

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
 Data File : BG022229.D
 Acq On : 8 Jun 2016 9:23
 Operator : UM/SJ
 Sample : H3425-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP16-LF2MW6

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	494	rBV	112563	221688	9.73%	2.727%
2	7.274	749	755	769	rBV	74637	151291	6.64%	1.861%
3	7.644	811	818	833	rBV	251805	517533	22.71%	6.367%
4	8.120	892	899	910	rBV	53614	106493	4.67%	1.310%
5	8.443	944	954	966	rBV	247164	501930	22.02%	6.175%
6	9.289	1090	1098	1110	rBV	169152	359694	15.78%	4.425%
7	10.935	1371	1378	1393	rBV	79567	183580	8.05%	2.259%
8	13.367	1785	1792	1807	rBV	725747	1239925	54.40%	15.254%
9	14.748	2018	2027	2040	rBV2	179258	325487	14.28%	4.004%
10	16.234	2273	2280	2296	rBV	516402	906667	39.78%	11.154%
11	17.492	2487	2494	2506	rBV2	231614	417811	18.33%	5.140%
12	20.083	2929	2935	2952	rBV	1511129	2279162	100.00%	28.039%
13	21.763	3214	3221	3236	rBV2	232857	467669	20.52%	5.754%
14	25.059	3768	3782	3796	rBV2	137310	449472	19.72%	5.530%

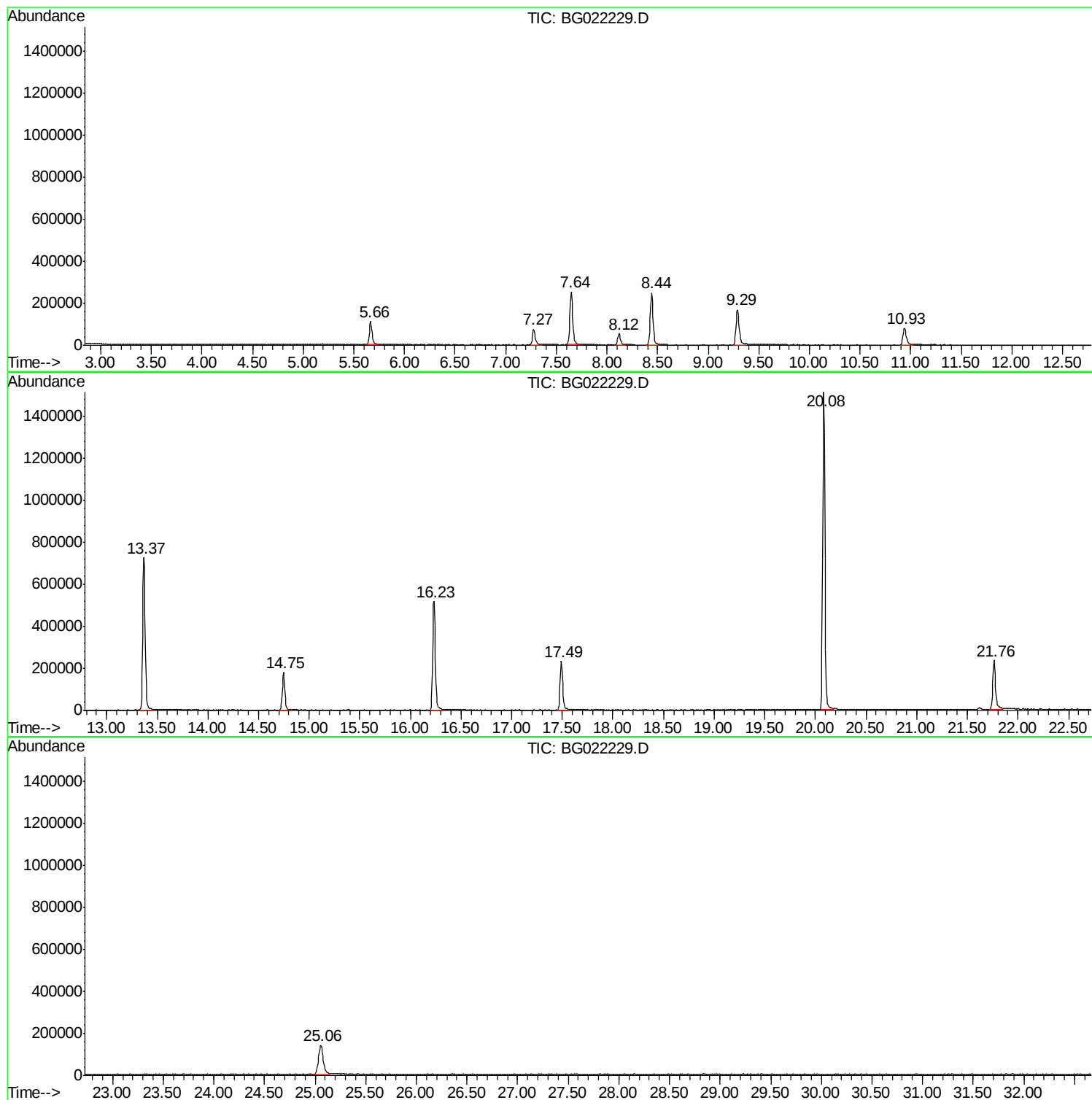
Sum of corrected areas: 8128402

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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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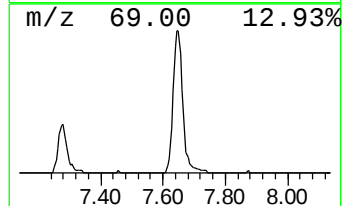
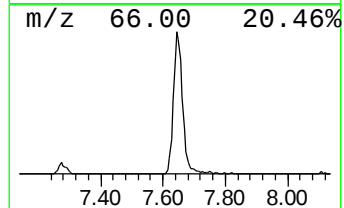
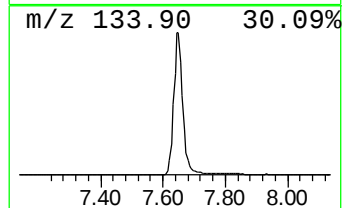
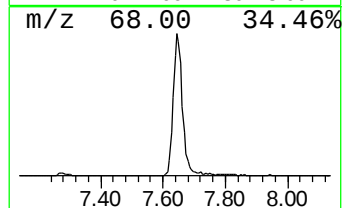
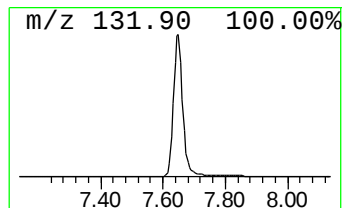
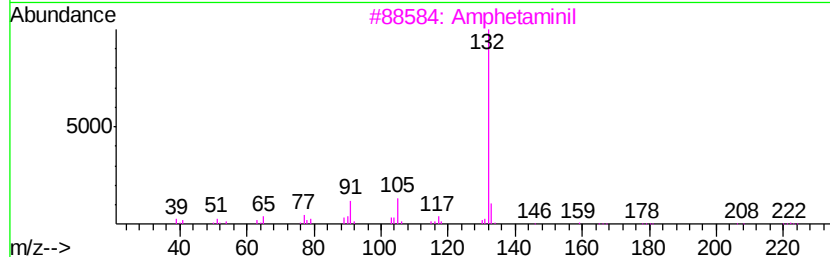
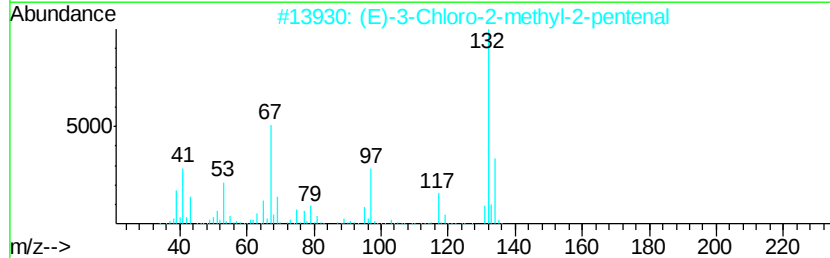
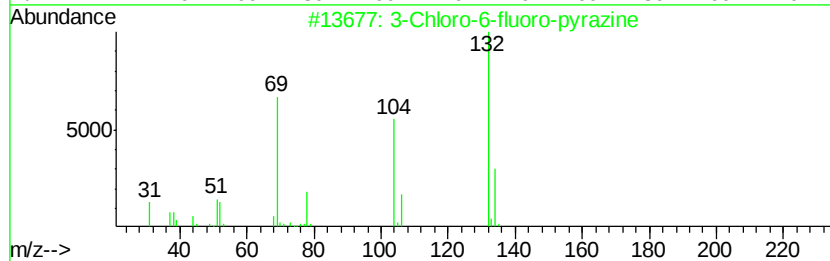
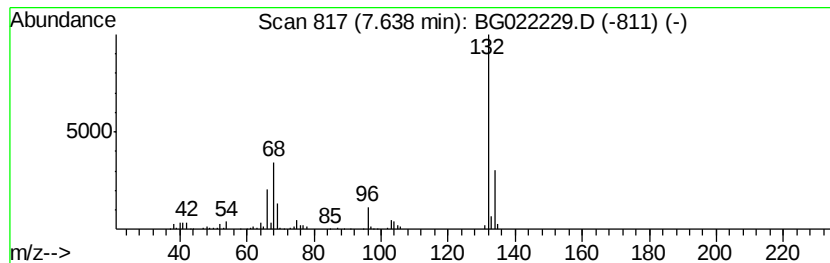
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Peak Number 1 unknown7.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	97.20 ng	517533	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	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.64	7.64	97.2	ng	517533	1	8.12	106493	20.0