

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092218\
 Data File : BG036912.D
 Acq On : 23 Sep 2018 3:57
 Operator : JU/SJ
 Sample : J5019-19
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 091918-TIPC-F-U-B

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-BG092018.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	5.150	312	317	325	rBV	30462	50719	1.14%	0.322%
2	5.615	390	396	414	rBV	219222	426793	9.59%	2.713%
3	7.207	658	667	684	rBV	131729	307978	6.92%	1.957%
4	7.577	722	730	748	rBV	416969	815563	18.33%	5.183%
5	8.047	801	810	825	rBV	114700	234938	5.28%	1.493%
6	8.370	856	865	887	rBV	580729	1121037	25.20%	7.125%
7	9.205	1000	1007	1027	rBV	487919	1049060	23.58%	6.668%
8	10.850	1277	1287	1300	rBV	162740	338514	7.61%	2.151%
9	13.294	1696	1703	1722	rBV	1508666	2540556	57.11%	16.147%
10	14.116	1836	1843	1852	rBV	28581	50619	1.14%	0.322%
11	14.669	1930	1937	1950	rBV2	307851	530589	11.93%	3.372%
12	16.155	2183	2190	2205	rBV	836944	1397888	31.42%	8.885%
13	17.418	2398	2405	2423	rBV2	431556	718591	16.15%	4.567%
14	20.021	2842	2848	2861	rBV	3069944	4448796	100.00%	28.275%
15	21.678	3124	3130	3140	rBV2	486166	847228	19.04%	5.385%
16	24.886	3666	3676	3688	rBV2	286799	854996	19.22%	5.434%

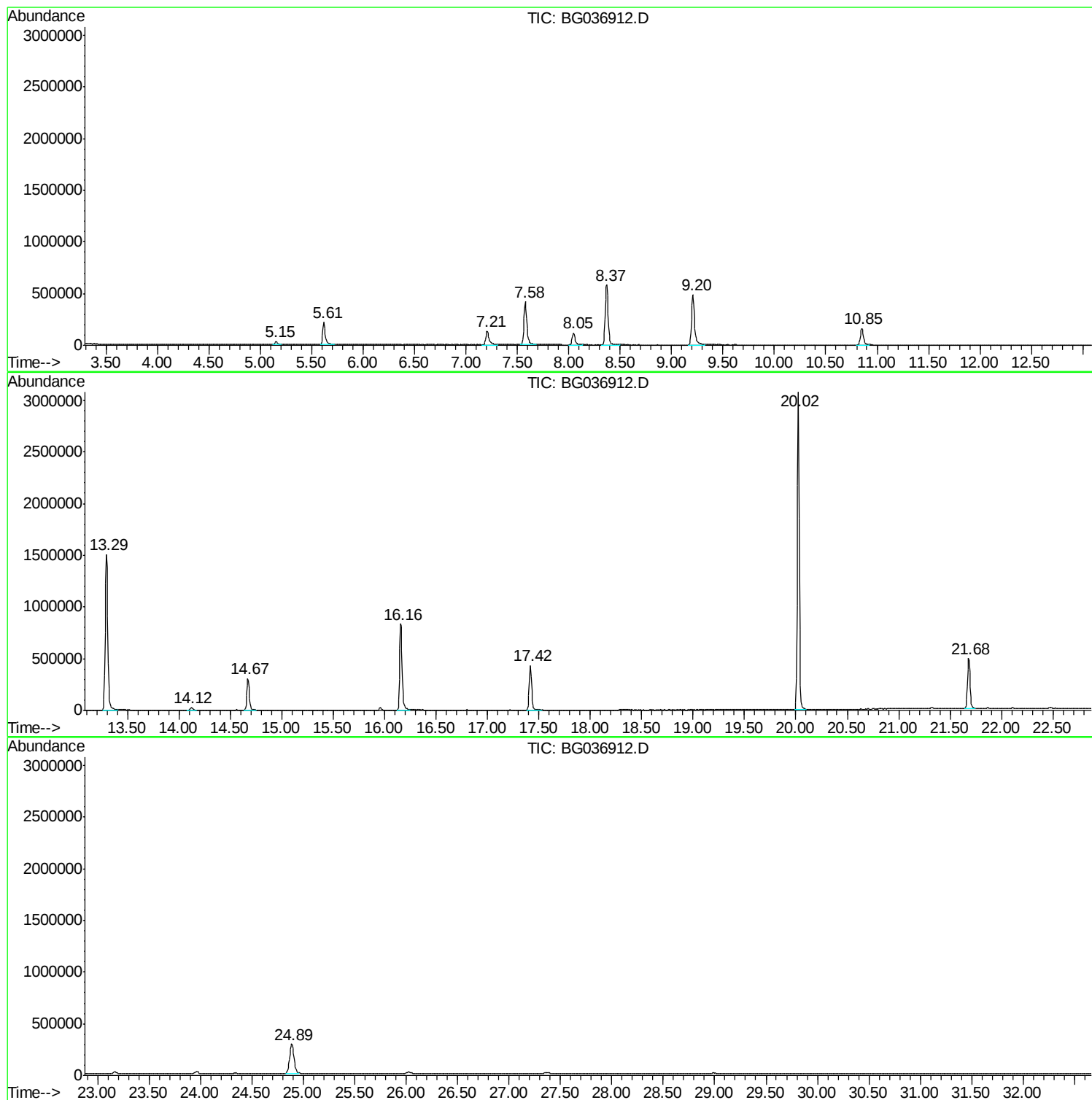
Sum of corrected areas: 15733865

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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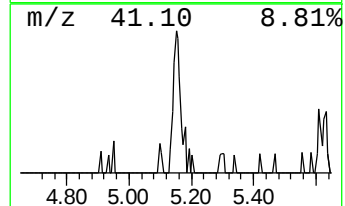
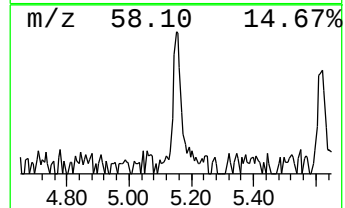
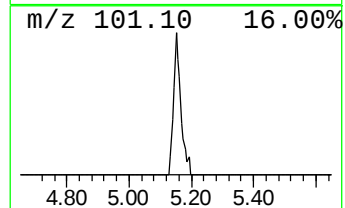
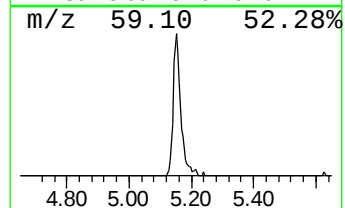
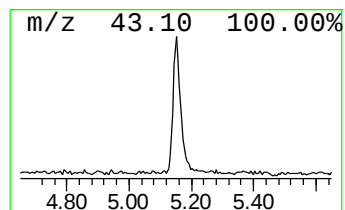
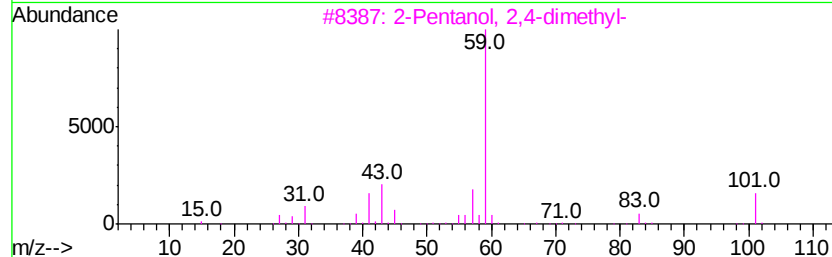
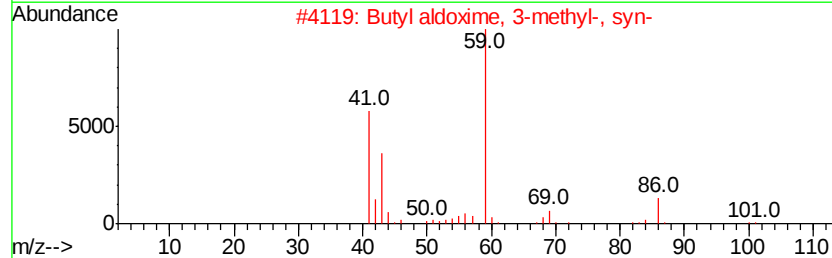
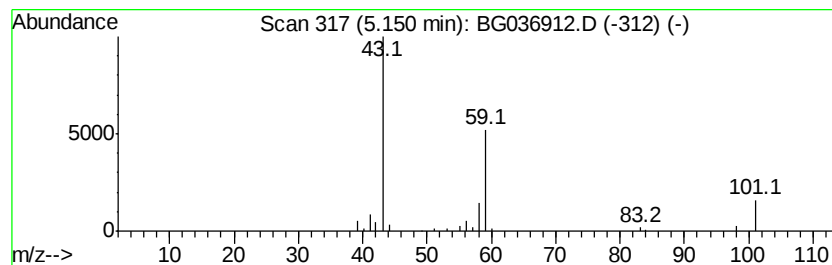
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	4.32 ng	50719	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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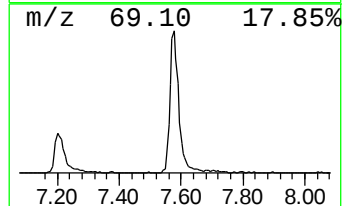
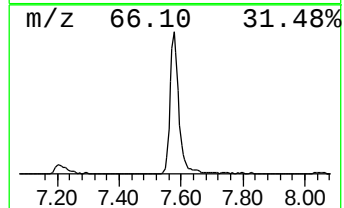
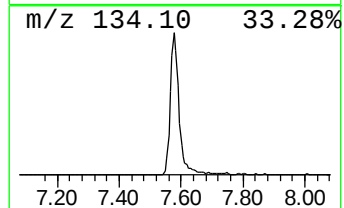
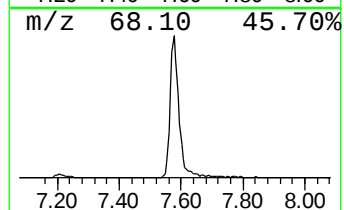
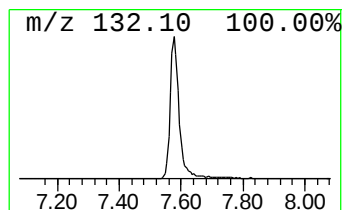
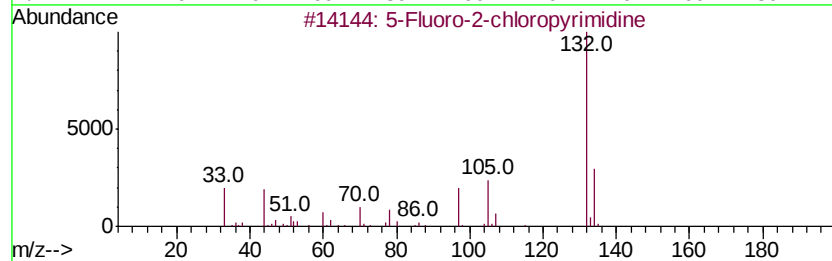
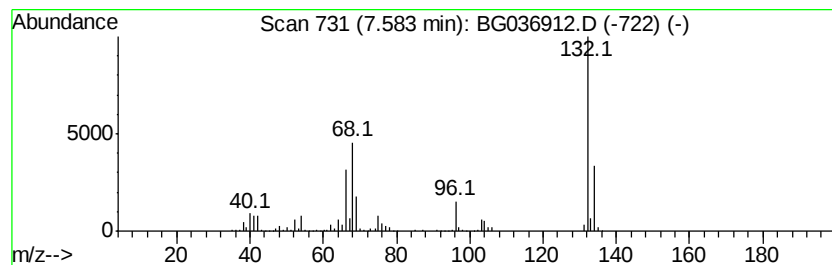
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	69.43 ng	815563	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	22
3			Tranvlcypromine-propionvl	189	C12H15NO	1000123-86-3	10
4			1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	10
5			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	4.3	ng	50719	1	8.04	234938	20.0
unknown7.58	7.58	69.4	ng	815563	1	8.04	234938	20.0