

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111915\
 Data File : BF083055.D
 Acq On : 20 Nov 2015 1:04
 Operator : UM/IZ
 Sample : PB86743BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB86743BL

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:\HPCHEM1\BNA_F\METHODS\8270-BF111915.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	4.853	238	242	252	rBV	1095098	1629087	45.41%	5.478%
2	5.299	279	281	284	rBV	2860415	2736729	76.28%	9.203%
3	6.350	370	373	375	rBV	3114408	3032890	84.54%	10.199%
4	6.465	381	383	386	rBV	2406662	2751607	76.70%	9.253%
5	6.693	401	403	406	rVB	536533	721518	20.11%	2.426%
6	6.853	415	417	420	rBV	2237365	2130747	59.39%	7.165%
7	7.265	450	453	455	rBV	2085166	1957393	54.56%	6.582%
8	7.985	514	516	519	rBV	1028355	950371	26.49%	3.196%
9	9.070	608	611	613	rBV	3737442	3524456	98.24%	11.852%
10	9.745	667	670	672	rBV	856015	1091022	30.41%	3.669%
11	10.522	736	738	741	rBV2	1645875	2130424	59.38%	7.164%
12	11.219	796	799	801	rBV	1442639	1240801	34.59%	4.173%
13	12.808	935	938	941	rBV	3869292	3587606	100.00%	12.064%
14	13.848	1027	1029	1031	rBV	1437954	1243551	34.66%	4.182%
15	15.231	1147	1150	1153	rBV	917116	1009107	28.13%	3.393%

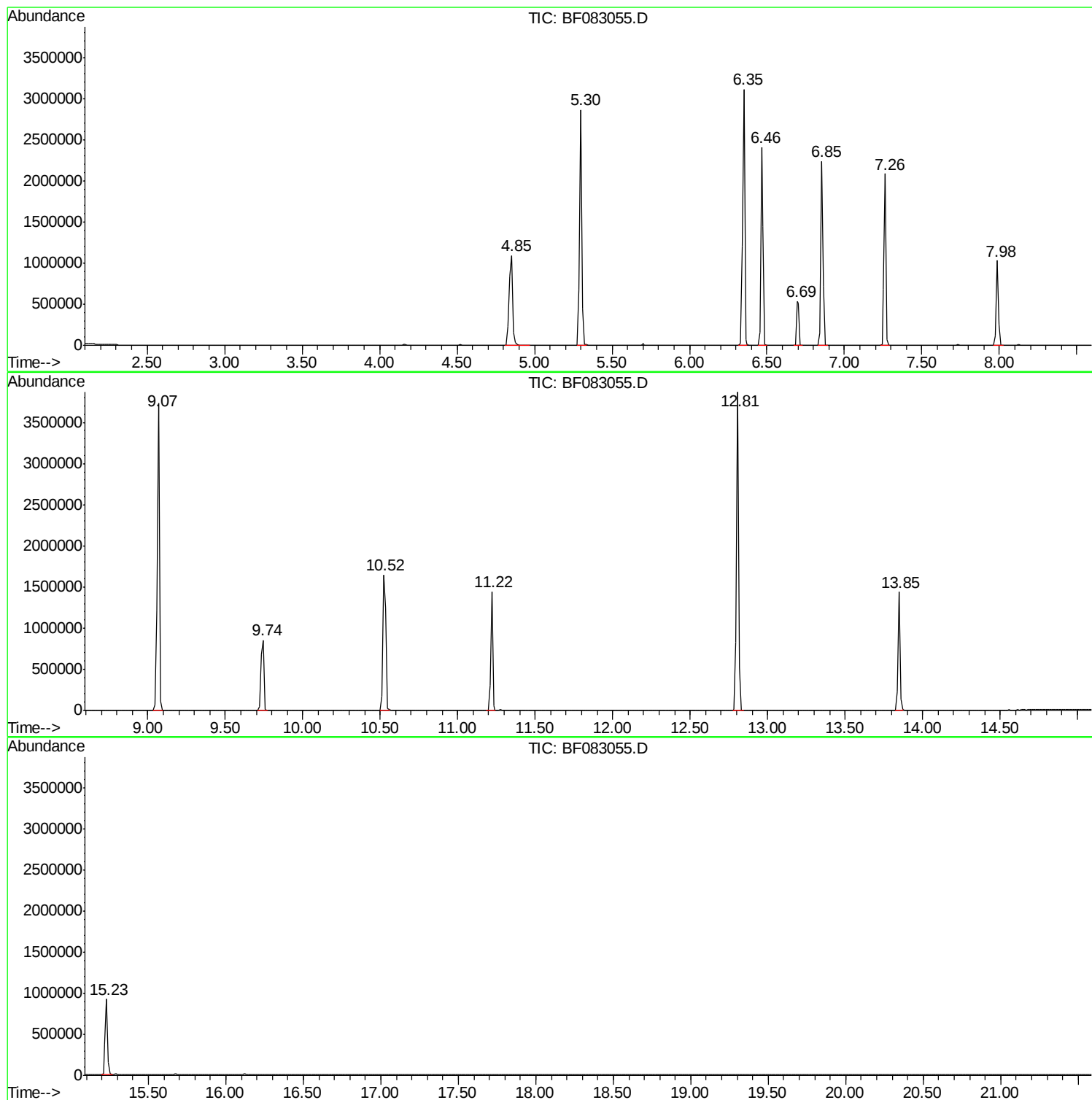
Sum of corrected areas: 29737309

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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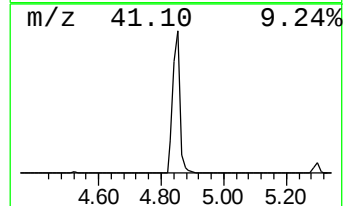
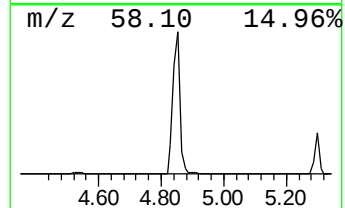
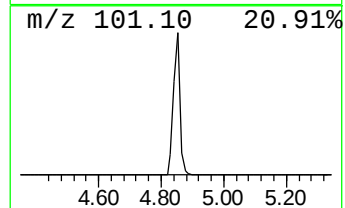
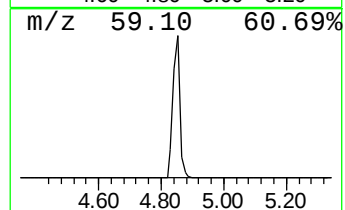
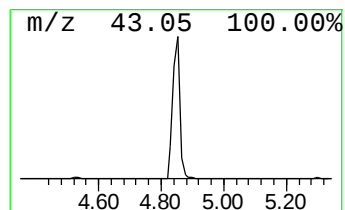
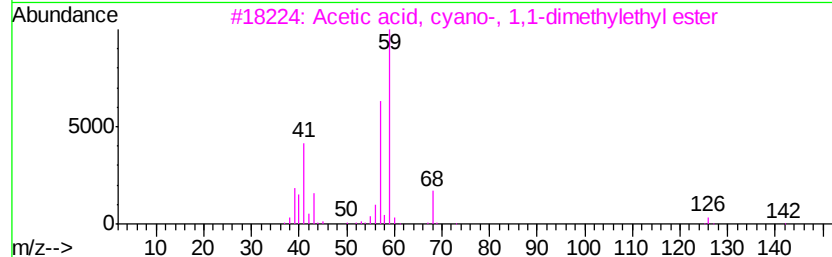
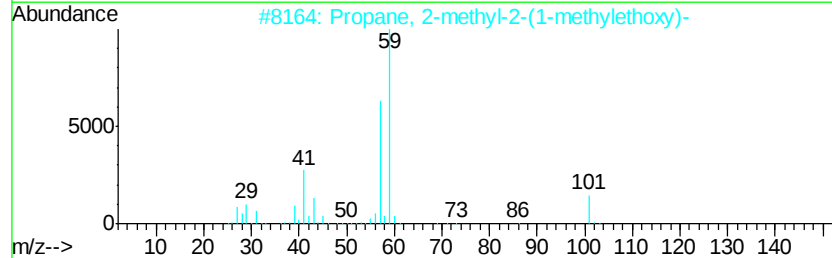
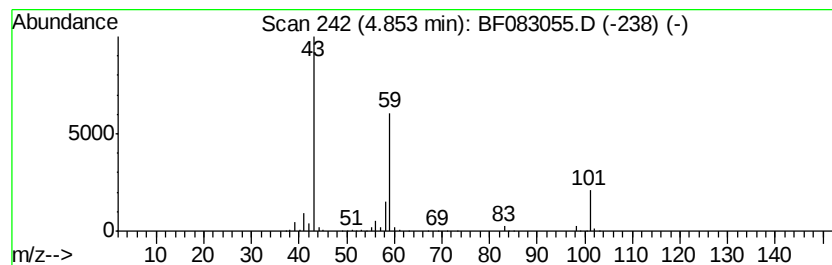
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.85	45.16 ng	1629090	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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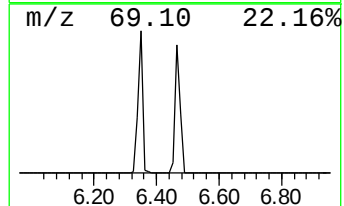
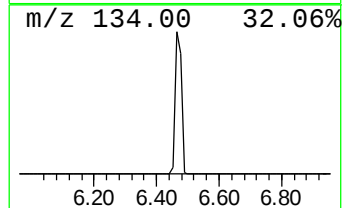
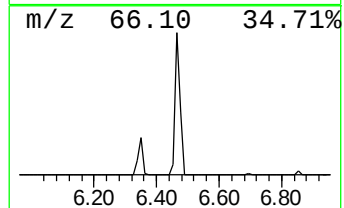
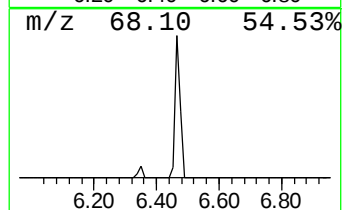
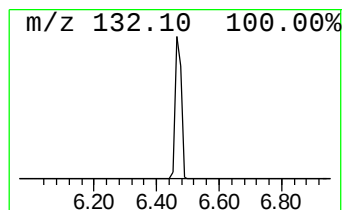
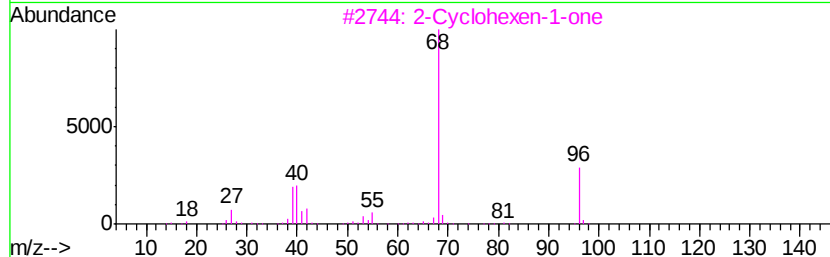
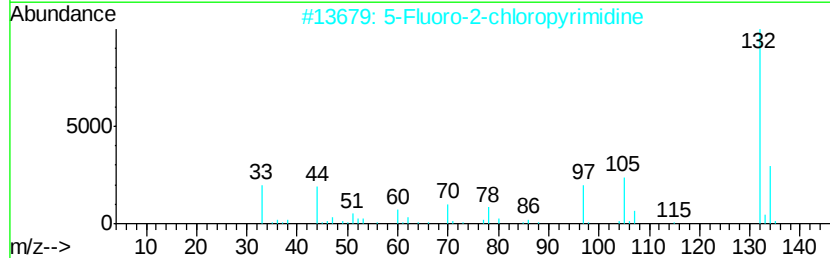
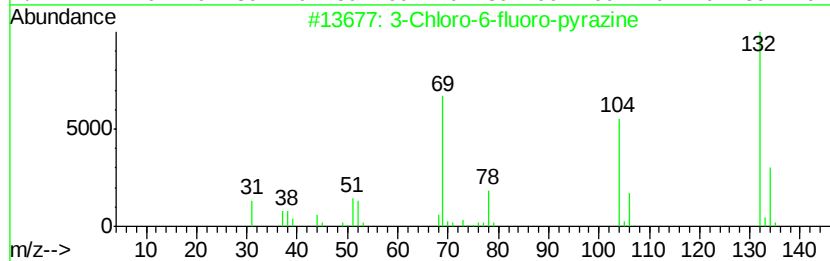
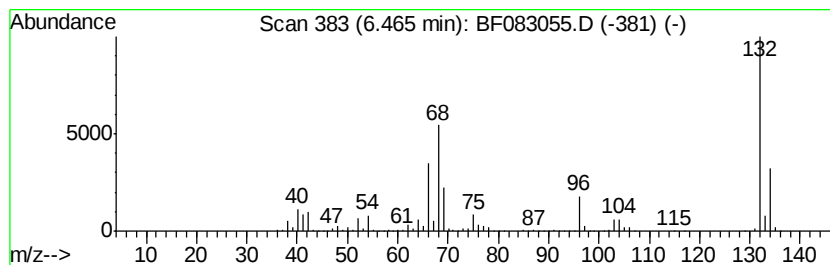
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Peak Number 2 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	76.27 ng	2751610	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.85	45.2	ng	1629090	1	6.70	721518	20.0
unknown6.46	6.46	76.3	ng	2751610	1	6.70	721518	20.0