

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050319\
 Data File : BG040727.D
 Acq On : 3 May 2019 16:36
 Operator : JU/SJ
 Sample : PB119443BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB119443BL

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-BG042319.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	3.153	6	11	21	rVB3	34955	67371	1.60%	0.379%
2	5.674	433	440	452	rBV	808623	1252323	29.83%	7.045%
3	7.284	706	714	723	rBV	852436	1446610	34.45%	8.138%
4	7.671	772	780	788	rBV	764863	1328627	31.64%	7.474%
5	8.147	855	861	868	rVB	143991	237939	5.67%	1.339%
6	8.470	907	916	926	rBV	542177	930318	22.16%	5.234%
7	9.322	1053	1061	1071	rBV	512977	887136	21.13%	4.991%
8	10.968	1334	1341	1349	rVB	178787	322342	7.68%	1.813%
9	13.400	1748	1755	1762	rBV	1336805	2051724	48.87%	11.542%
10	14.775	1983	1989	1998	rBV2	332807	519341	12.37%	2.922%
11	16.255	2234	2241	2255	rBV	1247671	2031319	48.38%	11.427%
12	17.519	2449	2456	2467	rVB2	536620	799770	19.05%	4.499%
13	20.110	2891	2897	2908	rBV	3057609	4198740	100.00%	23.620%
14	21.802	3178	3185	3200	rBV	555987	944551	22.50%	5.314%
15	25.163	3746	3757	3767	rBV2	239731	757947	18.05%	4.264%

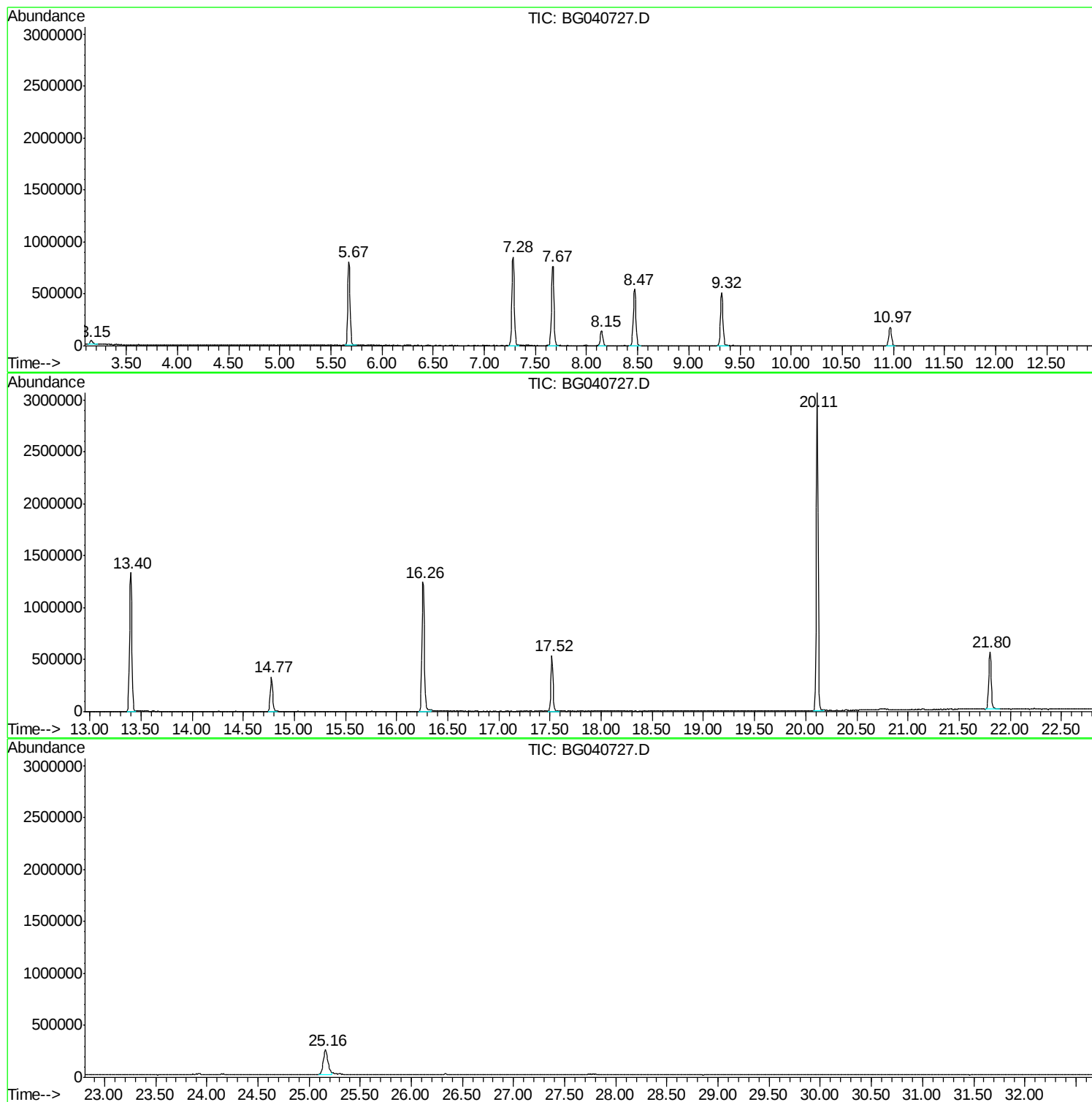
Sum of corrected areas: 17776058

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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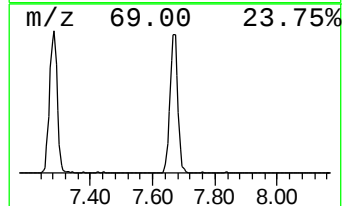
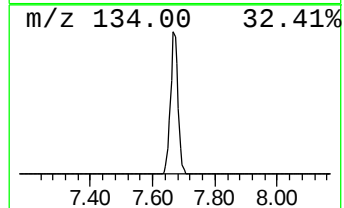
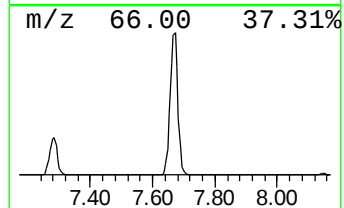
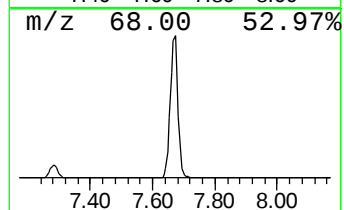
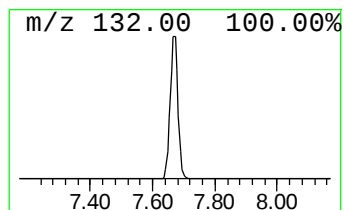
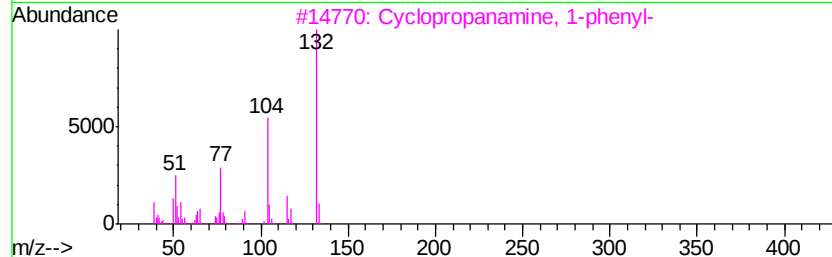
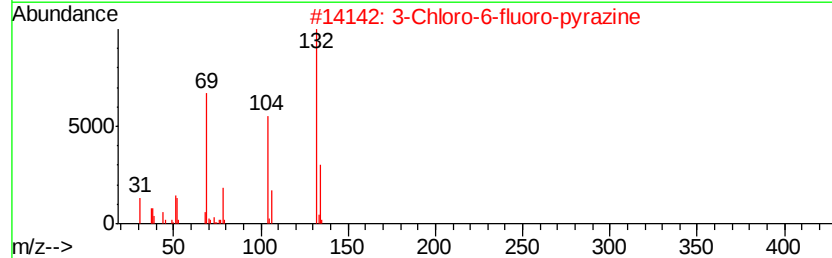
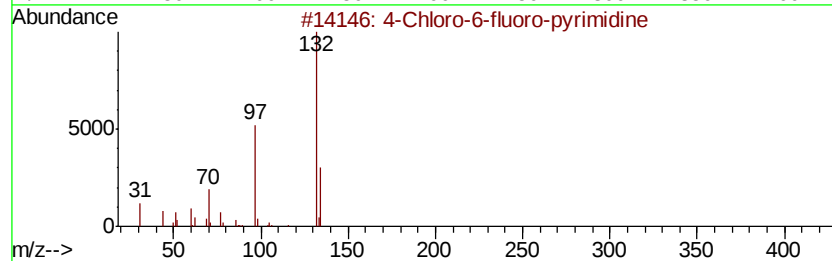
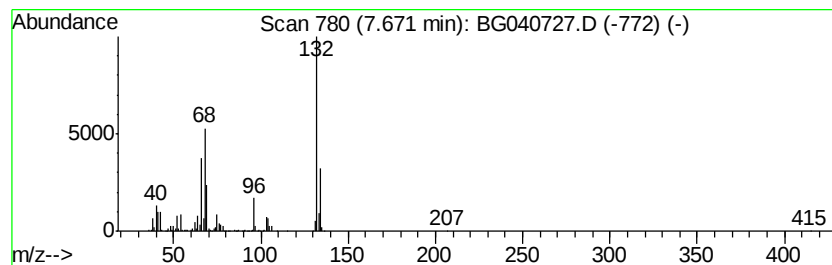
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Peak Number 2 unknown7.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	111.68 ng	1328630	1,4-Dichlorobenzene-d4	8.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3			Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
4			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5			Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.67	7.67	111.7	ng	1328630	1	8.15	237939	20.0