

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071818\
 Data File : BG035747.D
 Acq On : 19 Jul 2018 10:47
 Operator : JU/SJ
 Sample : PB111228BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111228BL

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.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	5.147	311	316	328	rBV3	32748	55481	1.30%	0.324%
2	5.622	390	397	412	rBV	615770	1054179	24.64%	6.150%
3	7.203	657	666	679	rBV	636077	1125346	26.31%	6.566%
4	7.573	722	729	739	rBV	628652	1089880	25.48%	6.359%
5	8.037	800	808	818	rBV	167792	286632	6.70%	1.672%
6	8.360	855	863	871	rBV	706624	1249182	29.20%	7.288%
7	9.200	998	1006	1020	rBV	531974	1021837	23.89%	5.962%
8	10.839	1277	1285	1295	rBV	168065	340376	7.96%	1.986%
9	13.283	1694	1701	1716	rBV	1633895	2533765	59.23%	14.783%
10	14.663	1928	1936	1948	rBV2	303079	525412	12.28%	3.065%
11	16.149	2182	2189	2209	rBV	546639	1026887	24.00%	5.991%
12	17.401	2394	2402	2419	rBV	444538	765275	17.89%	4.465%
13	20.009	2840	2846	2856	rBV	3155478	4278047	100.00%	24.960%
14	21.665	3119	3128	3143	rBV2	518218	948359	22.17%	5.533%
15	24.879	3663	3675	3696	rBV2	266628	839234	19.62%	4.896%

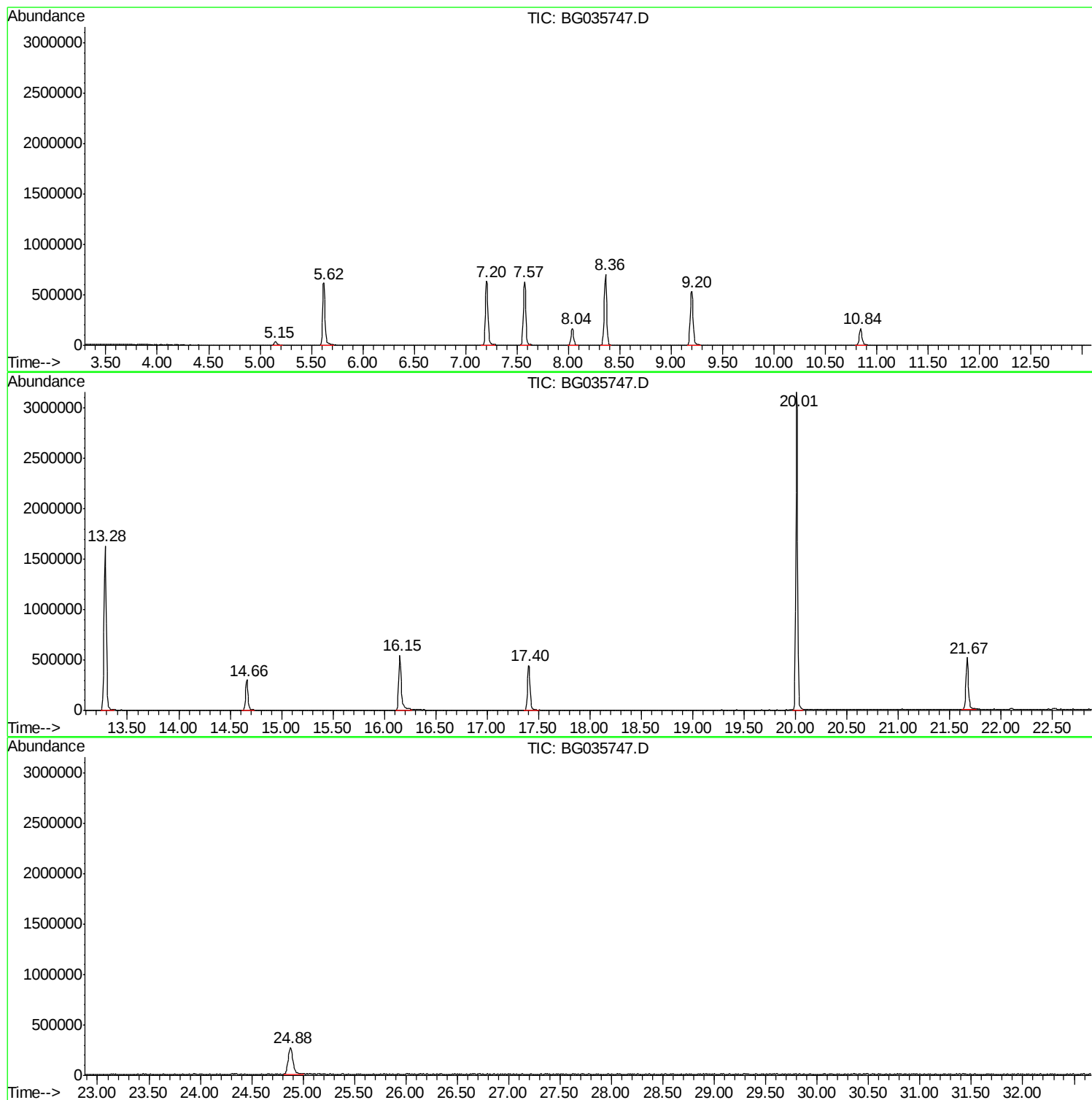
Sum of corrected areas: 17139892

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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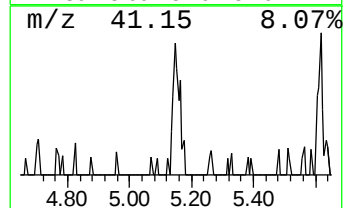
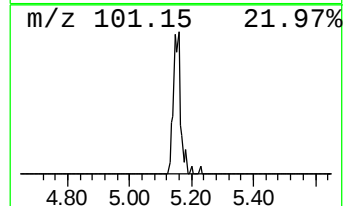
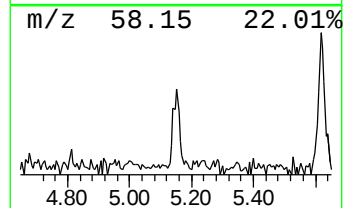
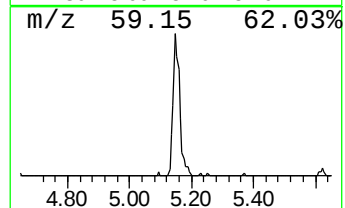
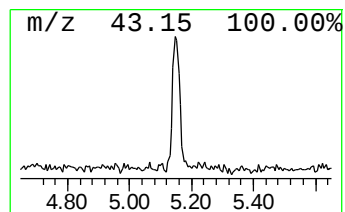
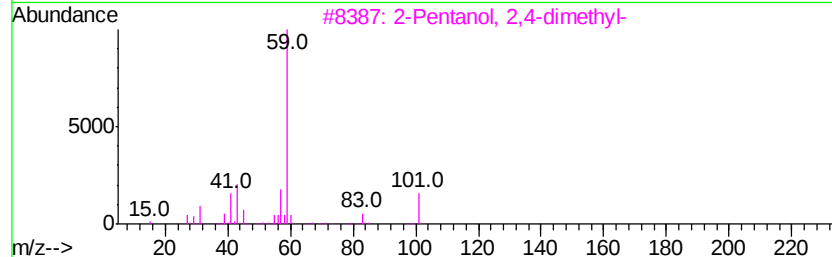
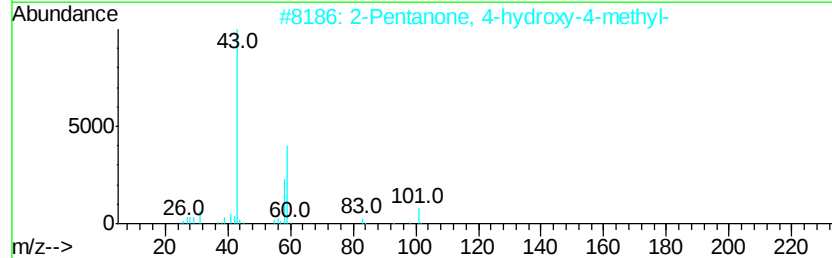
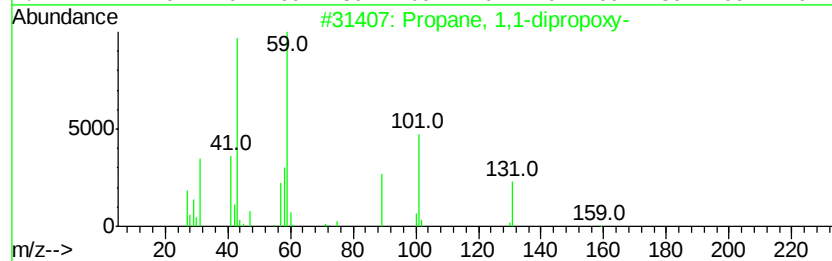
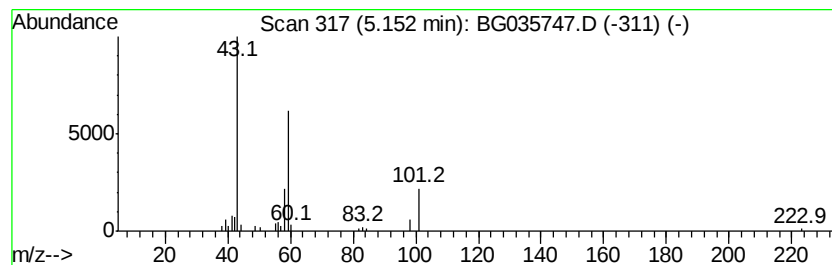
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.87 ng	55481	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	50
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	16
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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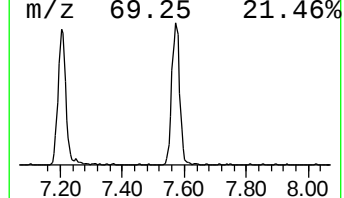
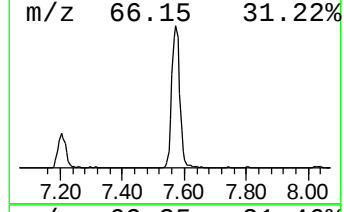
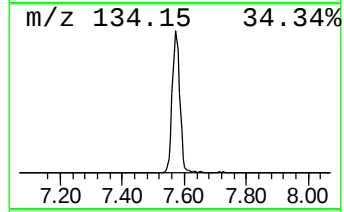
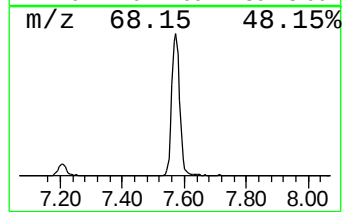
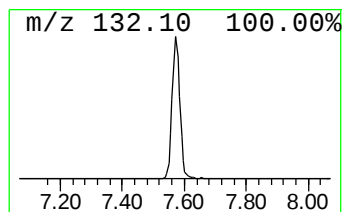
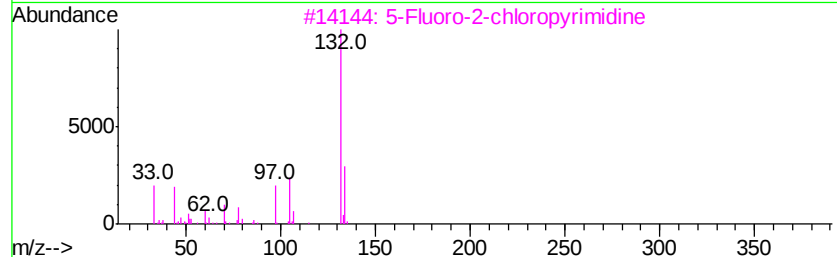
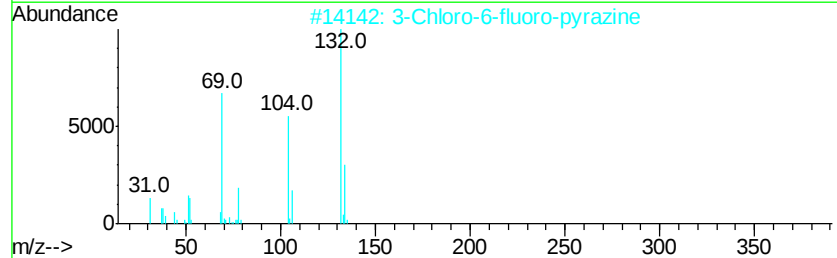
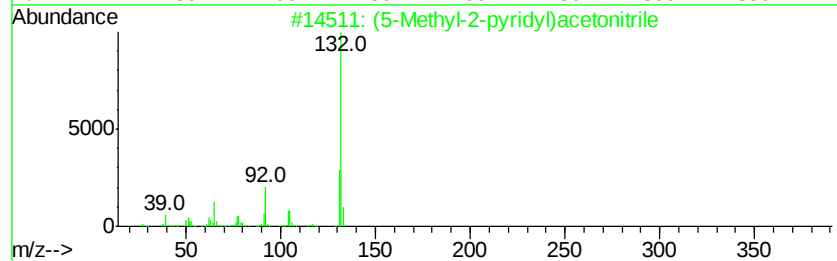
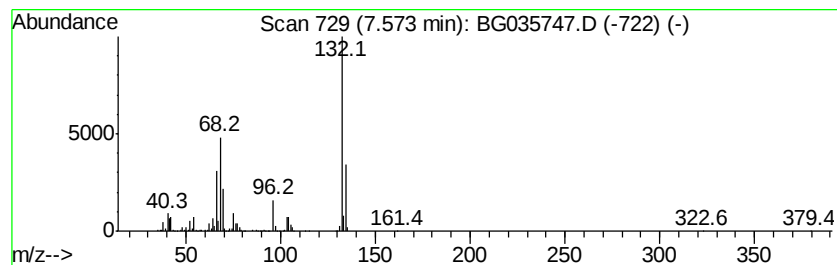
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Peak Number 2 unknown7.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	76.05 ng	1089880	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	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...	5.15	3.9	ng	55481	1	8.04	286632	20.0
unknown7.57	7.57	76.0	ng	1089880	1	8.04	286632	20.0