

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121817\
 Data File : BG031632.D
 Acq On : 18 Dec 2017 13:25
 Operator : SJ/JU
 Sample : I6909-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 121217-DBRI-E-D-S

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-BG121117.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	5.656	397	403	416	rBV	303472	564333	8.85%	2.366%
2	7.266	671	677	689	rBV	178699	371013	5.82%	1.555%
3	7.630	732	739	758	rBV	646329	1188025	18.62%	4.981%
4	8.095	811	818	827	rBV	225051	407115	6.38%	1.707%
5	8.424	866	874	886	rBV	934633	1739767	27.27%	7.294%
6	9.270	1008	1018	1039	rBV	666162	1346174	21.10%	5.644%
7	10.915	1290	1298	1315	rBV	291716	631132	9.89%	2.646%
8	13.341	1704	1711	1726	rBV	2319522	3953703	61.97%	16.575%
9	14.722	1937	1946	1965	rBV2	568223	1025172	16.07%	4.298%
10	16.209	2192	2199	2226	rBV	1105421	1948288	30.54%	8.168%
11	17.466	2406	2413	2429	rBV2	835012	1371271	21.49%	5.749%
12	20.057	2848	2854	2867	rBV	4654098	6380116	100.00%	26.747%
13	21.737	3131	3140	3156	rBV	885882	1619028	25.38%	6.787%
14	25.010	3686	3697	3712	rBV2	430381	1308121	20.50%	5.484%

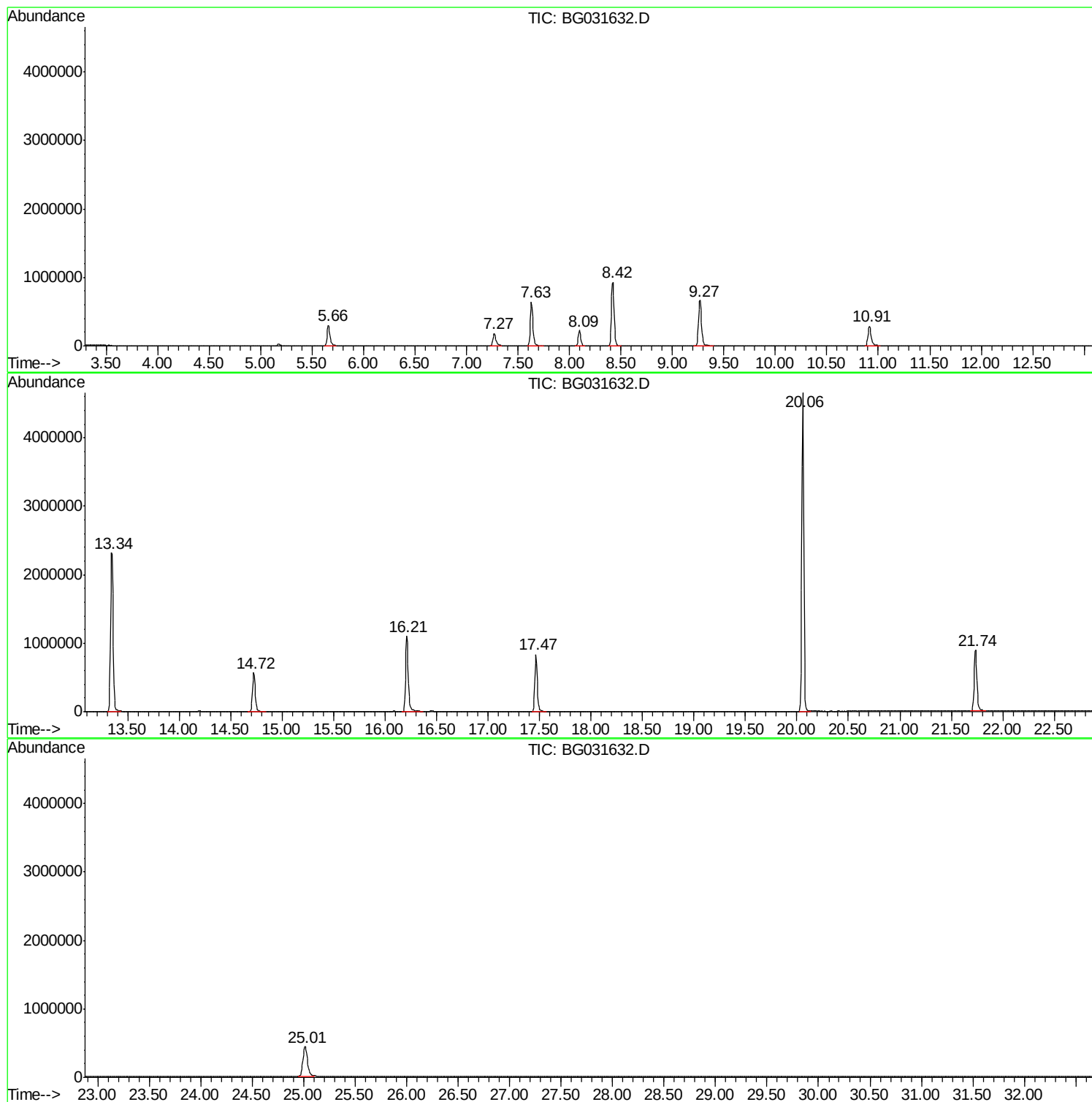
Sum of corrected areas: 23853258

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.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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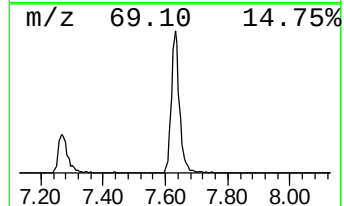
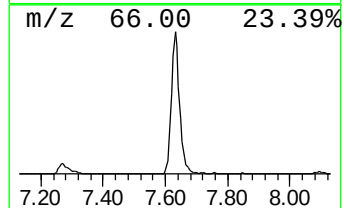
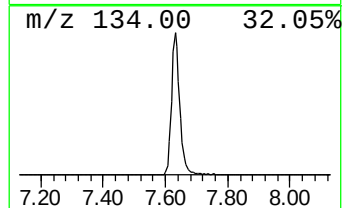
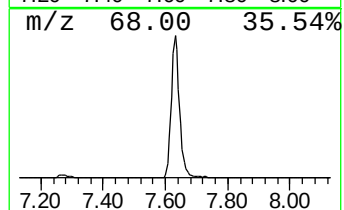
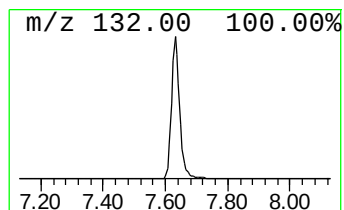
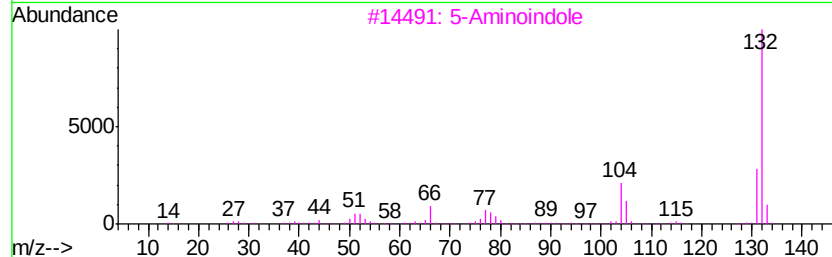
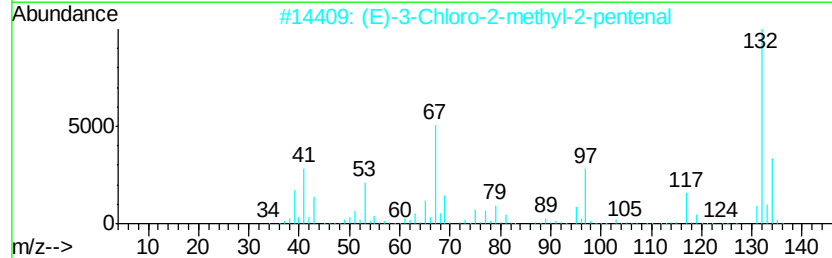
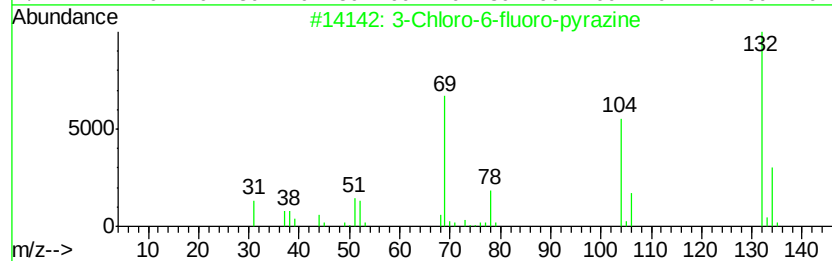
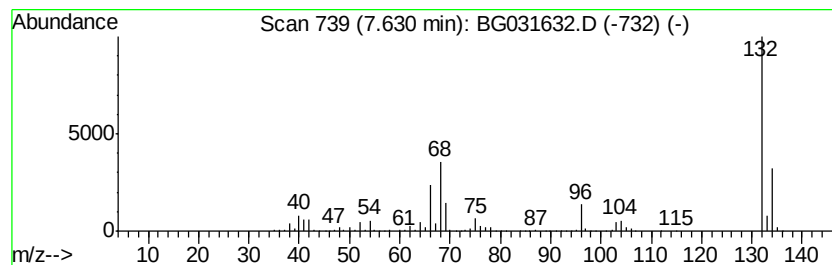
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Peak Number 1 unknown7.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	58.36 ng	1188030	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	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	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
unknown7.63	7.63	58.4	ng	1188030	1	8.10	407115	20.0