

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM042319\
 Data File : BM020058.D
 Acq On : 24 Apr 2019 05:28
 Operator : JU/SJ
 Sample : PB119111TB
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB119111TB

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.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	5.104	355	360	387	rBV	1413859	2426053	56.18%	9.822%
2	6.687	623	629	652	rBV	1257561	2393648	55.43%	9.691%
3	7.028	680	687	713	rBV	1624932	2715865	62.89%	10.996%
4	7.493	760	766	780	rVB	297662	495138	11.47%	2.005%
5	7.798	811	818	837	rBV	1203485	2037701	47.19%	8.250%
6	8.669	960	966	992	rBV	548029	1354216	31.36%	5.483%
7	10.281	1230	1240	1264	rBV	188568	549117	12.72%	2.223%
8	12.763	1656	1662	1686	rBV	1634826	3055897	70.77%	12.372%
9	14.169	1895	1901	1922	rBV2	304362	648733	15.02%	2.626%
10	15.680	2151	2158	2186	rBV	949932	2102892	48.70%	8.514%
11	16.939	2367	2372	2397	rBV	280252	693036	16.05%	2.806%
12	19.574	2814	2820	2833	rBV	3099169	4318358	100.00%	17.484%
13	21.156	3084	3089	3107	rBV	482610	960261	22.24%	3.888%
14	23.333	3453	3459	3473	rBV	397287	948690	21.97%	3.841%

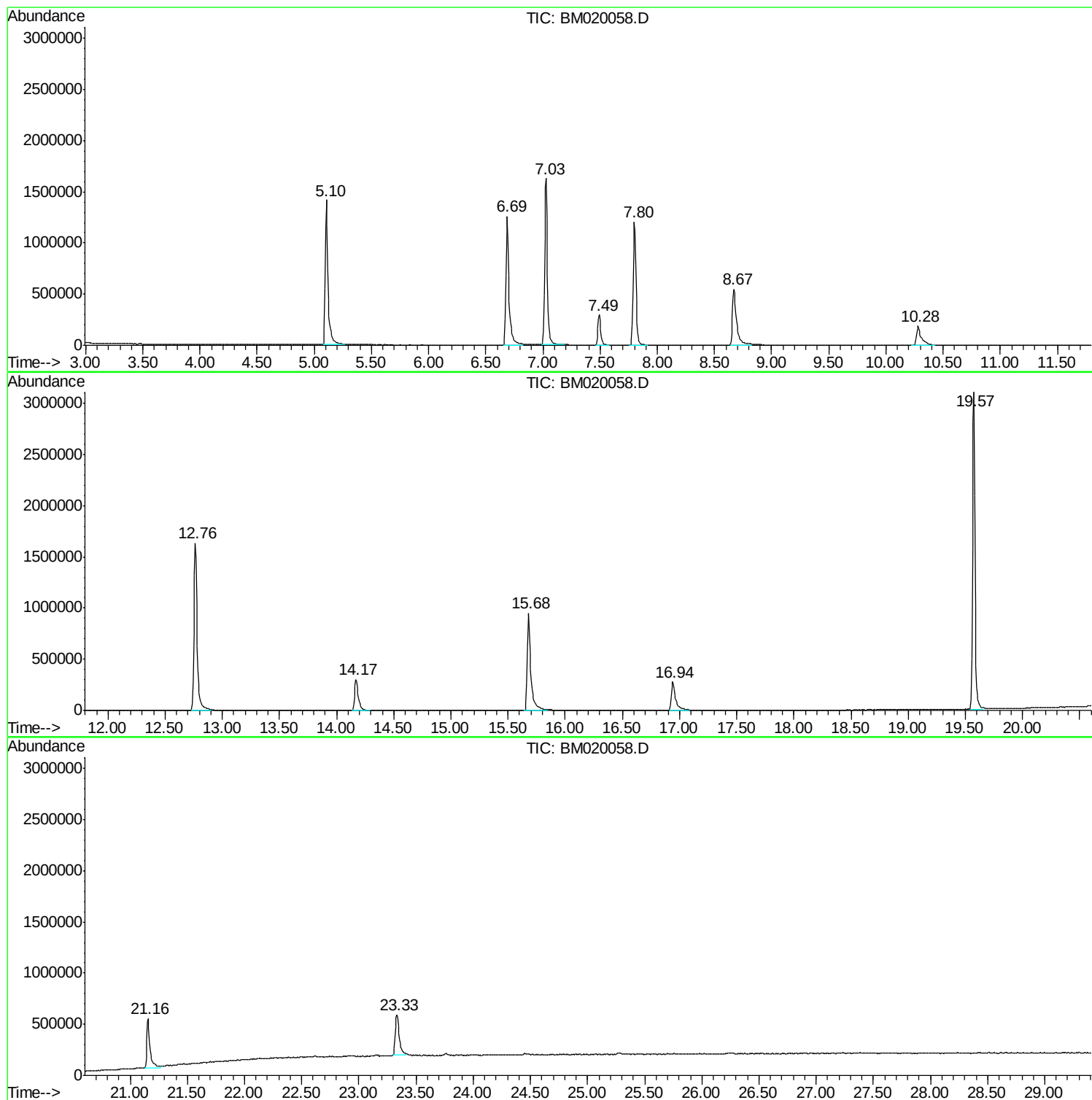
Sum of corrected areas: 24699605

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM041519.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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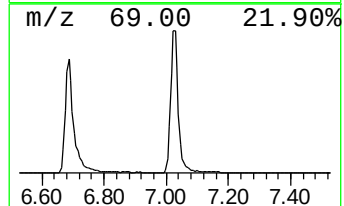
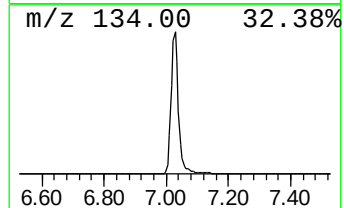
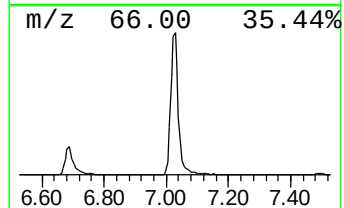
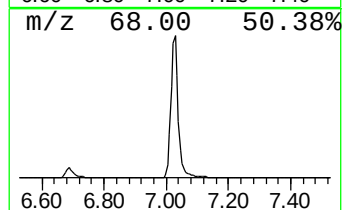
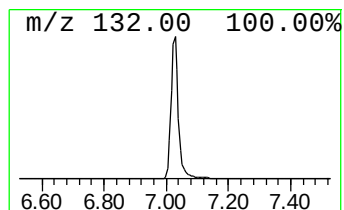
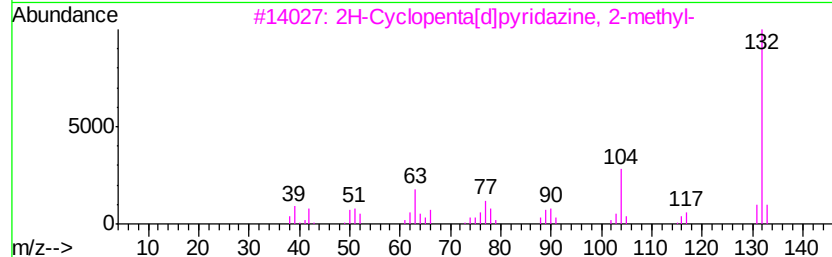
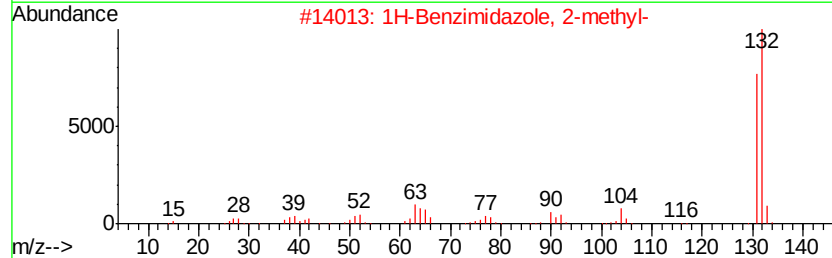
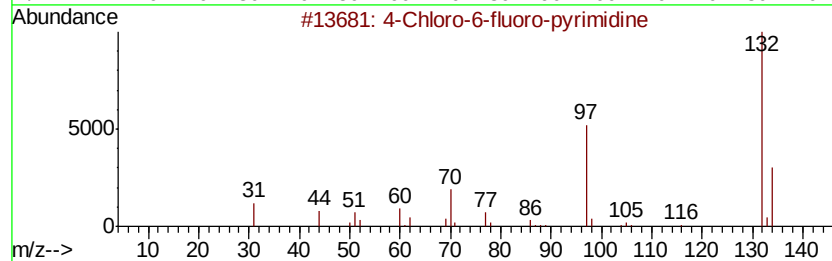
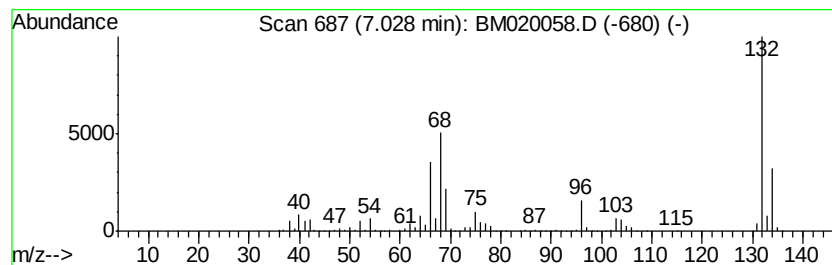
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.03	109.70 ng	2715870	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		2H-Cyclopentaf[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.03	7.03	109.7	ng	2715870	1	7.49	495138	20.0