

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060718\
 Data File : BG034955.D
 Acq On : 7 Jun 2018 16:47
 Operator : SJ/JU
 Sample : PB110027BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB110027BL

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-BG060718.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	5.186	318	323	332	rBV	48833	78951	2.30%	0.511%
2	5.650	393	402	412	rBV	715094	1110349	32.35%	7.191%
3	7.230	663	671	680	rBV	732129	1224678	35.69%	7.931%
4	7.612	728	736	744	rVB	684719	1148210	33.46%	7.436%
5	8.082	810	816	822	rBV	150057	252491	7.36%	1.635%
6	8.405	864	871	878	rBV	500924	864012	25.18%	5.595%
7	9.239	1003	1013	1023	rBV	402932	715693	20.85%	4.635%
8	10.884	1285	1293	1300	rVB	178813	324756	9.46%	2.103%
9	13.333	1703	1710	1717	rBV	1196431	1810152	52.75%	11.723%
10	14.702	1937	1943	1952	rVB3	318766	508295	14.81%	3.292%
11	16.188	2189	2196	2204	rBV	995435	1554044	45.28%	10.064%
12	17.446	2404	2410	2421	rVB2	483441	709676	20.68%	4.596%
13	20.060	2849	2855	2862	rBV	2645448	3431780	100.00%	22.224%
14	21.722	3131	3138	3145	rVB	540296	883479	25.74%	5.721%
15	24.965	3679	3690	3703	rBV2	285325	825084	24.04%	5.343%

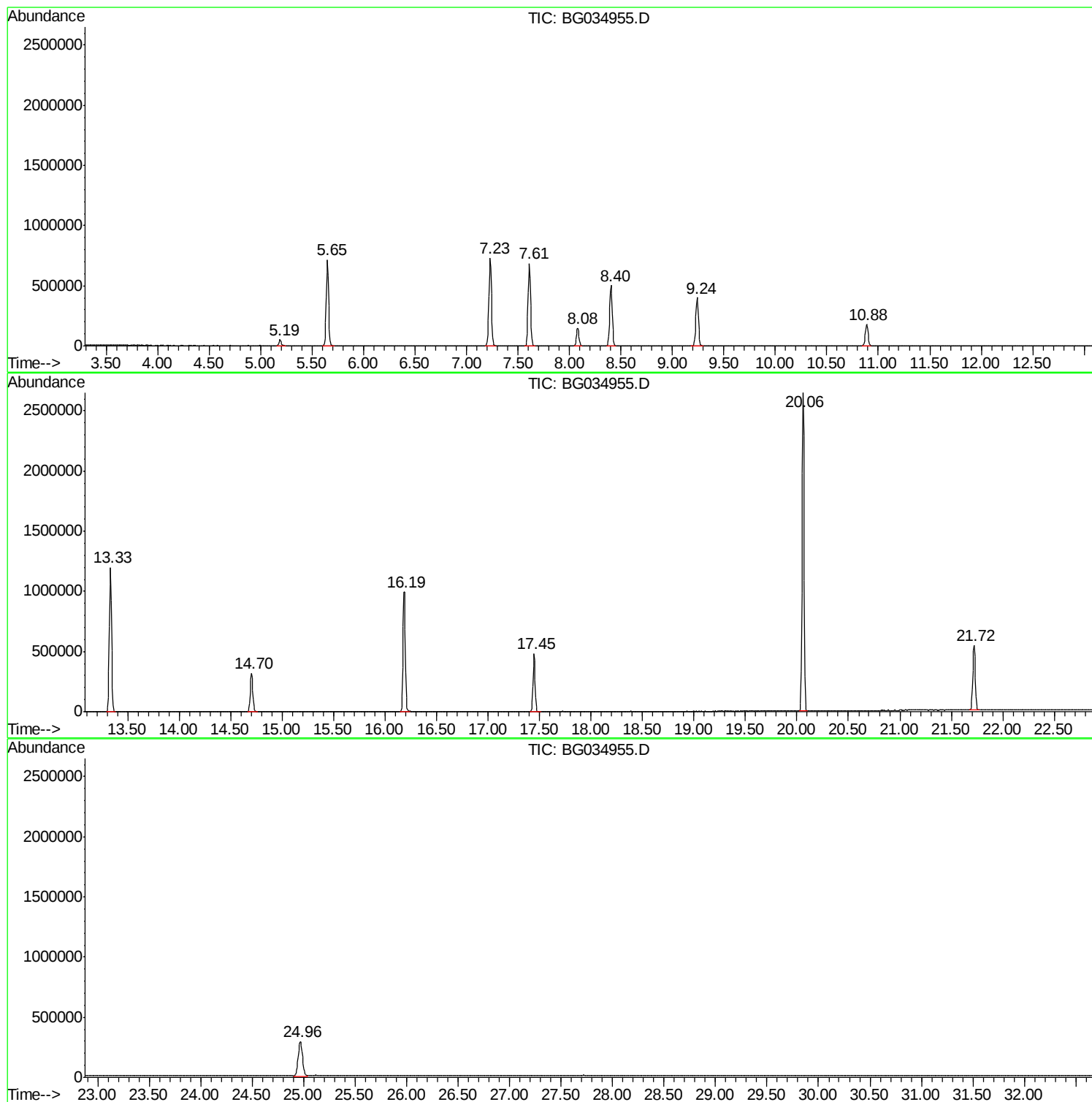
Sum of corrected areas: 15441650

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



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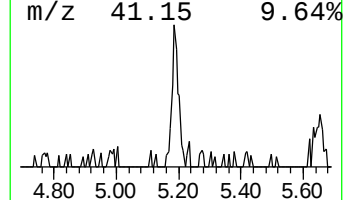
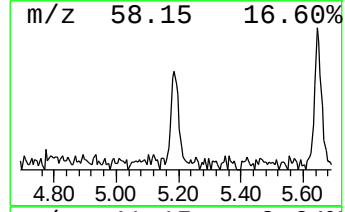
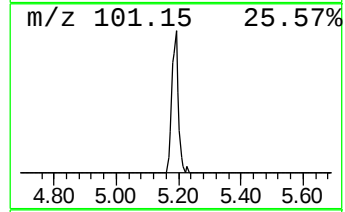
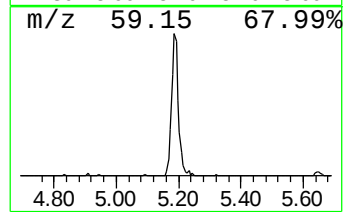
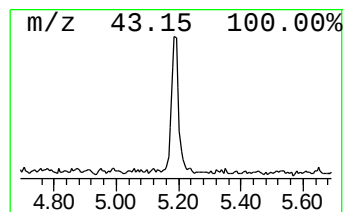
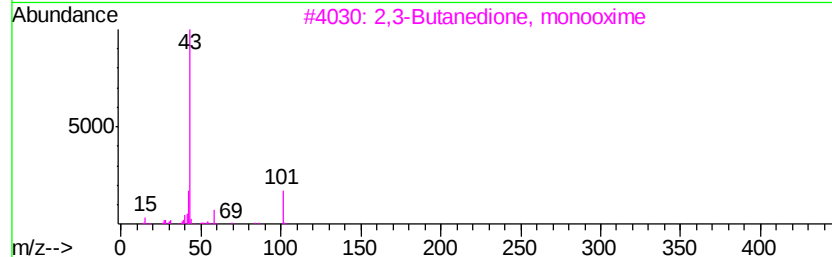
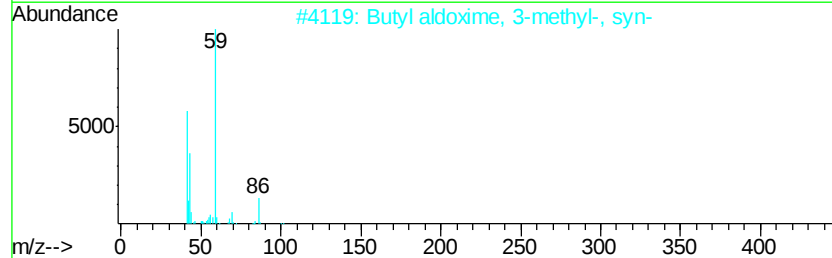
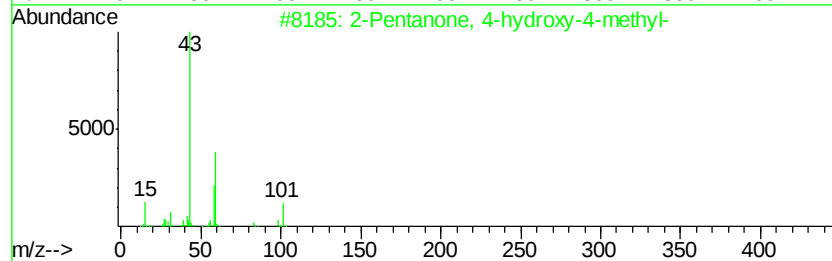
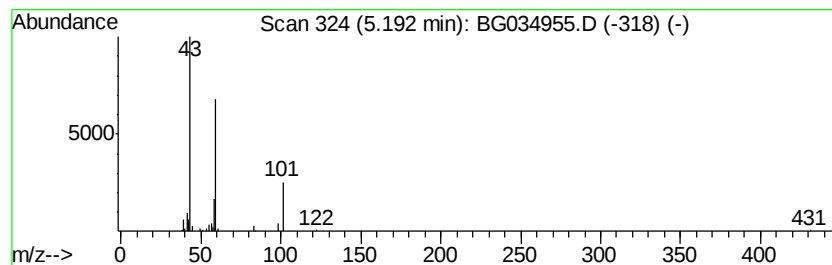
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	6.25 ng	78951	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	10
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Acetone	58	C3H6O	000067-64-1	9



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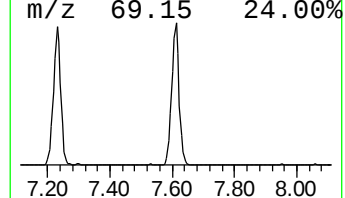
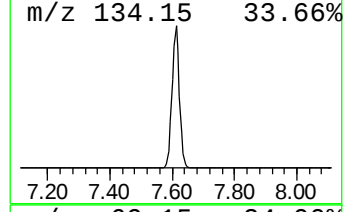
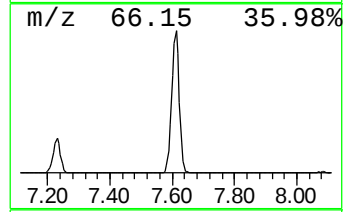
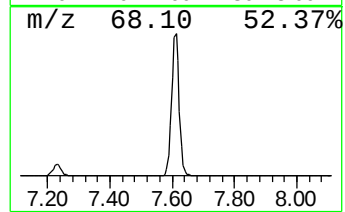
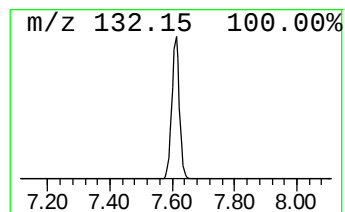
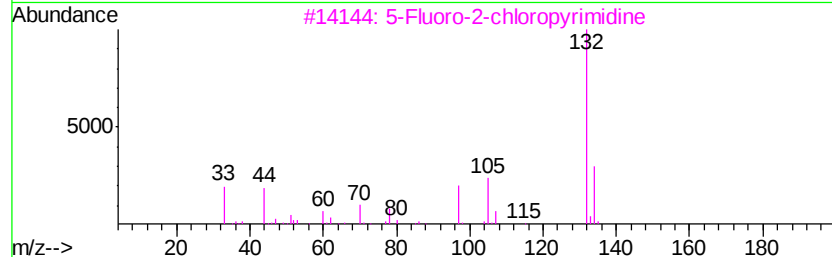
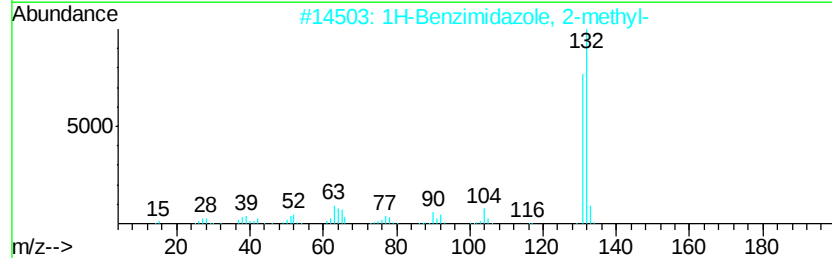
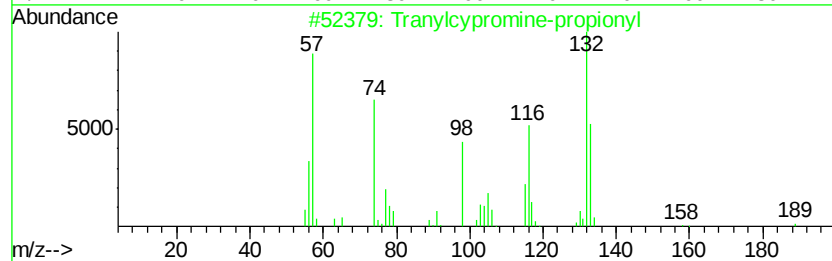
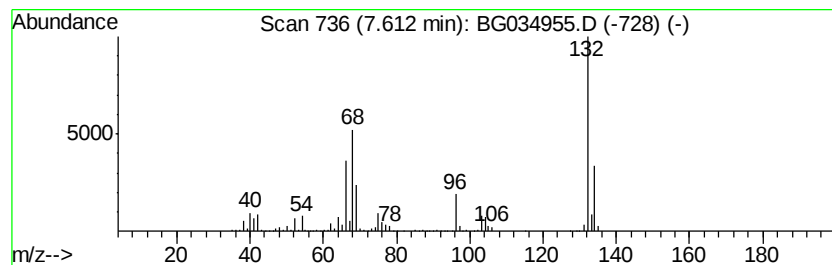
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	90.95 ng	1148210	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	6.3	ng	78951	1	8.08	252491	20.0
unknown7.61	7.61	91.0	ng	1148210	1	8.08	252491	20.0