

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110518\
 Data File : BG037798.D
 Acq On : 5 Nov 2018 11:56
 Operator : JU/SJ
 Sample : PB114465BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114465BL

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-BG101818.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	5.174	315	321	333	rBV	34532	59062	1.90%	0.404%
2	5.638	393	400	409	rBV	508094	830828	26.77%	5.684%
3	7.236	663	672	683	rBV	529221	956610	30.82%	6.545%
4	7.618	729	737	747	rBV	508125	899742	28.99%	6.156%
5	8.094	811	818	828	rVB	167201	297477	9.59%	2.035%
6	8.417	865	873	884	rBV	568504	1023372	32.98%	7.002%
7	9.269	1011	1018	1030	rBV	473425	877806	28.29%	6.006%
8	10.914	1291	1298	1309	rVB	245134	464369	14.96%	3.177%
9	13.352	1705	1713	1722	rBV	1389298	2301201	74.15%	15.745%
10	14.727	1937	1947	1960	rBV3	423805	713618	22.99%	4.883%
11	16.214	2192	2200	2218	rBV2	500296	841395	27.11%	5.757%
12	17.477	2408	2415	2425	rBV2	508653	824523	26.57%	5.641%
13	20.074	2851	2857	2869	rBV	2342450	3103407	100.00%	21.233%
14	21.760	3137	3144	3158	rBV	424812	827559	26.67%	5.662%
15	25.086	3697	3710	3724	rBV2	153305	594780	19.17%	4.069%

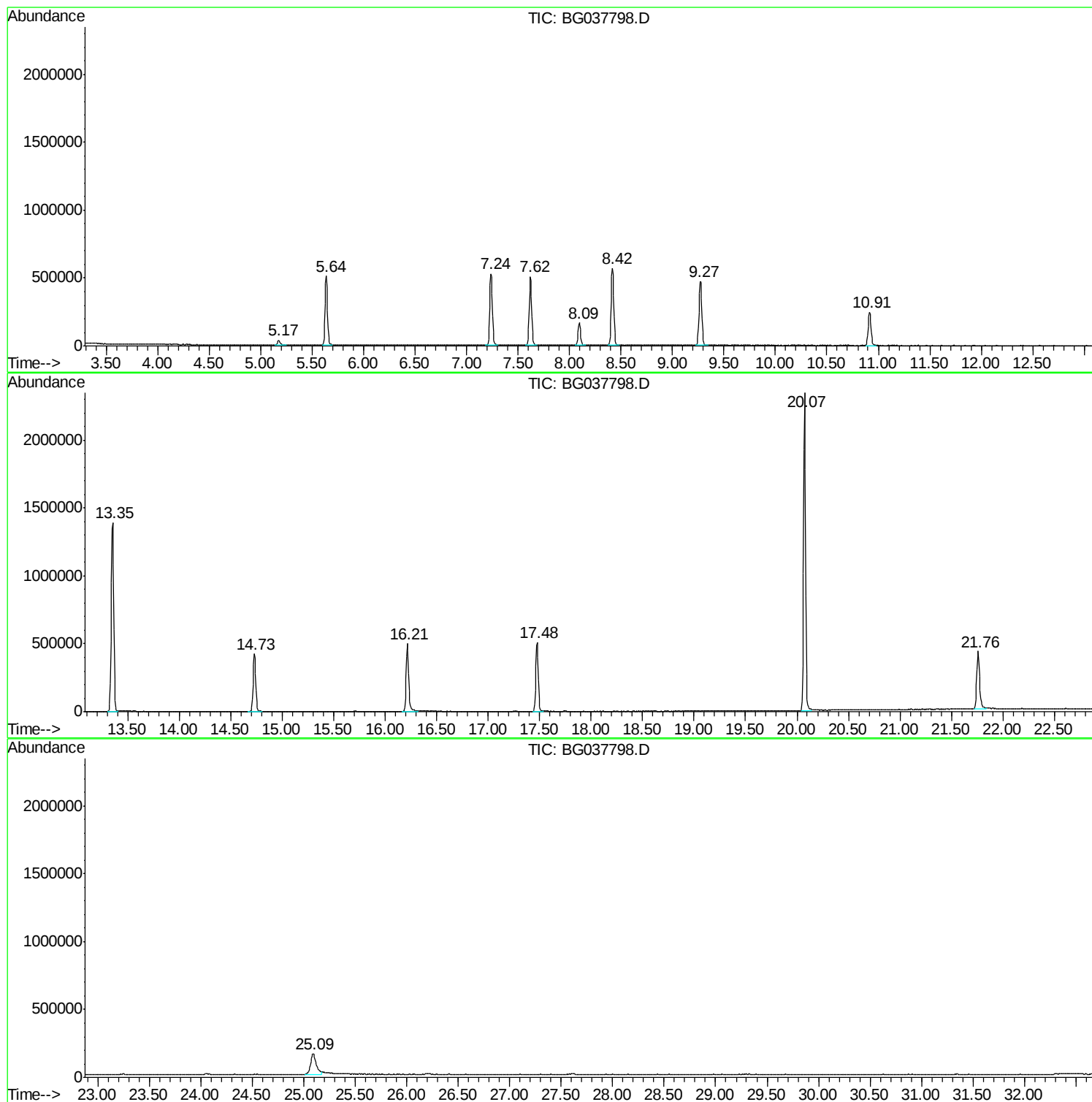
Sum of corrected areas: 14615749

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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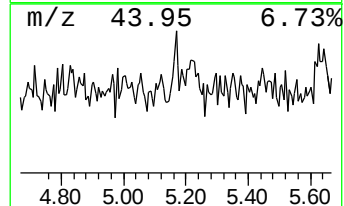
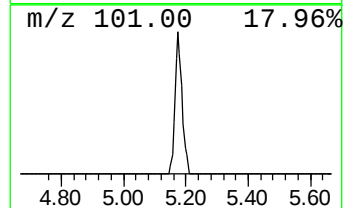
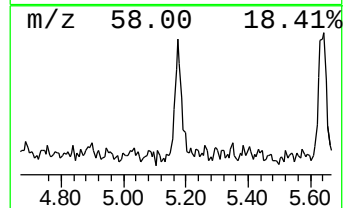
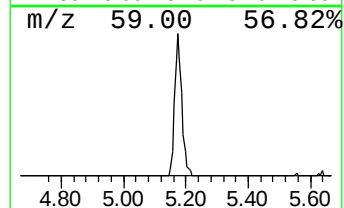
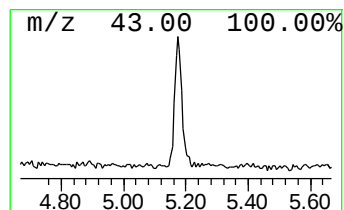
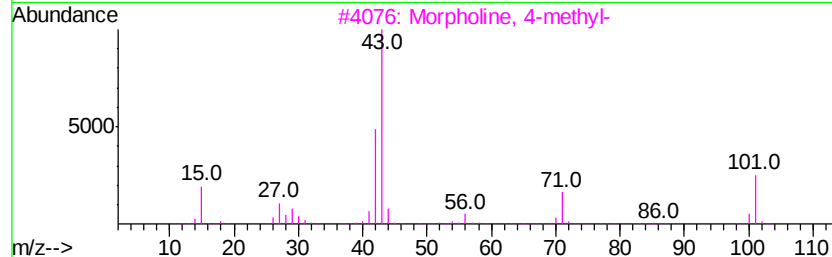
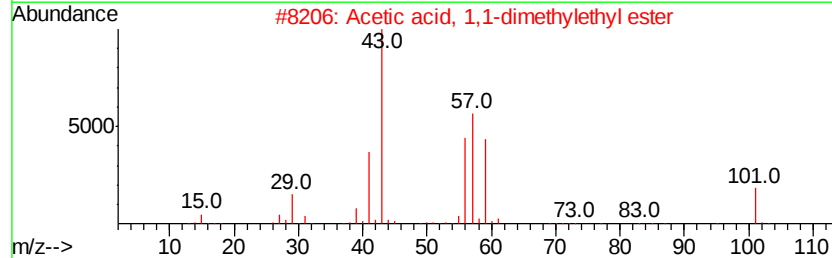
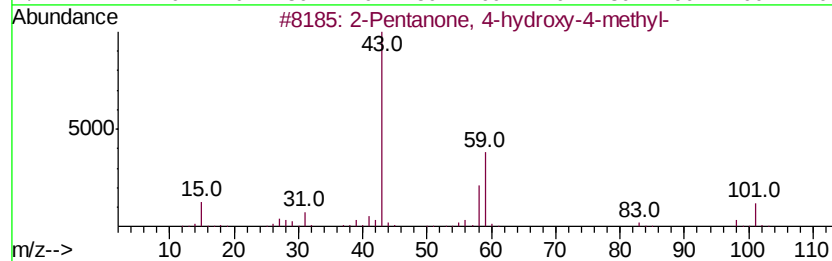
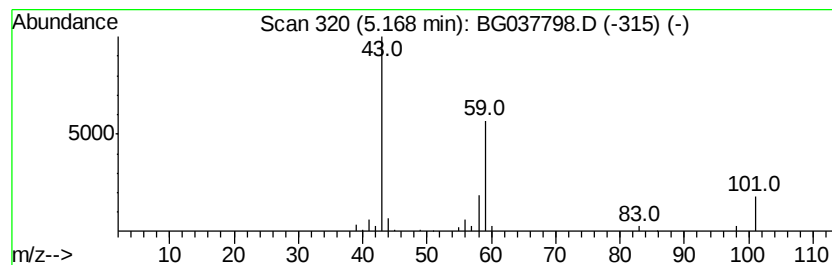
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.97 ng	59062	1,4-Dichlorobenzene-d4	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	33
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Acetone	58	C3H6O	000067-64-1	9



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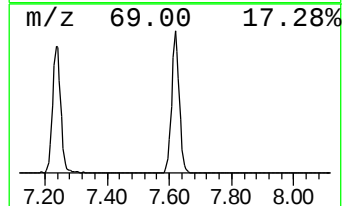
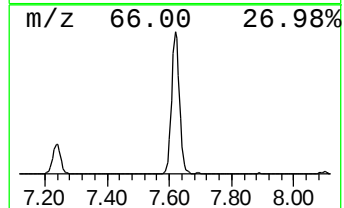
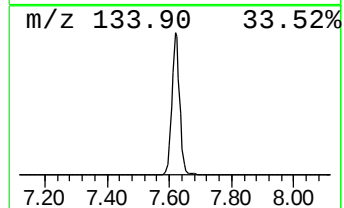
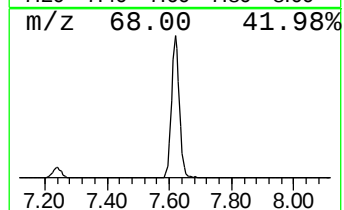
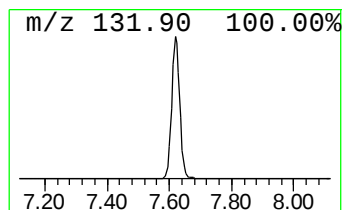
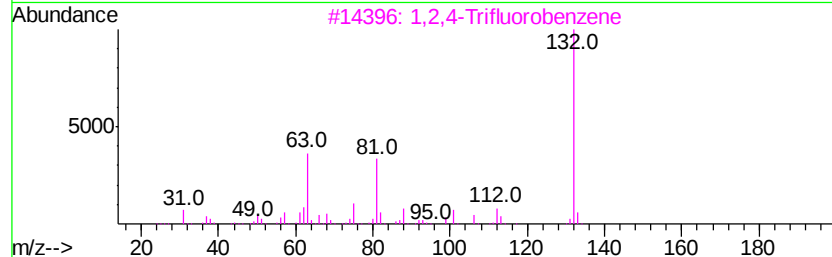
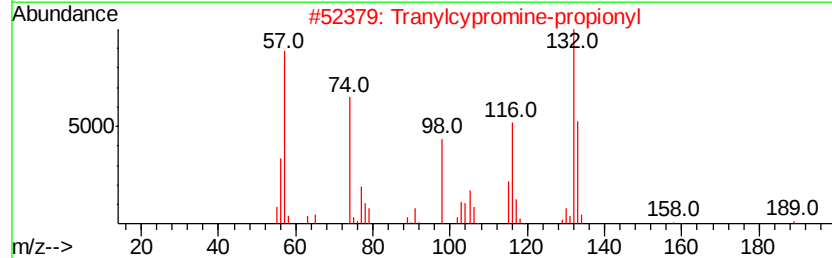
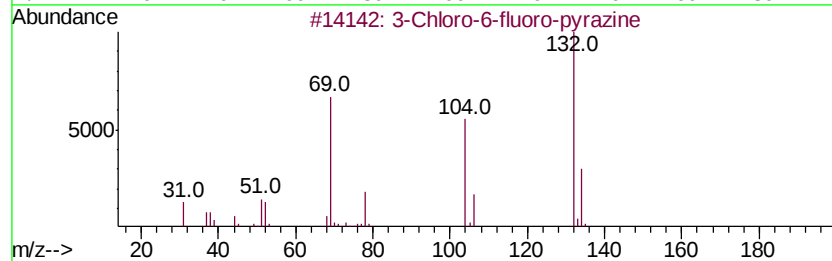
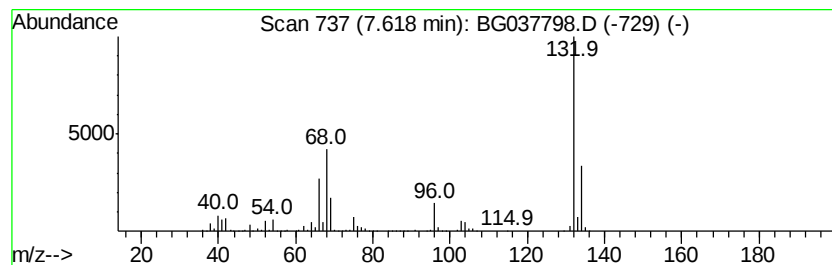
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Peak Number 2 unknown7.62 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	60.49 ng	899742	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	4.0	ng	59062	1	8.09	297477	20.0
unknown7.62	7.62	60.5	ng	899742	1	8.09	297477	20.0