

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121818\
 Data File : BG038702.D
 Acq On : 18 Dec 2018 19:59
 Operator : JU/SJ
 Sample : PB115802BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115802BL

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-BG121218.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	5.138	344	349	356	rBV	15056	25882	1.81%	0.382%
2	5.608	422	429	441	rBV	289290	474847	33.23%	7.000%
3	7.206	693	701	711	rBV	289814	529333	37.04%	7.803%
4	7.582	758	765	775	rBV	276059	497850	34.84%	7.339%
5	8.058	839	846	856	rVB	74097	133744	9.36%	1.972%
6	8.375	893	900	910	rVB	211146	387767	27.13%	5.716%
7	9.233	1037	1046	1057	rBV	169594	315669	22.09%	4.654%
8	10.872	1316	1325	1334	rBV	93793	178496	12.49%	2.631%
9	13.311	1733	1740	1752	rBV	487645	785064	54.93%	11.573%
10	14.691	1969	1975	1990	rVB2	163569	272600	19.07%	4.019%
11	16.184	2223	2229	2240	rBV2	350698	583721	40.84%	8.605%
12	17.441	2436	2443	2452	rBV	240701	367625	25.72%	5.419%
13	20.044	2880	2886	2896	rBV	1059236	1429120	100.00%	21.068%
14	21.718	3164	3171	3184	rBV2	240778	421396	29.49%	6.212%
15	25.014	3721	3732	3744	rBV2	129784	380286	26.61%	5.606%

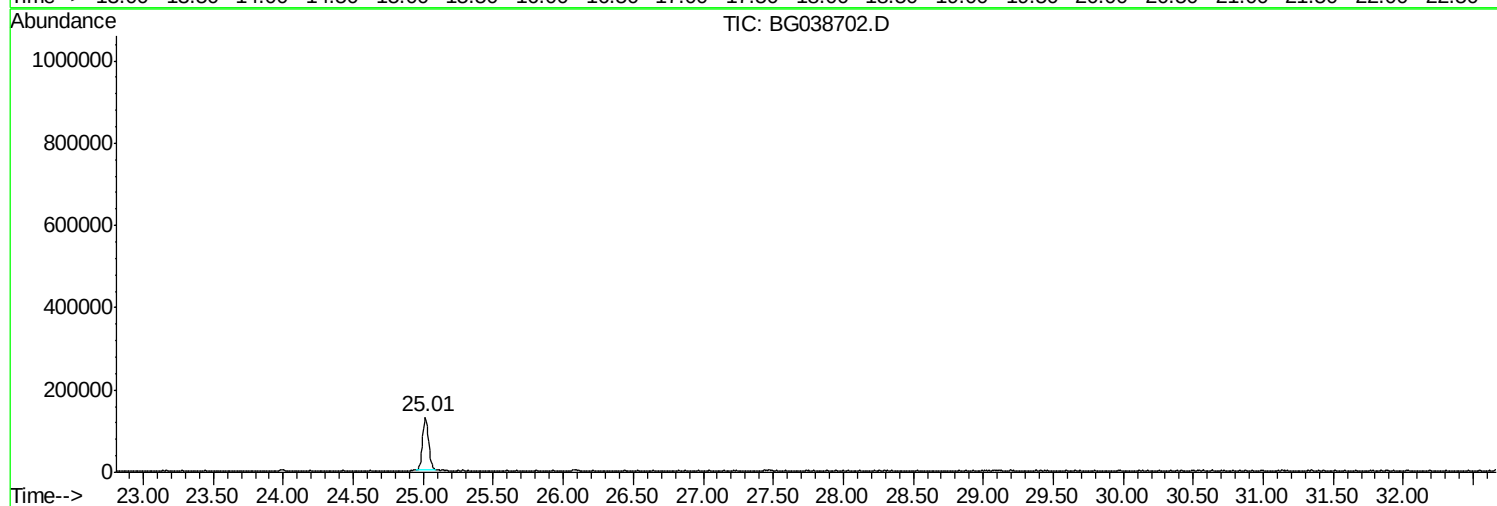
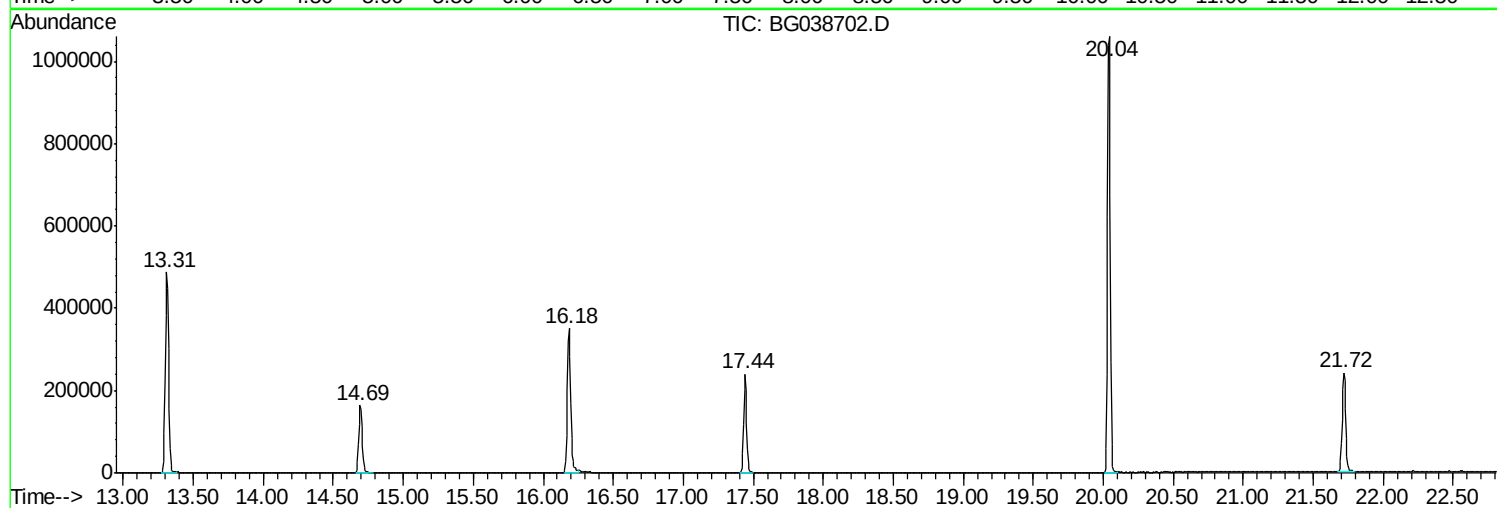
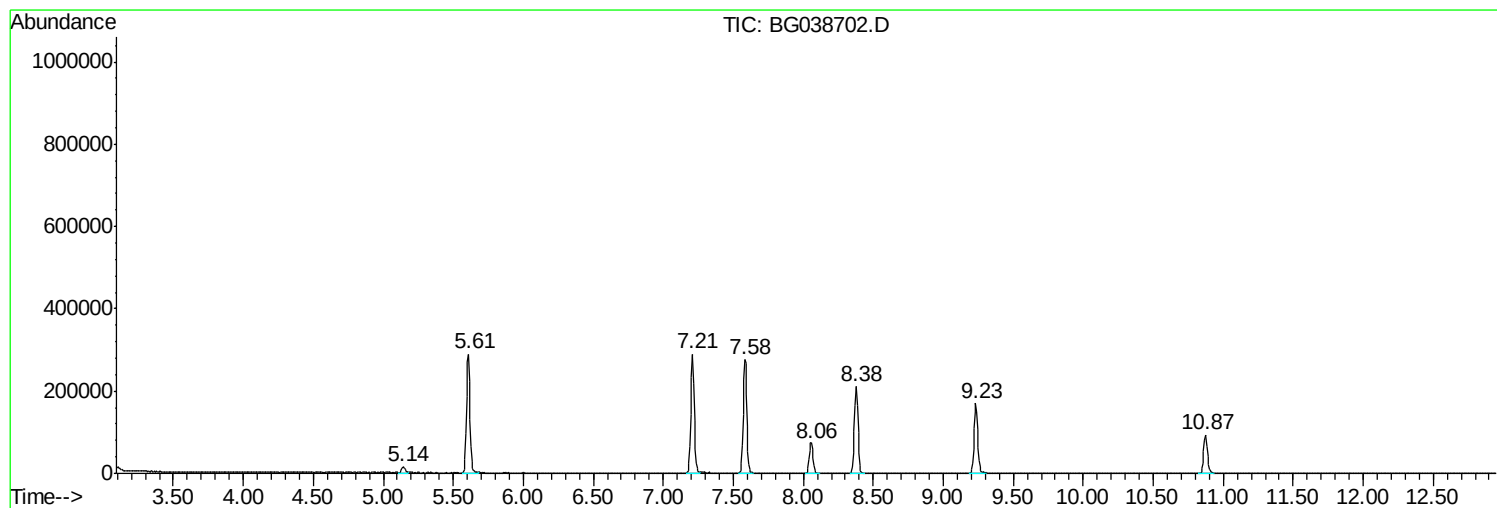
Sum of corrected areas: 6783400

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.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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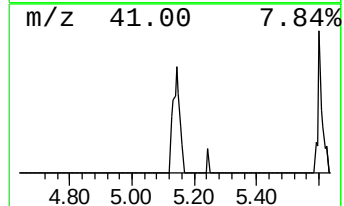
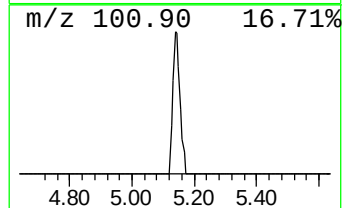
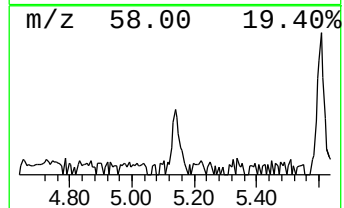
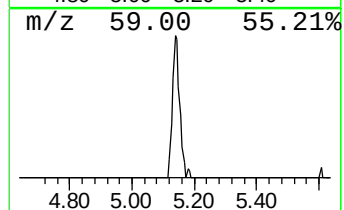
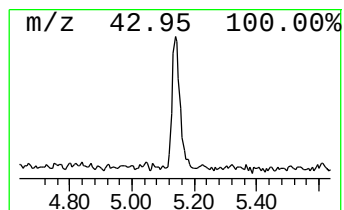
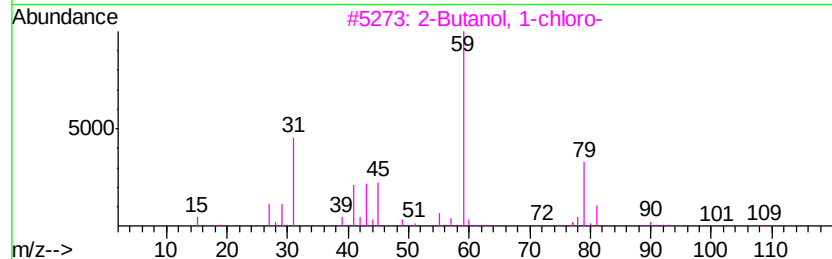
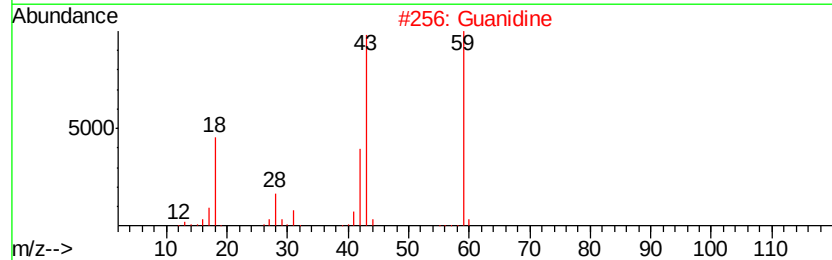
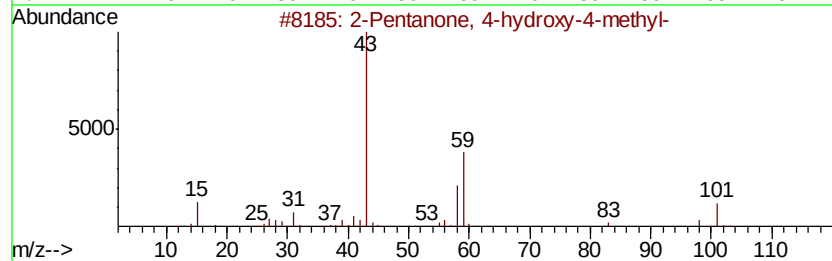
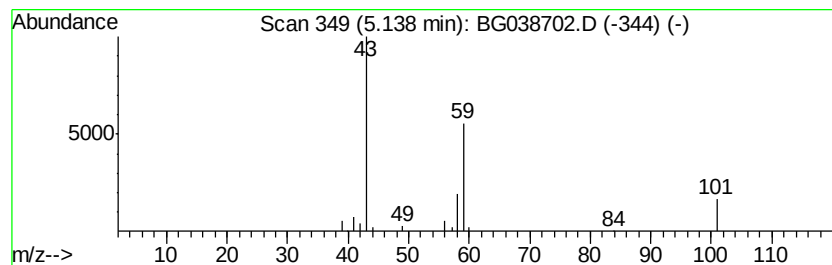
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	3.87 ng	25882	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		Guanidine	59	CH5N3	000113-00-8	9
3		2-Butanol, 1-chloro-	108	C4H9ClO	001873-25-2	9
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	9



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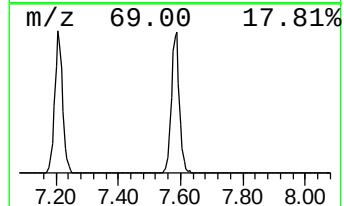
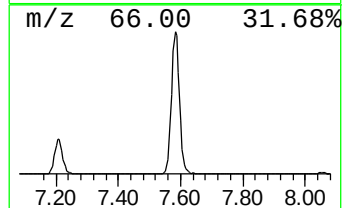
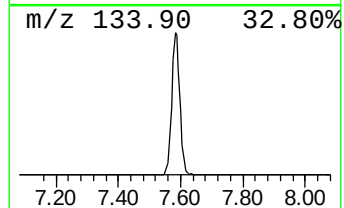
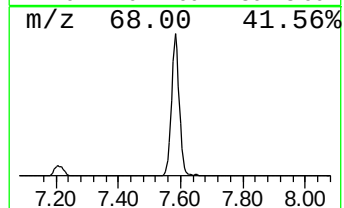
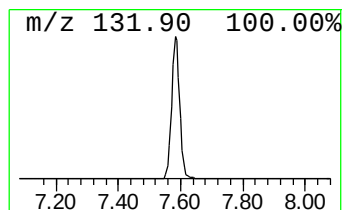
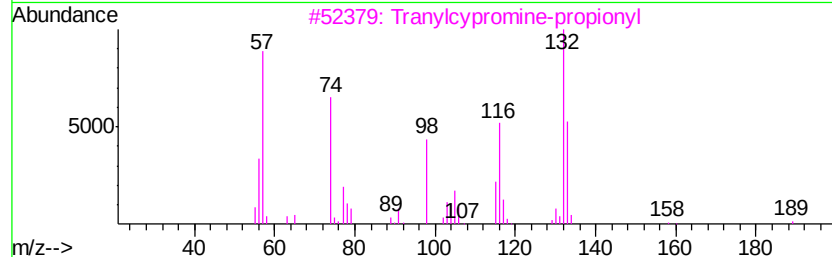
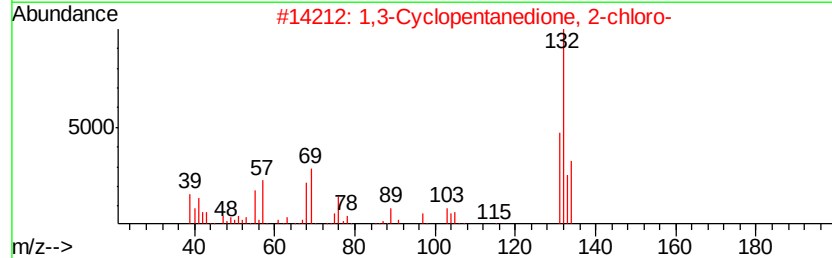
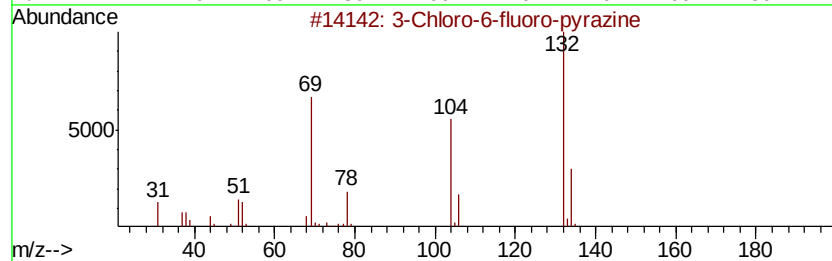
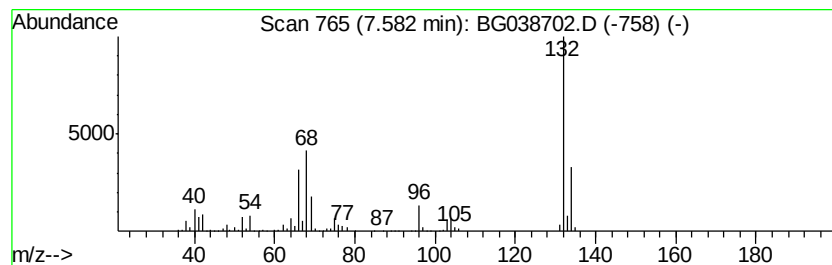
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	74.45 ng	497850	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		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	23
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Aminoindole	132	C8H8N2	005192-03-0	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	3.9	ng	25882	1	8.05	133744	20.0
unknown7.58	7.58	74.5	ng	497850	1	8.05	133744	20.0