

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091118\
 Data File : BG036672.D
 Acq On : 12 Sep 2018 1:19
 Operator : JU/SJ
 Sample : PB112760BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112760BL

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	328	rBV	126371	186183	3.52%	0.865%
2	5.621	390	397	410	rBV	818926	1296161	24.48%	6.024%
3	7.208	659	667	677	rBV	848264	1446059	27.31%	6.720%
4	7.584	723	731	743	rBV	780858	1340043	25.31%	6.228%
5	8.054	804	811	819	rVB	206693	342569	6.47%	1.592%
6	8.377	857	866	874	rBV	849815	1458044	27.54%	6.776%
7	9.211	1000	1008	1017	rBV	711129	1273382	24.05%	5.918%
8	10.856	1279	1288	1297	rBV	282846	501170	9.47%	2.329%
9	13.300	1697	1704	1717	rBV	2025171	3130097	59.12%	14.546%
10	14.675	1931	1938	1947	rBV2	471701	741198	14.00%	3.445%
11	16.162	2184	2191	2201	rBV	924497	1470034	27.76%	6.832%
12	17.419	2398	2405	2415	rBV2	656669	987708	18.65%	4.590%
13	20.028	2842	2849	2871	rBV	3897841	5294825	100.00%	24.606%
14	21.685	3123	3131	3147	rBV	640386	1157659	21.86%	5.380%
15	24.893	3663	3677	3692	rBV2	243382	892897	16.86%	4.150%

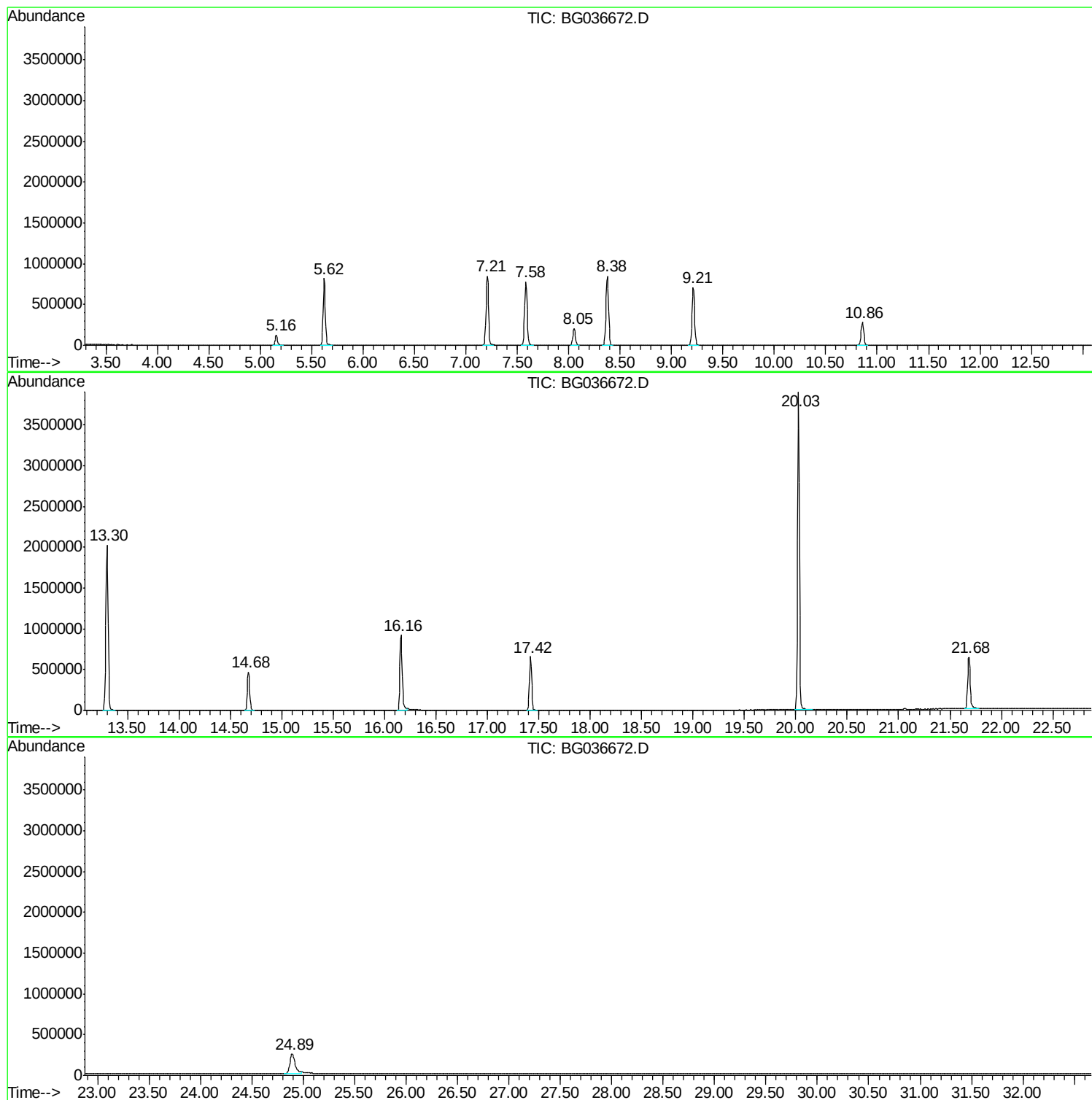
Sum of corrected areas: 21518029

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



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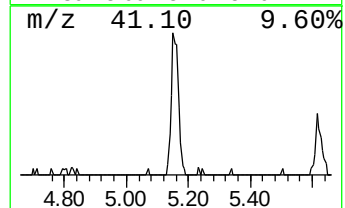
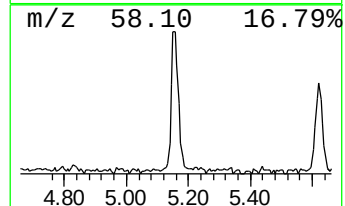
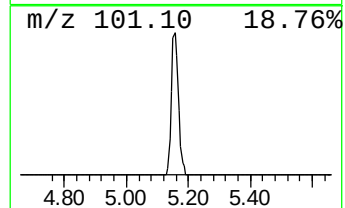
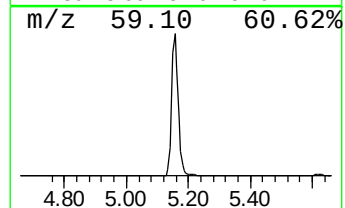
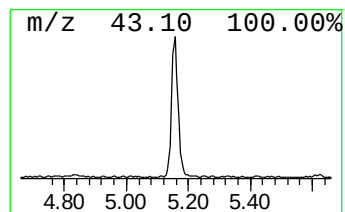
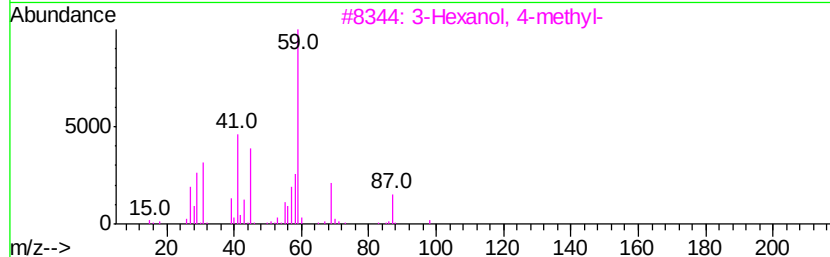
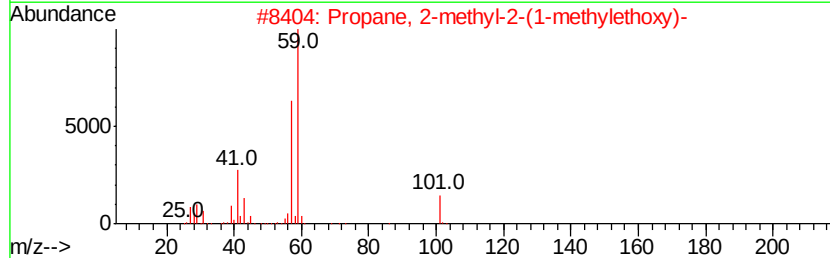
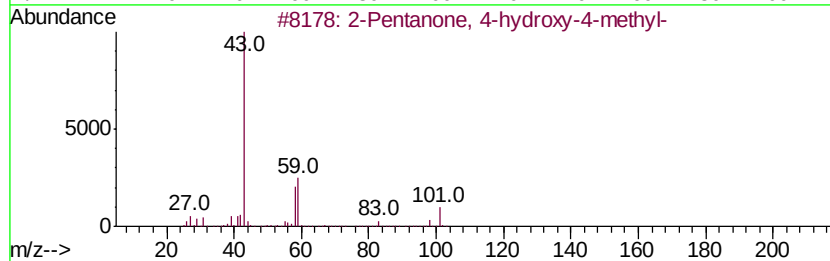
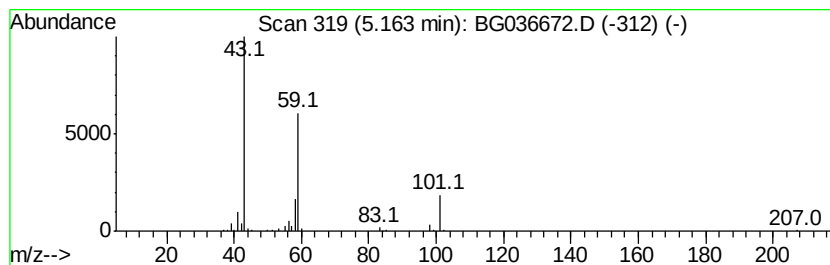
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	10.87 ng	186183	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			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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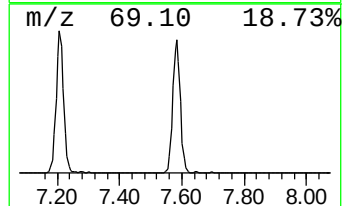
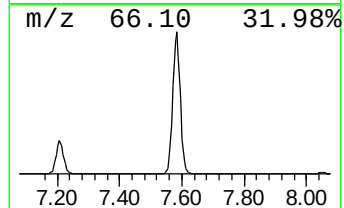
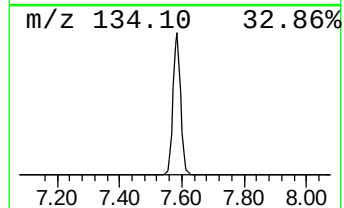
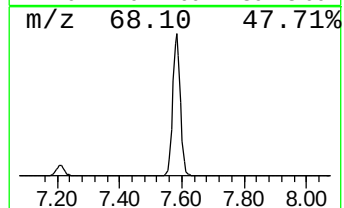
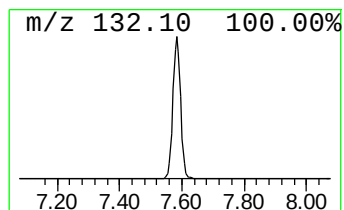
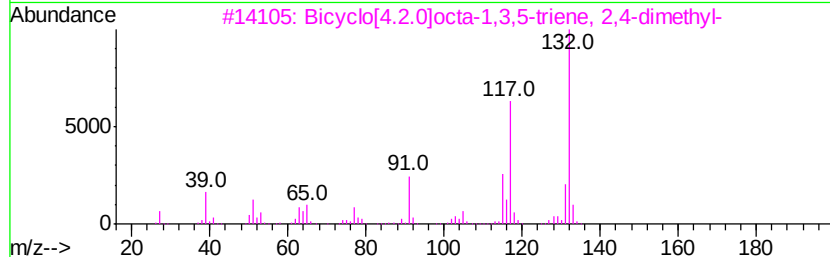
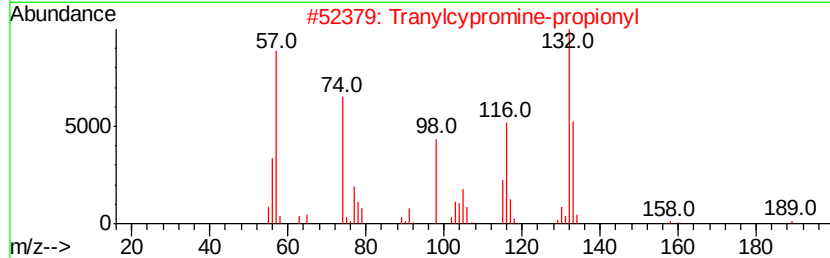
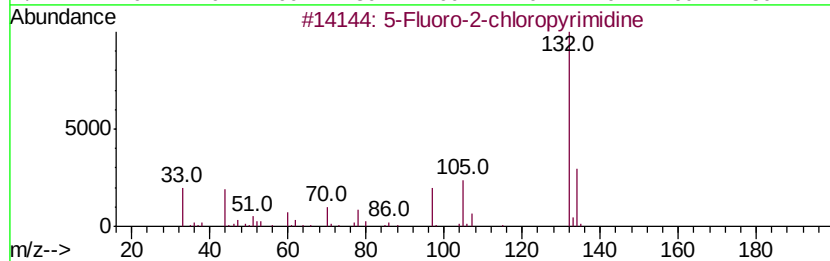
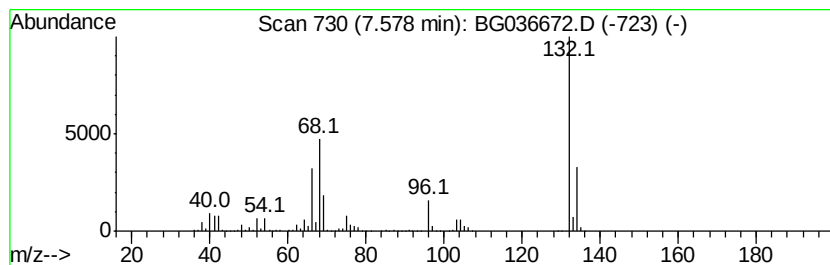
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	78.23 ng	1340040	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	10.9	ng	186183	1	8.05	342569	20.0
unknown7.58	7.58	78.2	ng	1340040	1	8.05	342569	20.0