

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010919\
 Data File : BG039026.D
 Acq On : 9 Jan 2019 18:31
 Operator : JU/SJ
 Sample : PB116250BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116250BL

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-BG010919.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	5.108	340	344	354	rVB	13639	21224	1.17%	0.304%
2	5.578	418	424	437	rBV	248808	414844	22.82%	5.938%
3	7.188	690	698	709	rBV	265476	469468	25.82%	6.719%
4	7.558	752	761	771	rBV	252958	448415	24.66%	6.418%
5	8.029	835	841	849	rVB	59258	100752	5.54%	1.442%
6	8.352	887	896	904	rBV	193691	340734	18.74%	4.877%
7	9.209	1032	1042	1053	rBV	149524	282568	15.54%	4.044%
8	10.849	1313	1321	1331	rBV	73916	138709	7.63%	1.985%
9	13.287	1729	1736	1751	rBV	490739	784661	43.16%	11.231%
10	14.674	1965	1972	1985	rVB2	134902	232004	12.76%	3.321%
11	16.160	2218	2225	2234	rBV	427895	677695	37.28%	9.700%
12	17.423	2433	2440	2450	rBV	221989	347735	19.13%	4.977%
13	20.020	2876	2882	2890	rBV	1329442	1818039	100.00%	26.021%
14	21.695	3160	3167	3178	rBV2	273277	471546	25.94%	6.749%
15	24.979	3715	3726	3738	rVB2	152049	438442	24.12%	6.275%

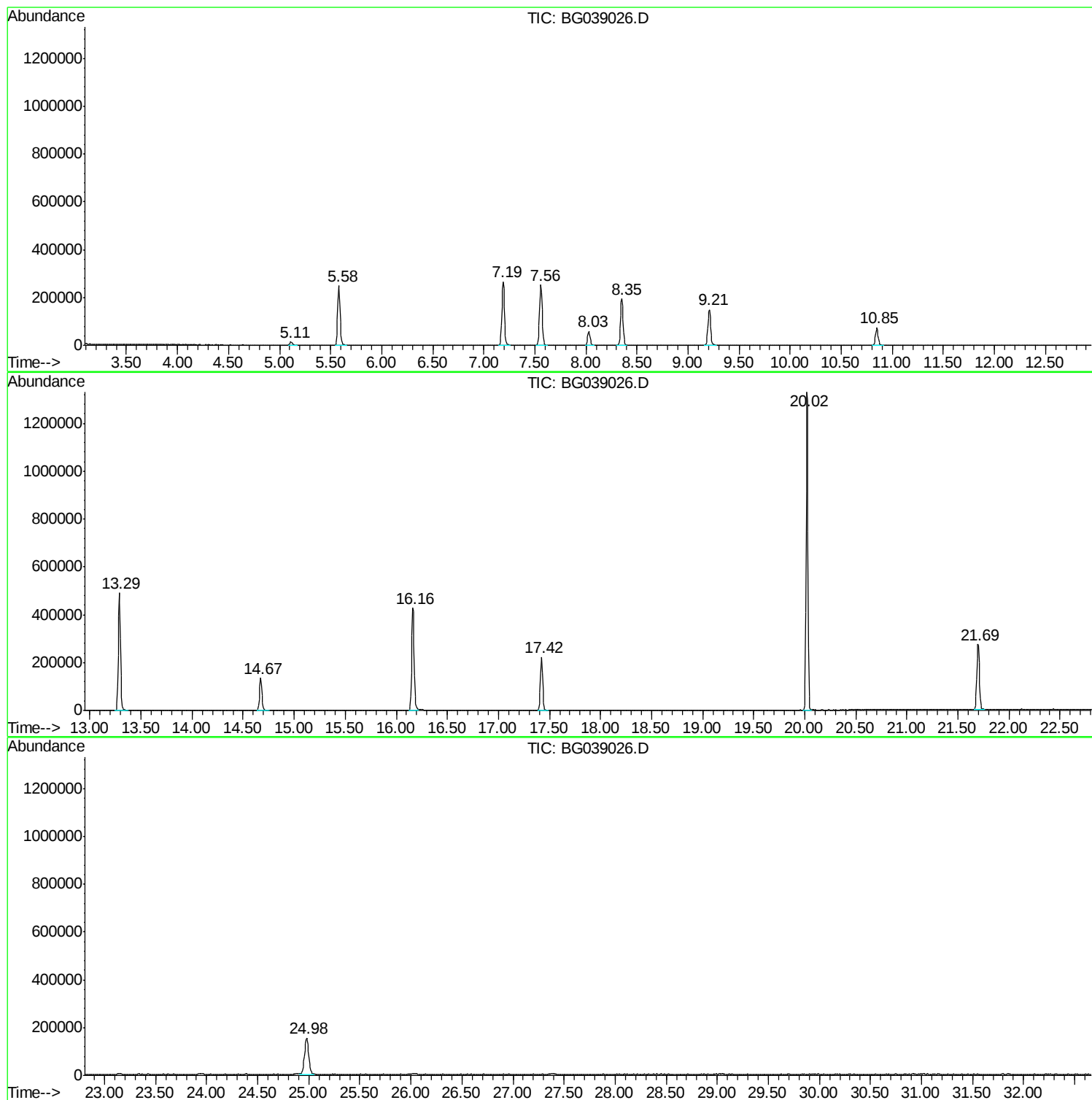
Sum of corrected areas: 6986836

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



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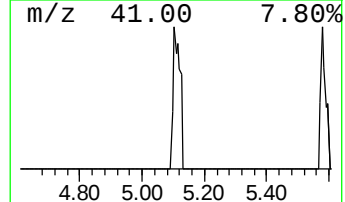
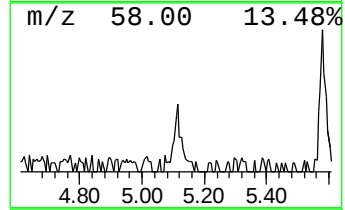
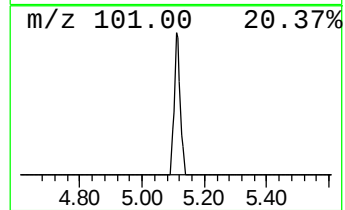
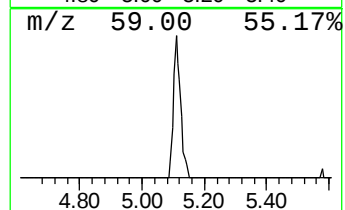
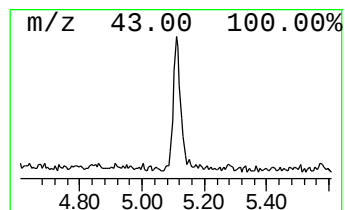
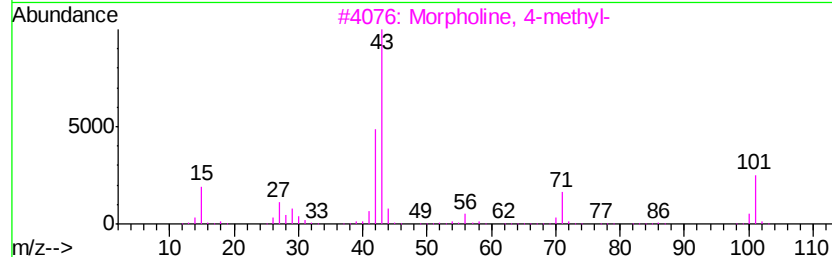
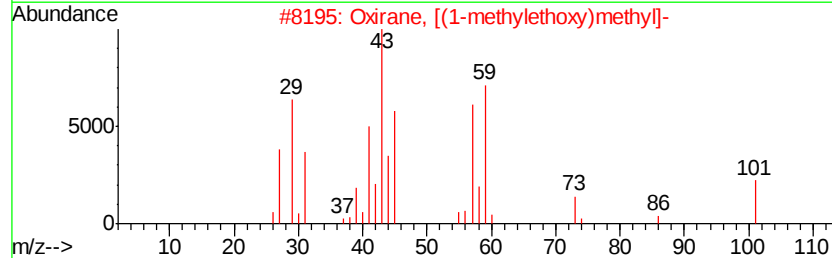
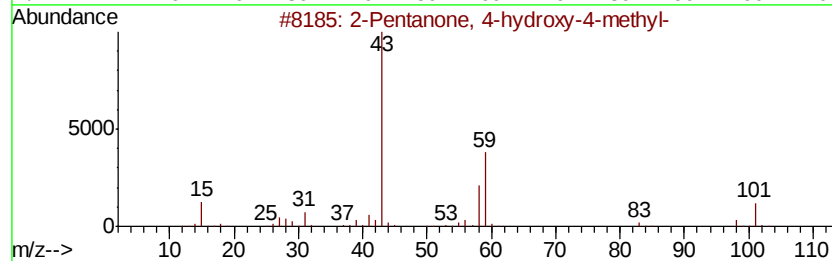
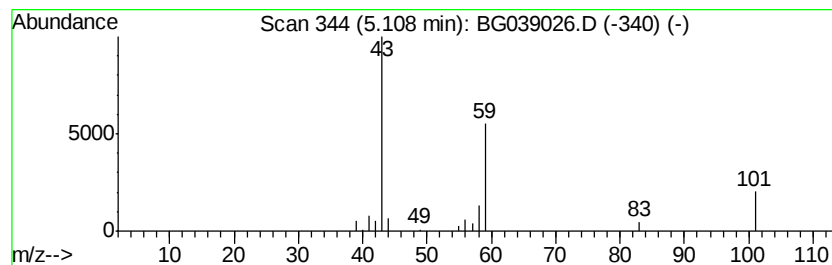
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	4.21 ng	21224	1,4-Dichlorobenzene-d4	8.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	40
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			1,3,4-Oxadiazole, 2-(acetyloxy)-...	172	C7H12N2O3	066398-57-0	9
5			Ethanol, 1-(methylenecyclopropyl)-	98	C6H10O	1000152-74-6	9



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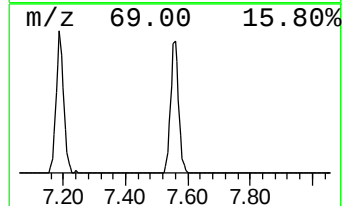
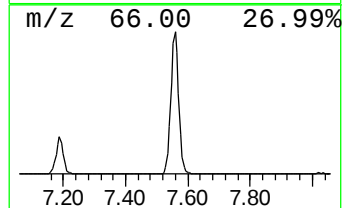
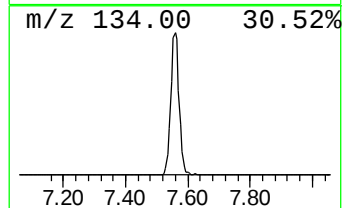
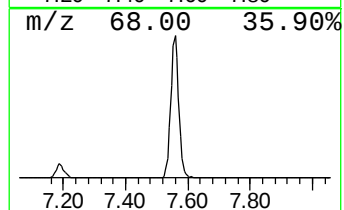
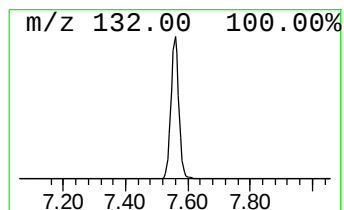
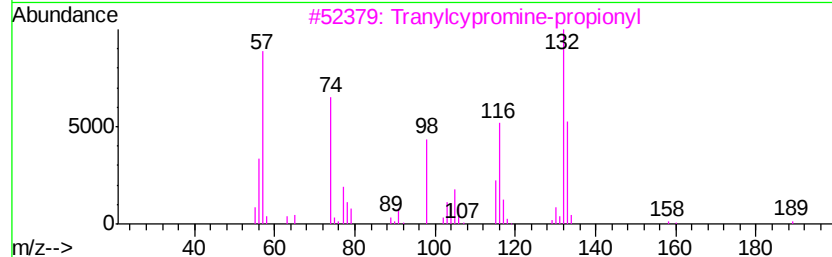
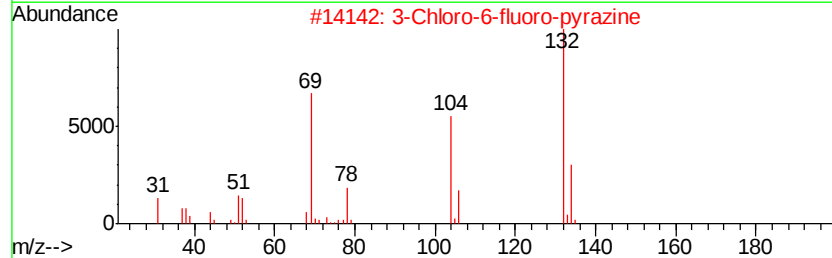
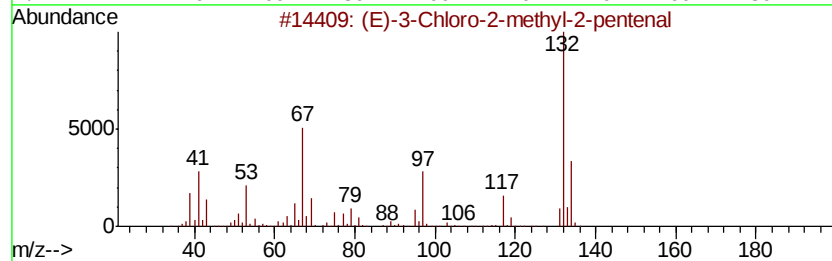
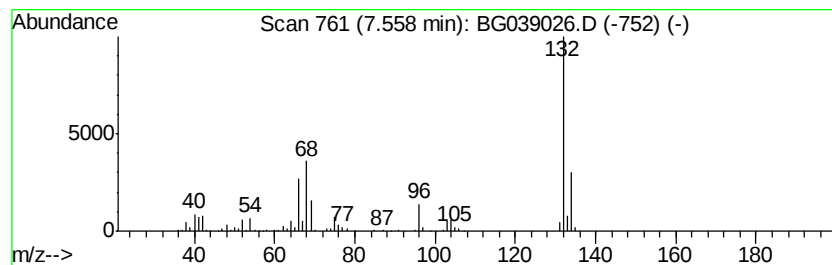
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	89.01 ng	448415	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Tranvlcypropamine-propionyl	189	C12H15NO	1000123-86-3	16
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	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
2-Pentanone, 4-hy...	5.11	4.2	ng	21224	1	8.03	100752	20.0
unknown7.56	7.56	89.0	ng	448415	1	8.03	100752	20.0