

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF071218\
 Data File : BF107337.D
 Acq On : 12 Jul 2018 14:30
 Operator : JU/SJ
 Sample : PB110949BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110949BL

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	5.172	478	482	502	rBB	124113	173575	10.44%	1.439%
2	5.578	547	551	573	rBV	1199054	1177343	70.78%	9.759%
3	6.578	716	721	739	rBV	1142767	1180860	71.00%	9.788%
4	6.707	739	743	746	rBV	1295643	1179962	70.94%	9.780%
5	6.925	777	780	787	rBB	344593	299749	18.02%	2.485%
6	7.084	803	807	820	rBB	1075242	973829	58.55%	8.072%
7	7.495	872	877	897	rBV	783678	795533	47.83%	6.594%
8	8.213	994	999	1018	rBB	395880	394209	23.70%	3.267%
9	9.289	1176	1182	1185	rBV	1528849	1505079	90.49%	12.475%
10	9.978	1294	1299	1309	rVB	471510	466261	28.03%	3.865%
11	10.778	1431	1435	1453	rBV	981232	959108	57.66%	7.950%
12	11.472	1550	1553	1563	rBV	491832	509370	30.62%	4.222%
13	13.060	1819	1823	1826	rBV	1660339	1663297	100.00%	13.786%
14	14.136	2002	2006	2012	rBV	437156	417410	25.10%	3.460%
15	15.677	2263	2268	2276	rBV	248285	369146	22.19%	3.060%

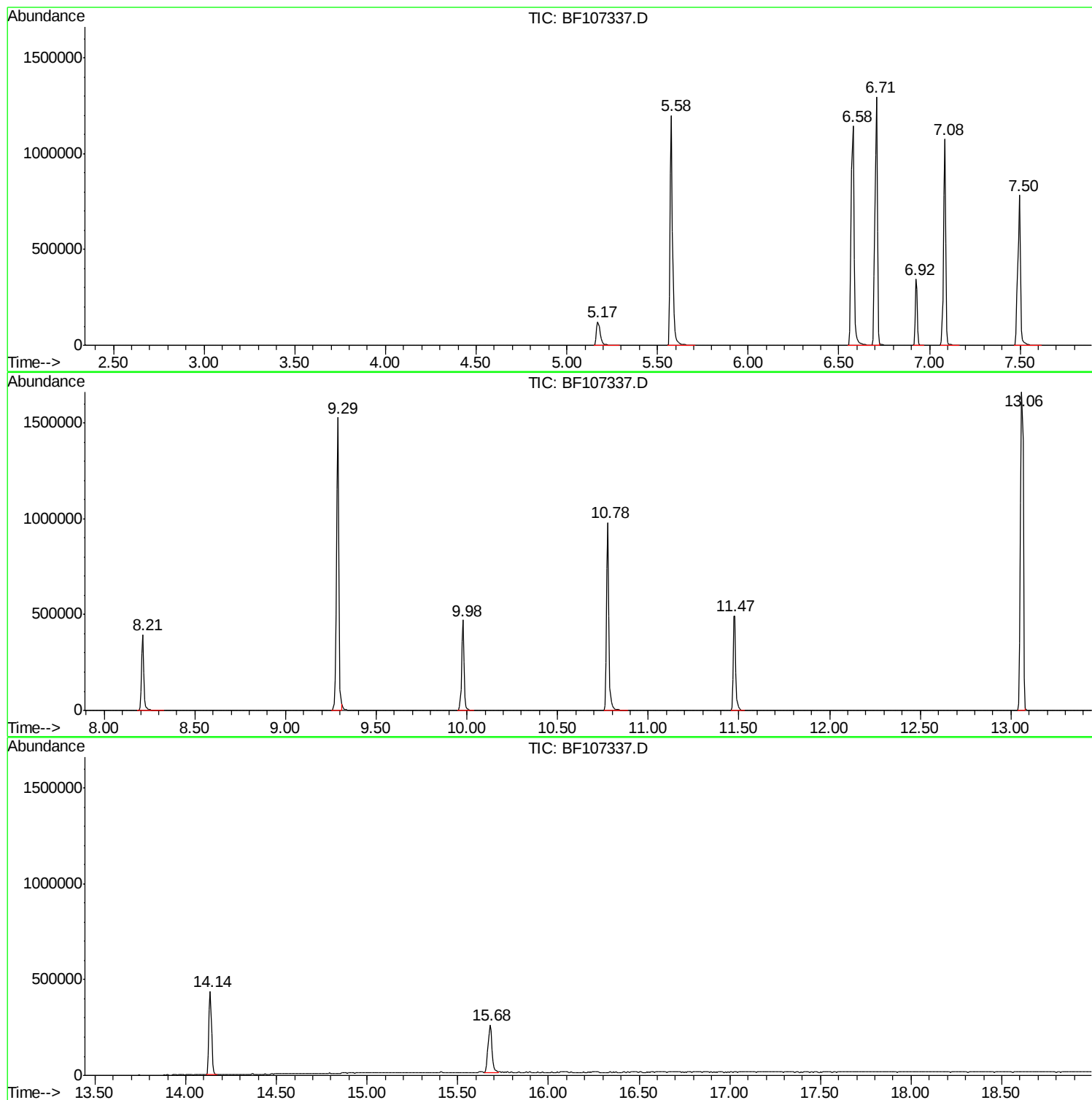
Sum of corrected areas: 12064731

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



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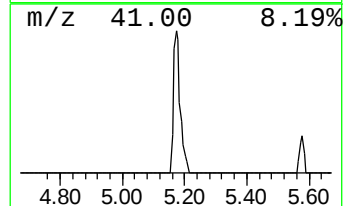
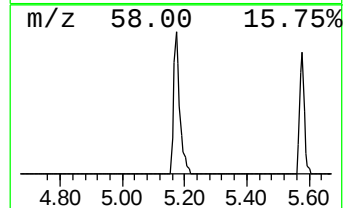
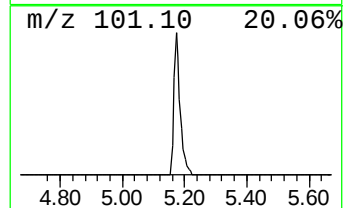
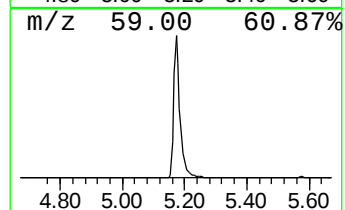
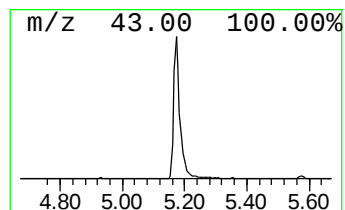
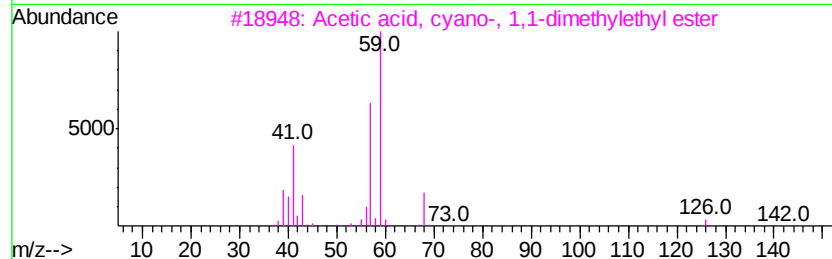
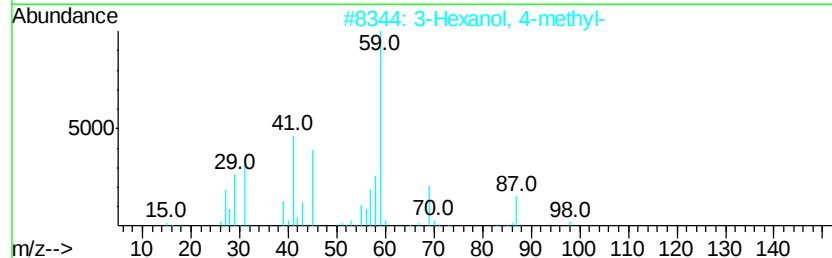
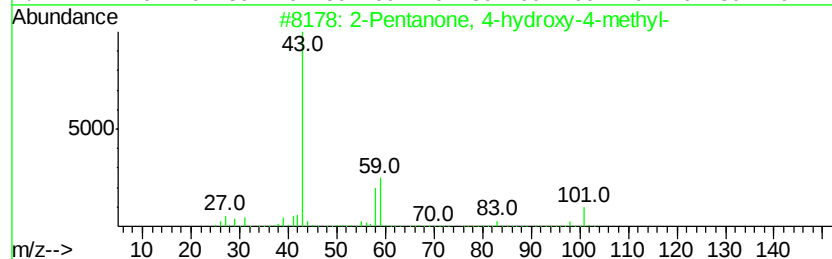
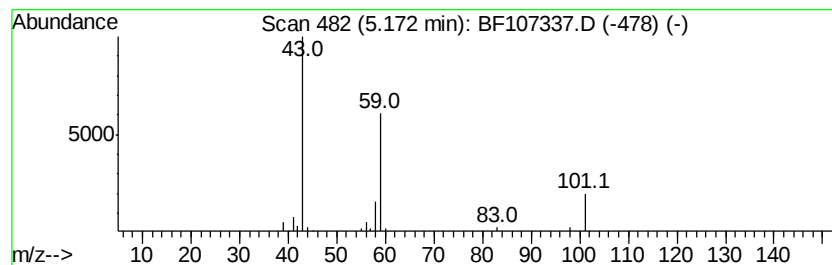
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	11.58 ng	173575	1,4-Dichlorobenzene-d4	6.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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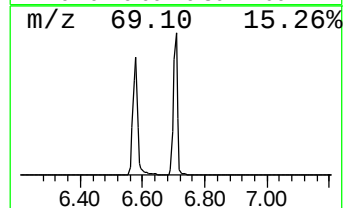
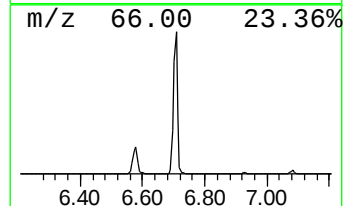
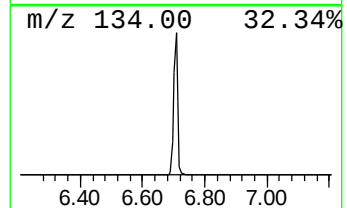
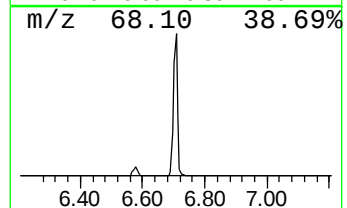
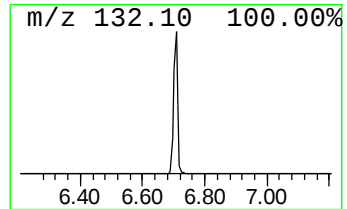
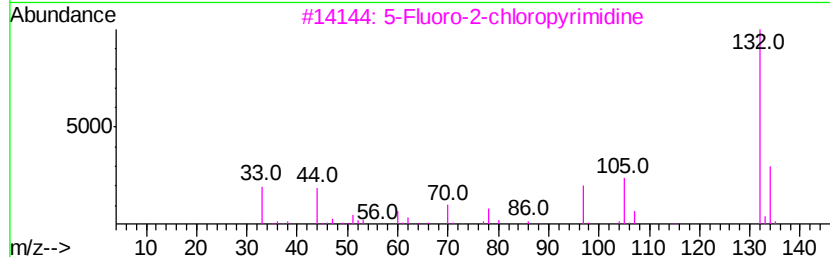
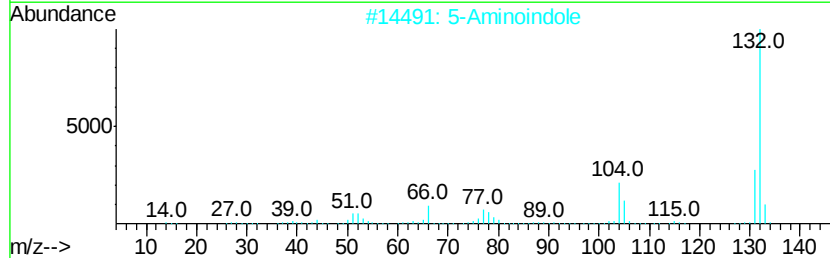
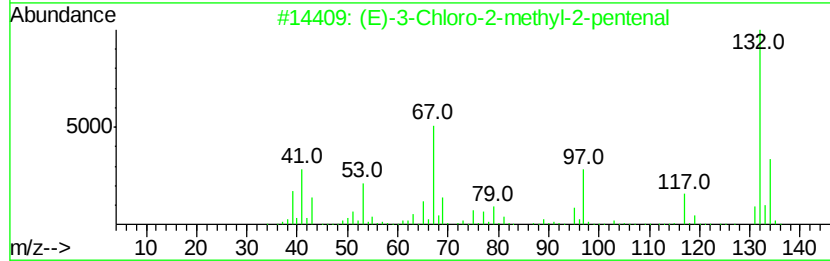
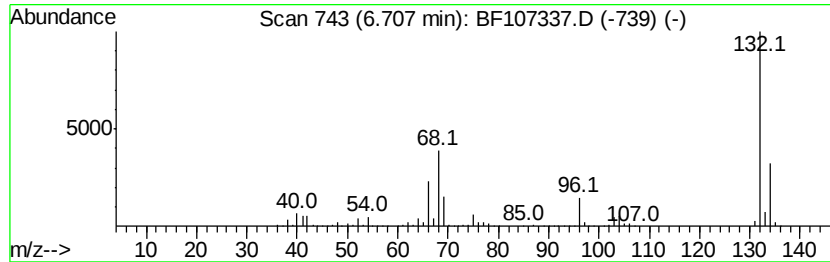
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Peak Number 2 unknown6.71 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	78.73 ng	1179960	1,4-Dichlorobenzene-d4	6.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	9
5			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	11.6	ng	173575	1	6.92	299749	20.0
unknown6.71	6.71	78.7	ng	1179960	1	6.92	299749	20.0