

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082318\
 Data File : BG036365.D
 Acq On : 23 Aug 2018 5:42
 Operator : JU/SJ
 Sample : PB112219BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112219BL

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-BG082318.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	5.175	316	321	331	rVB	51212	82165	1.74%	0.414%
2	5.639	393	400	409	rBV	796647	1265393	26.77%	6.376%
3	7.219	660	669	682	rBV	799816	1379174	29.18%	6.949%
4	7.601	727	734	742	rVB	759722	1271802	26.91%	6.408%
5	8.071	807	814	821	rVB	179342	302061	6.39%	1.522%
6	8.395	861	869	878	rVB	829706	1416906	29.98%	7.139%
7	9.229	1002	1011	1020	rBV	661439	1193666	25.26%	6.014%
8	10.880	1285	1292	1300	rVB	241721	418897	8.86%	2.111%
9	13.324	1699	1708	1715	rBV	1895045	2871901	60.77%	14.470%
10	14.699	1934	1942	1950	rBV2	367662	602396	12.75%	3.035%
11	16.180	2185	2194	2209	rBV	953954	1422538	30.10%	7.167%
12	17.437	2402	2408	2418	rBV2	536068	809810	17.13%	4.080%
13	20.051	2847	2853	2863	rVB	3523012	4726098	100.00%	23.812%
14	21.708	3127	3135	3143	rBV	649964	1064494	22.52%	5.363%
15	24.928	3672	3683	3696	rBV3	368807	1020273	21.59%	5.141%

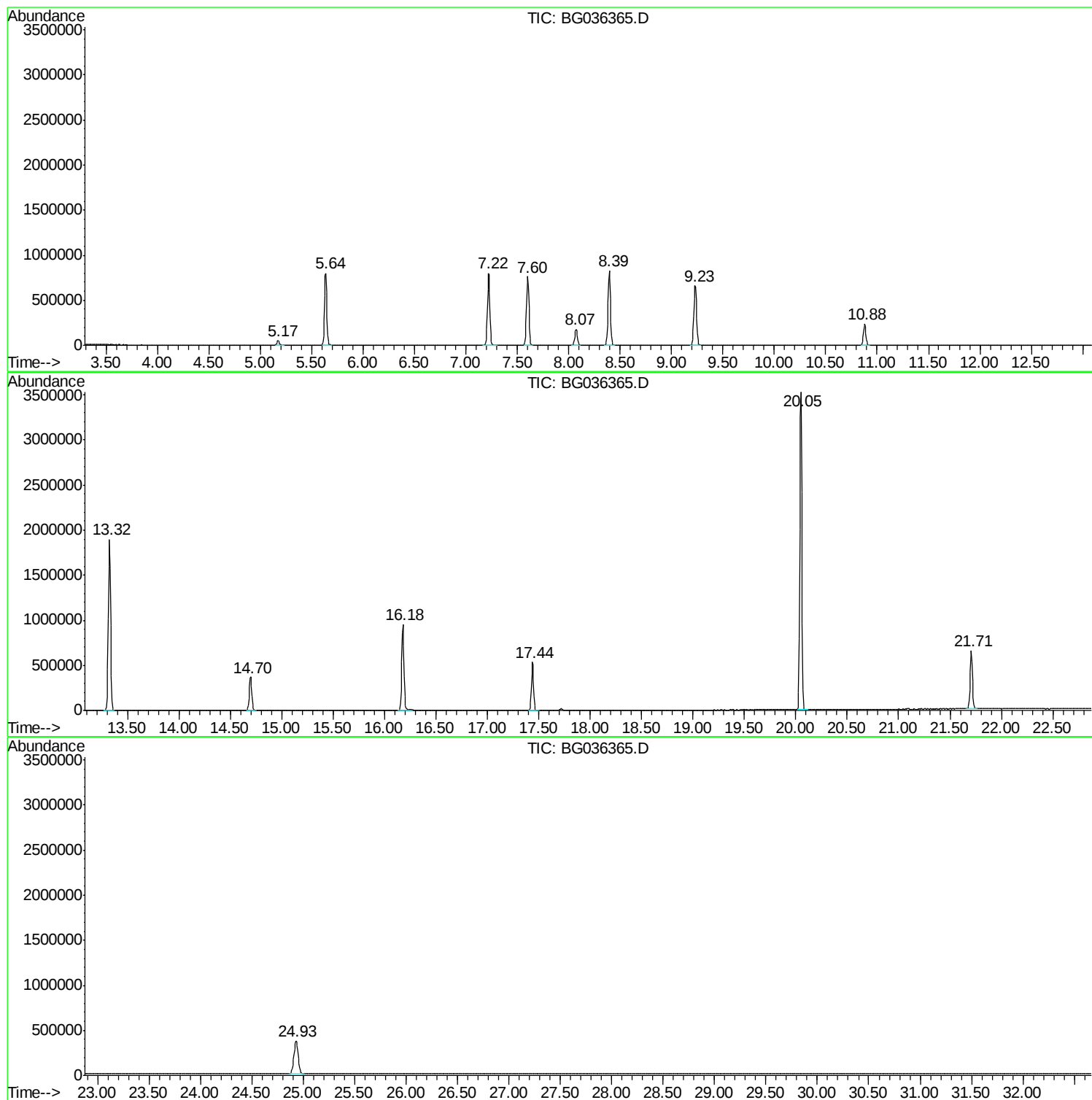
Sum of corrected areas: 19847574

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



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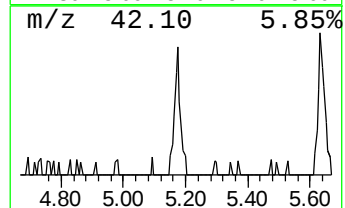
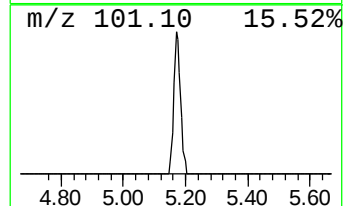
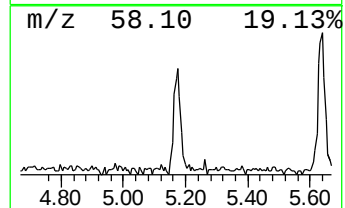
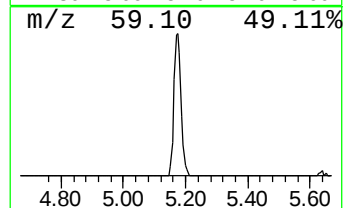
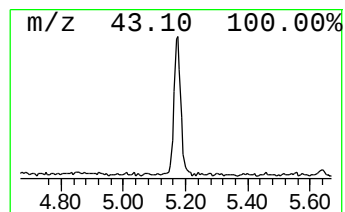
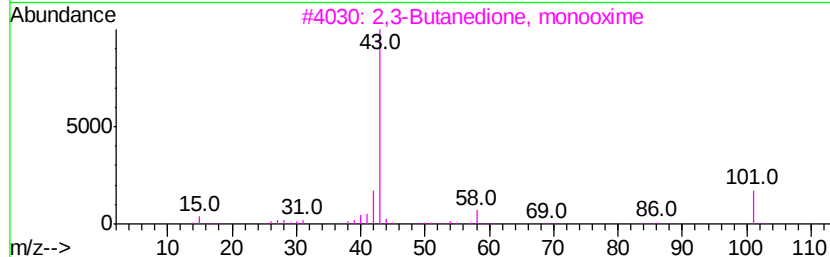
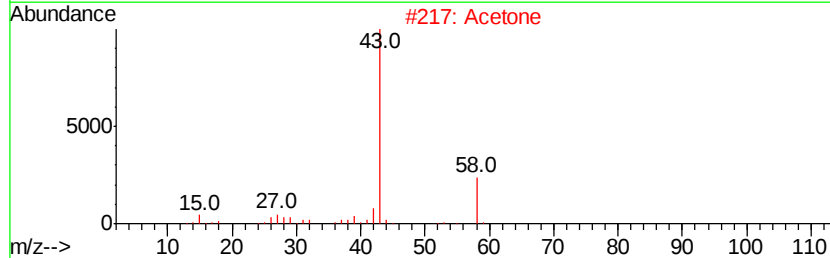
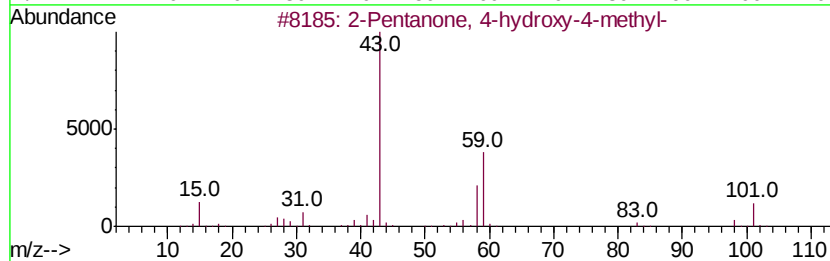
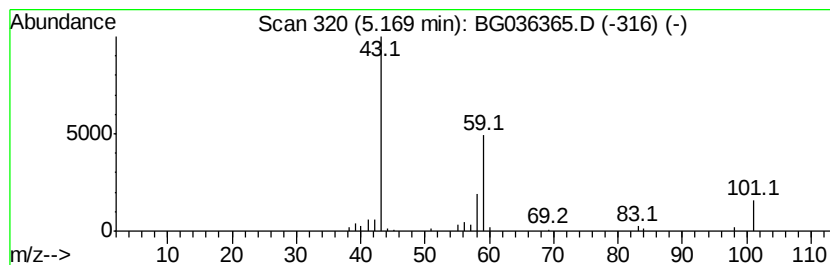
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	5.44 ng	82165	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			Acetone	58	C3H6O	000067-64-1	16
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			Propylamine	59	C3H9N	000107-10-8	5



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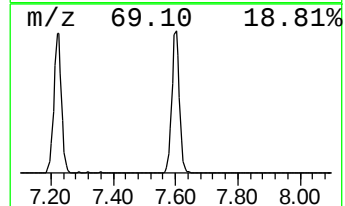
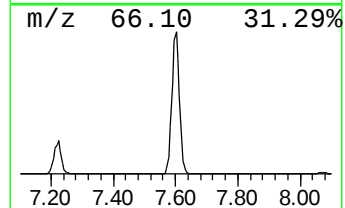
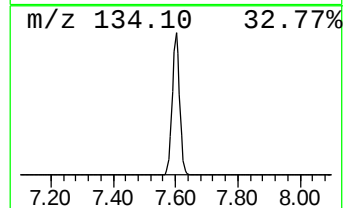
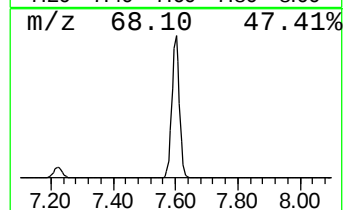
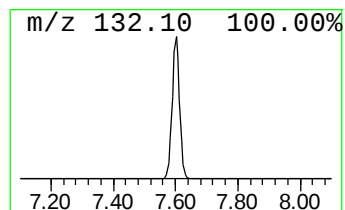
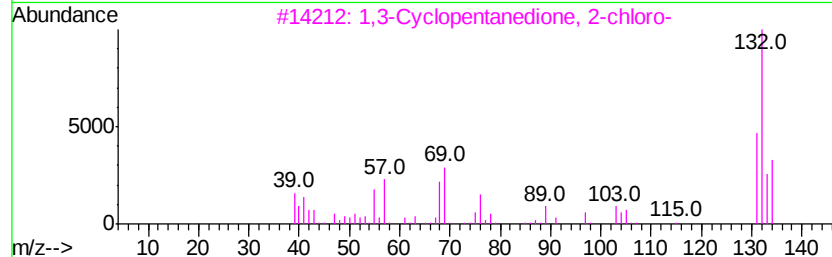
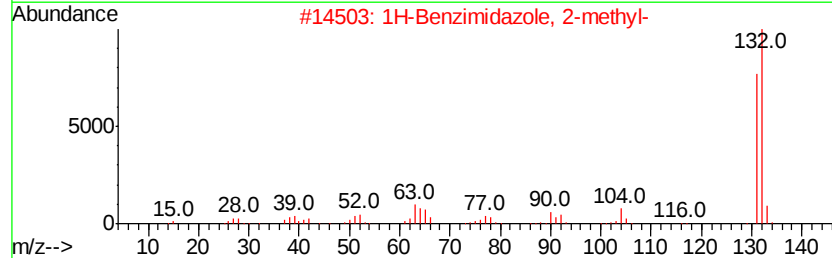
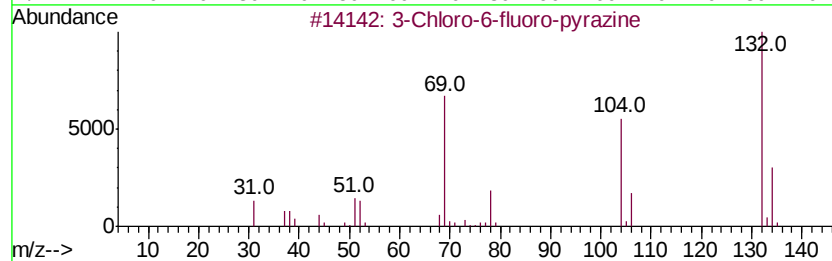
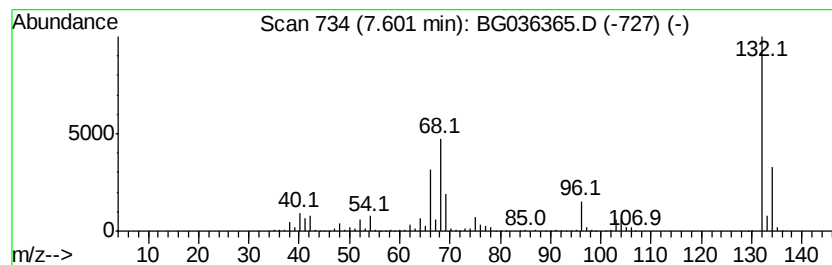
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Peak Number 2 unknown7.60 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	84.21 ng	1271800	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	5.4	ng	82165	1	8.07	302061	20.0
unknown7.60	7.60	84.2	ng	1271800	1	8.07	302061	20.0