

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121918\
 Data File : BG038731.D
 Acq On : 19 Dec 2018 17:45
 Operator : JU/SJ
 Sample : PB115831BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115831BL

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	358	rBV	19664	33236	1.83%	0.387%
2	5.608	422	429	440	rBV	353543	589853	32.48%	6.866%
3	7.212	694	702	712	rBV	363880	653602	35.99%	7.608%
4	7.582	758	765	776	rBV	345468	620648	34.17%	7.225%
5	8.058	837	846	854	rBV	93165	163113	8.98%	1.899%
6	8.381	893	901	909	rBV	267249	480257	26.44%	5.591%
7	9.233	1037	1046	1061	rBV	204756	394400	21.72%	4.591%
8	10.872	1318	1325	1336	rBV	119686	220507	12.14%	2.567%
9	13.311	1732	1740	1754	rBV	604217	983556	54.15%	11.449%
10	14.697	1969	1976	1986	rBV2	194519	335580	18.48%	3.906%
11	16.184	2222	2229	2247	rBV	469727	781352	43.02%	9.096%
12	17.441	2436	2443	2456	rBV	287609	457516	25.19%	5.326%
13	20.044	2880	2886	2897	rVB	1380191	1816190	100.00%	21.142%
14	21.718	3162	3171	3180	rBV	339178	551394	30.36%	6.419%
15	25.020	3721	3733	3748	rVB2	162081	509317	28.04%	5.929%

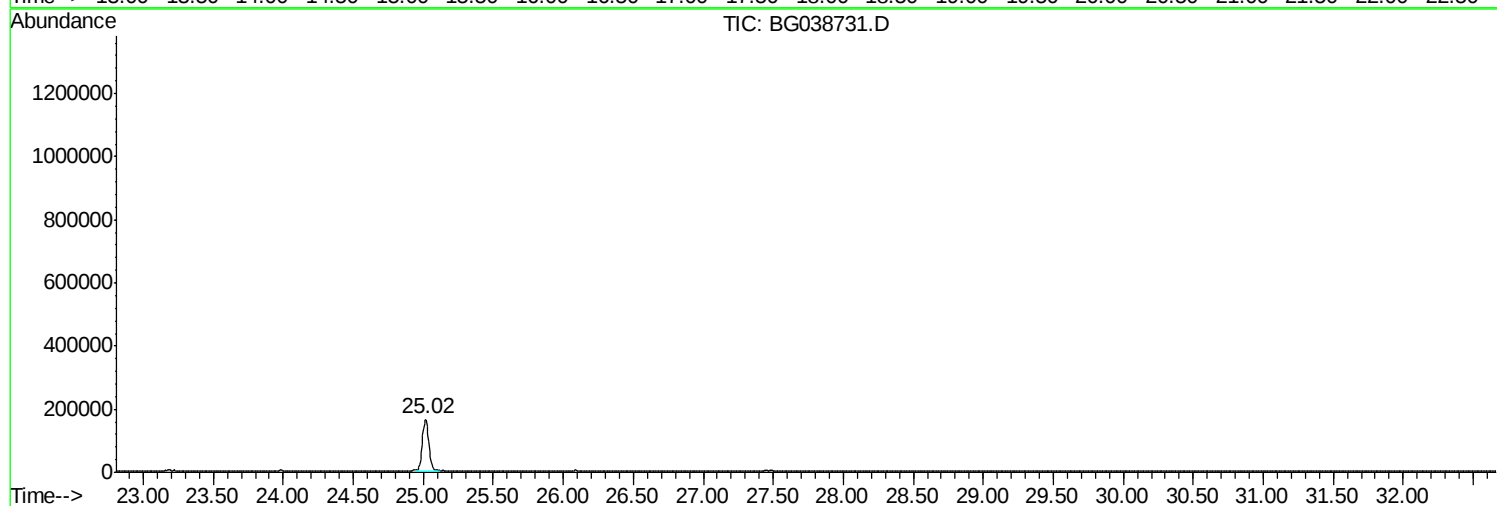
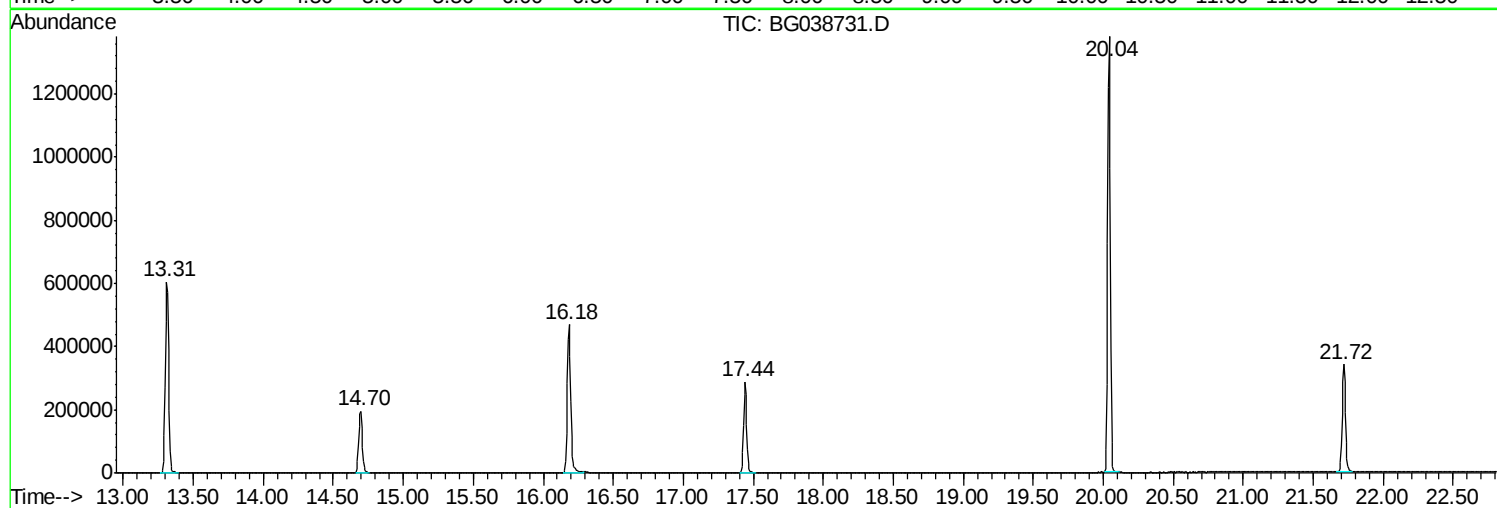
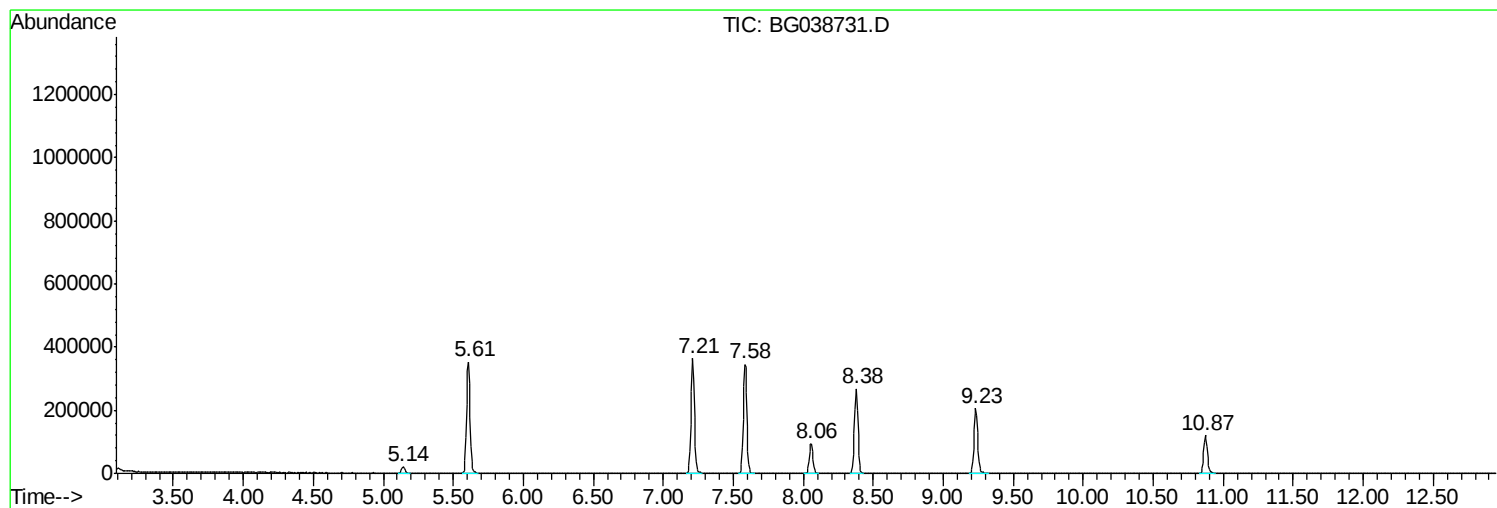
Sum of corrected areas: 8590521

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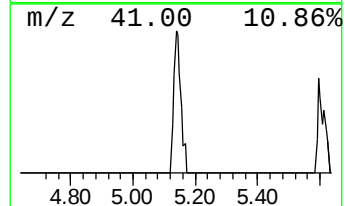
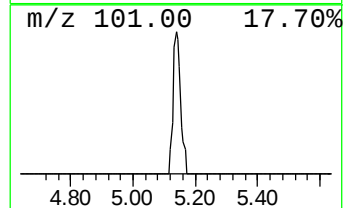
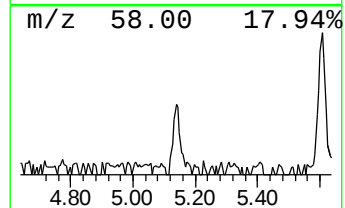
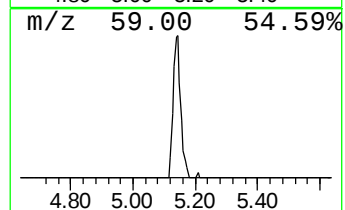
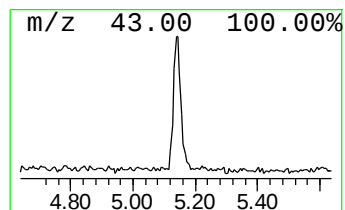
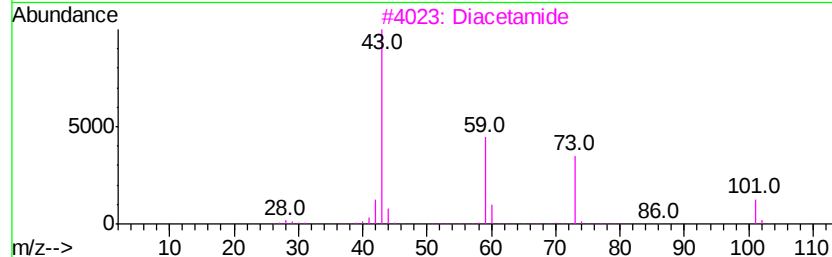
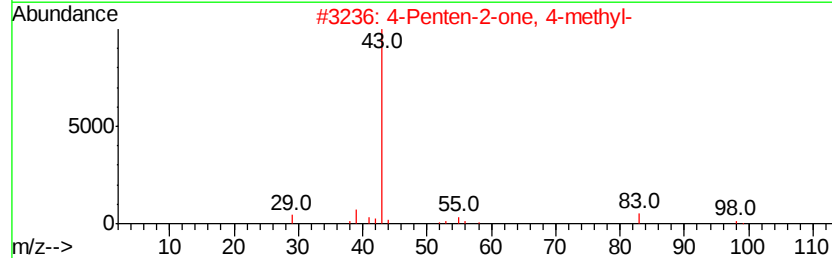
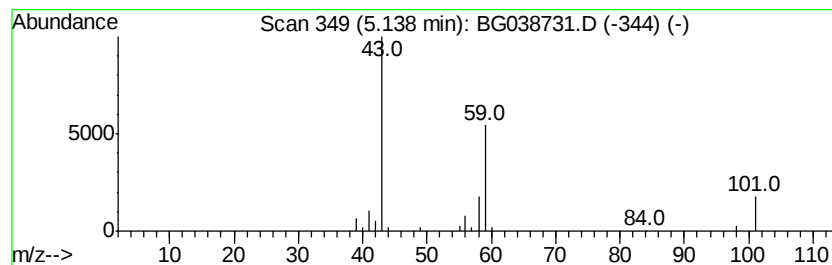
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.08 ng	33236	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9		
3	Diacetamide	101	C4H7NO2	000625-77-4	9		
4	Acetone	58	C3H6O	000067-64-1	9		
5	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9		



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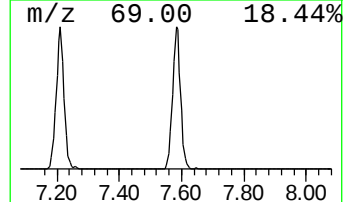
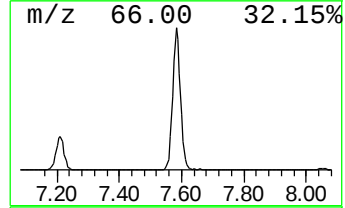
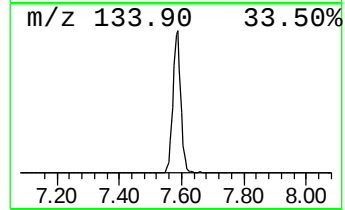
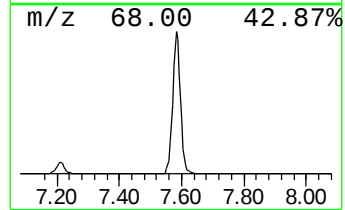
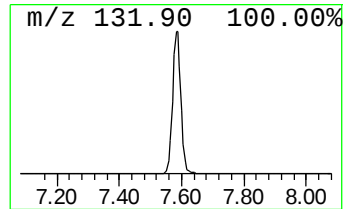
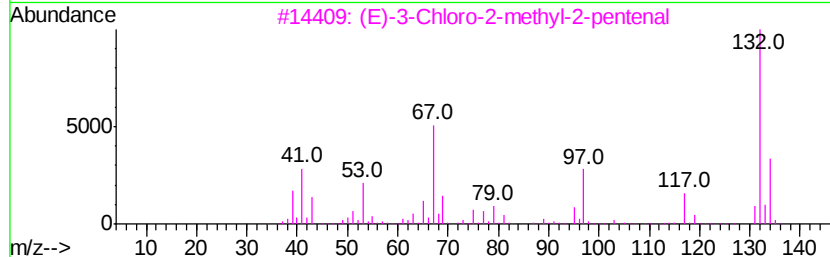
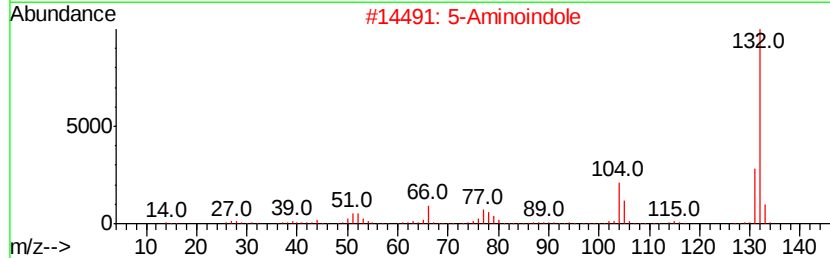
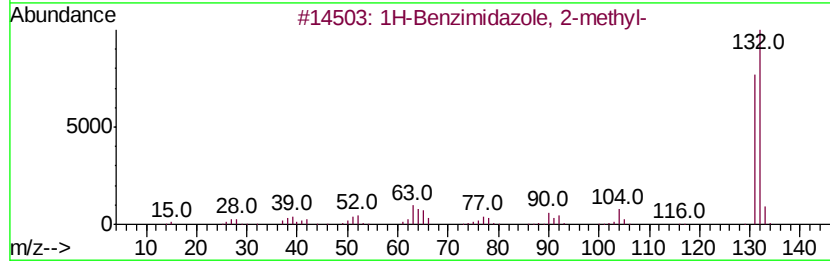
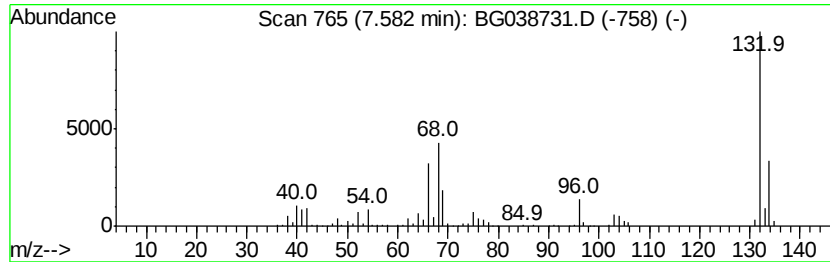
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	76.10 ng	620648	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		1-Isoquinolinemethanol, .alpha.-...	191	C12H17NO	114608-92-3	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	4.1	ng	33236	1	8.06	163113	20.0
unknown7.58	7.58	76.1	ng	620648	1	8.06	163113	20.0