

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072718\
 Data File : BG035940.D
 Acq On : 27 Jul 2018 16:44
 Operator : JU/SJ
 Sample : J4192-03
 Misc :
 ALS Vial : 12 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:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

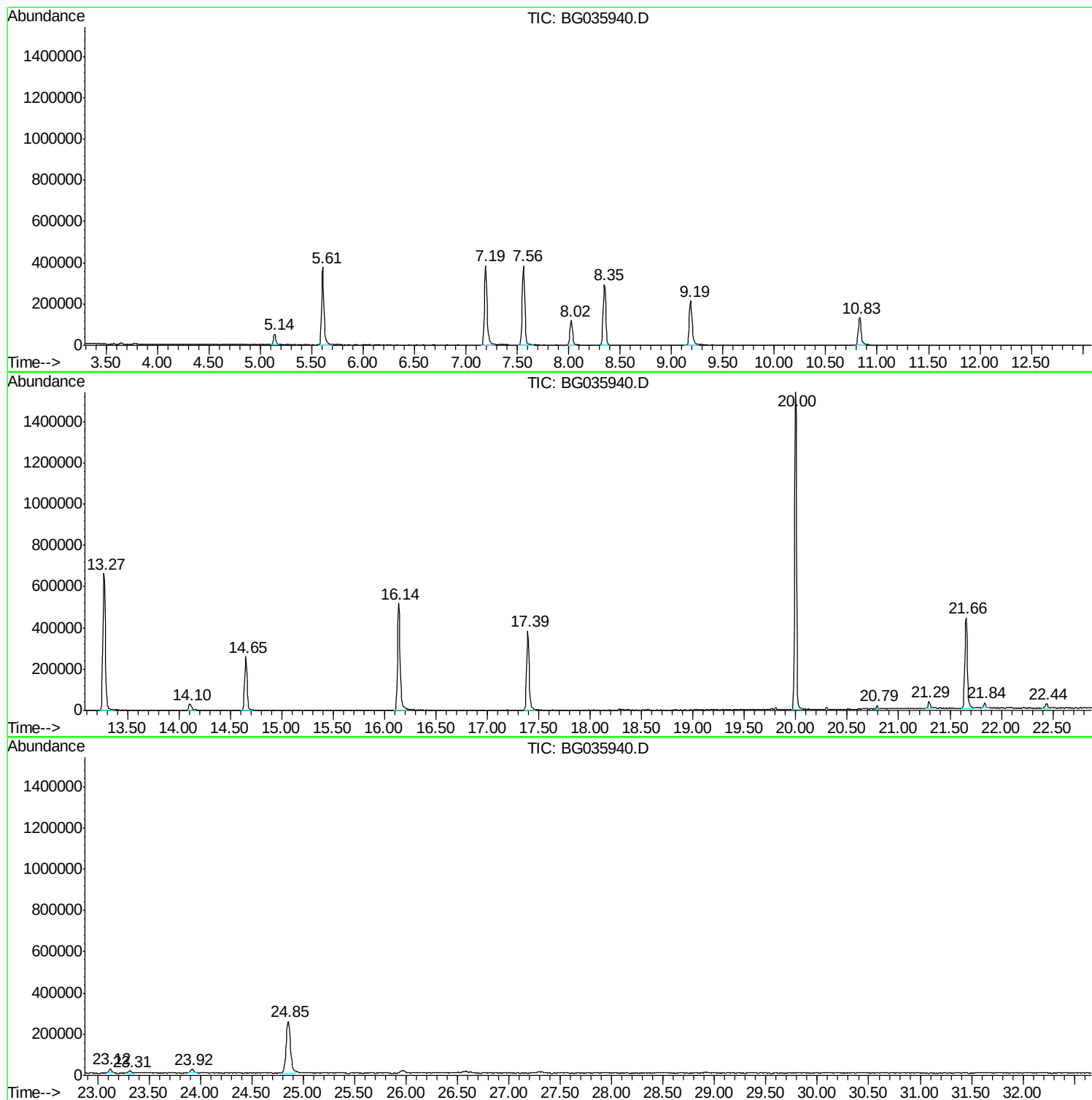
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.139	309	315	321	rBV2	53974	83754	4.09%	0.791%
2	5.609	388	395	411	rBV	377050	646190	31.53%	6.100%
3	7.195	656	665	679	rBV	384657	741014	36.16%	6.995%
4	7.559	719	727	743	rBV	385250	686635	33.50%	6.481%
5	8.023	800	806	814	rBV	120444	209465	10.22%	1.977%
6	8.346	854	861	869	rVB2	293696	529289	25.82%	4.996%
7	9.186	994	1004	1017	rBV	215237	434671	21.21%	4.103%
8	10.831	1276	1284	1294	rBV	134644	269313	13.14%	2.542%
9	13.269	1692	1699	1715	rBV	664434	1116756	54.49%	10.541%
10	14.103	1837	1841	1850	rBV2	29460	61898	3.02%	0.584%
11	14.650	1926	1934	1944	rBV2	262060	440359	21.49%	4.157%
12	16.136	2180	2187	2201	rBV2	519526	894003	43.62%	8.439%
13	17.393	2394	2401	2413	rBV	384411	643813	31.41%	6.077%
14	19.995	2838	2844	2854	rBV	1539958	2049525	100.00%	19.346%
15	20.788	2975	2979	2983	rBV3	18459	24339	1.19%	0.230%
16	21.293	3059	3065	3070	rBV3	36545	60103	2.93%	0.567%
17	21.658	3119	3127	3140	rBV	440460	773969	37.76%	7.306%
18	21.840	3153	3158	3164	rVB2	24861	34766	1.70%	0.328%
19	22.439	3253	3260	3265	rBV4	23857	42084	2.05%	0.397%
20	23.120	3371	3376	3381	rBV5	20886	34233	1.67%	0.323%
21	23.314	3404	3409	3415	rVB5	12149	26874	1.31%	0.254%
22	23.919	3506	3512	3519	rBV8	18957	39906	1.95%	0.377%
23	24.853	3658	3671	3682	rBV	253778	751177	36.65%	7.090%

Sum of corrected areas: 10594136

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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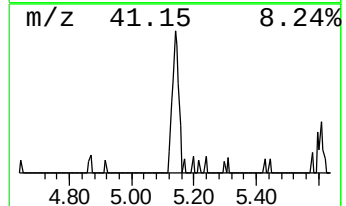
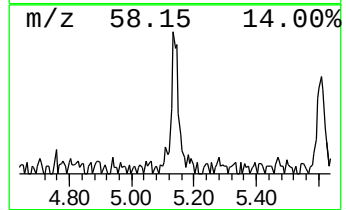
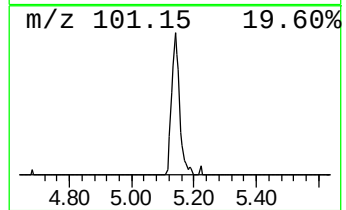
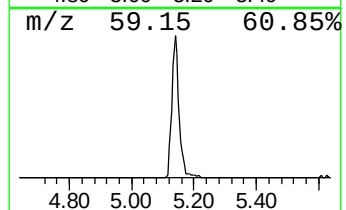
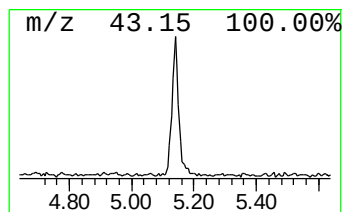
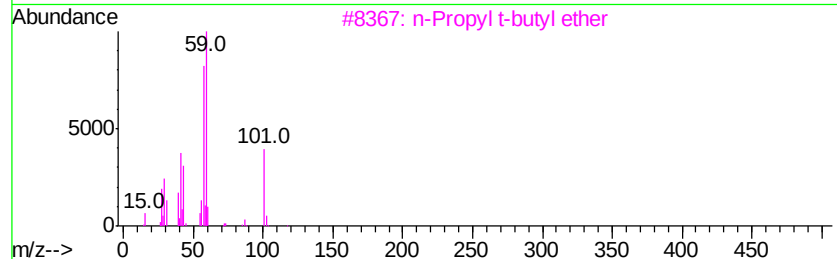
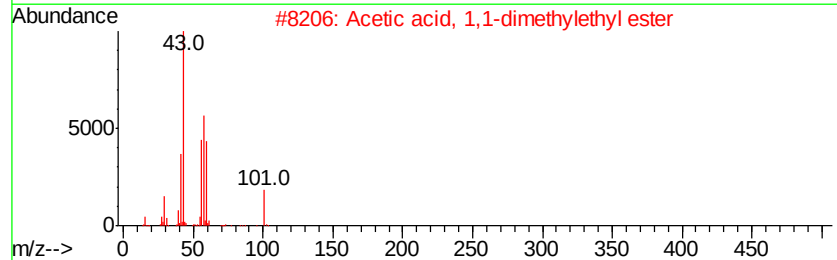
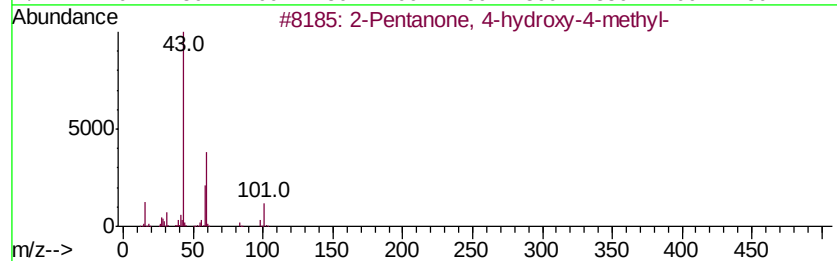
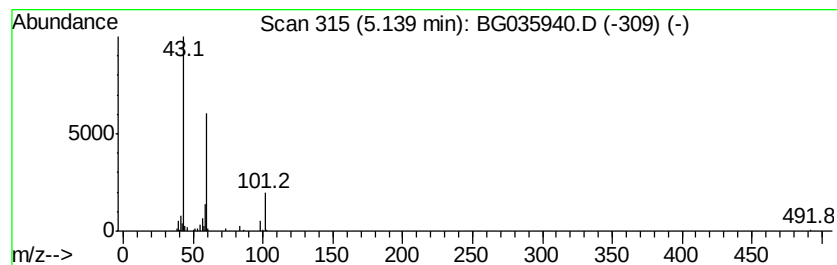
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	8.00 ng	83754	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	45
3			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
4			5-Hexen-2-one	98	C6H10O	000109-49-9	10
5			Guanidine	59	CH5N3	000113-00-8	9



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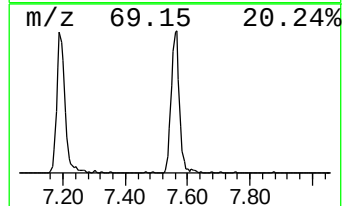
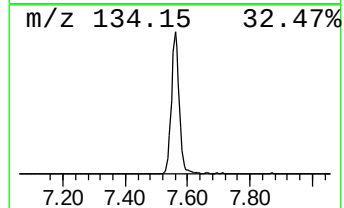
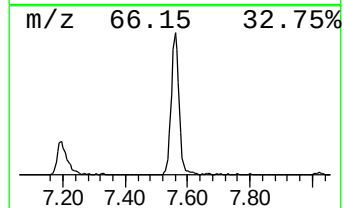
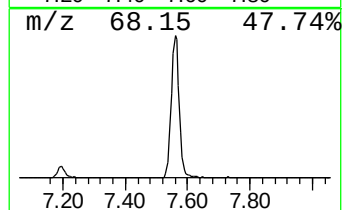
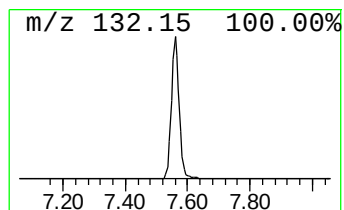
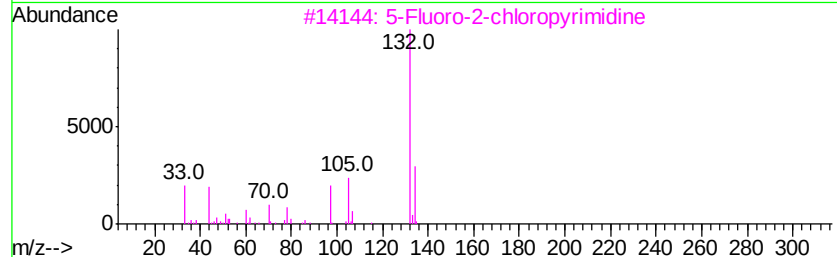
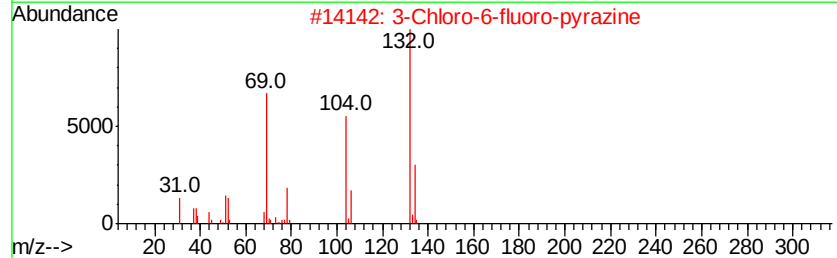
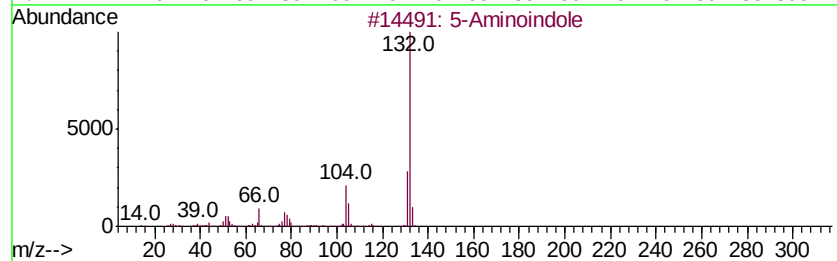
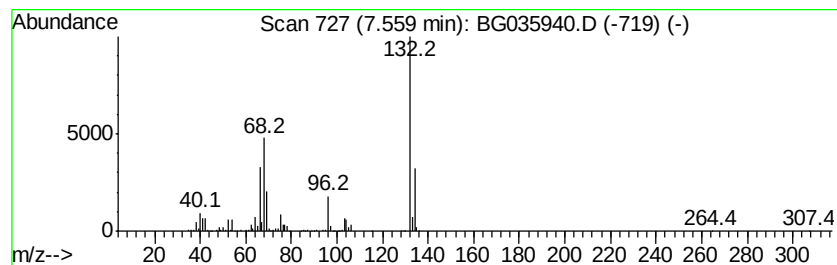
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	65.56 ng	686635	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	8.0	ng	83754	1	8.02	209465	20.0
unknown7.56	7.56	65.6	ng	686635	1	8.02	209465	20.0