

Data Path : Z:\HPCHEM1\BNA G\DATA\BG052918\
 Data File : BG034779.D
 Acq On : 29 May 2018 16:23
 Operator : SJ/JU
 Sample : PB109739BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

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-BG051618.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

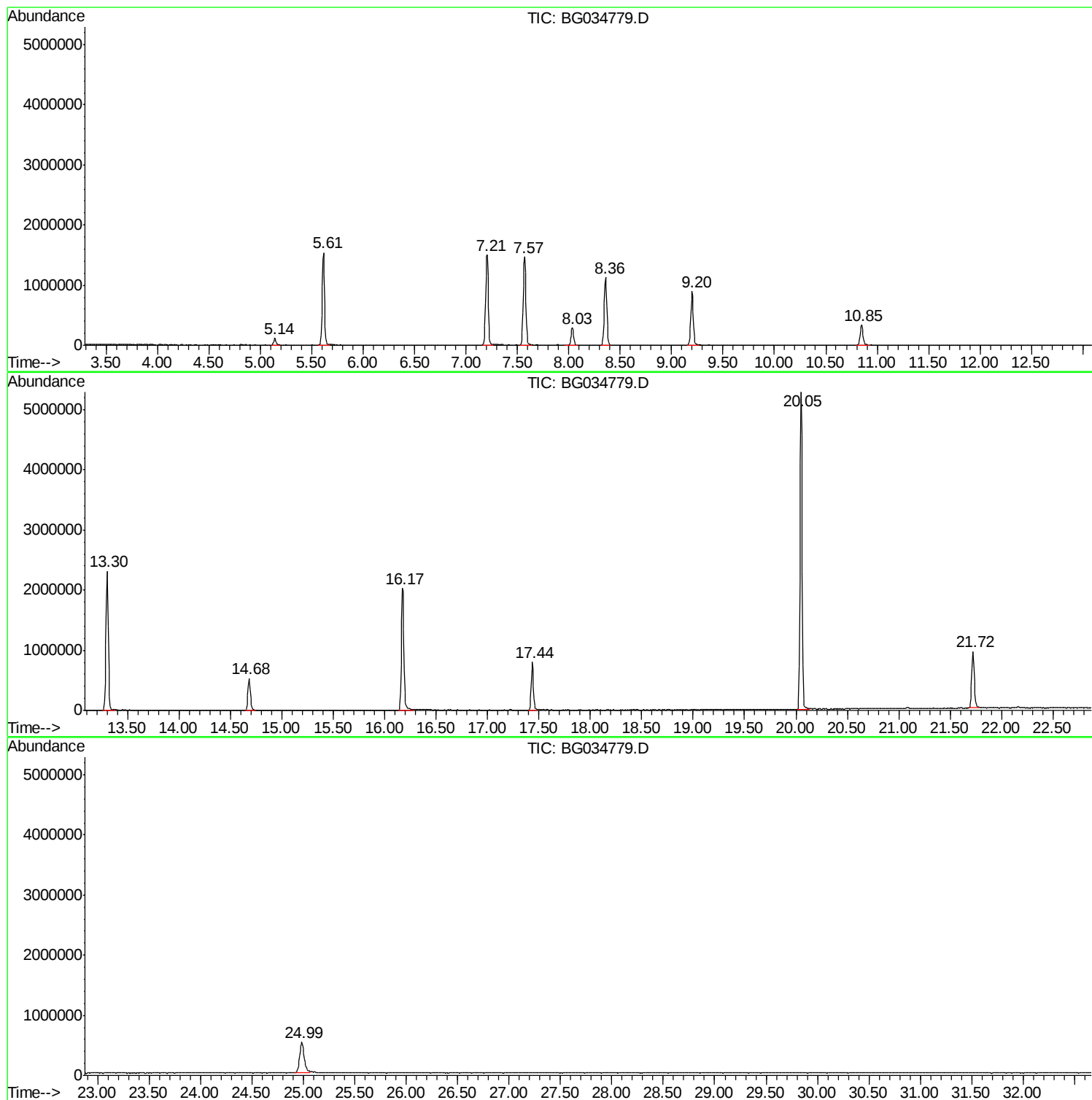
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.138	310	315	322	rVB	116919	175812	2.44%	0.547%
2	5.614	388	396	406	rBV	1539258	2480064	34.44%	7.710%
3	7.206	659	667	677	rBV	1499892	2683526	37.27%	8.343%
4	7.571	721	729	740	rBV	1464747	2521988	35.02%	7.841%
5	8.035	796	808	816	rBV	287020	501865	6.97%	1.560%
6	8.358	855	863	870	rBV	1127128	1947101	27.04%	6.053%
7	9.198	998	1006	1017	rBV	891752	1586200	22.03%	4.931%
8	10.849	1278	1287	1300	rBV	331920	623349	8.66%	1.938%
9	13.299	1697	1704	1717	rBV	2311982	3696319	51.33%	11.491%
10	14.680	1933	1939	1948	rBV2	526774	882699	12.26%	2.744%
11	16.172	2186	2193	2211	rBV	2023578	3314152	46.02%	10.303%
12	17.436	2401	2408	2416	rBV2	794885	1261420	17.52%	3.922%
13	20.050	2847	2853	2867	rBV	5278147	7201038	100.00%	22.387%
14	21.725	3131	3138	3146	rBV	942786	1649786	22.91%	5.129%
15	24.986	3681	3693	3706	rBV2	512872	1640524	22.78%	5.100%

Sum of corrected areas: 32165843

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG051618.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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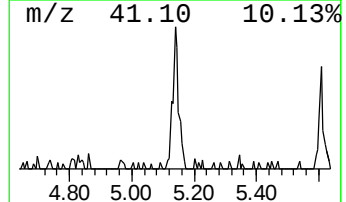
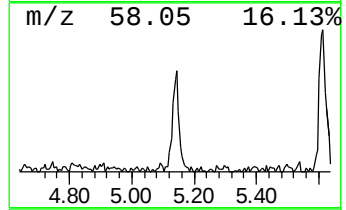
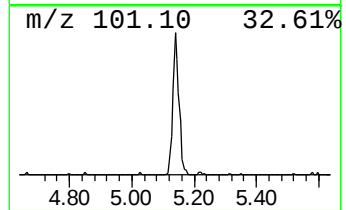
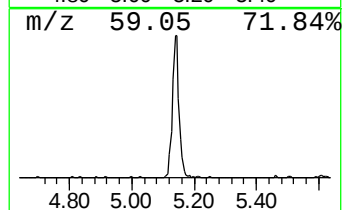
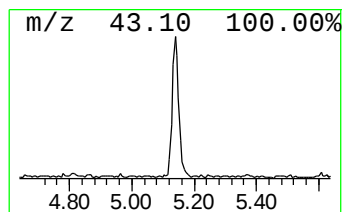
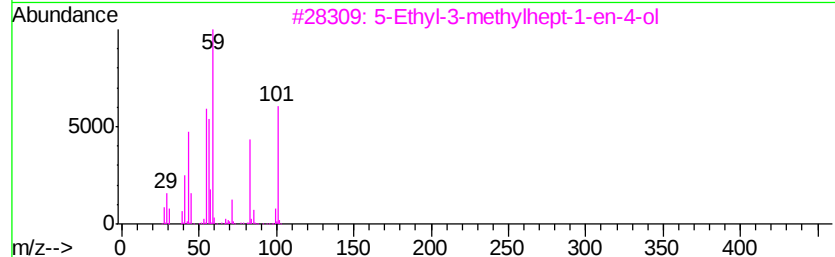
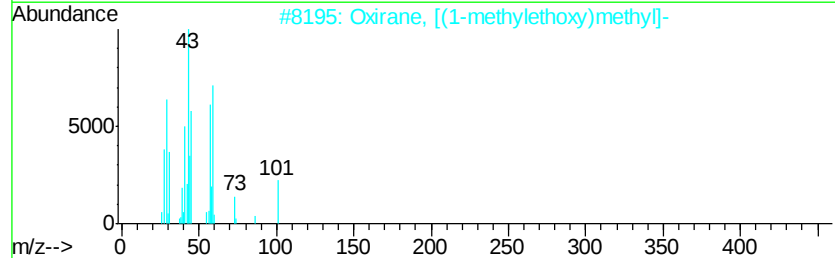
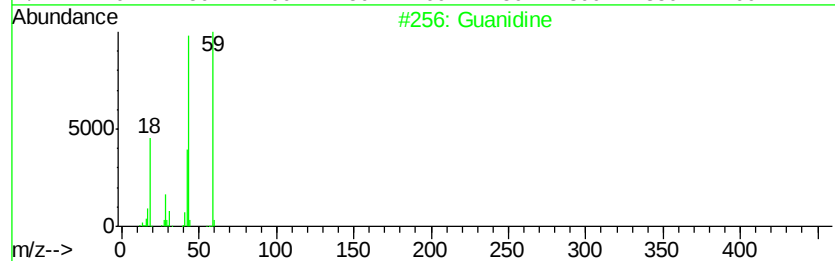
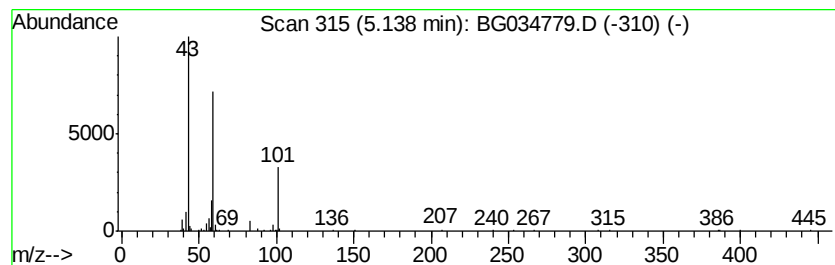
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Peak Number 1 unknown5.14 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	7.01 ng	175812	1,4-Dichlorobenzene-d4	8.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Guanidine	59	CH5N3	000113-00-8	43
2			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	38
3			5-Ethyl-3-methylhept-1-en-4-ol	156	C10H20O	286424-80-4	23
4			Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	17
5			2,3-Dimethyloxirane-2-carboxylic...	130	C6H10O3	1000306-64-4	17



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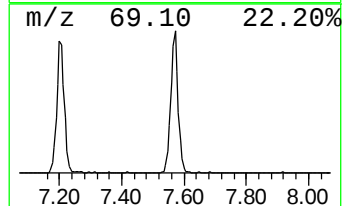
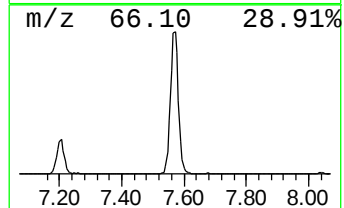
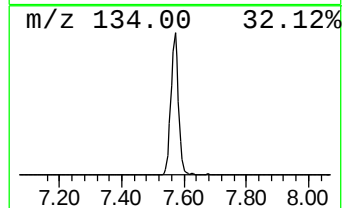
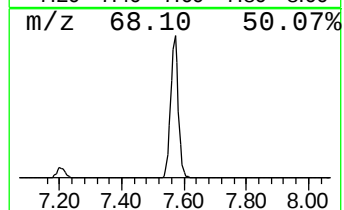
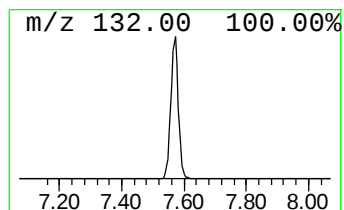
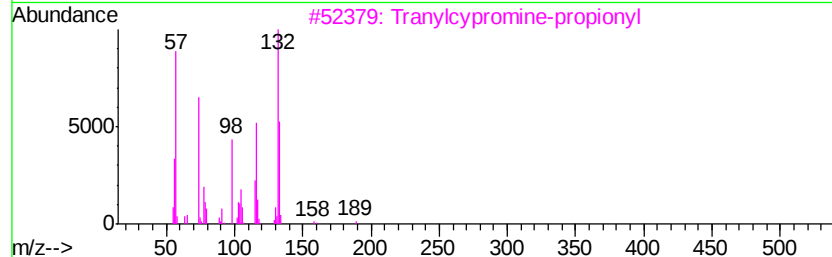
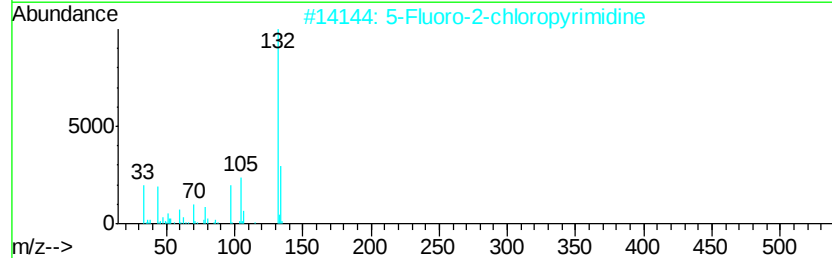
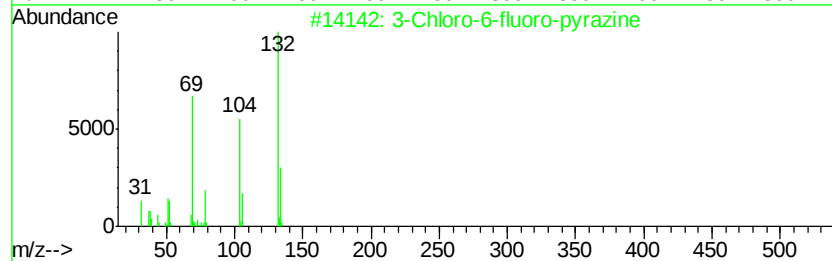
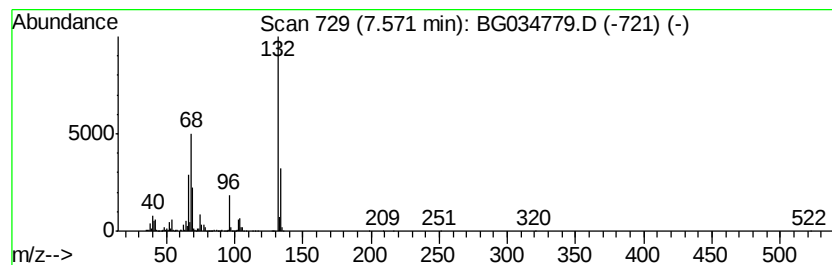
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	100.51 ng	2521990	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.14	5.14	7.0	ng	175812	1	8.03	501865	20.0
unknown7.57	7.57	100.5	ng	2521990	1	8.03	501865	20.0