

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090415\
Data File : BF081431.D
Acq On : 4 Sep 2015 21:25
Operator : UM/IZ
Sample : PB85409BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB85409BL

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-BF090115.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	4.410	244	247	259	rBV	515487	846877	47.39%	5.550%
2	4.913	288	291	303	rBV	1536632	1579684	88.40%	10.352%
3	6.021	386	388	391	rBV	1230809	1508065	84.39%	9.883%
4	6.124	395	397	399	rBV	1813198	1559541	87.27%	10.220%
5	6.353	415	417	420	rBV	464631	437398	24.48%	2.866%
6	6.513	428	431	433	rBV	1409493	1229589	68.81%	8.058%
7	6.936	465	468	470	rBV	1045999	1114630	62.37%	7.304%
8	7.644	527	530	533	rBV	651605	578241	32.36%	3.789%
9	8.730	622	625	627	rBV	1917773	1787038	100.00%	11.711%
10	9.382	680	682	685	rBV	596632	594957	33.29%	3.899%
11	10.170	748	751	753	rBV	960883	985708	55.16%	6.460%
12	10.845	808	810	813	rBV2	549797	612004	34.25%	4.011%
13	12.434	946	949	952	rBV	1659142	1570991	87.91%	10.295%
14	13.451	1036	1038	1041	rBV	433205	509038	28.49%	3.336%
15	14.765	1150	1153	1155	rBV	306607	345781	19.35%	2.266%

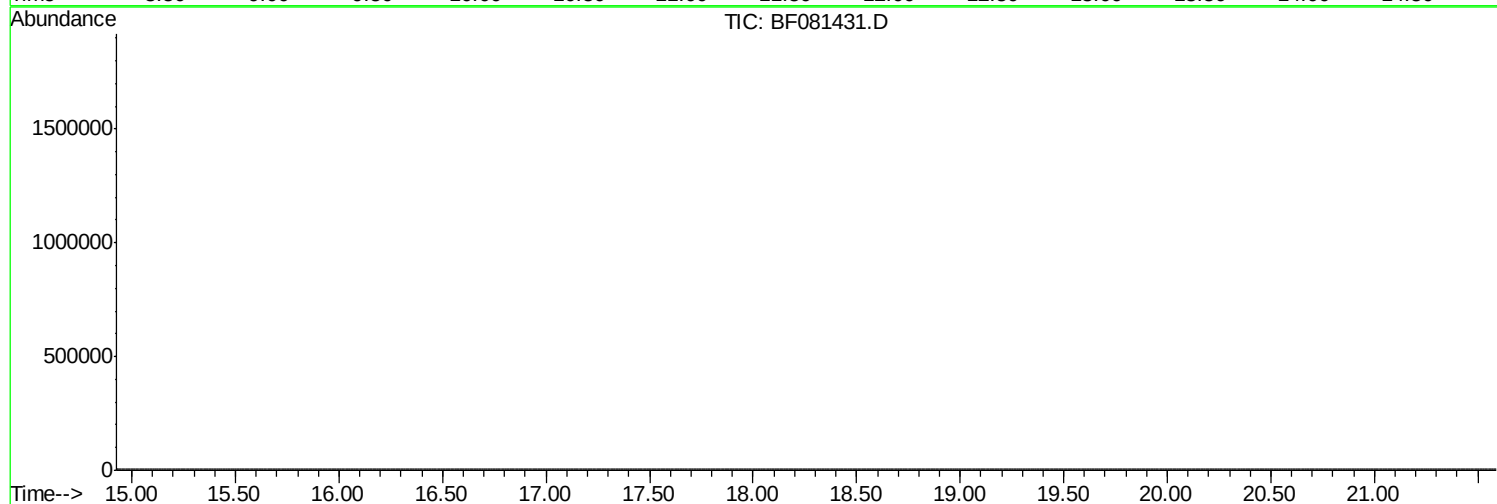
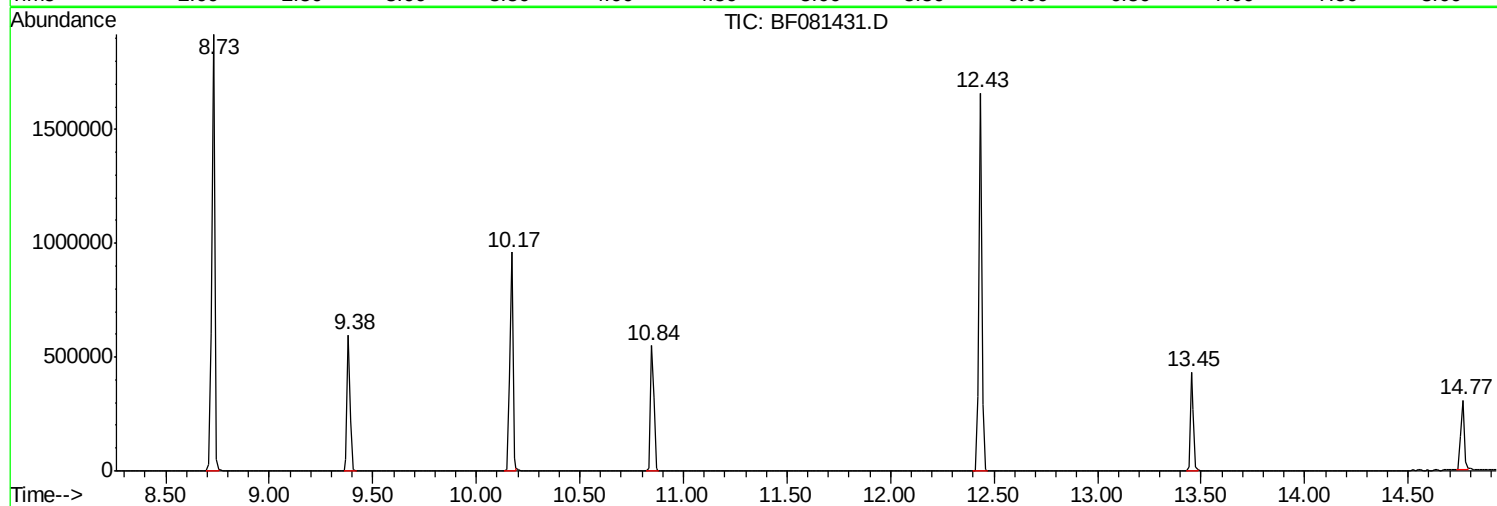
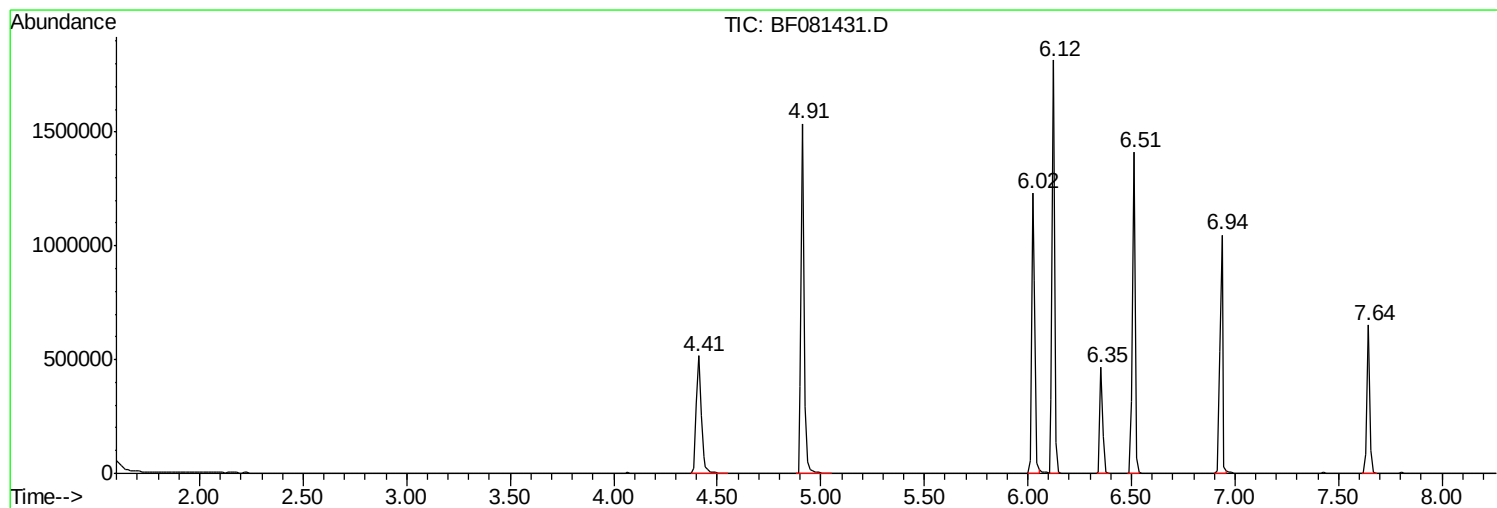
Sum of corrected areas: 15259542

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.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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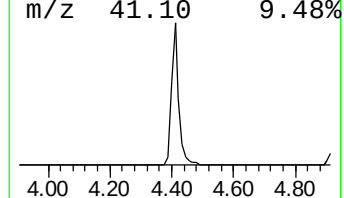
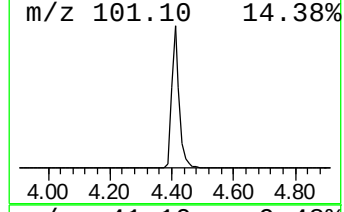
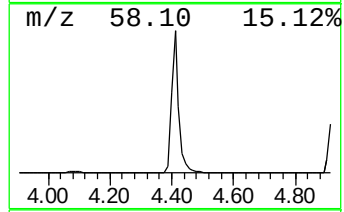
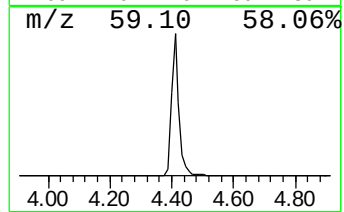
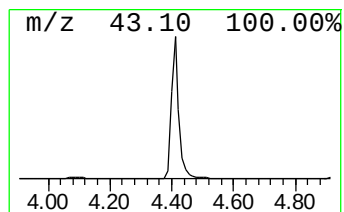
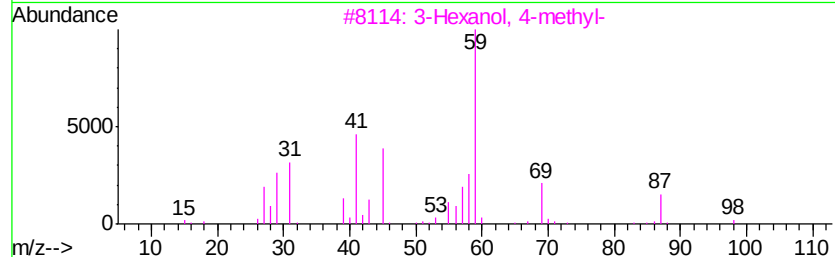
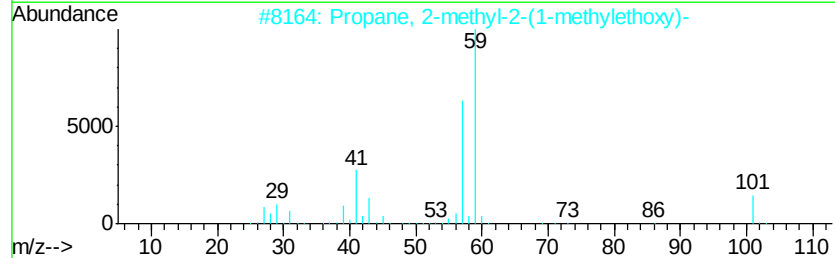
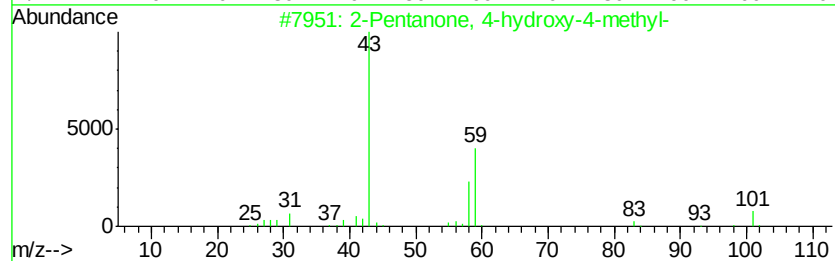
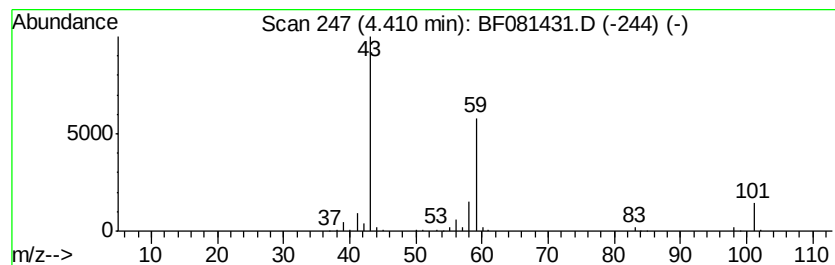
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	38.72 ng	846877	1,4-Dichlorobenzene-d4	6.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-			116	C6H12O2	000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...			116	C7H16O	017348-59-3	36
3	3-Hexanol, 4-methyl-			116	C7H16O	000615-29-2	33
4	2-Hexanol, 2-methyl-			116	C7H16O	000625-23-0	33
5	Acetic acid, cyano-, 1,1-dimethy...			141	C7H11NO2	001116-98-9	32



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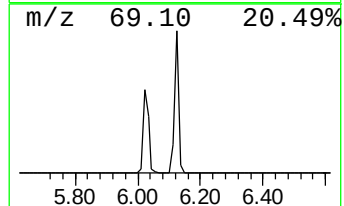
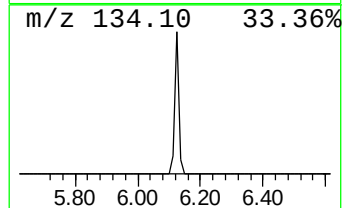
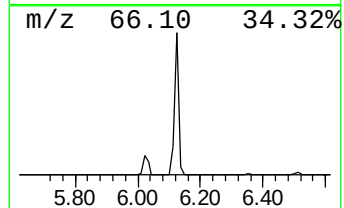
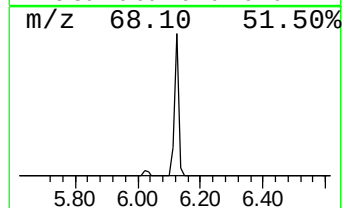
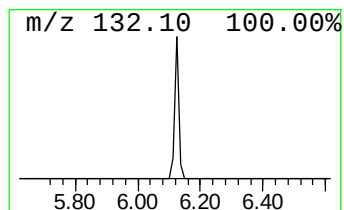
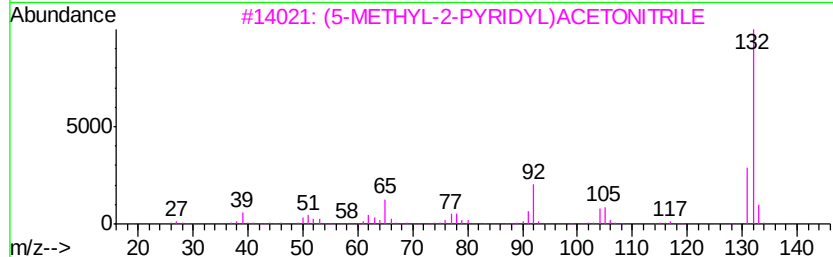
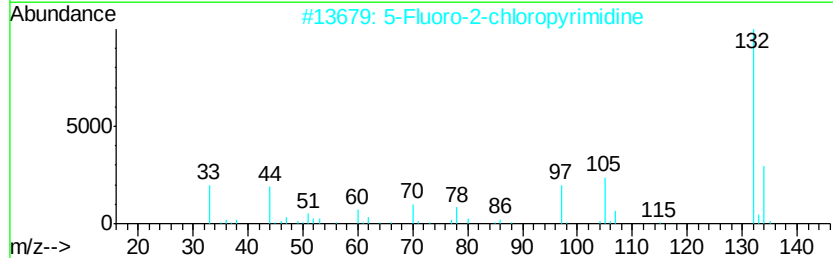
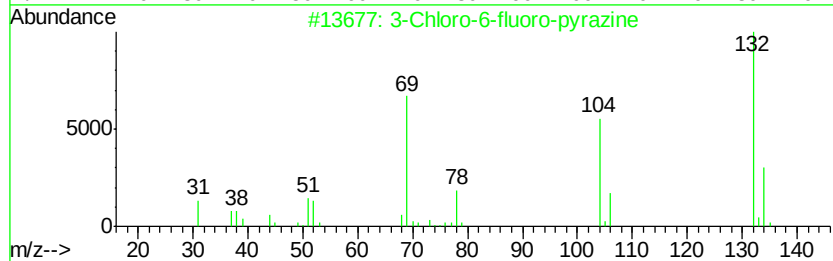
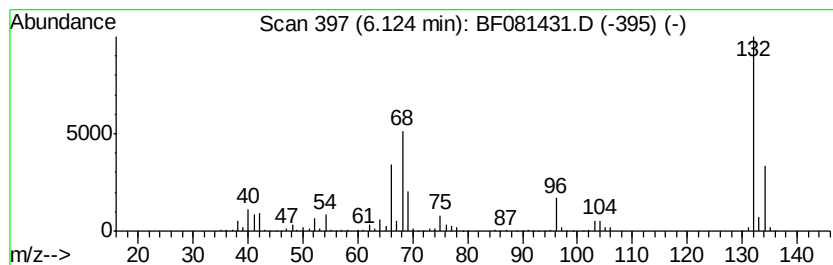
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Peak Number 2 unknown6.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	71.31 ng	1559540	1,4-Dichlorobenzene-d4	6.35

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	27
3		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
4		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
5		Tranlylcypromine-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...	4.41	38.7	ng	846877	1	6.35	437398	20.0
unknown6.12	6.12	71.3	ng	1559540	1	6.35	437398	20.0