

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073018\
 Data File : BG035974.D
 Acq On : 30 Jul 2018 13:44
 Operator : JU/SJ
 Sample : J4196-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 072518-DBRI-E-D-B

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_G\METHODS\8270-BG071218.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	5.606	387	394	407	rBV	184337	330933	9.11%	2.299%
2	7.192	655	664	675	rBV	132615	261922	7.21%	1.820%
3	7.556	719	726	739	rBV	385072	687939	18.93%	4.780%
4	8.020	799	805	814	rBV	140535	245848	6.76%	1.708%
5	8.343	851	860	872	rBV	553326	988871	27.21%	6.871%
6	9.183	993	1003	1014	rBV	455635	865966	23.83%	6.017%
7	10.828	1274	1283	1297	rBV	160782	335106	9.22%	2.328%
8	13.272	1691	1699	1710	rBV	1413446	2260985	62.21%	15.710%
9	14.100	1835	1840	1849	rBV	43614	87622	2.41%	0.609%
10	14.646	1926	1933	1944	rBV2	329166	574151	15.80%	3.989%
11	15.933	2147	2152	2157	rBV2	64206	88253	2.43%	0.613%
12	16.133	2178	2186	2202	rBV	694289	1183570	32.56%	8.224%
13	17.390	2393	2400	2408	rBV2	481744	807656	22.22%	5.612%
14	19.998	2838	2844	2857	rBV	2826429	3634487	100.00%	25.253%
15	21.302	3061	3066	3074	rBV6	30738	79258	2.18%	0.551%
16	21.655	3119	3126	3138	rBV	547200	942653	25.94%	6.550%
17	23.123	3370	3376	3381	rBV7	24952	49969	1.37%	0.347%
18	23.904	3506	3509	3516	rVB8	18680	40117	1.10%	0.279%
19	24.850	3659	3670	3681	rBV	303814	886784	24.40%	6.162%
20	25.960	3852	3859	3861	rBV8	18905	39930	1.10%	0.277%

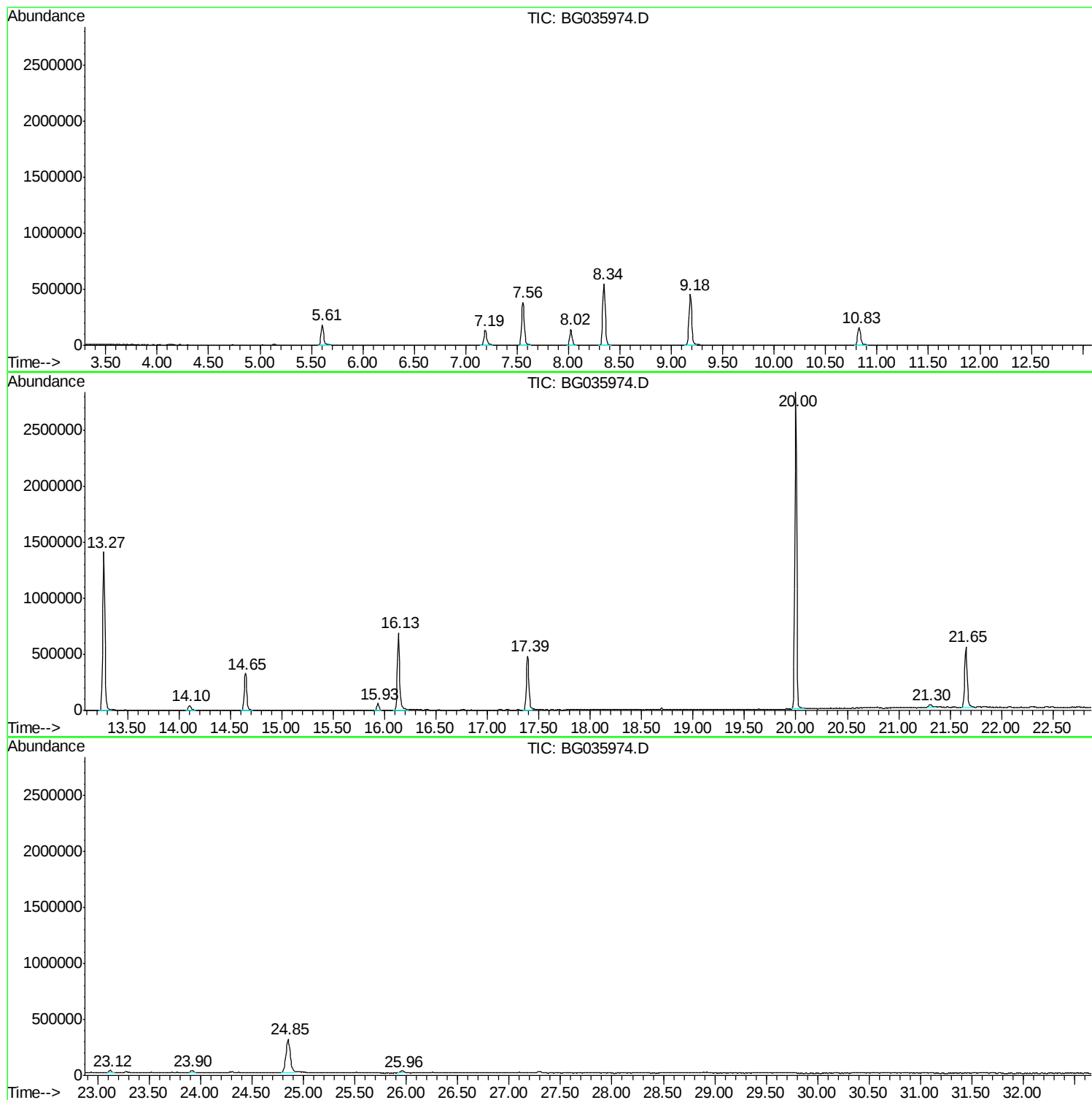
Sum of corrected areas: 14392020

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.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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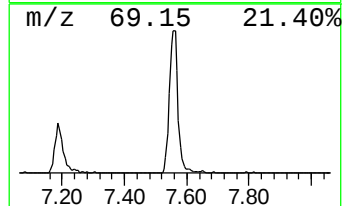
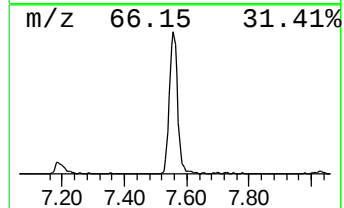
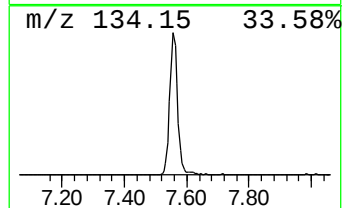
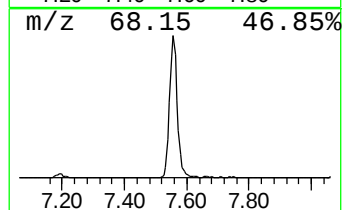
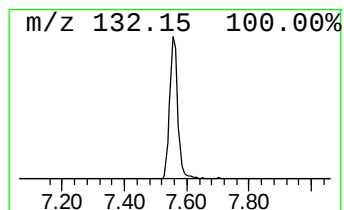
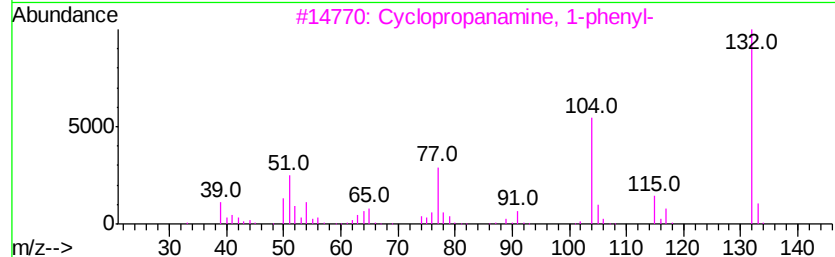
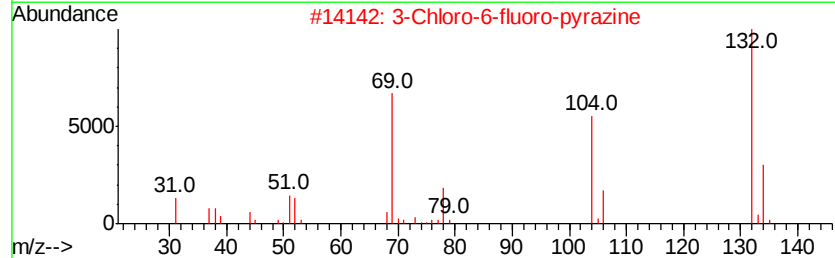
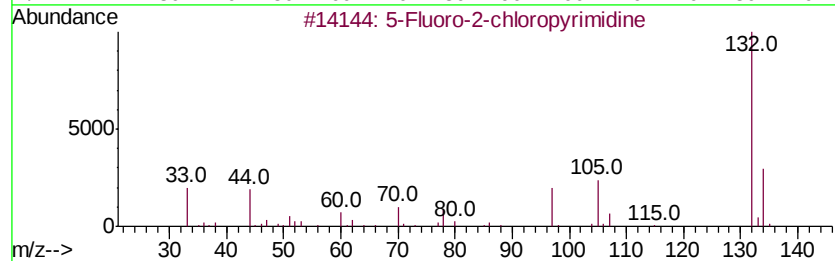
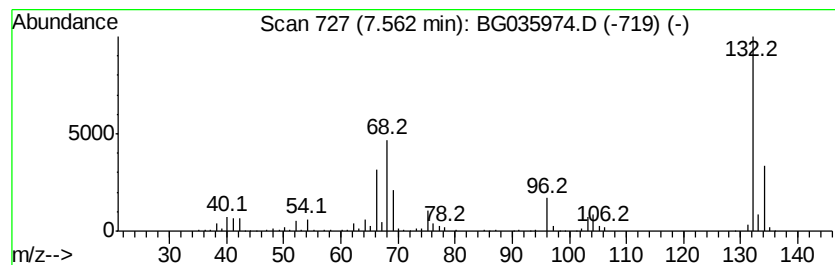
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
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Peak Number 1 unknown7.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	55.96 ng	687939	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		5-Aminoindole	132	C8H8N2	005192-03-0	10



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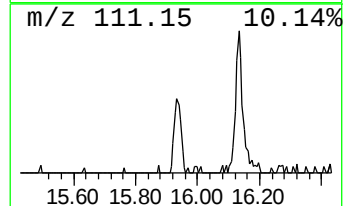
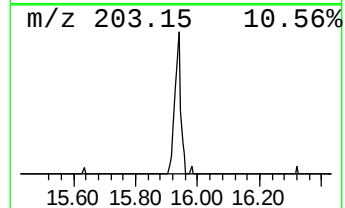
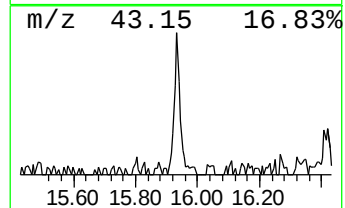
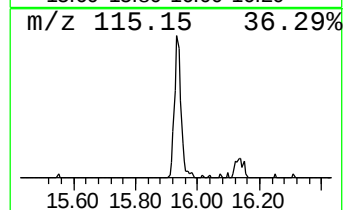
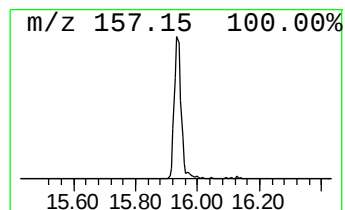
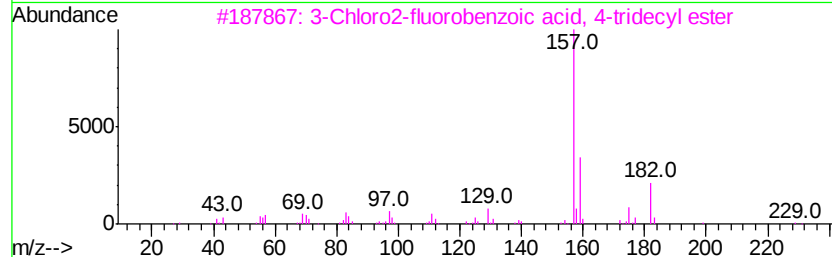
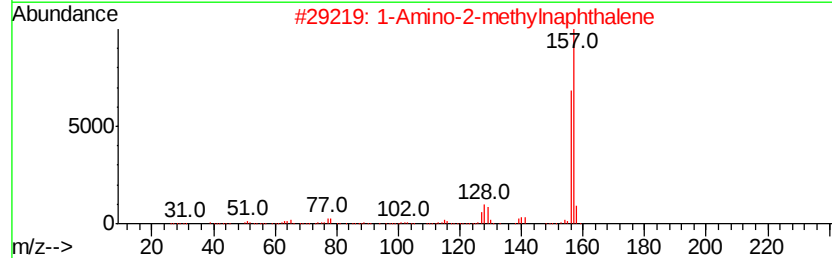
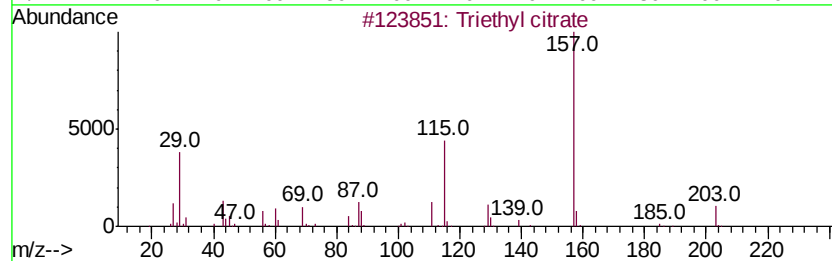
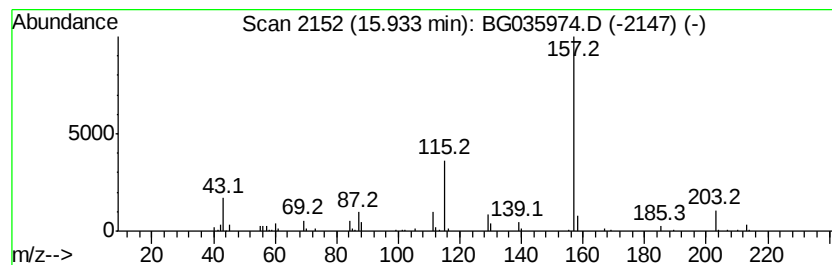
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Peak Number 3 Triethyl citrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	3.07 ng	88253	Acenaphthene-d10	14.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triethyl citrate	276	C12H20O7	000077-93-0	83
2		1-Amino-2-methylnaphthalene	157	C11H11N	002246-44-8	50
3		3-Chloro2-fluorobenzoic acid, 4-...	356	C20H30ClF02	1000338-64-4	40
4		Adipic acid, isobutyl 3-(2-metho...	358	C20H38O5	1000324-28-7	36
5		3-Chloro2-fluorobenzoic acid, 4-...	384	C22H34ClF02	1000338-65-1	33



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.56	7.56	56.0	ng	687939	1	8.02	245848	20.0
Triethyl citrate	15.93	3.1	ng	88253	3	14.65	574151	20.0