

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091218\
 Data File : BG036683.D
 Acq On : 12 Sep 2018 21:40
 Operator : JU/SJ
 Sample : PB112784BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112784BL

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-BG082318.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	5.157	312	318	325	rBV	47025	69645	1.49%	0.314%
2	5.621	390	397	410	rBV	1120410	1781754	38.11%	8.028%
3	7.207	659	667	677	rBV	1143197	1966954	42.07%	8.862%
4	7.583	723	731	745	rBV	1058378	1822105	38.97%	8.210%
5	8.053	802	811	818	rBV	209578	353485	7.56%	1.593%
6	8.376	858	866	874	rBV	773821	1350655	28.89%	6.086%
7	9.211	1000	1008	1020	rBV	658446	1189248	25.43%	5.358%
8	10.856	1281	1288	1299	rVB	299736	529570	11.33%	2.386%
9	13.300	1697	1704	1715	rBV	1908527	2923551	62.53%	13.172%
10	14.675	1931	1938	1946	rBV2	499595	782975	16.75%	3.528%
11	16.161	2184	2191	2210	rBV	1082012	1801388	38.53%	8.116%
12	17.419	2398	2405	2417	rBV2	676921	1021369	21.84%	4.602%
13	20.027	2842	2849	2864	rBV	3417891	4675799	100.00%	21.067%
14	21.684	3124	3131	3149	rBV	601062	1126334	24.09%	5.075%
15	24.892	3667	3677	3694	rBV3	204589	799746	17.10%	3.603%

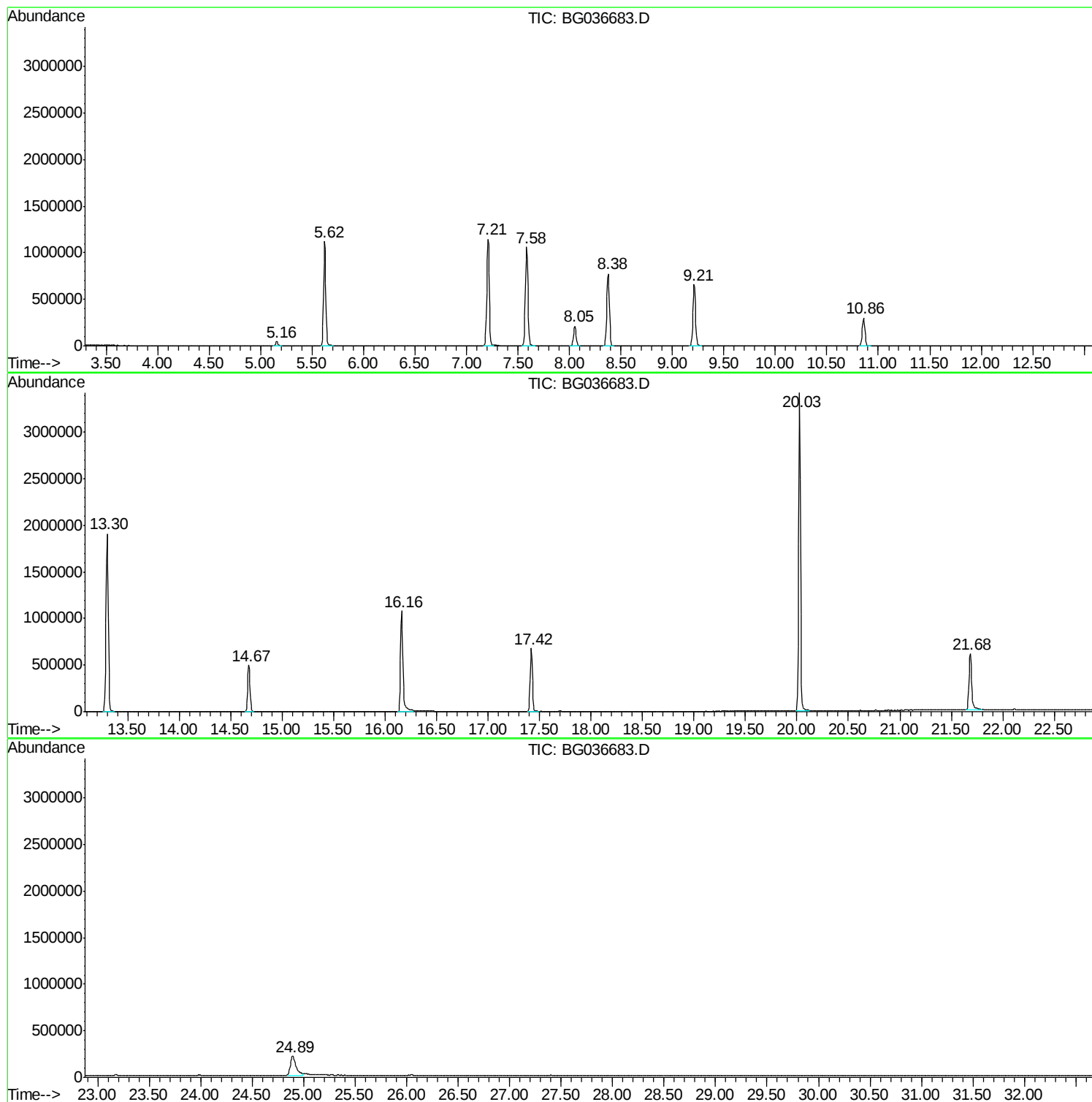
Sum of corrected areas: 22194578

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



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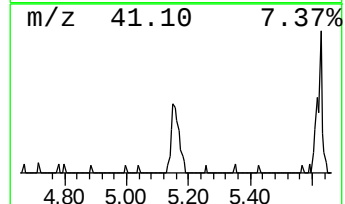
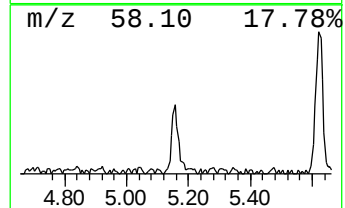
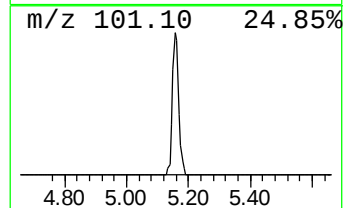
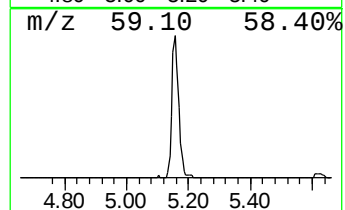
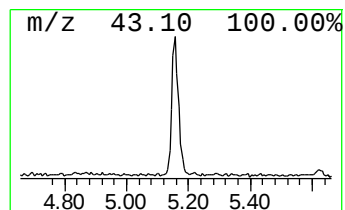
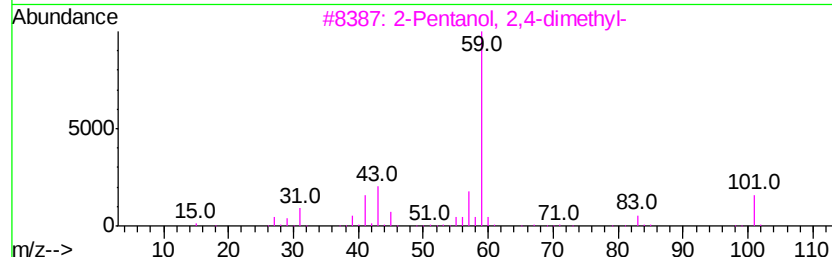
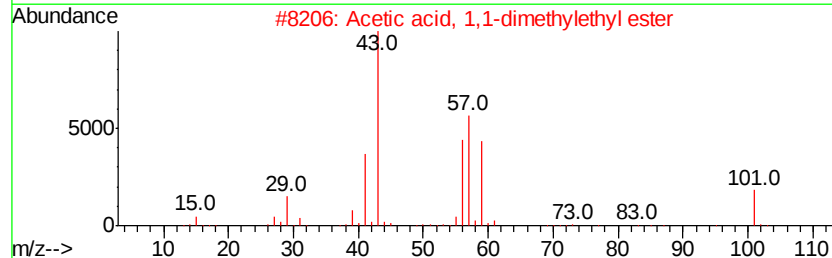
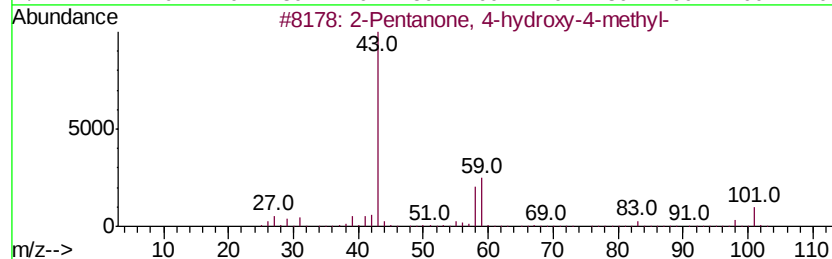
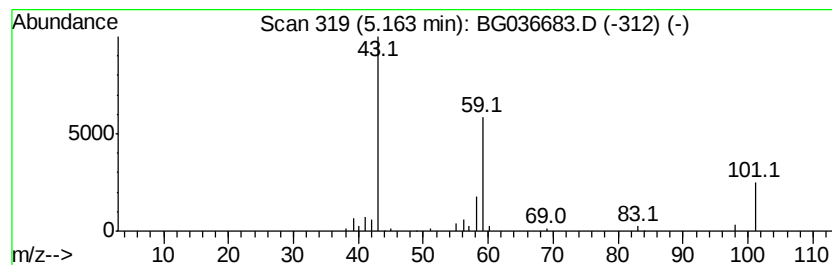
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	3.94 ng	69645	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	45
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	28
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	23
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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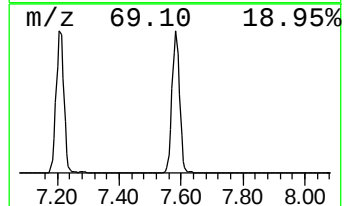
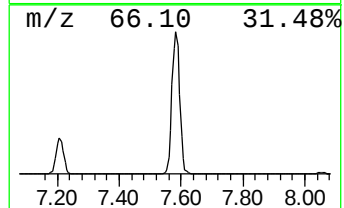
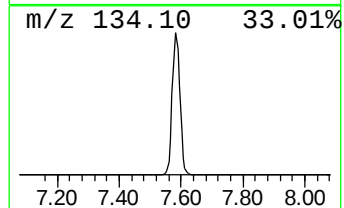
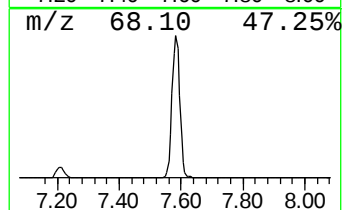
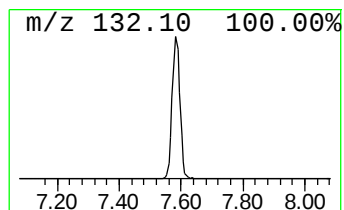
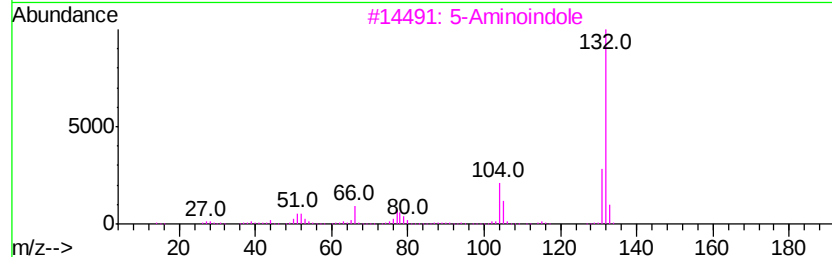
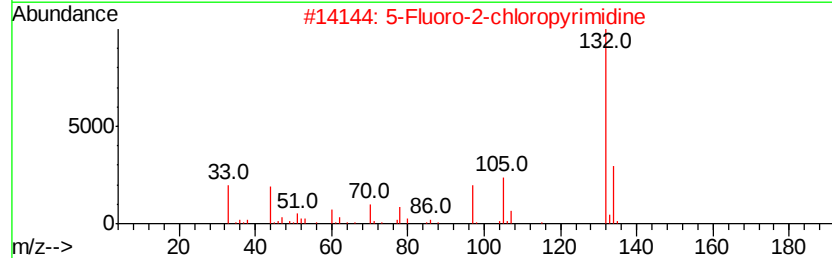
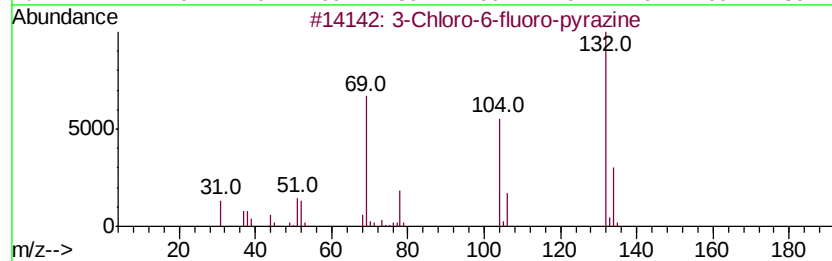
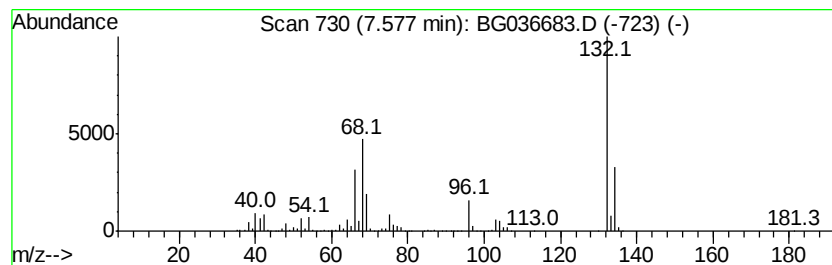
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	103.09 ng	1822110	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	3.9	ng	69645	1	8.05	353485	20.0
unknown7.58	7.58	103.1	ng	1822110	1	8.05	353485	20.0