

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012017\
 Data File : BG025539.D
 Acq On : 20 Jan 2017 15:20
 Operator : UM/SJ
 Sample : PB96092BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB96092BL

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:\HPCHEM1\BNA_G\METHODS\8270-BG122116.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	5.235	400	408	416	rBV	658631	1081754	13.19%	2.981%
2	5.717	483	490	505	rBV	1524327	2624879	32.01%	7.233%
3	7.339	758	766	778	rBV	1566644	2840008	34.63%	7.826%
4	7.709	821	829	839	rBV	1513906	2704371	32.98%	7.452%
5	8.179	903	909	917	rVB2	268397	481248	5.87%	1.326%
6	8.502	956	964	973	rBV	1113918	2043601	24.92%	5.631%
7	9.366	1102	1111	1122	rBV	889081	1769940	21.58%	4.877%
8	11.011	1384	1391	1401	rBV	287780	614999	7.50%	1.695%
9	13.431	1795	1803	1812	rBV	2435309	4025707	49.09%	11.093%
10	14.812	2031	2038	2054	rBV	513993	935213	11.41%	2.577%
11	16.299	2284	2291	2307	rBV2	2248415	3863425	47.12%	10.646%
12	17.556	2499	2505	2514	rBV2	719220	1208453	14.74%	3.330%
13	20.135	2938	2944	2952	rBV	5987789	8199929	100.00%	22.595%
14	21.845	3228	3235	3245	rBV	1020491	1913920	23.34%	5.274%
15	25.223	3800	3810	3833	rVB2	617714	1982876	24.18%	5.464%

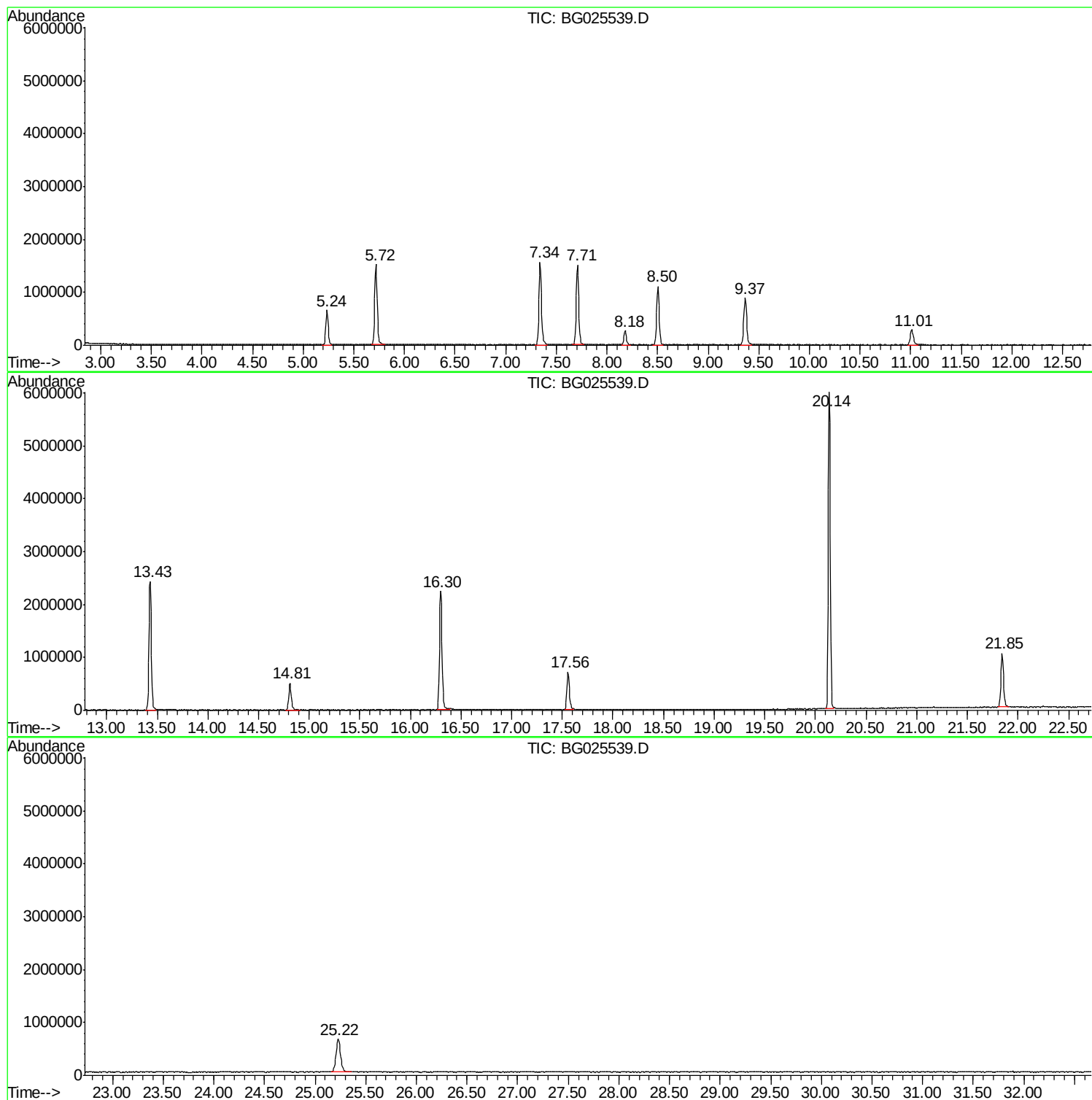
Sum of corrected areas: 36290323

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.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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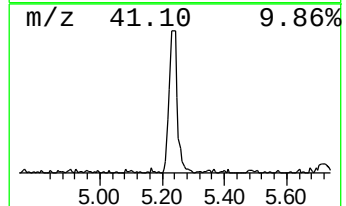
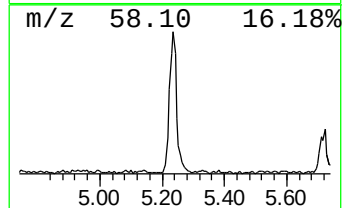
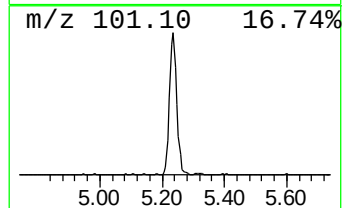
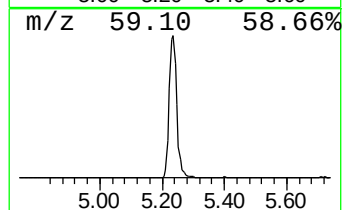
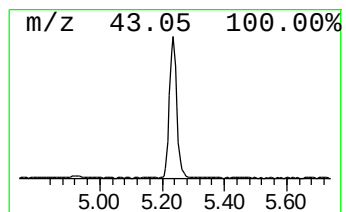
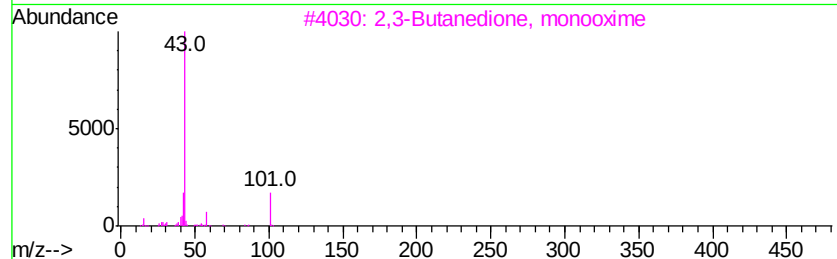
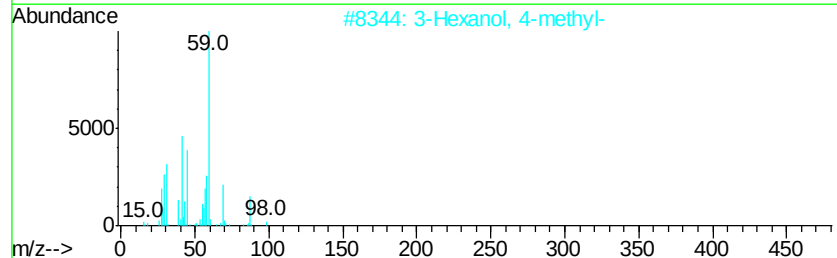
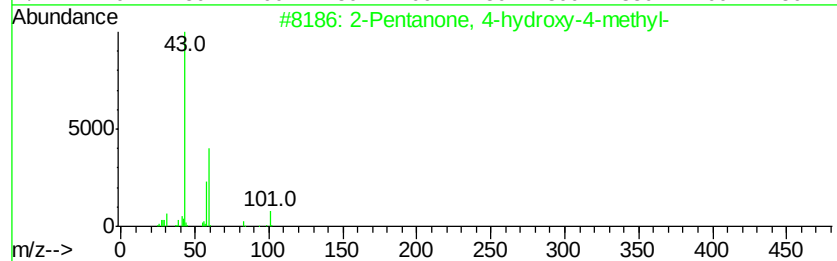
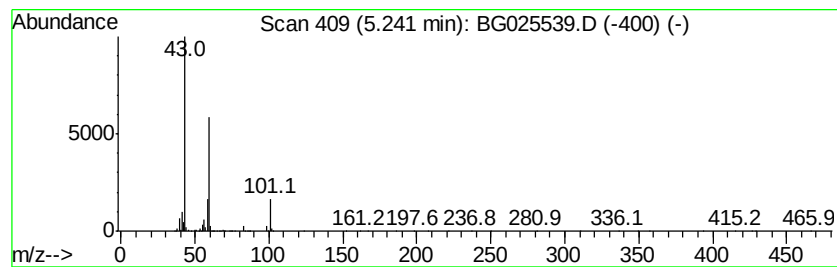
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	44.96 ng	1081750	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4		Oxirane, tetramethyl-	100	C6H12O	005076-20-0	9
5		4-Methoxy-1-pentene	100	C6H12O	098386-09-5	9



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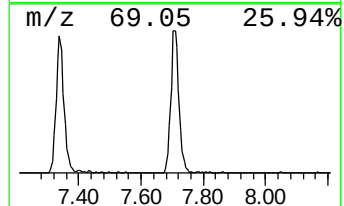
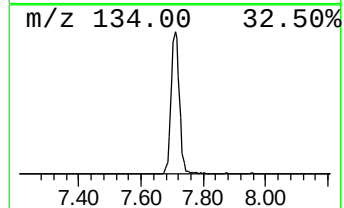
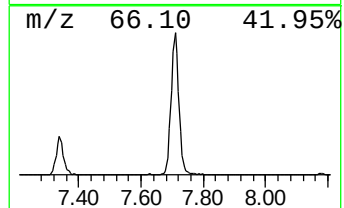
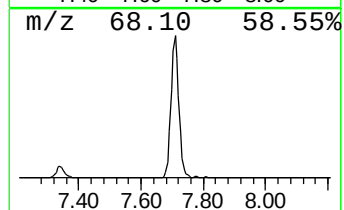
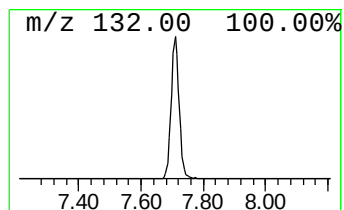
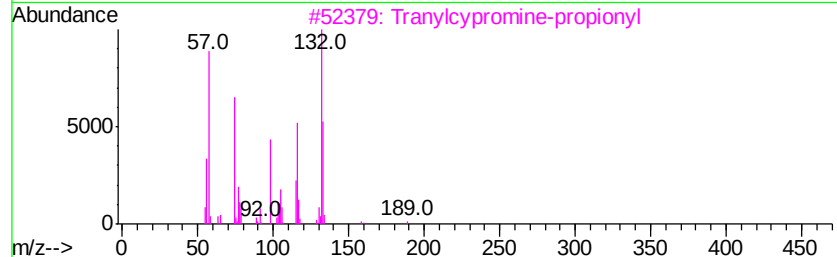
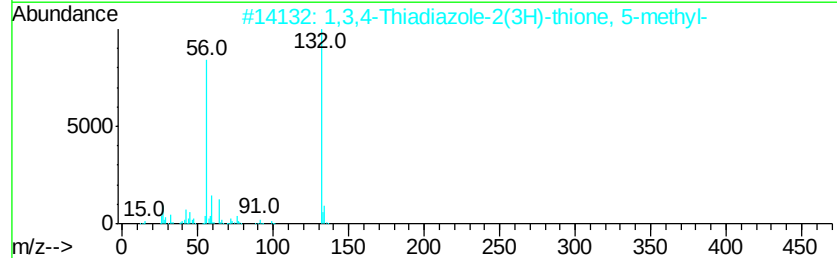
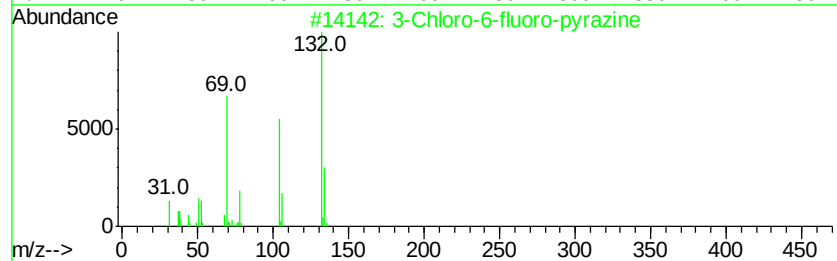
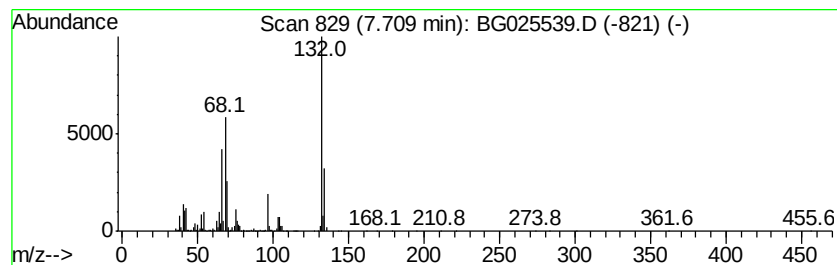
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Peak Number 2 unknown7.71 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.71	112.39 ng	2704370	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
2		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	45.0	ng	1081750	1	8.18	481248	20.0
unknown7.71	7.71	112.4	ng	2704370	1	8.18	481248	20.0