

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062818\
 Data File : BF106929.D
 Acq On : 28 Jun 2018 19:08
 Operator : JU/SJ
 Sample : PB110673BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110673BL

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	5.207	484	488	501	rBV	169983	229920	8.49%	1.237%
2	5.601	550	555	558	rBV	1858617	1856948	68.59%	9.994%
3	6.595	719	724	728	rBV	1595645	1900588	70.20%	10.228%
4	6.731	742	747	750	rBV	1897329	1934287	71.44%	10.410%
5	6.954	781	785	788	rBV	427886	387432	14.31%	2.085%
6	7.107	807	811	815	rBV	1649797	1589205	58.70%	8.553%
7	7.519	875	881	884	rBV	1198024	1271089	46.95%	6.841%
8	8.236	998	1003	1015	rBV	474580	490453	18.11%	2.639%
9	9.307	1180	1185	1189	rBV	2280088	2381178	87.95%	12.815%
10	9.995	1297	1302	1312	rBB	585512	577623	21.33%	3.109%
11	10.789	1432	1437	1440	rBV	1635677	1614962	59.65%	8.691%
12	11.483	1551	1555	1569	rBB	641976	635464	23.47%	3.420%
13	13.071	1820	1825	1828	rBV	2506402	2707451	100.00%	14.571%
14	14.136	2002	2006	2015	rBV	620068	570342	21.07%	3.069%
15	15.671	2261	2267	2277	rVB	306505	434539	16.05%	2.339%

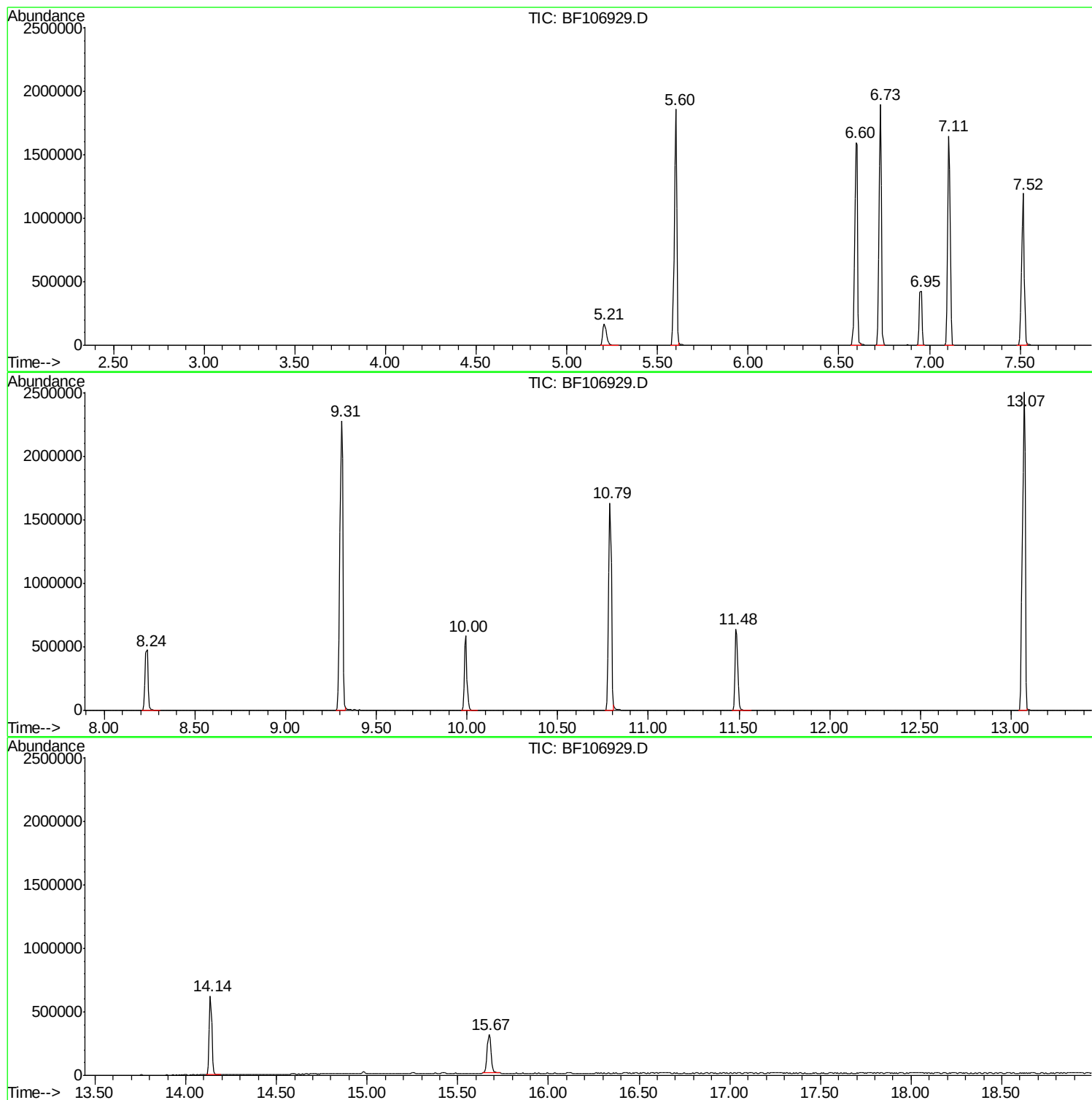
Sum of corrected areas: 18581481

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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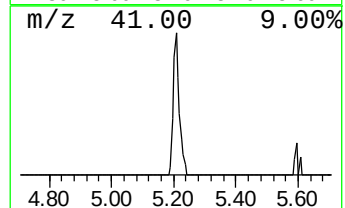
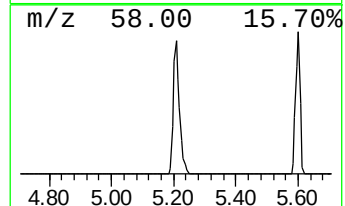
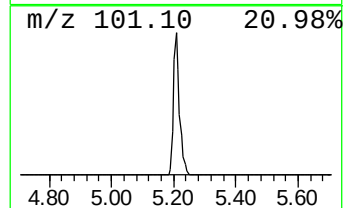
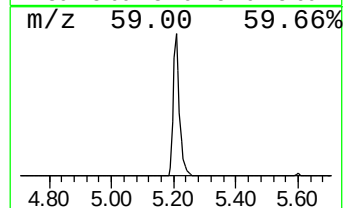
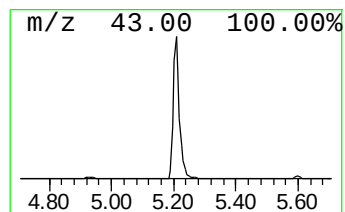
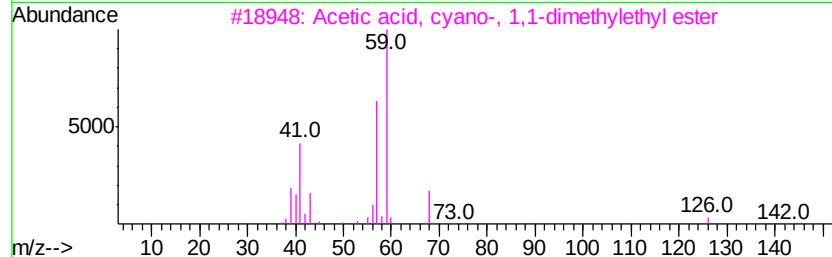
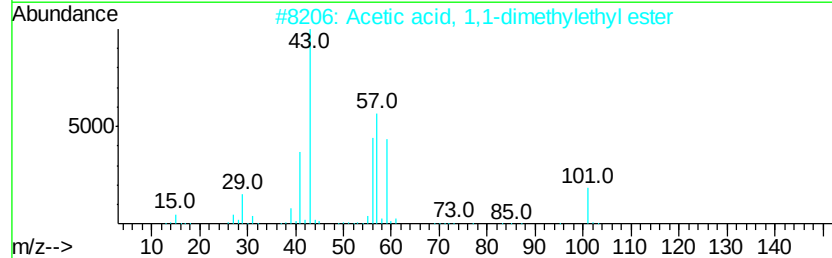
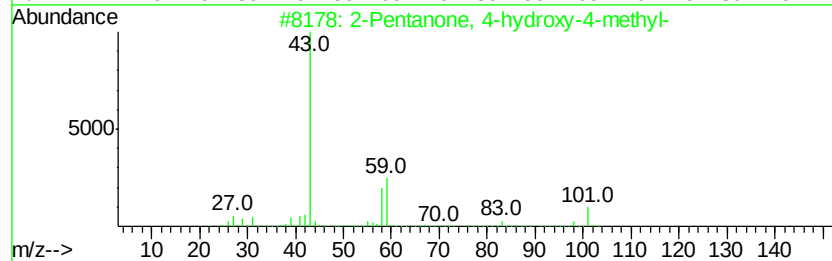
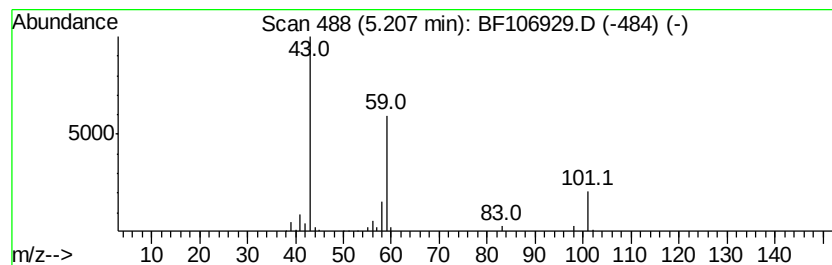
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	11.87 ng	229920	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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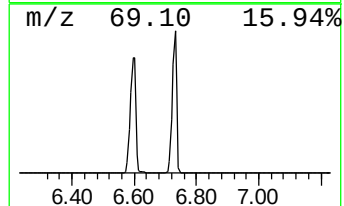
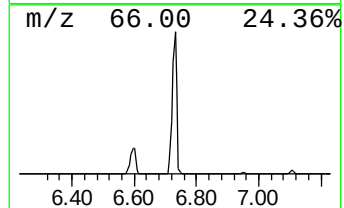
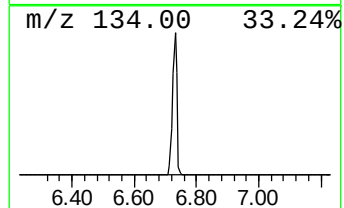
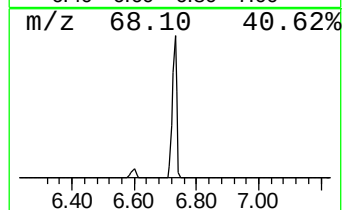
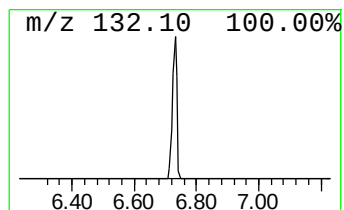
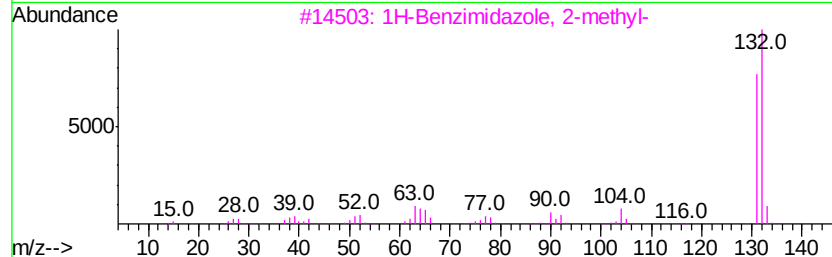
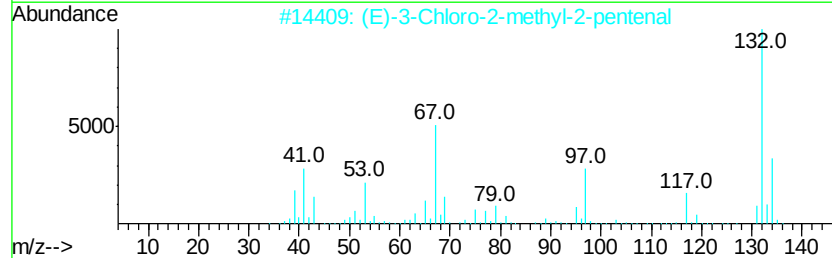
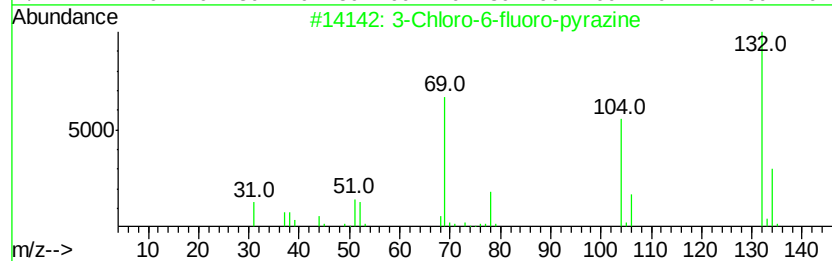
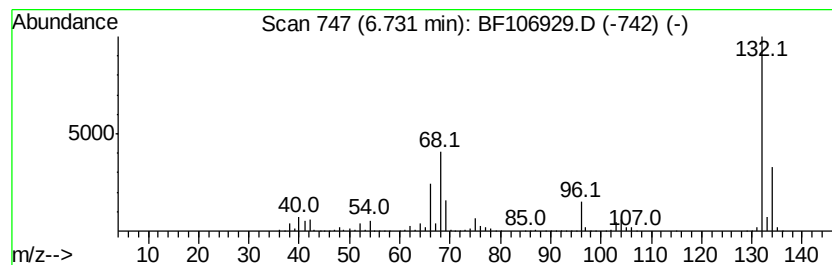
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	99.85 ng	1934290	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
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...	5.21	11.9	ng	229920	1	6.95	387432	20.0
unknown6.73	6.73	99.8	ng	1934290	1	6.95	387432	20.0