

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060716\
 Data File : BG022218.D
 Acq On : 8 Jun 2016 00:08
 Operator : UM/SJ
 Sample : H3399-23
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DUP-1

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-BG060616.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	5.664	474	481	497	rBV	113391	225671	10.50%	2.847%
2	7.274	749	755	772	rBV	75439	160348	7.46%	2.023%
3	7.650	809	819	836	rBV	250609	521376	24.25%	6.578%
4	8.120	892	899	912	rVB	54921	107478	5.00%	1.356%
5	8.443	946	954	963	rBV	247504	502058	23.35%	6.334%
6	9.289	1090	1098	1113	rBV	167591	371017	17.26%	4.681%
7	10.940	1371	1379	1392	rBV	82185	183004	8.51%	2.309%
8	13.367	1785	1792	1810	rBV	711931	1244050	57.86%	15.695%
9	14.747	2020	2027	2040	rBV2	182039	323102	15.03%	4.076%
10	16.234	2273	2280	2297	rBV	489258	819383	38.11%	10.337%
11	17.491	2487	2494	2506	rBV2	230421	413308	19.22%	5.214%
12	20.082	2929	2935	2948	rBV	1460803	2150106	100.00%	27.125%
13	21.763	3214	3221	3233	rBV	239675	460360	21.41%	5.808%
14	25.053	3770	3781	3794	rBV2	138530	445301	20.71%	5.618%

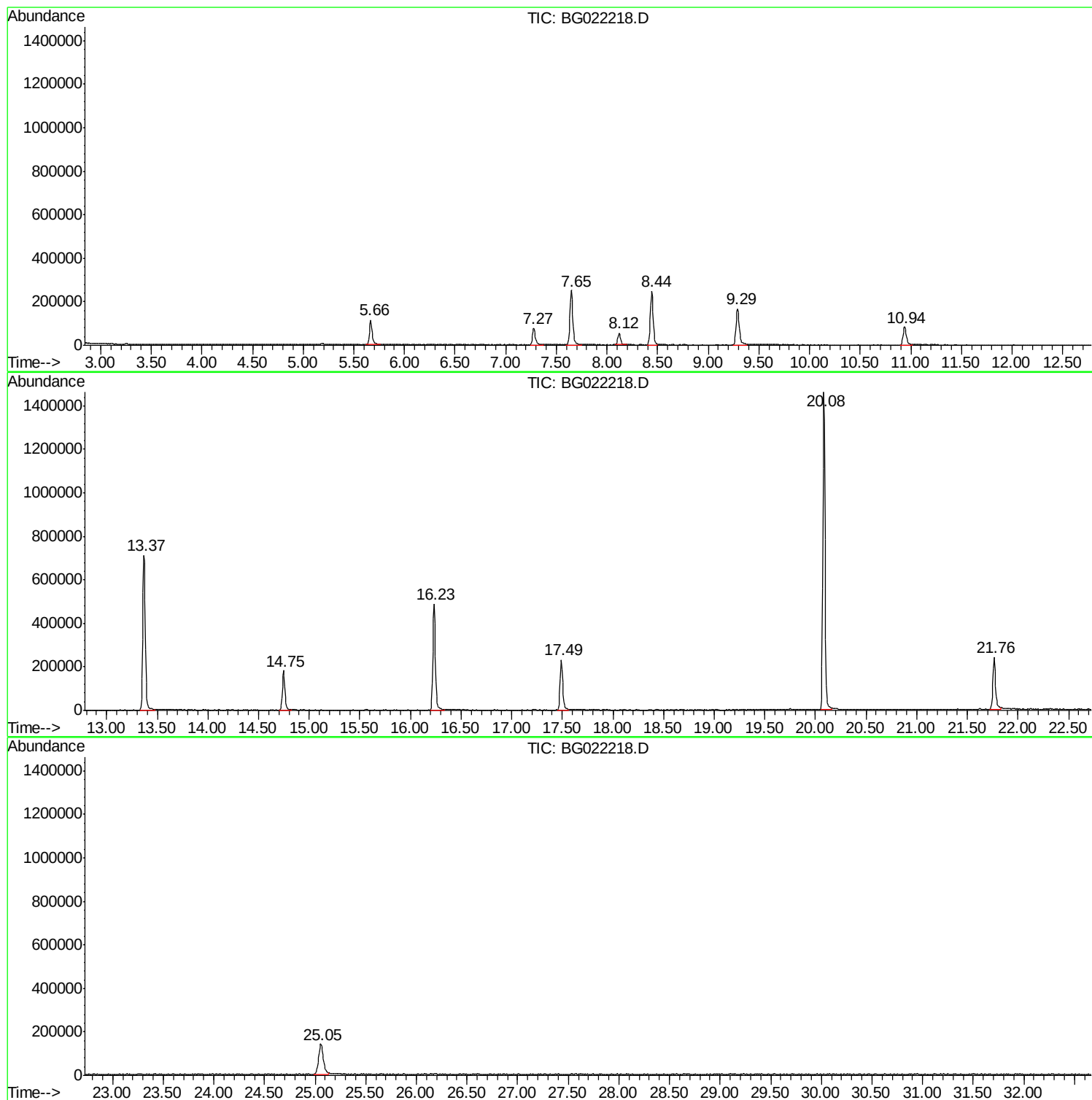
Sum of corrected areas: 7926562

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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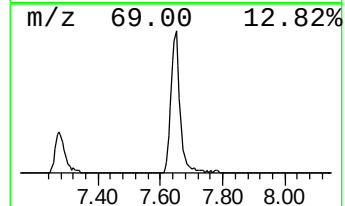
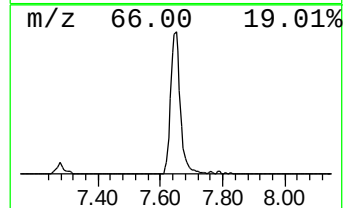
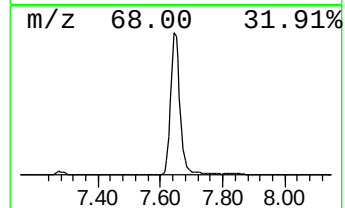
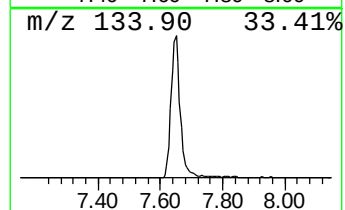
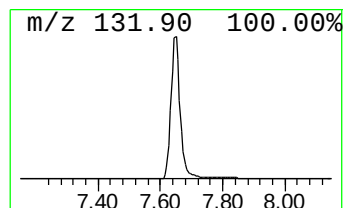
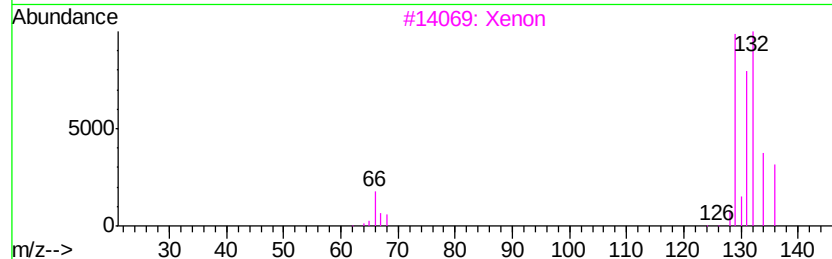
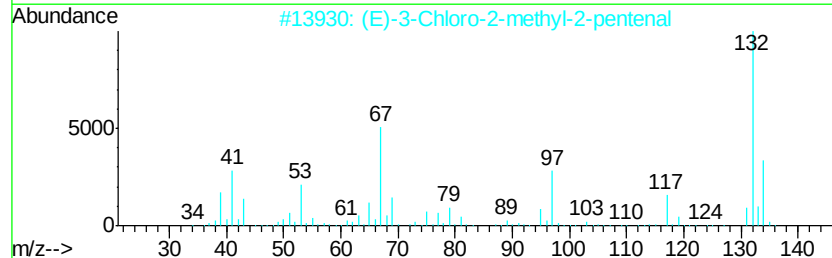
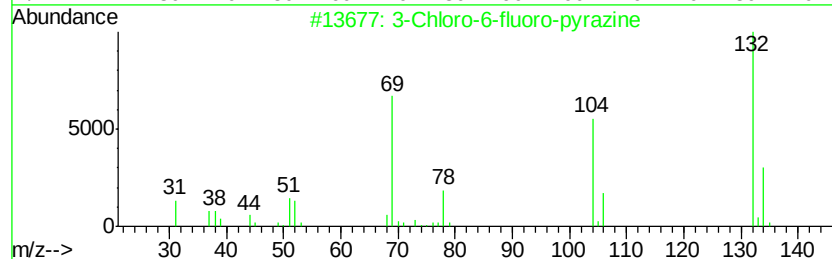
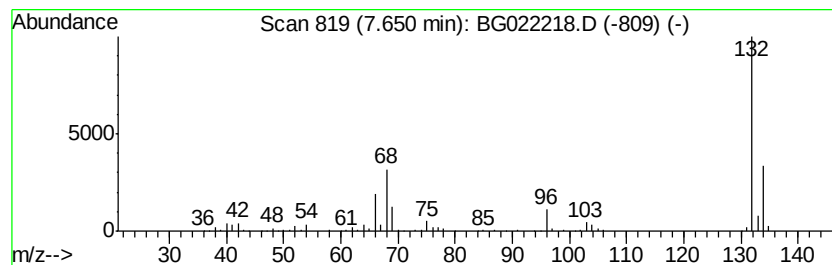
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Peak Number 1 unknown7.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	97.02 ng	521376	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	40
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	38
3		Xenon	132	Xe	007440-63-3	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.65	7.65	97.0	ng	521376	1	8.12	107478	20.0