

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG030419\
 Data File : BG039737.D
 Acq On : 4 Mar 2019 18:50
 Operator : JU/SJ
 Sample : PB117513BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB117513BL

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-BG030419.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	3.211	14	21	28	rBV2	25365	48676	1.53%	0.327%
2	5.702	437	445	455	rBV	688627	1089871	34.26%	7.322%
3	7.300	710	717	726	rBV	707680	1209721	38.02%	8.127%
4	7.681	775	782	791	rBV	637219	1117590	35.13%	7.508%
5	8.157	856	863	870	rBV	138240	234431	7.37%	1.575%
6	8.480	910	918	927	rVB	475347	810123	25.46%	5.443%
7	9.332	1055	1063	1074	rBV	390650	695148	21.85%	4.670%
8	10.977	1336	1343	1351	rBV	192414	344268	10.82%	2.313%
9	13.403	1749	1756	1766	rVB	1175363	1790227	56.27%	12.027%
10	14.784	1984	1991	2000	rVB2	320669	518062	16.28%	3.480%
11	16.270	2237	2244	2259	rBV	890178	1354461	42.57%	9.100%
12	17.527	2450	2458	2470	rVB	484079	723958	22.75%	4.864%
13	20.123	2893	2900	2908	rBV	2344215	3181543	100.00%	21.374%
14	21.815	3181	3188	3196	rBV2	537887	905365	28.46%	6.082%
15	25.187	3748	3762	3775	rBV3	285080	861515	27.08%	5.788%

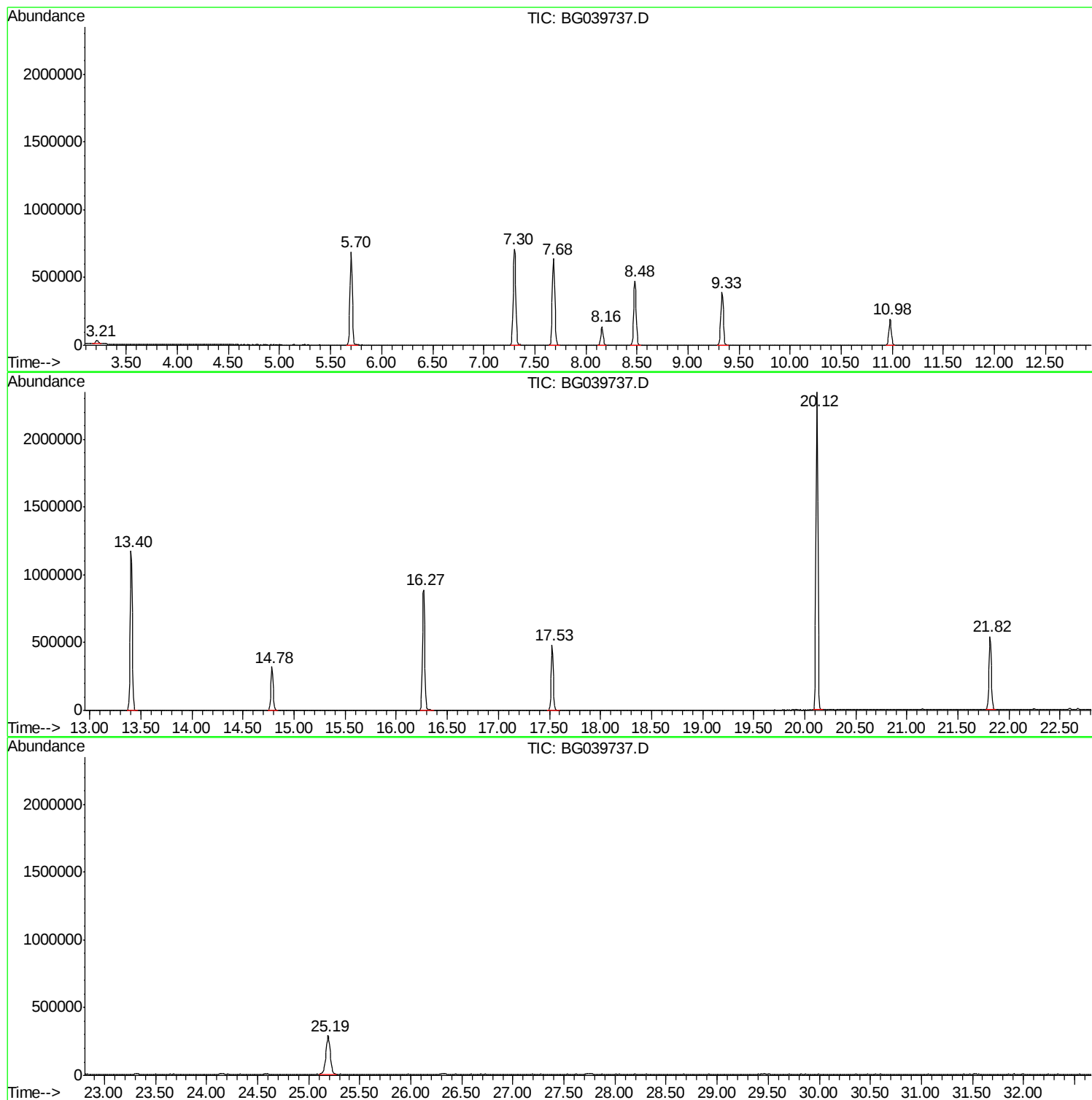
Sum of corrected areas: 14884959

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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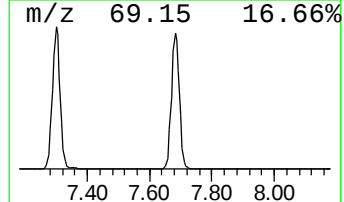
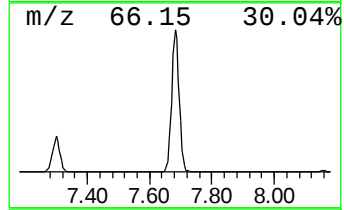
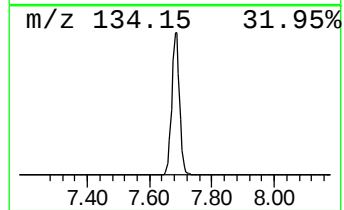
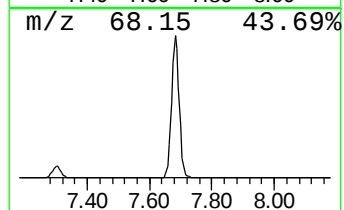
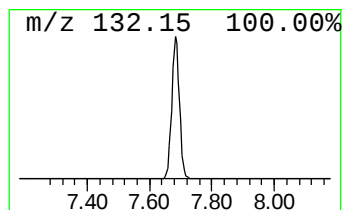
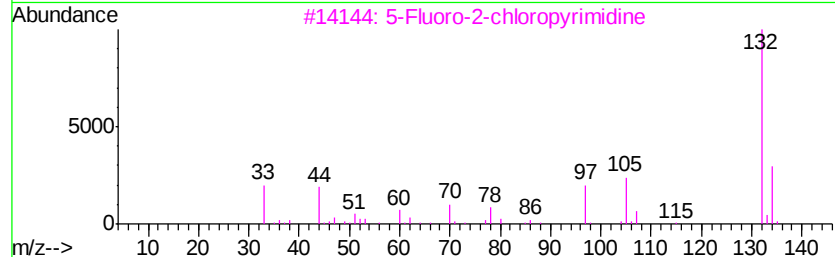
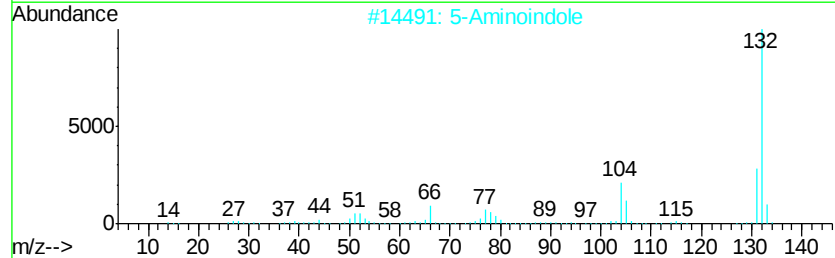
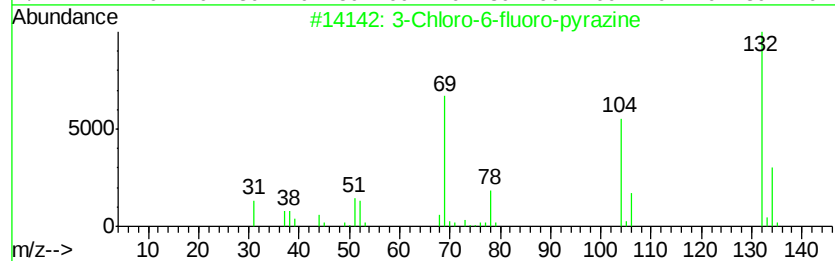
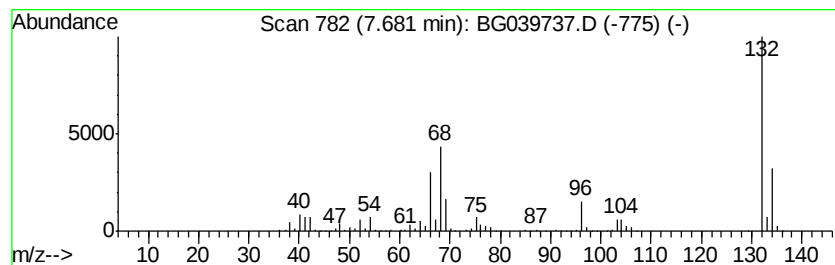
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	95.34 ng	1117590	1,4-Dichlorobenzene-d4	8.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			5-Aminoindole	132	C8H8N2	005192-03-0	22
3			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5			Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.68	7.68	95.3	ng	1117590	1	8.16	234431	20.0