

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
 Data File : BG038945.D
 Acq On : 5 Jan 2019 1:02
 Operator : JU/SJ
 Sample : PB116172BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116172BL

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-BG121218.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	5.130	341	347	357	rVB2	22774	39095	1.66%	0.401%
2	5.600	421	427	442	rBV	423898	699609	29.68%	7.170%
3	7.209	689	701	714	rBV	436205	773387	32.81%	7.926%
4	7.574	756	763	776	rBV	414300	747470	31.71%	7.661%
5	8.044	835	843	851	rVB	74282	124772	5.29%	1.279%
6	8.367	890	898	908	rVB	319901	562389	23.86%	5.764%
7	9.225	1036	1044	1054	rBV	250836	463164	19.65%	4.747%
8	10.864	1316	1323	1331	rBV	88811	167337	7.10%	1.715%
9	13.302	1731	1738	1749	rBV	729762	1200433	50.92%	12.303%
10	14.683	1966	1973	1984	rBV2	161964	267850	11.36%	2.745%
11	16.175	2220	2227	2247	rBV	663685	1111581	47.15%	11.392%
12	17.433	2434	2441	2458	rBV	251754	394751	16.74%	4.046%
13	20.036	2878	2884	2895	rBV	1749535	2357459	100.00%	24.161%
14	21.710	3163	3169	3180	rBV	247181	440035	18.67%	4.510%
15	25.006	3719	3730	3745	rBV	131847	407916	17.30%	4.181%

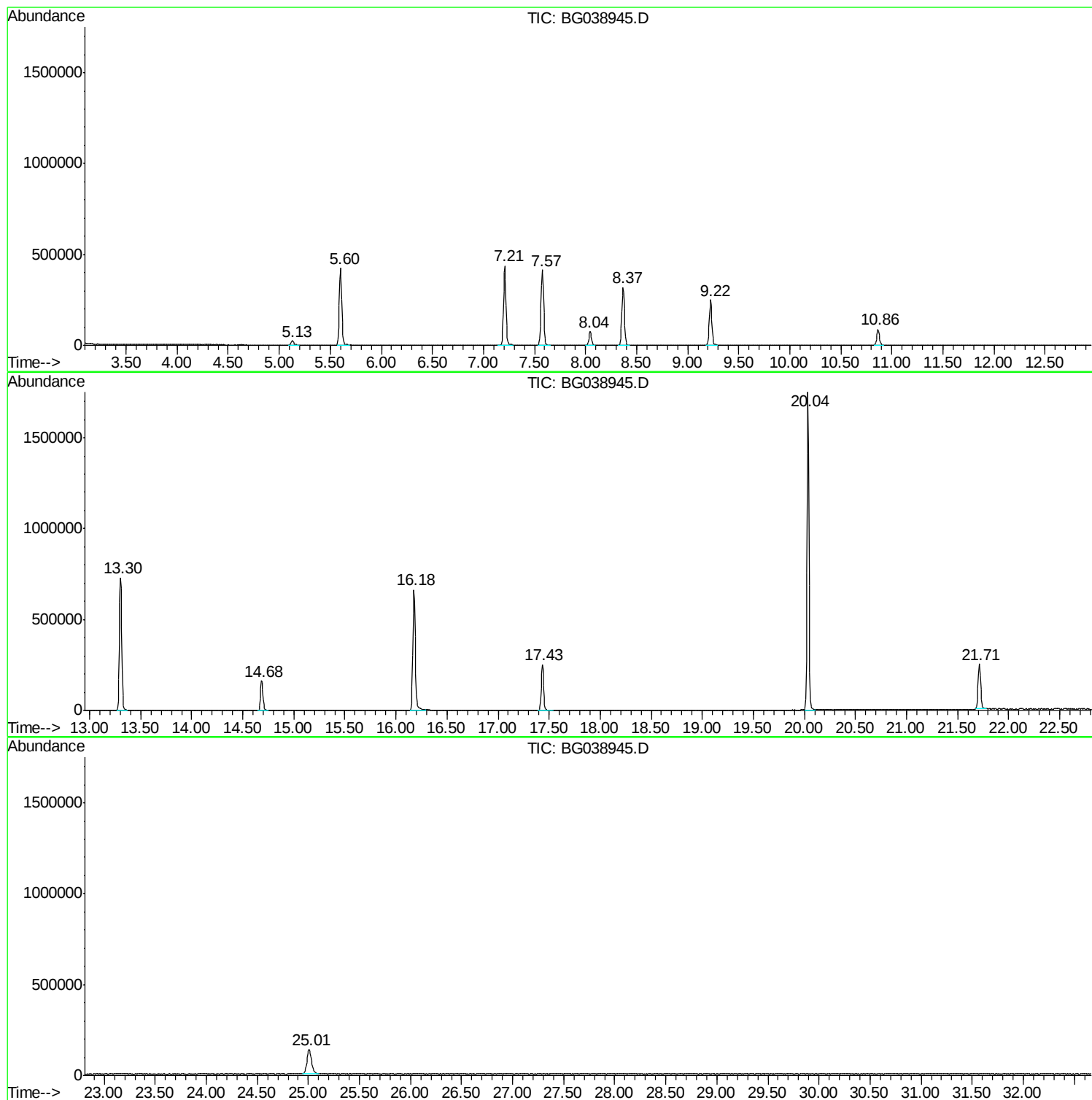
Sum of corrected areas: 9757248

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



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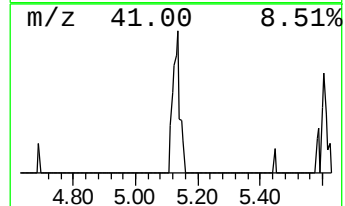
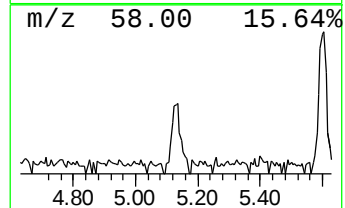
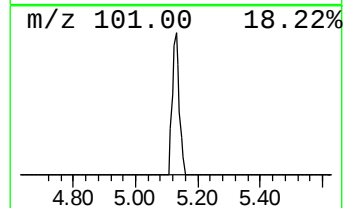
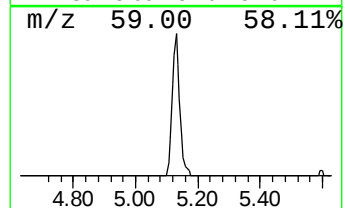
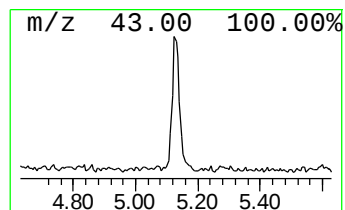
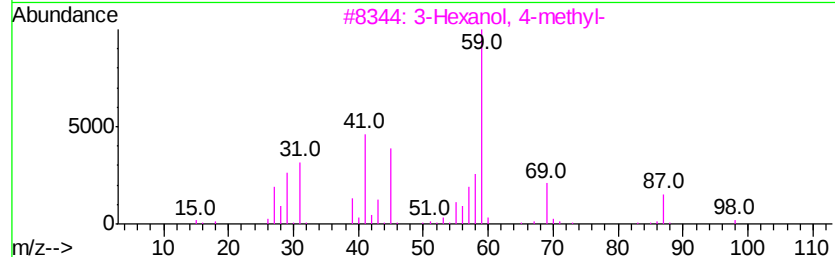
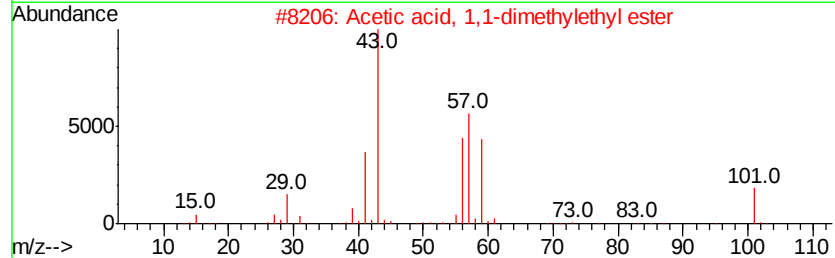
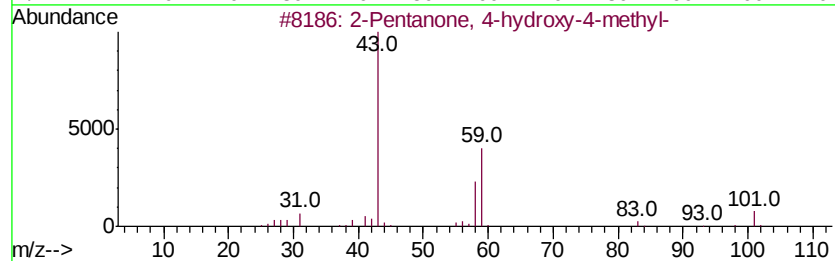
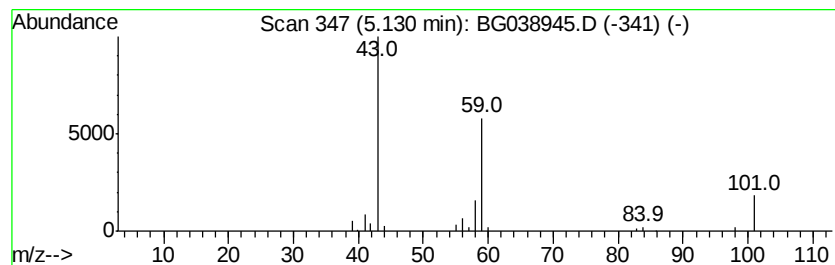
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	6.27 ng	39095	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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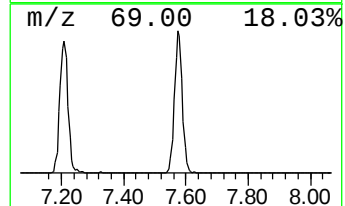
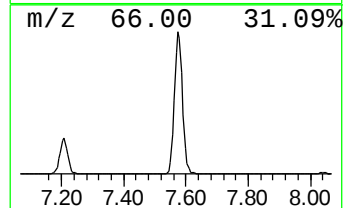
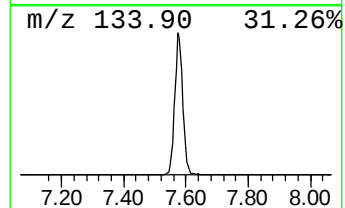
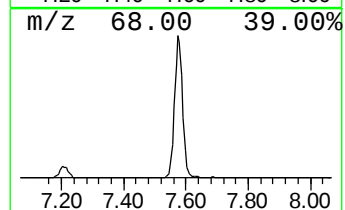
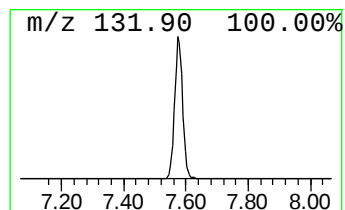
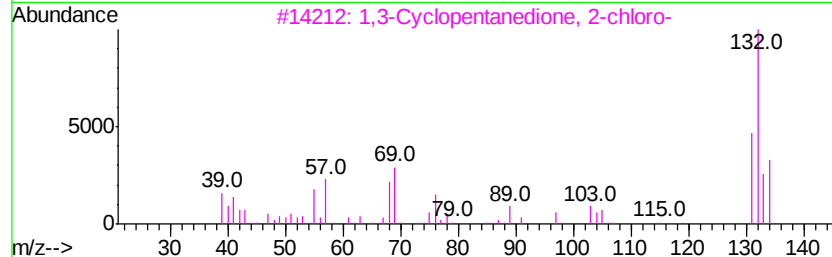
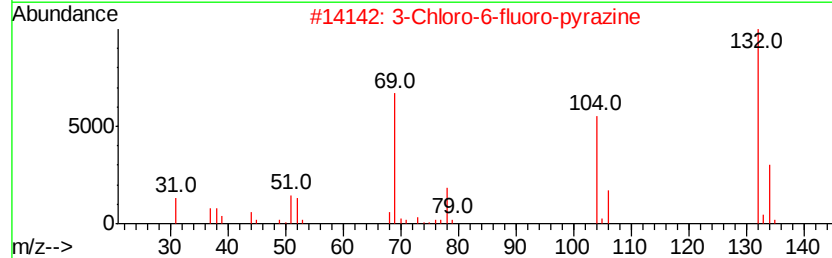
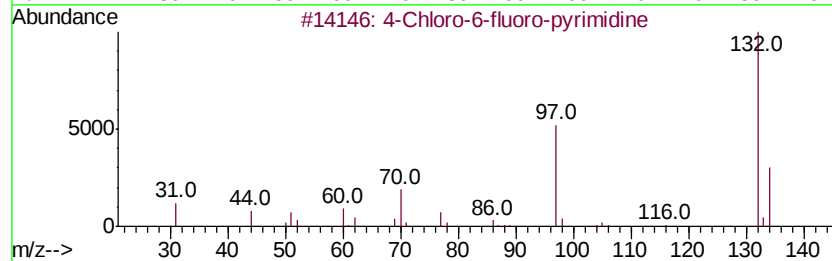
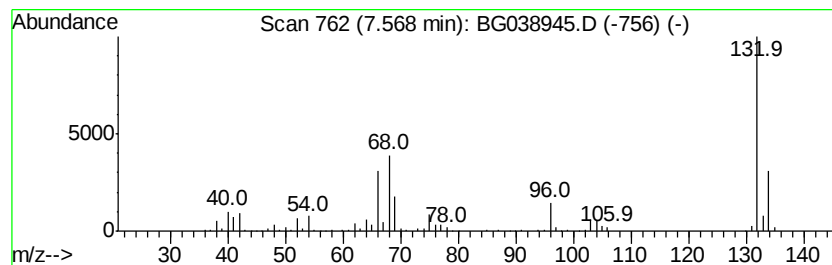
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	119.81 ng	747470	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
4		Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	6.3	ng	39095	1	8.04	124772	20.0
unknown7.57	7.57	119.8	ng	747470	1	8.04	124772	20.0