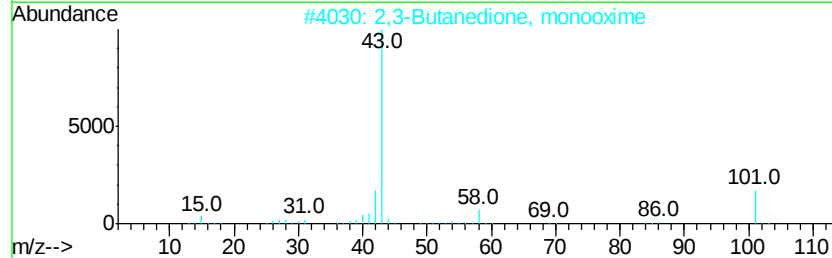
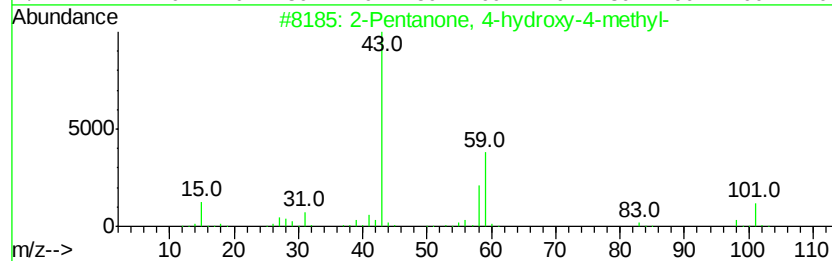
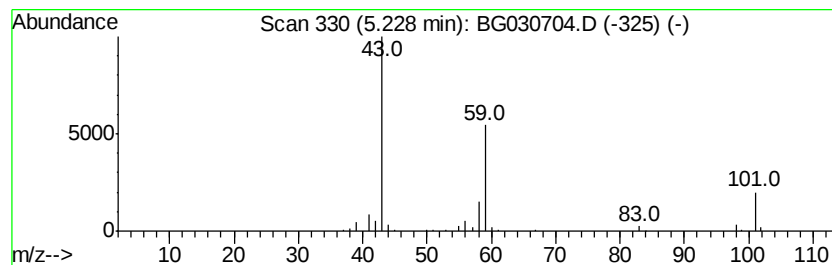
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	10.75 ng	181541	1,4-Dichlorobenzene-d4	8.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3			Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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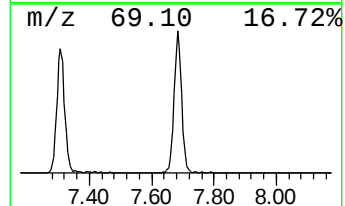
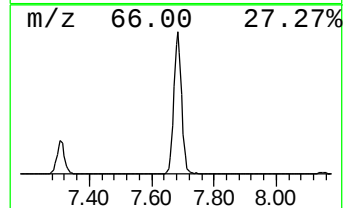
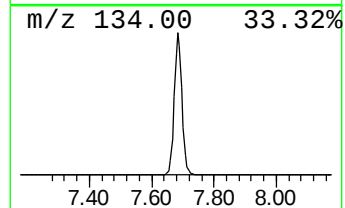
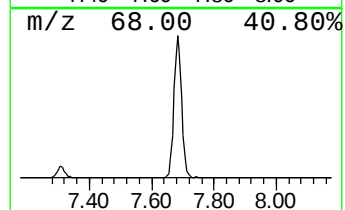
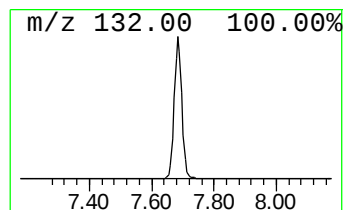
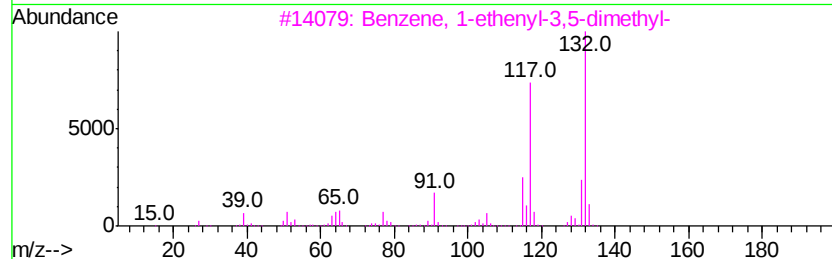
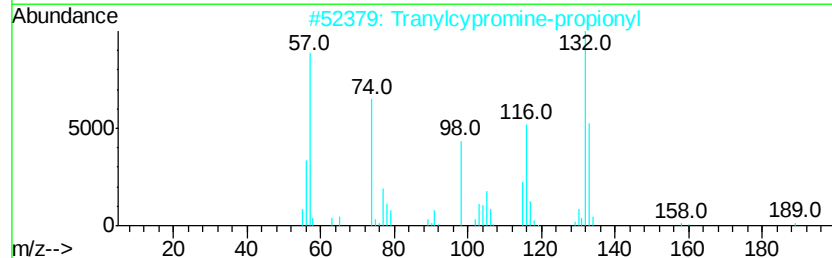
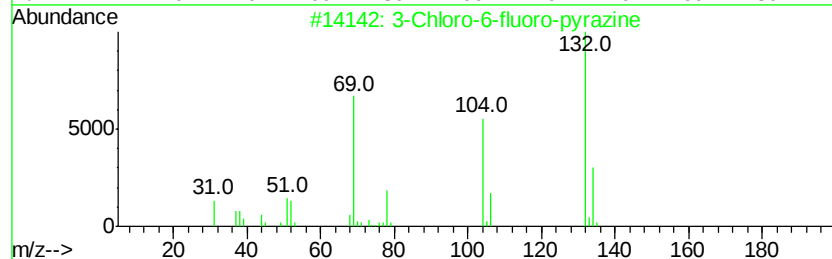
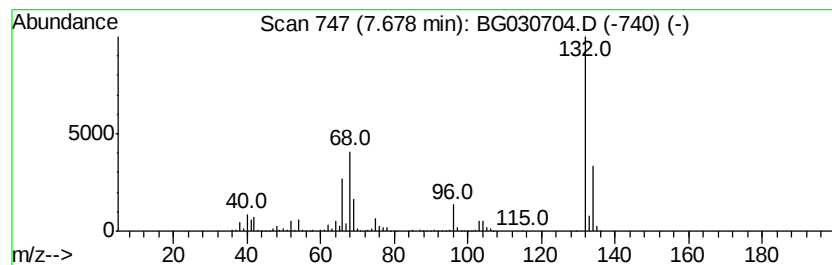
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	73.18 ng	1236130	1,4-Dichlorobenzene-d4	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.23	10.8	ng	181541	1	8.15	337834	20.0
unknown7.68	7.68	73.2	ng	1236130	1	8.15	337834	20.0