

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
 Data File : BG037644.D
 Acq On : 30 Oct 2018 4:55
 Operator : JU/SJ
 Sample : PB114248BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114248BL

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	5.181	314	322	331	rBV	29504	52688	1.85%	0.336%
2	5.645	394	401	412	rVB	650796	1081853	37.91%	6.895%
3	7.243	664	673	683	rBV	705650	1233535	43.23%	7.862%
4	7.625	729	738	750	rBV	645805	1150280	40.31%	7.331%
5	8.101	811	819	828	rVB	186281	318591	11.16%	2.031%
6	8.424	866	874	883	rBV	477258	851977	29.86%	5.430%
7	9.276	1012	1019	1030	rBV	397467	733533	25.71%	4.675%
8	10.921	1292	1299	1311	rVB	264469	486850	17.06%	3.103%
9	13.353	1706	1713	1724	rBV	1128021	1890888	66.27%	12.052%
10	14.734	1939	1948	1957	rBV2	464846	762523	26.72%	4.860%
11	16.221	2194	2201	2218	rBV	838824	1344589	47.12%	8.570%
12	17.484	2409	2416	2427	rBV	628690	982829	34.44%	6.264%
13	20.081	2852	2858	2870	rBV	2186023	2853506	100.00%	18.187%
14	21.767	3138	3145	3156	rBV	593396	1031965	36.16%	6.577%
15	25.098	3698	3712	3728	rBV3	253616	914270	32.04%	5.827%

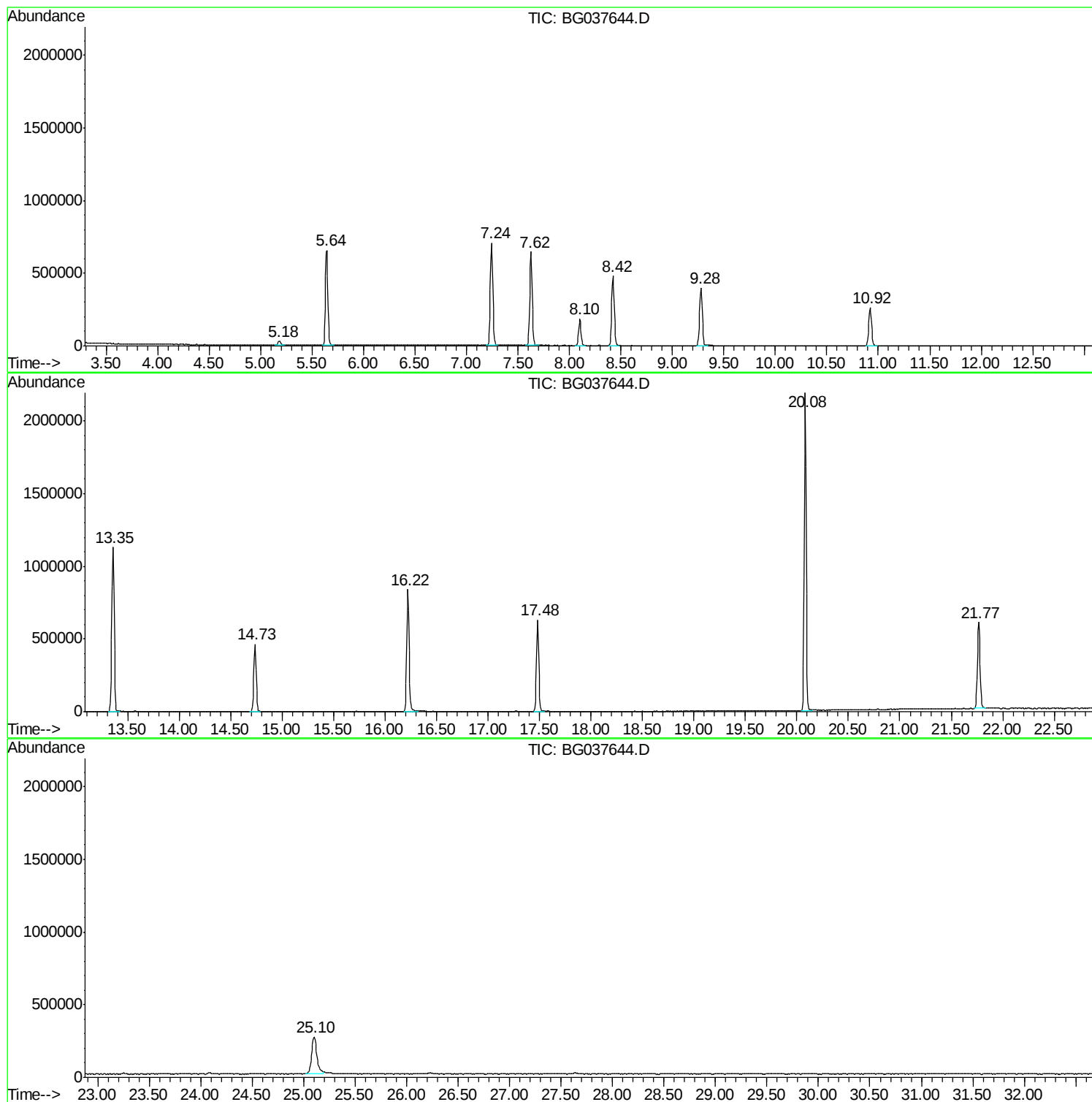
Sum of corrected areas: 15689877

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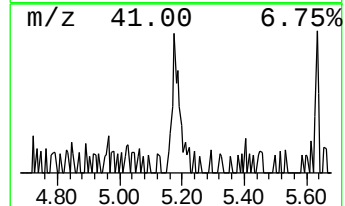
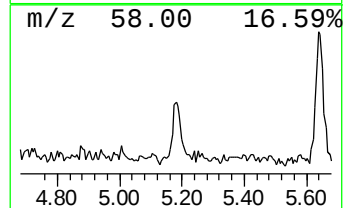
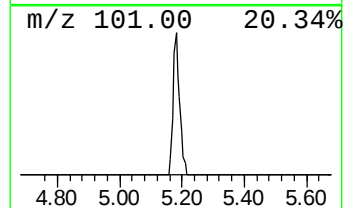
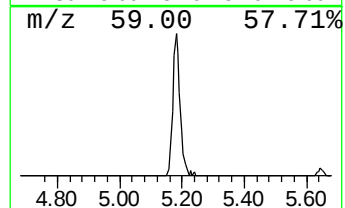
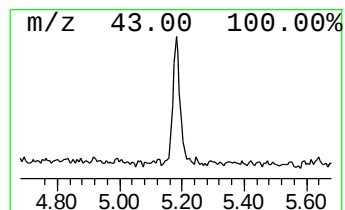
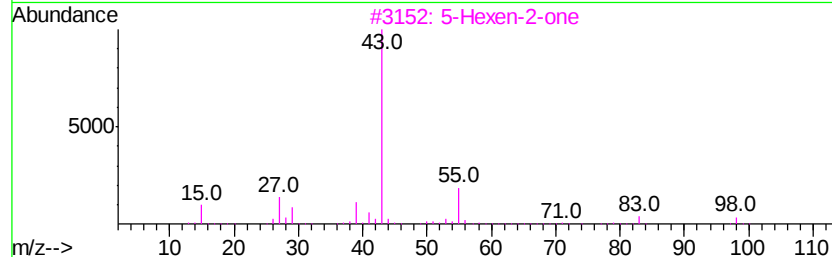
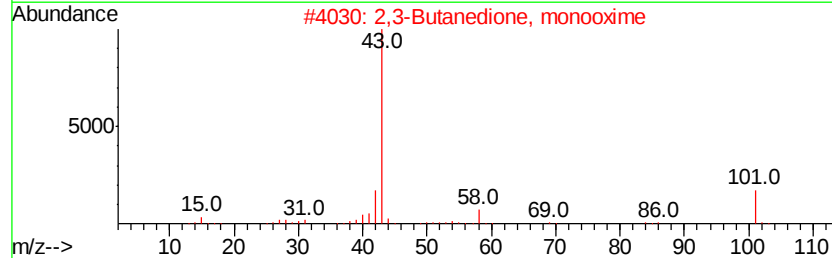
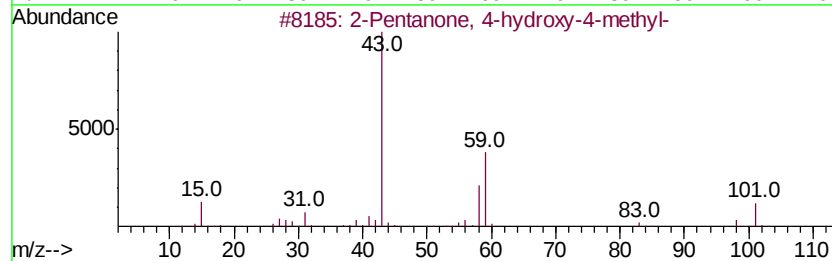
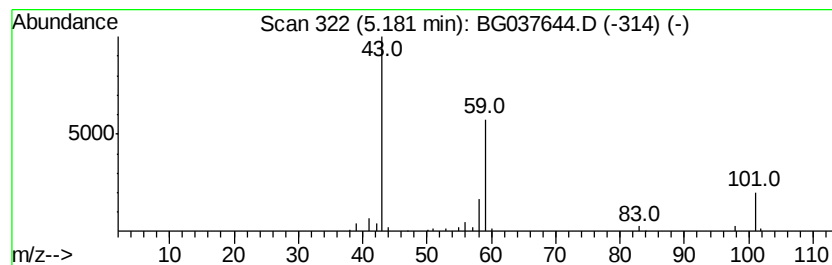
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.31 ng	52688	1,4-Dichlorobenzene-d4	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			5-Hexen-2-one	98	C6H10O	000109-49-9	10
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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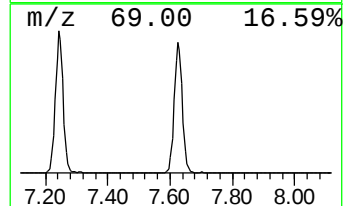
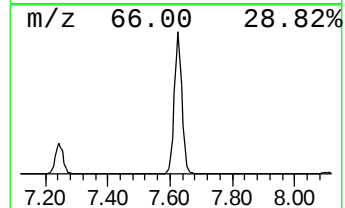
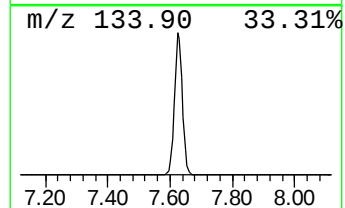
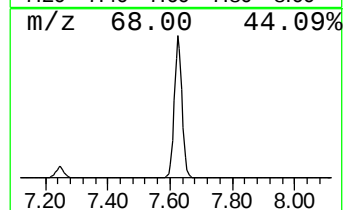
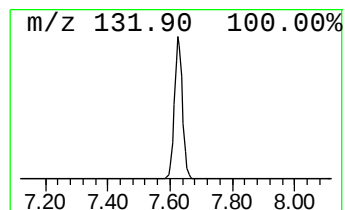
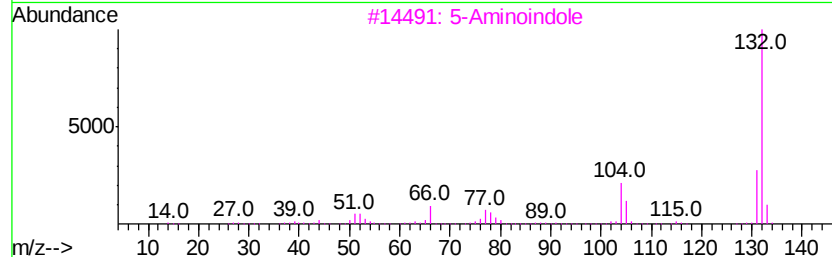
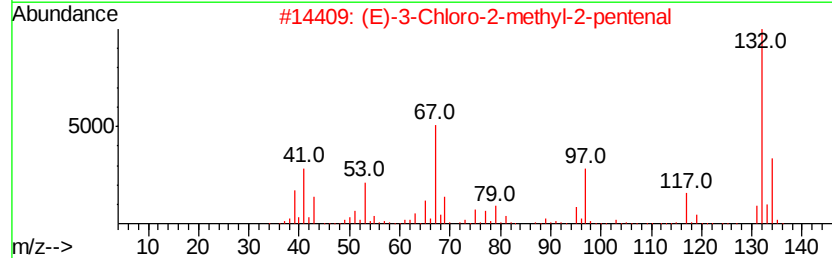
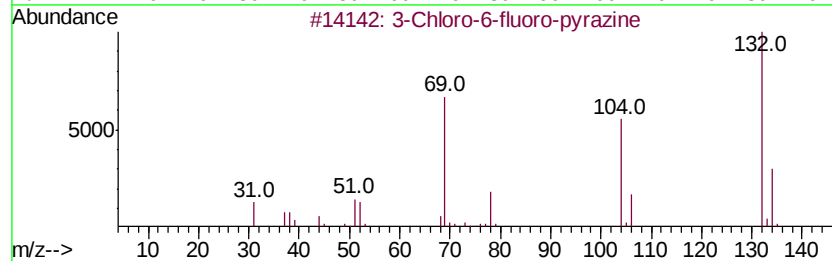
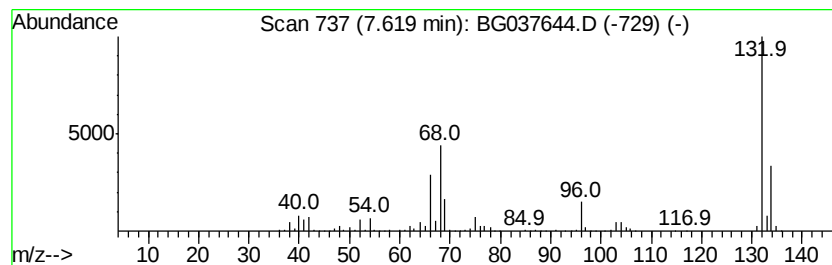
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	72.21 ng	1150280	1,4-Dichlorobenzene-d4	8.10

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Amphetaminil	250	C17H18N2	017590-01-1	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	3.3	ng	52688	1	8.10	318591	20.0
unknown7.62	7.62	72.2	ng	1150280	1	8.10	318591	20.0