

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF071118\
 Data File : BF107328.D
 Acq On : 12 Jul 2018 4:27
 Operator : JU/SJ
 Sample : PB111003BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB111003BL

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_F\METHODS\8270-BF061918.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	5.166	478	481	492	rBV	58159	69609	4.71%	0.711%
2	5.578	547	551	571	rBV	834685	855388	57.85%	8.735%
3	6.578	717	721	739	rBV	861308	856996	57.96%	8.751%
4	6.707	739	743	757	rVB	988989	855658	57.87%	8.738%
5	6.931	777	781	787	rBB	343497	290427	19.64%	2.966%
6	7.084	803	807	811	rBV	1038066	1003608	67.88%	10.249%
7	7.495	873	877	892	rBV	731896	768526	51.98%	7.848%
8	8.213	995	999	1012	rBV	372505	344248	23.28%	3.515%
9	9.289	1177	1182	1192	rBV	1509513	1404011	94.96%	14.337%
10	9.977	1295	1299	1308	rBB	400481	366312	24.78%	3.741%
11	10.777	1431	1435	1450	rBB	573345	523099	35.38%	5.342%
12	11.477	1550	1554	1563	rBV	439168	381980	25.83%	3.901%
13	13.060	1819	1823	1827	rBV	1485611	1478549	100.00%	15.099%
14	14.136	2002	2006	2010	rBV	368684	348624	23.58%	3.560%
15	15.671	2262	2267	2274	rVB	178822	245630	16.61%	2.508%

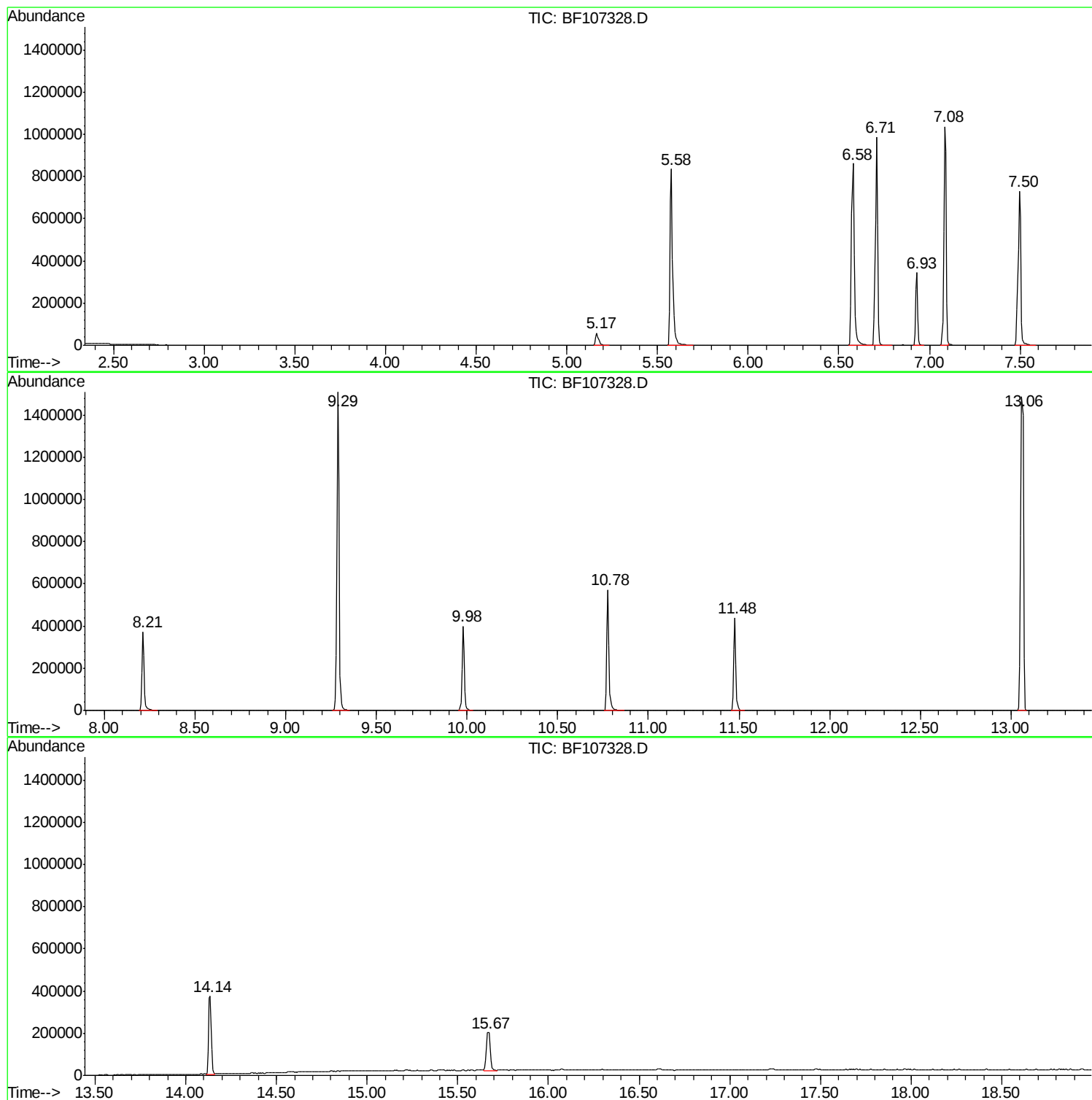
Sum of corrected areas: 9792665

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



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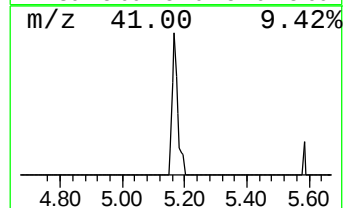
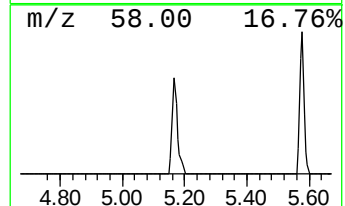
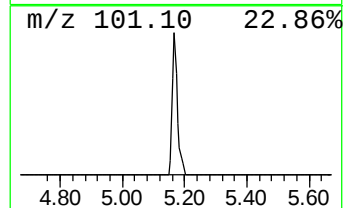
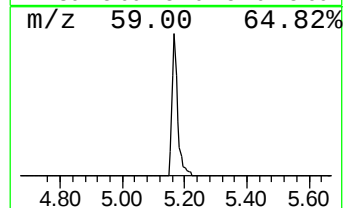
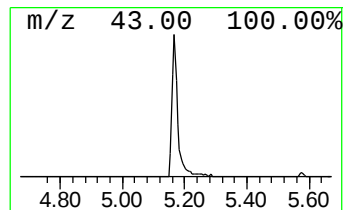
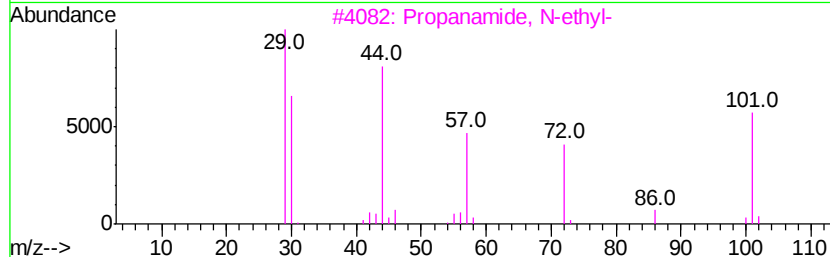
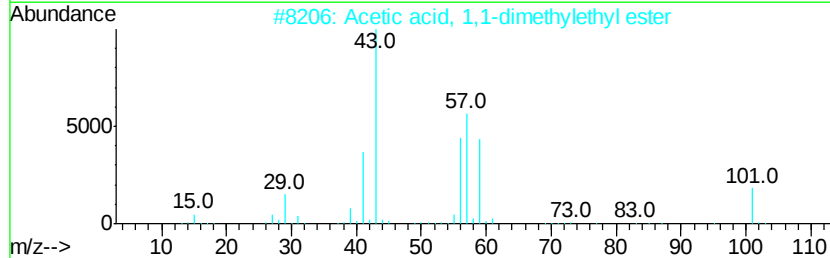
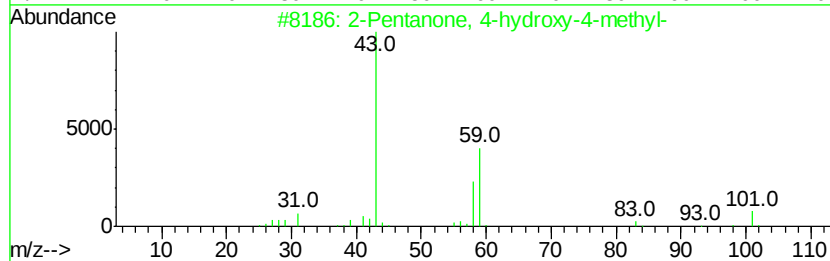
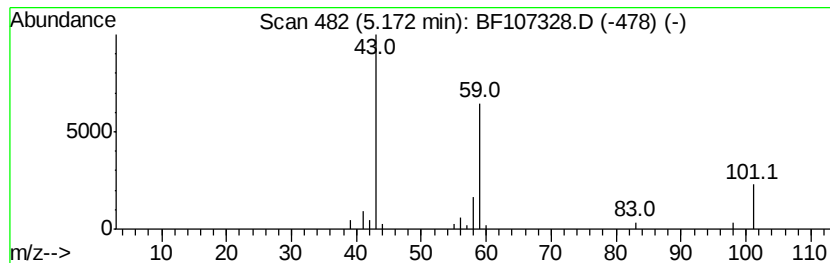
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.79 ng	69609	1,4-Dichlorobenzene-d4	6.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propanamide, N-ethyl-	101	C5H11NO	005129-72-6	9
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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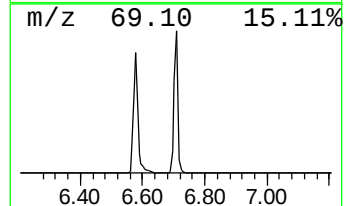
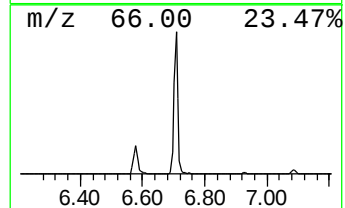
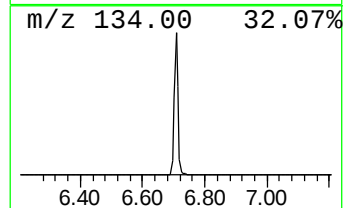
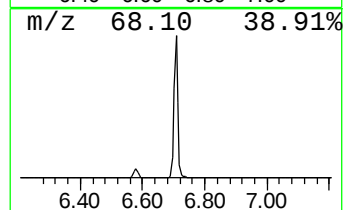
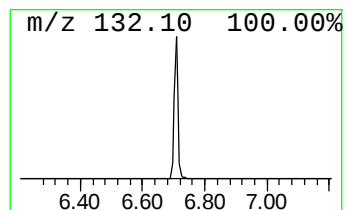
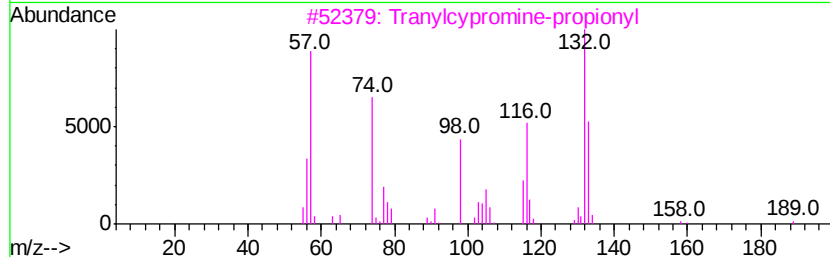
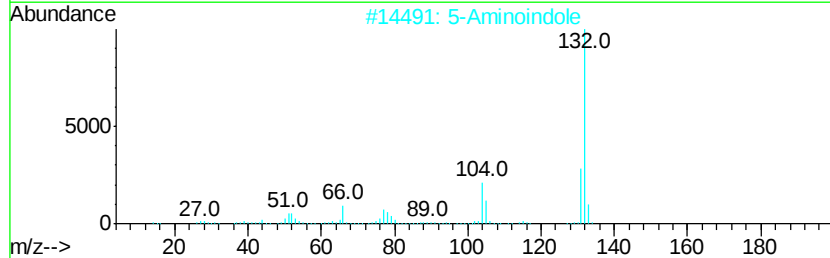
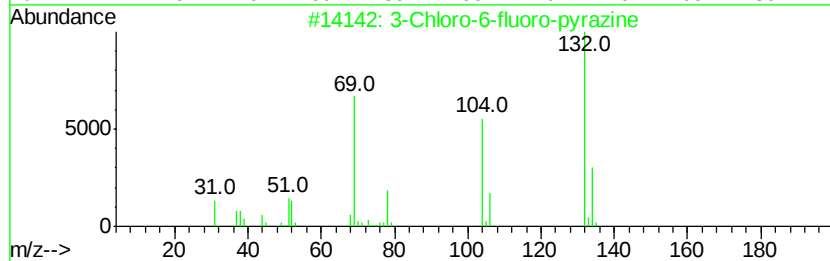
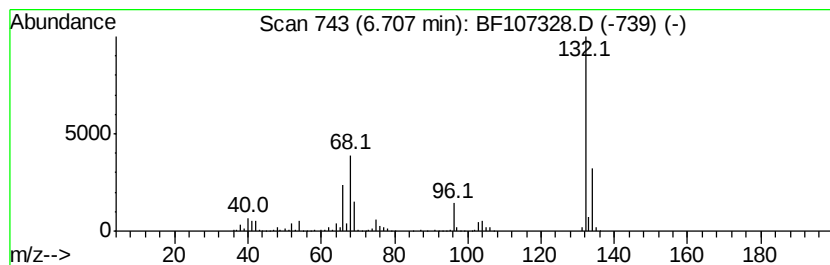
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Peak Number 2 unknown6.71 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	58.92 ng	855658	1,4-Dichlorobenzene-d4	6.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	4.8	ng	69609	1	6.93	290427	20.0
unknown6.71	6.71	58.9	ng	855658	1	6.93	290427	20.0