

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061218\
 Data File : BF106367.D
 Acq On : 13 Jun 2018 1:53
 Operator : JU/SJ
 Sample : PB110114BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110114BL

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-BF061118.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	5.219	485	490	497	rBV	265951	332331	8.85%	1.132%
2	5.619	555	558	577	rVB	2975143	3434147	91.48%	11.699%
3	6.619	724	728	747	rBV	2453139	3549346	94.55%	12.092%
4	6.760	747	752	755	rBV	3372531	3149178	83.89%	10.728%
5	6.972	785	788	792	rBV	1003009	823343	21.93%	2.805%
6	7.131	811	815	818	rBV	2975353	2651958	70.65%	9.034%
7	7.537	879	884	887	rBV	2201728	2070473	55.16%	7.054%
8	8.254	1002	1006	1010	rBV	1141898	922522	24.57%	3.143%
9	9.331	1184	1189	1192	rBV	3703982	3328572	88.67%	11.340%
10	10.013	1301	1305	1309	rBV	1035858	895022	23.84%	3.049%
11	10.807	1436	1440	1459	rBV	1834487	1789441	47.67%	6.096%
12	11.507	1555	1559	1562	rBV	1073588	926419	24.68%	3.156%
13	13.089	1824	1828	1832	rBV	3611303	3753915	100.00%	12.789%
14	14.154	2005	2009	2012	rBV	1097377	1002318	26.70%	3.415%
15	15.689	2265	2270	2277	rVB	567337	724763	19.31%	2.469%

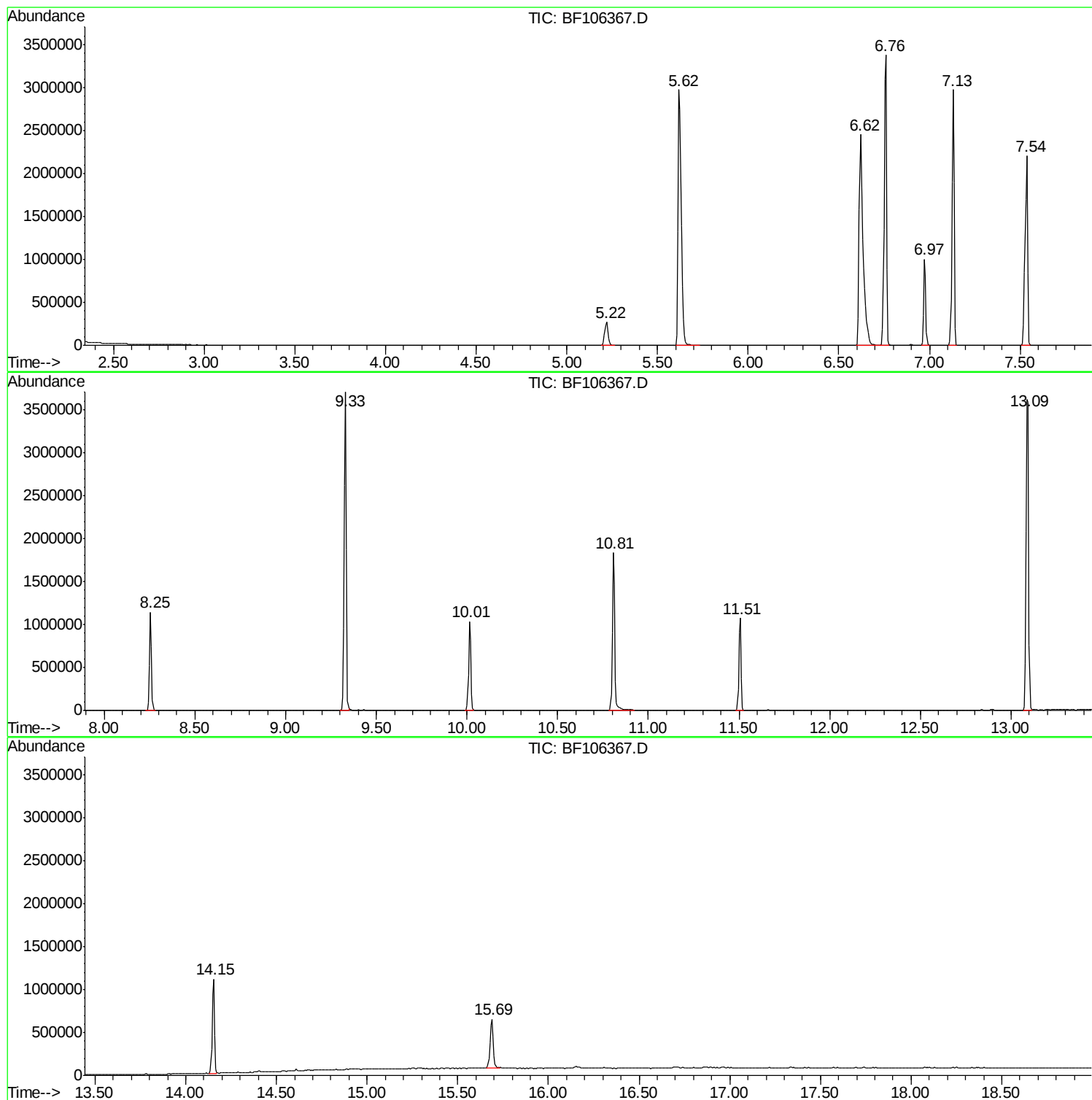
Sum of corrected areas: 29353748

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



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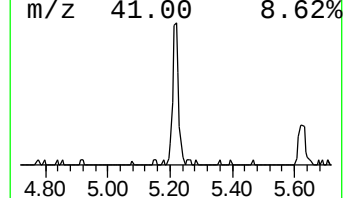
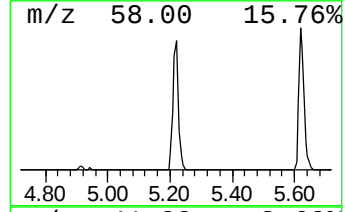
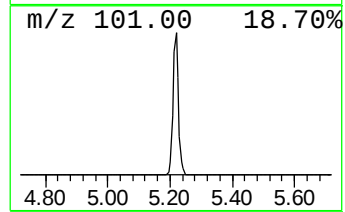
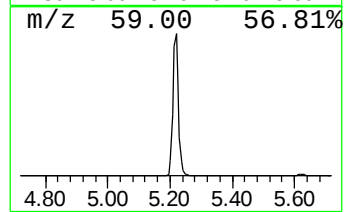
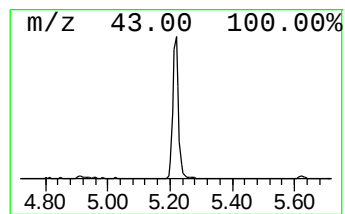
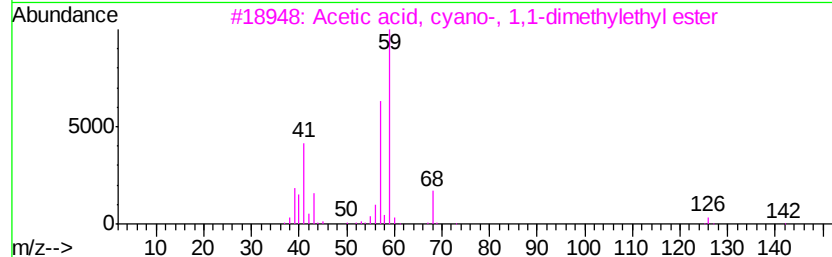
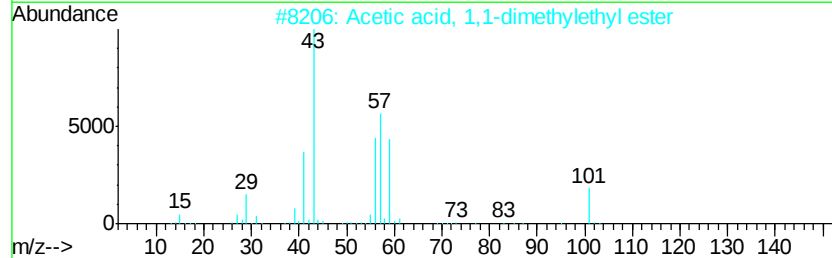
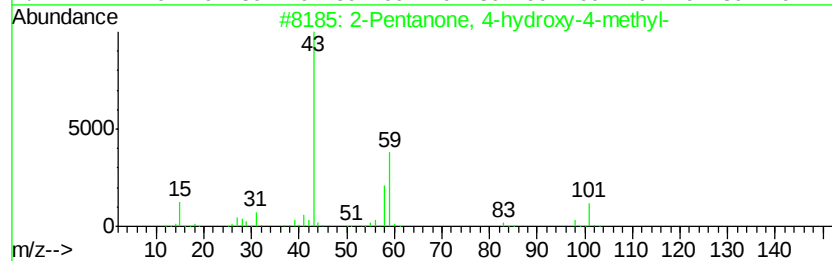
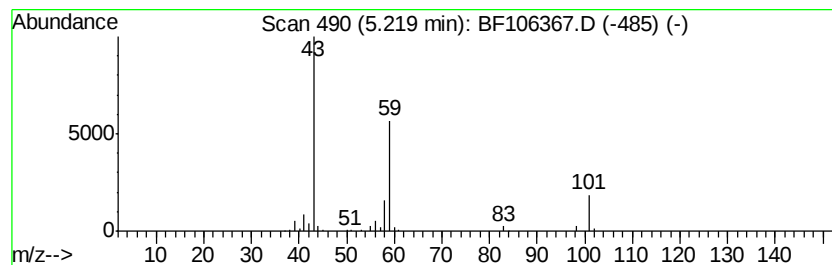
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	8.07 ng	332331	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	37
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33



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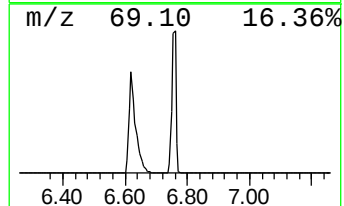
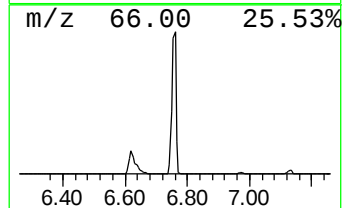
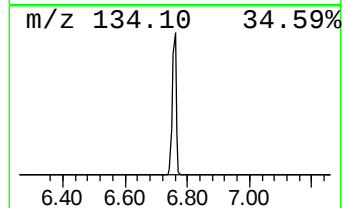
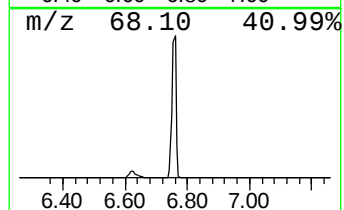
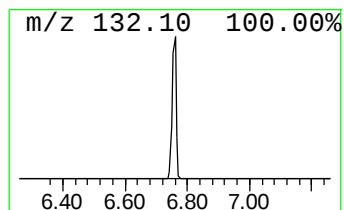
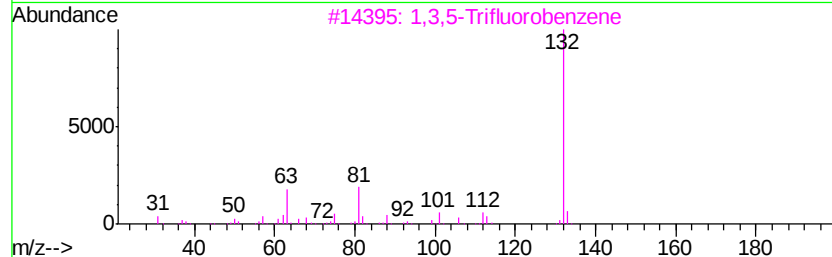
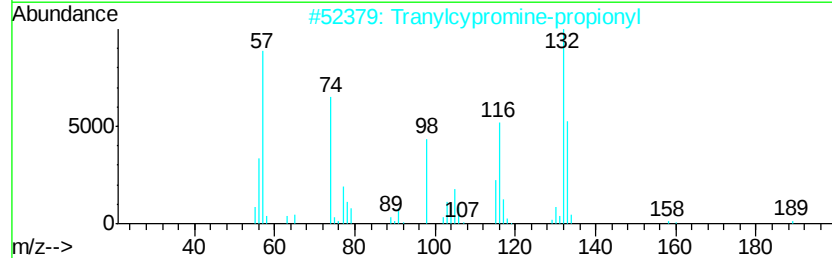
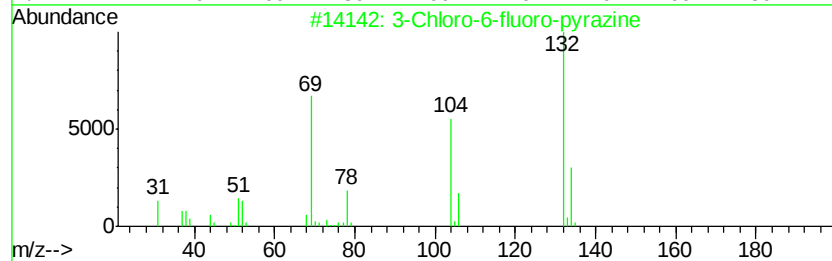
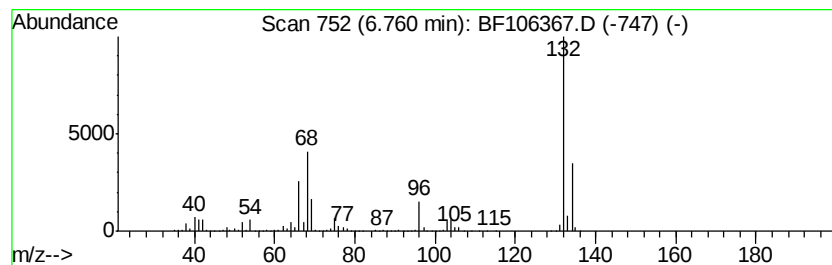
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Peak Number 2 unknown6.76 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	76.50 ng	3149180	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	8.1	ng	332331	1	6.97	823343	20.0
unknown6.76	6.76	76.5	ng	3149180	1	6.97	823343	20.0