

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036648.D
 Acq On : 11 Sep 2018 6:03
 Operator : JU/SJ
 Sample : J4777-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PL-05-090718-D

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	5.163	307	319	327	rBV	145263	229924	5.57%	1.166%
2	5.627	390	398	412	rBV	882416	1419862	34.40%	7.201%
3	7.213	659	668	682	rBV	812343	1452397	35.19%	7.366%
4	7.583	721	731	741	rBV	930959	1575090	38.16%	7.989%
5	8.053	803	811	819	rVB	202623	339250	8.22%	1.721%
6	8.377	858	866	875	rVB	709481	1202678	29.14%	6.100%
7	9.217	1000	1009	1019	rBV	596285	1075465	26.06%	5.455%
8	10.862	1281	1289	1302	rVB	265846	490472	11.88%	2.488%
9	13.300	1697	1704	1713	rBV	1629539	2544652	61.65%	12.906%
10	14.681	1930	1939	1949	rBV2	445608	734651	17.80%	3.726%
11	16.162	2182	2191	2211	rBV	1029365	1749654	42.39%	8.874%
12	17.419	2398	2405	2420	rBV2	641871	968032	23.45%	4.910%
13	20.028	2843	2849	2863	rBV	3091170	4127612	100.00%	20.934%
14	21.684	3125	3131	3147	rBV2	554783	1050660	25.45%	5.329%
15	24.898	3668	3678	3694	rBV	202002	756455	18.33%	3.837%

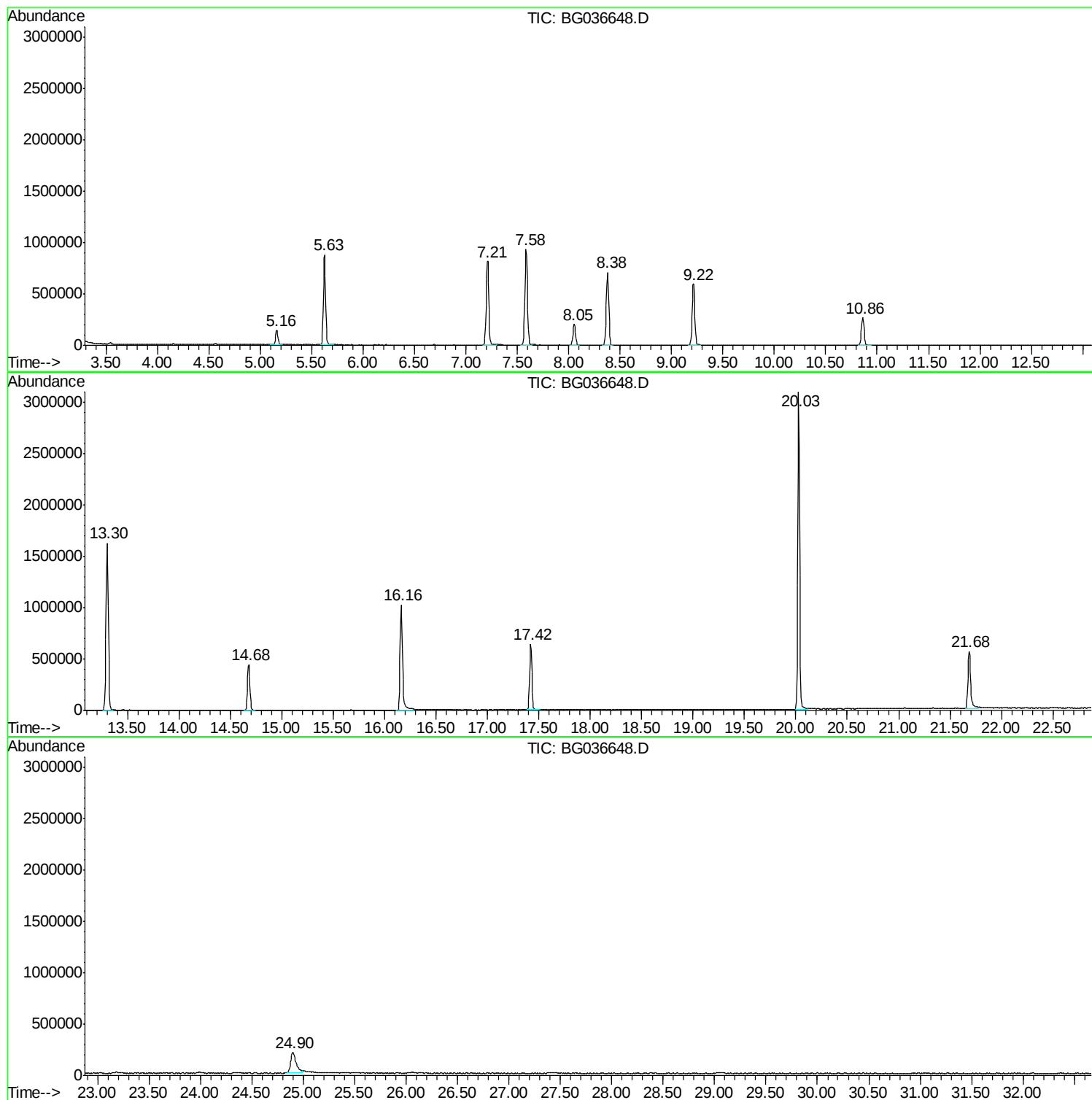
Sum of corrected areas: 19716854

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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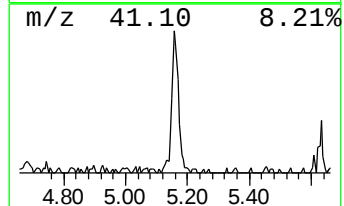
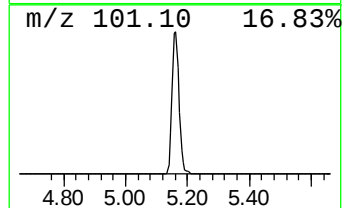
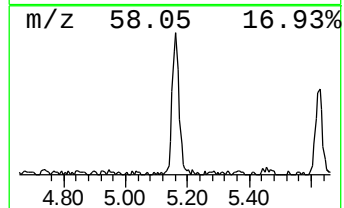
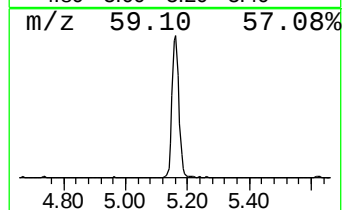
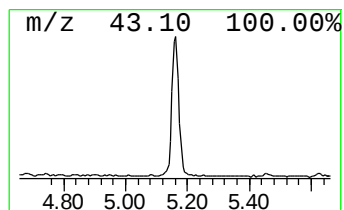
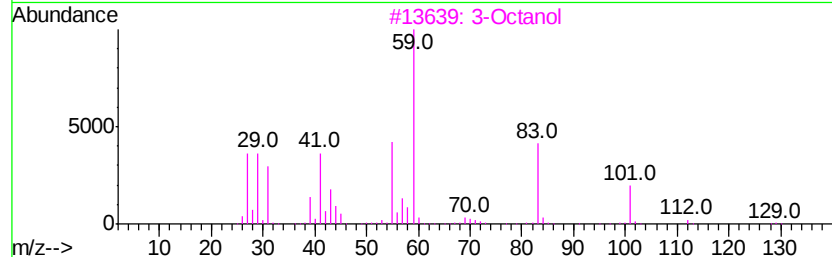
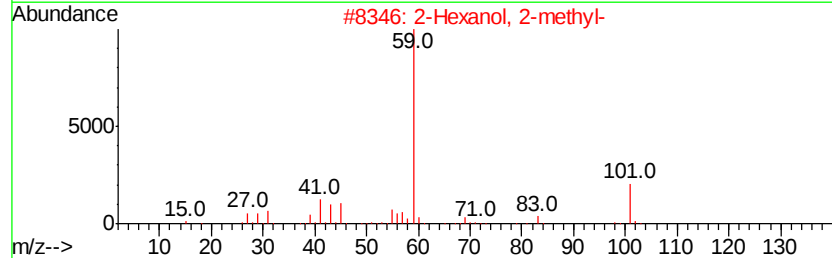
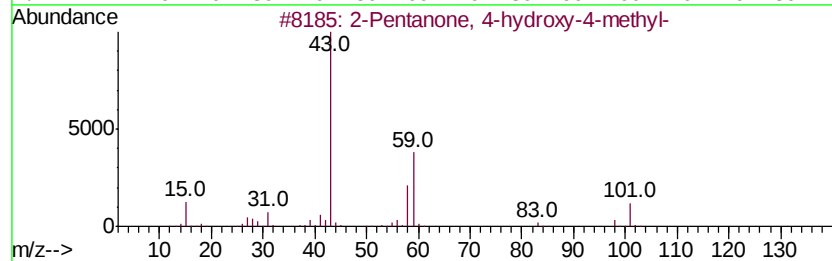
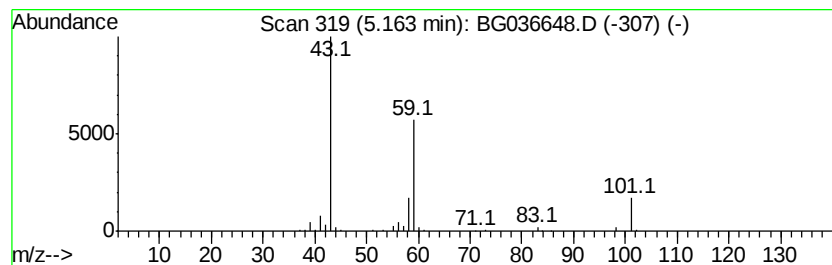
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	13.55 ng	229924	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	53
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3			3-Octanol	130	C8H18O	000589-98-0	33
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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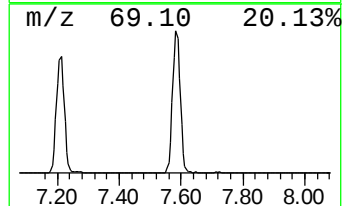
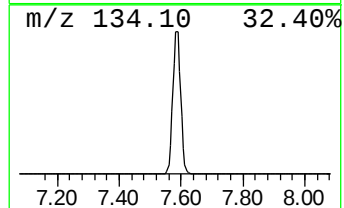
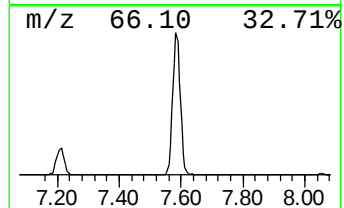
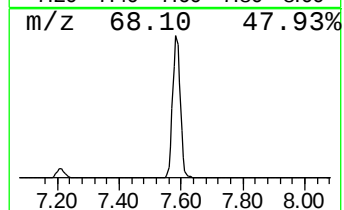
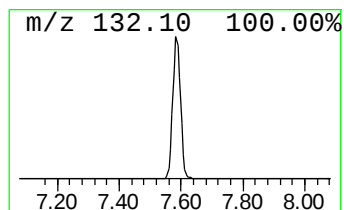
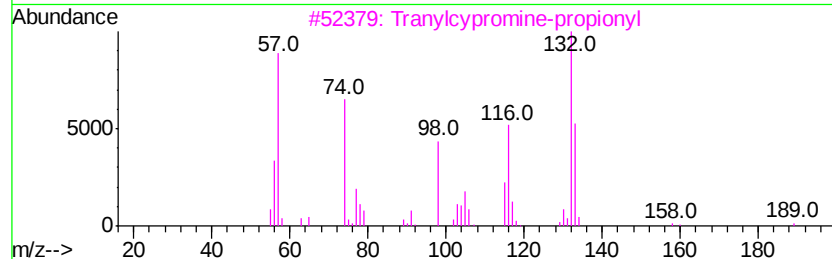
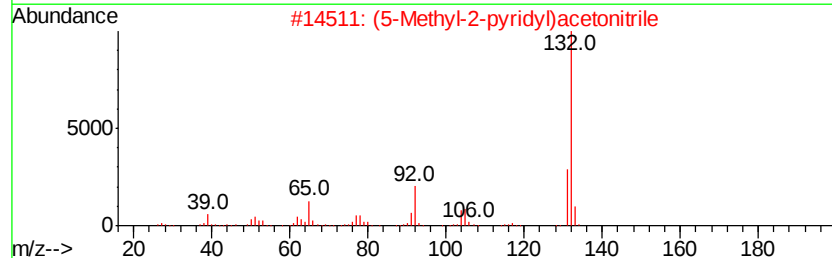
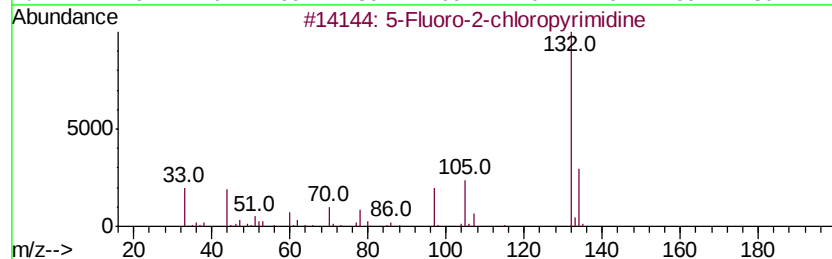
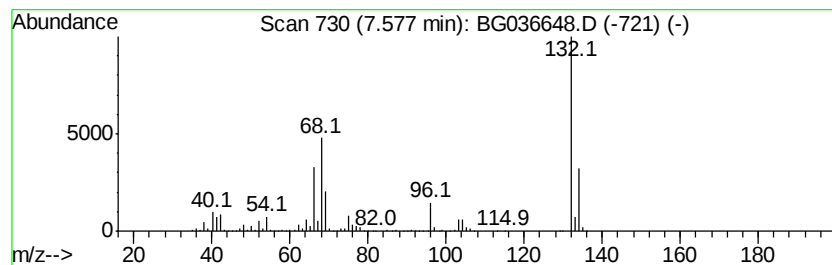
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	92.86 ng	1575090	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		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3		Tranvlcypramine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	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	13.6	ng	229924	1	8.05	339250	20.0
unknown7.58	7.58	92.9	ng	1575090	1	8.05	339250	20.0