

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071218\
 Data File : BG035605.D
 Acq On : 12 Jul 2018 21:49
 Operator : JU/SJ
 Sample : PB110987BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB110987BL

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	5.167	314	320	328	rBV	113683	175804	3.59%	0.808%
2	5.637	391	400	410	rBV	982200	1564784	31.98%	7.190%
3	7.223	662	670	682	rBV	1012889	1753875	35.85%	8.058%
4	7.593	725	733	744	rBV	950398	1616537	33.04%	7.427%
5	8.057	804	812	819	rBV2	191454	327640	6.70%	1.505%
6	8.380	857	867	875	rBV2	684665	1173609	23.99%	5.392%
7	9.220	1001	1010	1020	rBV	575803	1021455	20.88%	4.693%
8	10.859	1281	1289	1296	rBV	248647	437376	8.94%	2.010%
9	13.303	1696	1705	1720	rBV	1617200	2558648	52.30%	11.756%
10	14.678	1932	1939	1947	rBV2	427834	692630	14.16%	3.182%
11	16.164	2185	2192	2211	rBV	1420152	2200371	44.97%	10.110%
12	17.421	2397	2406	2415	rBV2	651735	1019731	20.84%	4.685%
13	20.023	2842	2849	2856	rBV	3661976	4892527	100.00%	22.479%
14	21.686	3124	3132	3143	rBV	735080	1242762	25.40%	5.710%
15	24.911	3669	3681	3692	rBV	374722	1086719	22.21%	4.993%

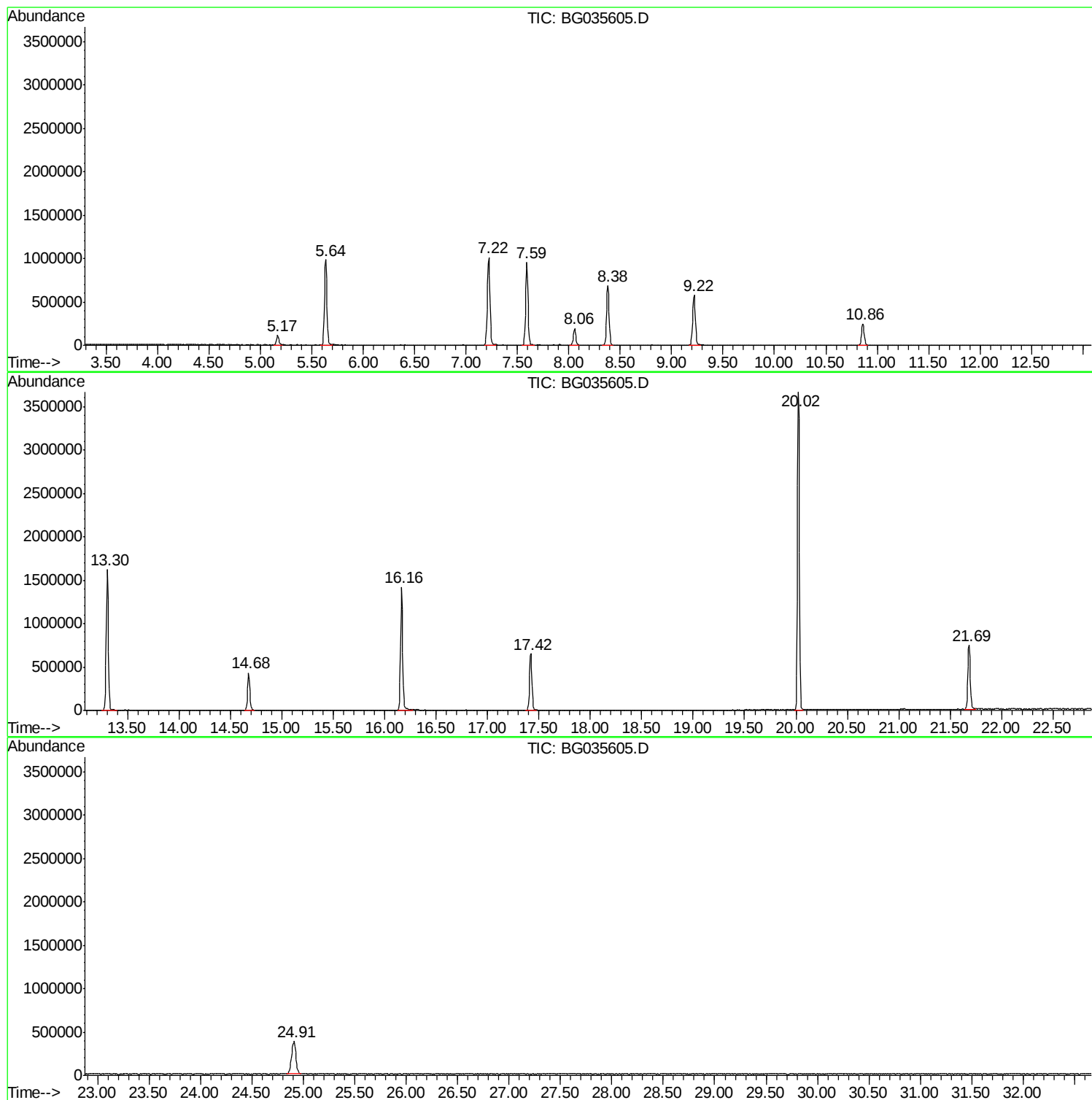
Sum of corrected areas: 21764468

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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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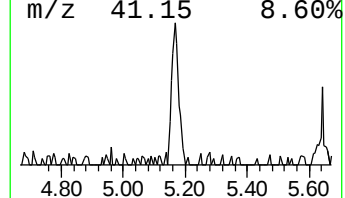
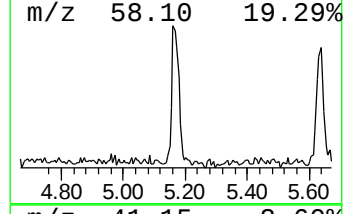
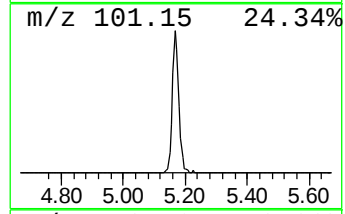
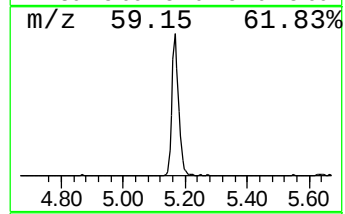
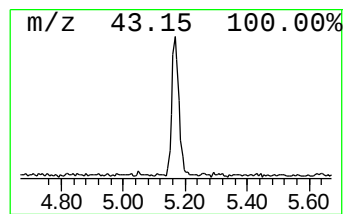
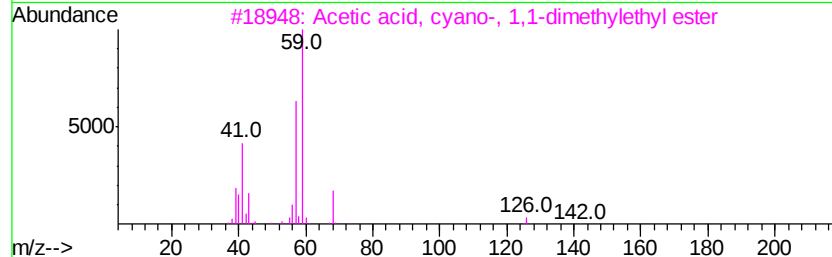
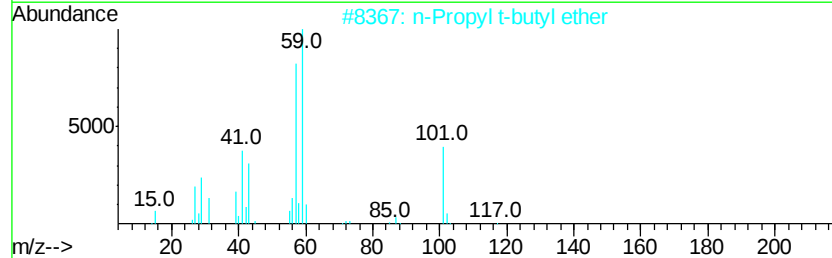
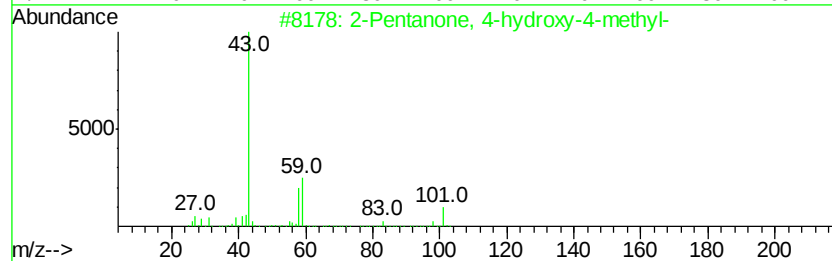
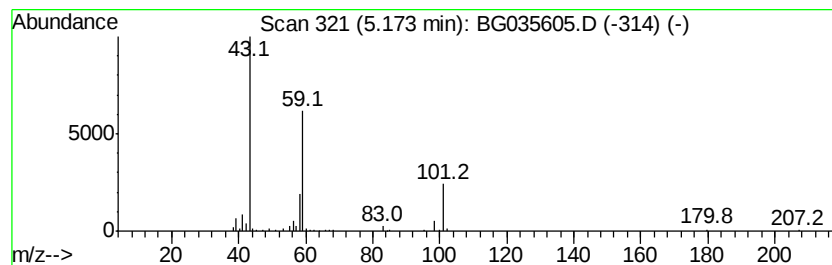
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	10.73 ng	175804	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
4			Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	28
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	16



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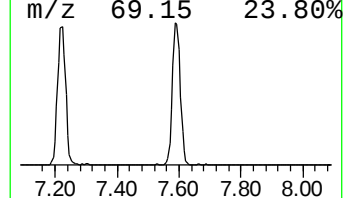
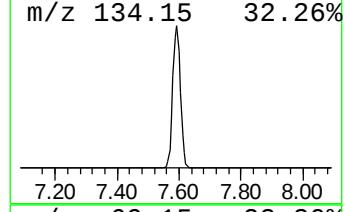
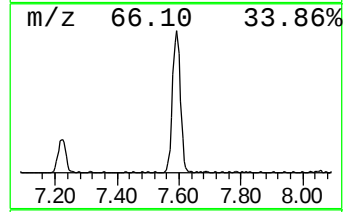
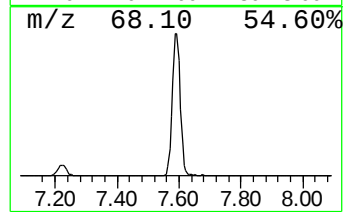
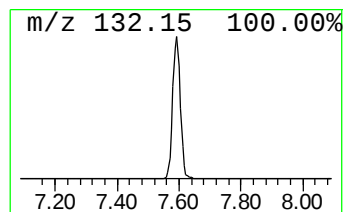
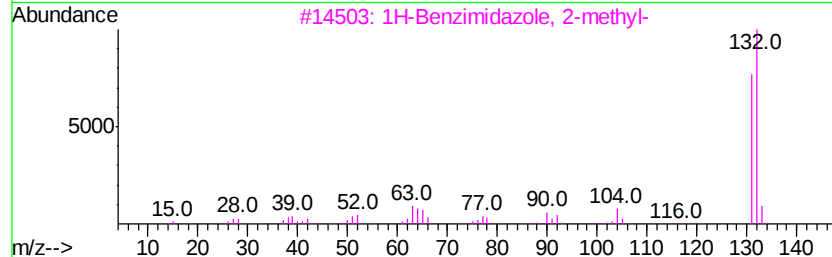
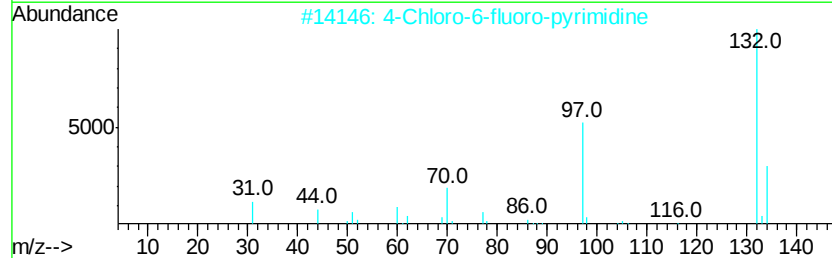
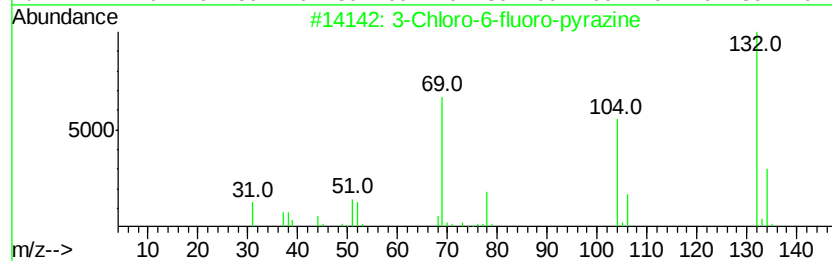
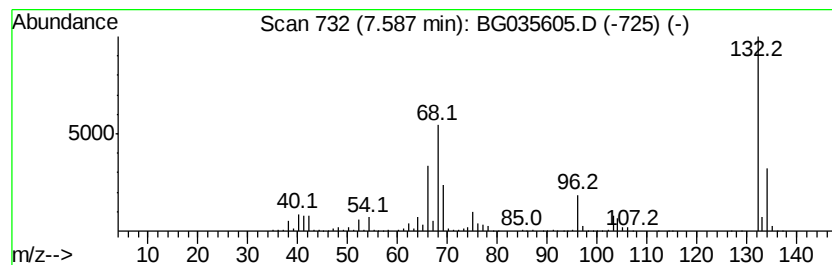
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	98.68 ng	1616540	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	10.7	ng	175804	1	8.06	327640	20.0
unknown7.59	7.59	98.7	ng	1616540	1	8.06	327640	20.0