

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
 Data File : BG038059.D
 Acq On : 17 Nov 2018 7:45
 Operator : JU/SJ
 Sample : PB114894BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114894BL

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-BG111318.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	5.168	315	320	326	rVB	37114	60704	1.94%	0.376%
2	5.626	391	398	407	rBV	787720	1325391	42.46%	8.206%
3	7.230	663	671	684	rBV	807471	1454453	46.59%	9.005%
4	7.612	728	736	749	rVB	774473	1375998	44.08%	8.519%
5	8.082	810	816	824	rVB	150050	256251	8.21%	1.586%
6	8.405	864	871	879	rVB	572359	1040138	33.32%	6.440%
7	9.263	1007	1017	1026	rBV	491110	913265	29.26%	5.654%
8	10.902	1288	1296	1306	rBV	199791	373276	11.96%	2.311%
9	13.341	1704	1711	1718	rBV	1320314	2106675	67.49%	13.043%
10	14.721	1939	1946	1954	rBV2	318997	525156	16.82%	3.251%
11	16.208	2192	2199	2212	rBV4	942716	1479842	47.41%	9.162%
12	17.465	2403	2413	2421	rBV2	414997	661313	21.19%	4.094%
13	20.068	2849	2856	2862	rBV	2300802	3121567	100.00%	19.326%
14	21.748	3135	3142	3152	rBV	435375	744931	23.86%	4.612%
15	25.068	3692	3707	3721	rBV2	230698	713093	22.84%	4.415%

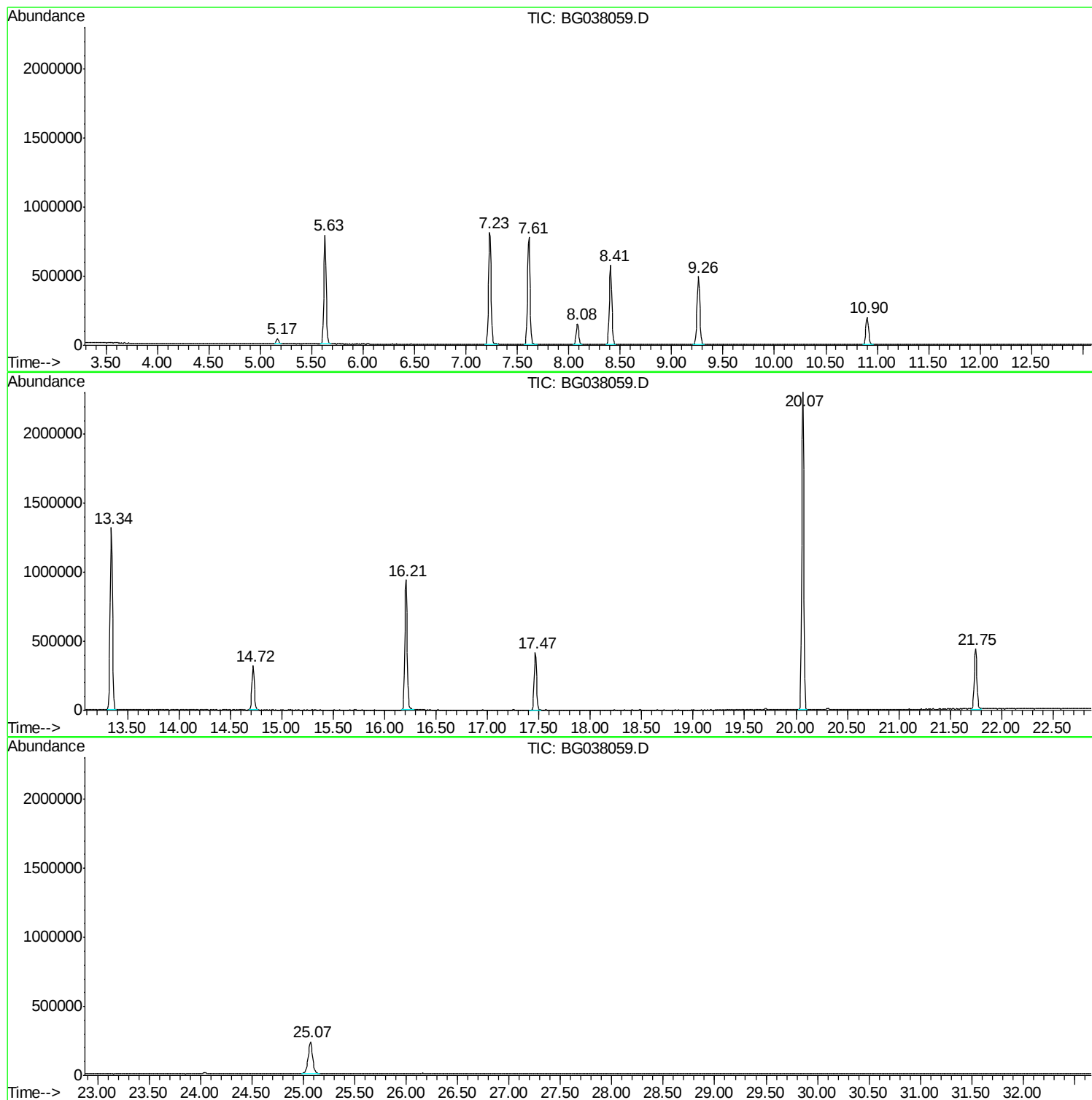
Sum of corrected areas: 16152053

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



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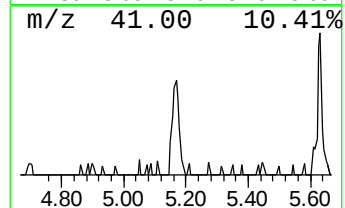
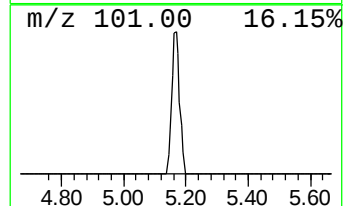
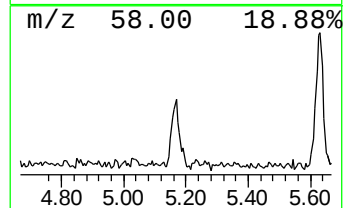
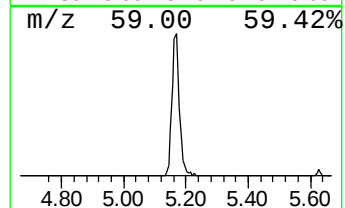
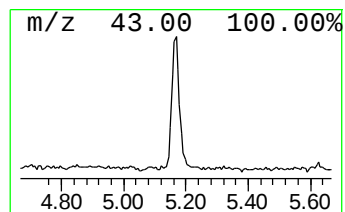
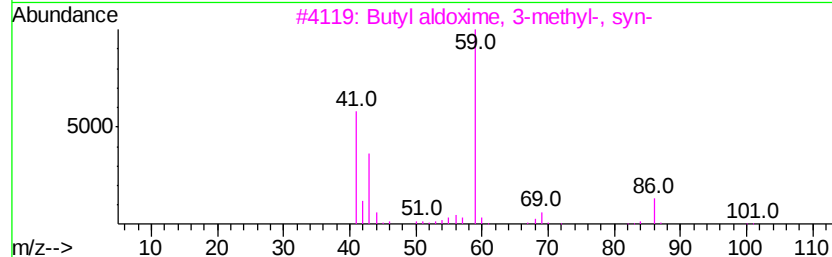
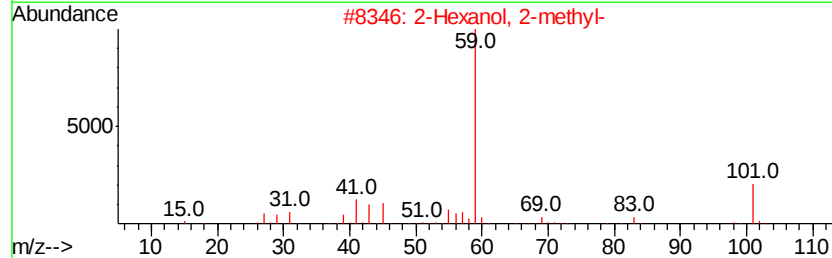
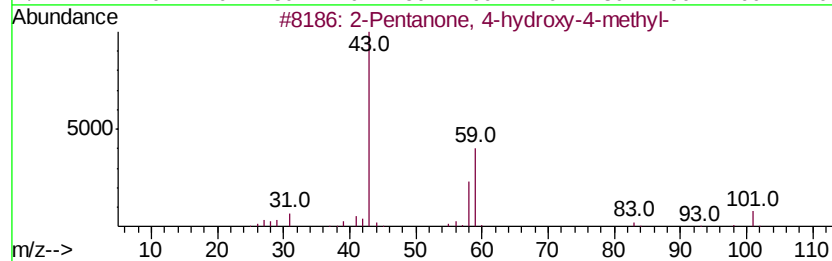
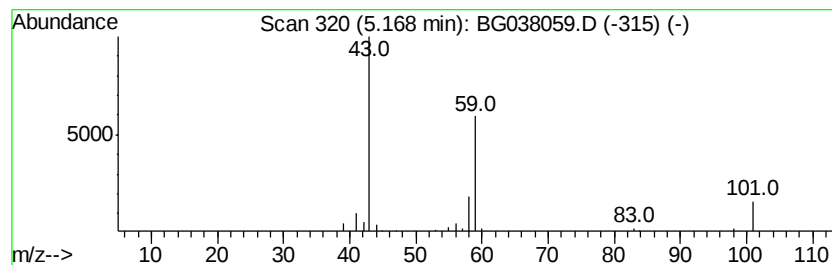
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.74 ng	60704	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			3-Hexanol	102	C6H14O	000623-37-0	9



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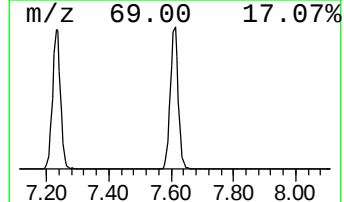
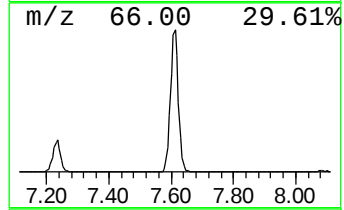
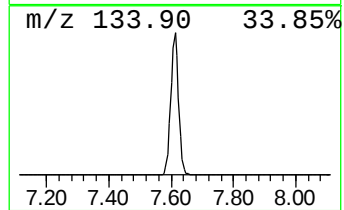
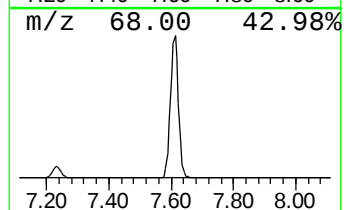
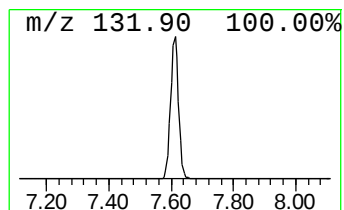
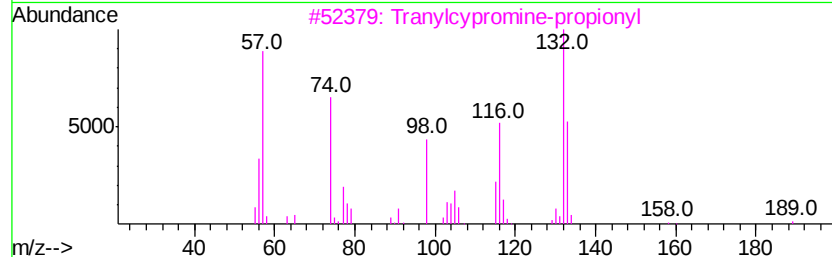
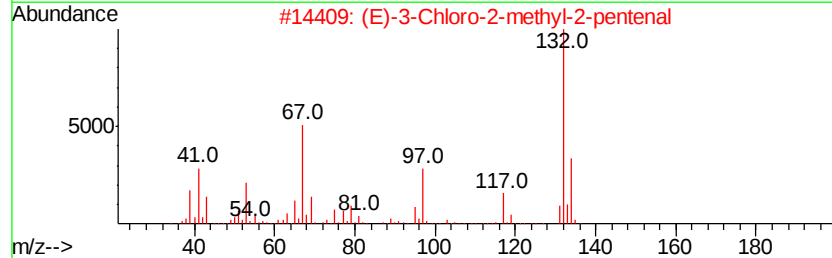
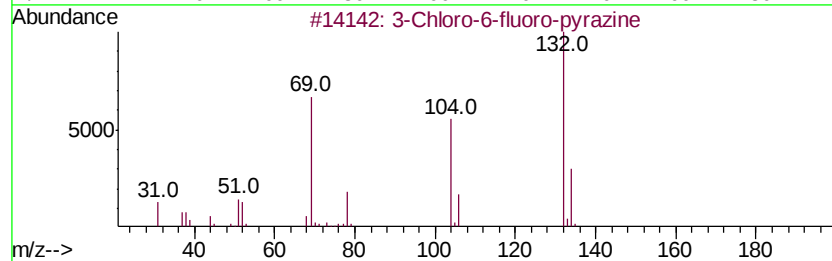
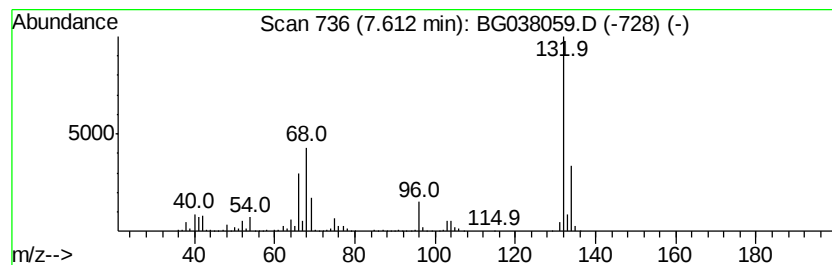
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	107.40 ng	1376000	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	4.7	ng	60704	1	8.08	256251	20.0
unknown7.61	7.61	107.4	ng	1376000	1	8.08	256251	20.0