

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG091217\
 Data File : BG028676.D
 Acq On : 12 Sep 2017 19:11
 Operator : SJ/JU
 Sample : PB102188BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB102188BL

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-BG091217.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	5.435	308	314	323	rVB2	46769	79943	2.46%	0.570%
2	5.905	387	394	411	rBV	605718	1023365	31.48%	7.291%
3	7.485	654	663	675	rBV	574334	1046633	32.19%	7.456%
4	7.867	718	728	737	rBV	560121	1012564	31.15%	7.214%
5	8.331	800	807	816	rVB	112176	200256	6.16%	1.427%
6	8.655	854	862	872	rBV	424731	764315	23.51%	5.445%
7	9.501	998	1006	1018	rBV	322232	621337	19.11%	4.426%
8	11.152	1279	1287	1296	rBV	141899	260374	8.01%	1.855%
9	13.561	1690	1697	1708	rBV	1036446	1604672	49.36%	11.432%
10	14.941	1925	1932	1942	rVB2	239850	411870	12.67%	2.934%
11	16.434	2179	2186	2203	rBV	979320	1577202	48.51%	11.236%
12	17.697	2393	2401	2410	rBV2	381330	640131	19.69%	4.560%
13	20.276	2834	2840	2851	rBV	2366432	3251022	100.00%	23.161%
14	22.027	3131	3138	3148	rBV	432043	778663	23.95%	5.547%
15	25.529	3724	3734	3748	rVB2	240297	764557	23.52%	5.447%

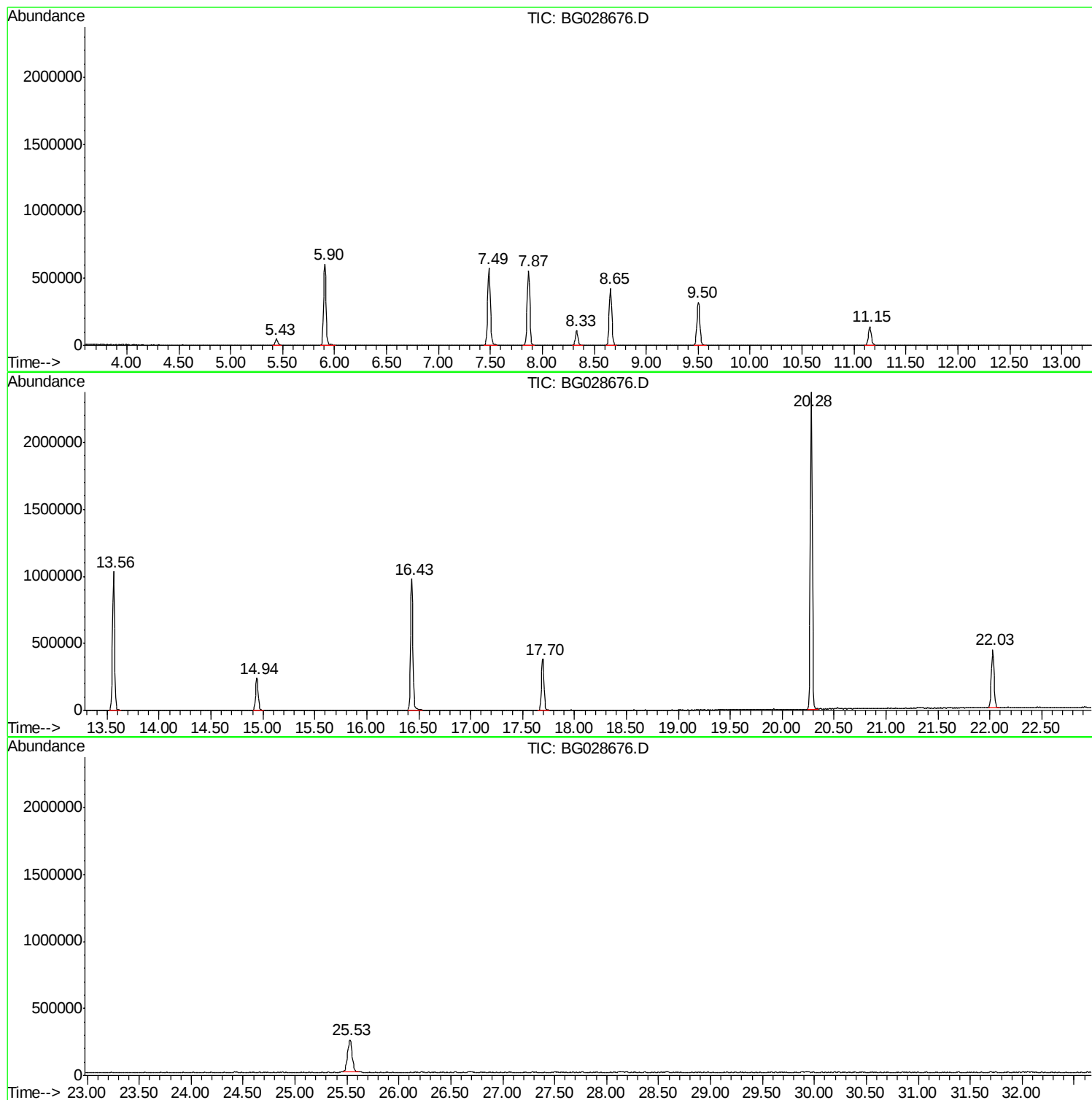
Sum of corrected areas: 14036904

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



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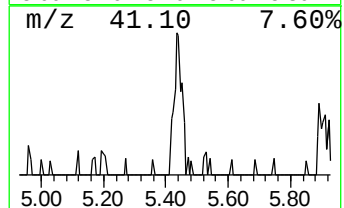
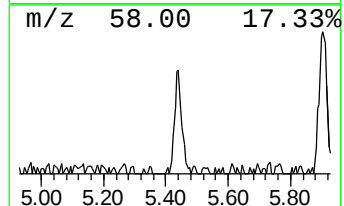
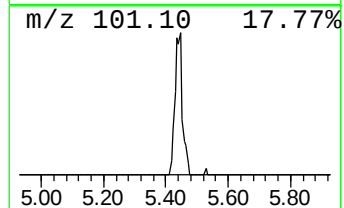
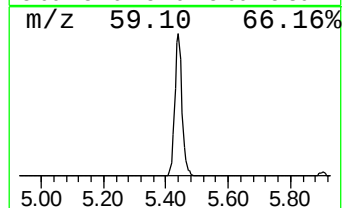
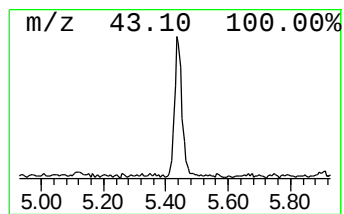
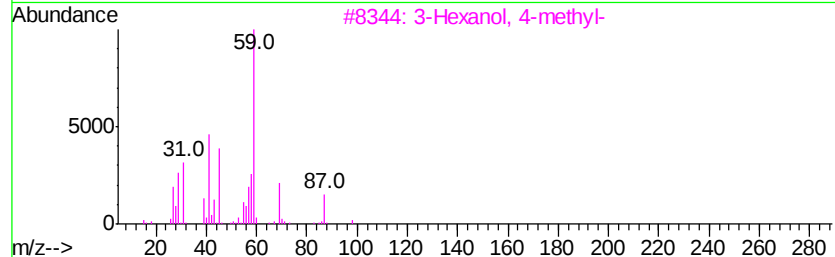
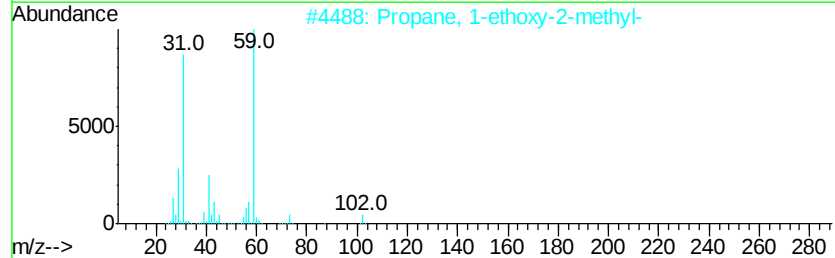
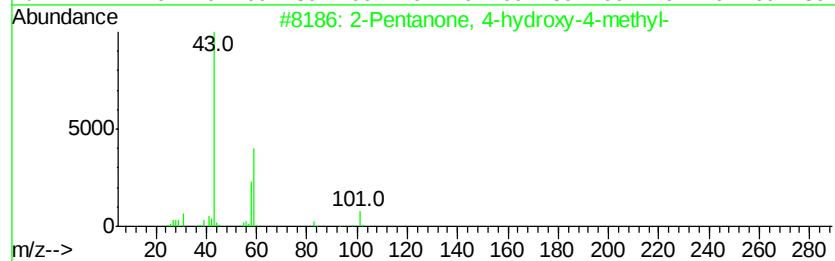
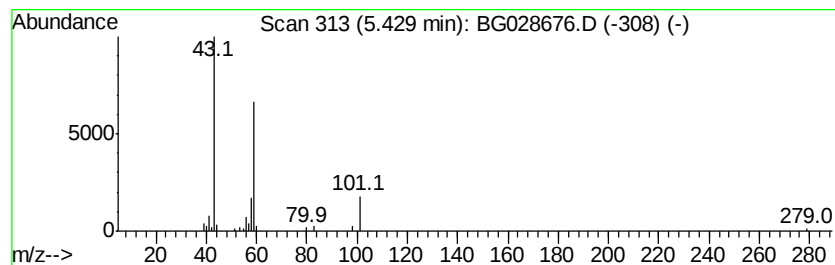
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.43	7.98 ng	79943	1,4-Dichlorobenzene-d4	8.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	43
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	38
4			Tetraolyme	222	C10H22O5	000143-24-8	28
5			2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	25



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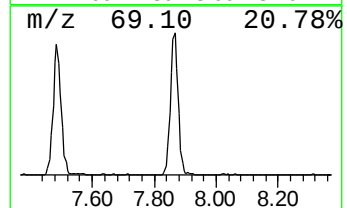
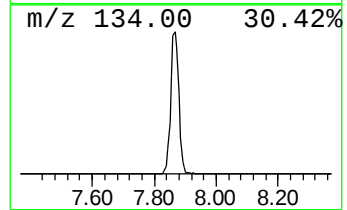
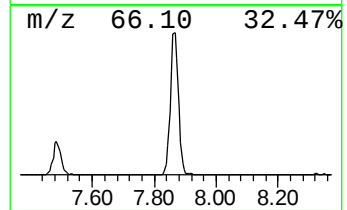
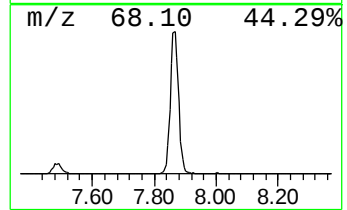
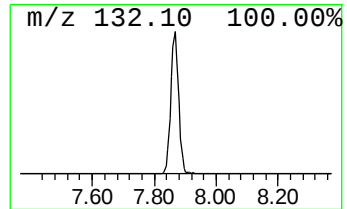
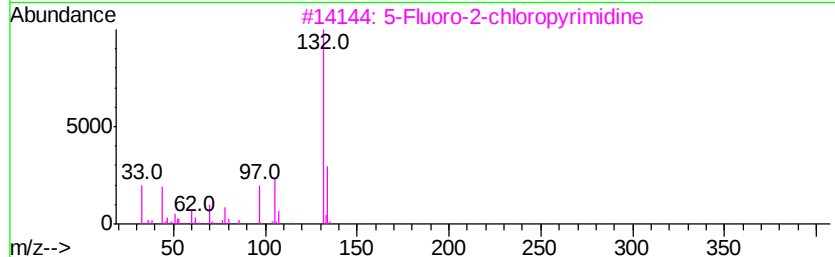
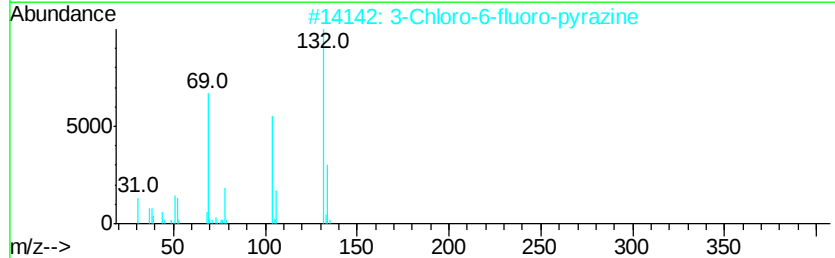
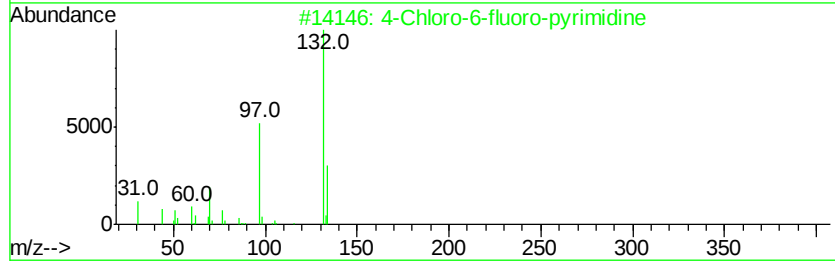
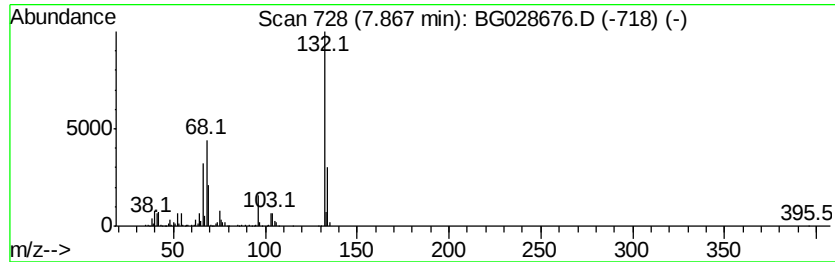
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Peak Number 2 unknown7.87 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	101.13 ng	1012560	1,4-Dichlorobenzene-d4	8.33

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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.43	8.0	ng	79943	1	8.33	200256	20.0
unknown7.87	7.87	101.1	ng	1012560	1	8.33	200256	20.0