

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090415\
Data File : BF081428.D
Acq On : 4 Sep 2015 19:52
Operator : UM/IZ
Sample : PB85375BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB85375BL

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	243	247	256	rBV	547927	942702	49.32%	5.876%
2	4.913	289	291	302	rBV	1628969	1727526	90.37%	10.769%
3	6.033	386	389	391	rBV	1344157	1816173	95.01%	11.321%
4	6.124	395	397	399	rBV	1861796	1627582	85.14%	10.146%
5	6.353	415	417	420	rBV	472266	442212	23.13%	2.757%
6	6.513	429	431	433	rBV	1499361	1319735	69.04%	8.227%
7	6.936	465	468	470	rBV	1146241	1195758	62.55%	7.454%
8	7.645	528	530	533	rBV	658798	573280	29.99%	3.574%
9	8.730	622	625	627	rBV	1997989	1911582	100.00%	11.916%
10	9.382	680	682	685	rBV	515884	539039	28.20%	3.360%
11	10.171	748	751	753	rBV	988730	987846	51.68%	6.158%
12	10.845	808	810	813	rBV	475357	567745	29.70%	3.539%
13	12.434	946	949	952	rBV	1558626	1555806	81.39%	9.698%
14	13.451	1036	1038	1041	rBV	402080	493384	25.81%	3.076%
15	14.765	1150	1153	1155	rBV	311596	341983	17.89%	2.132%

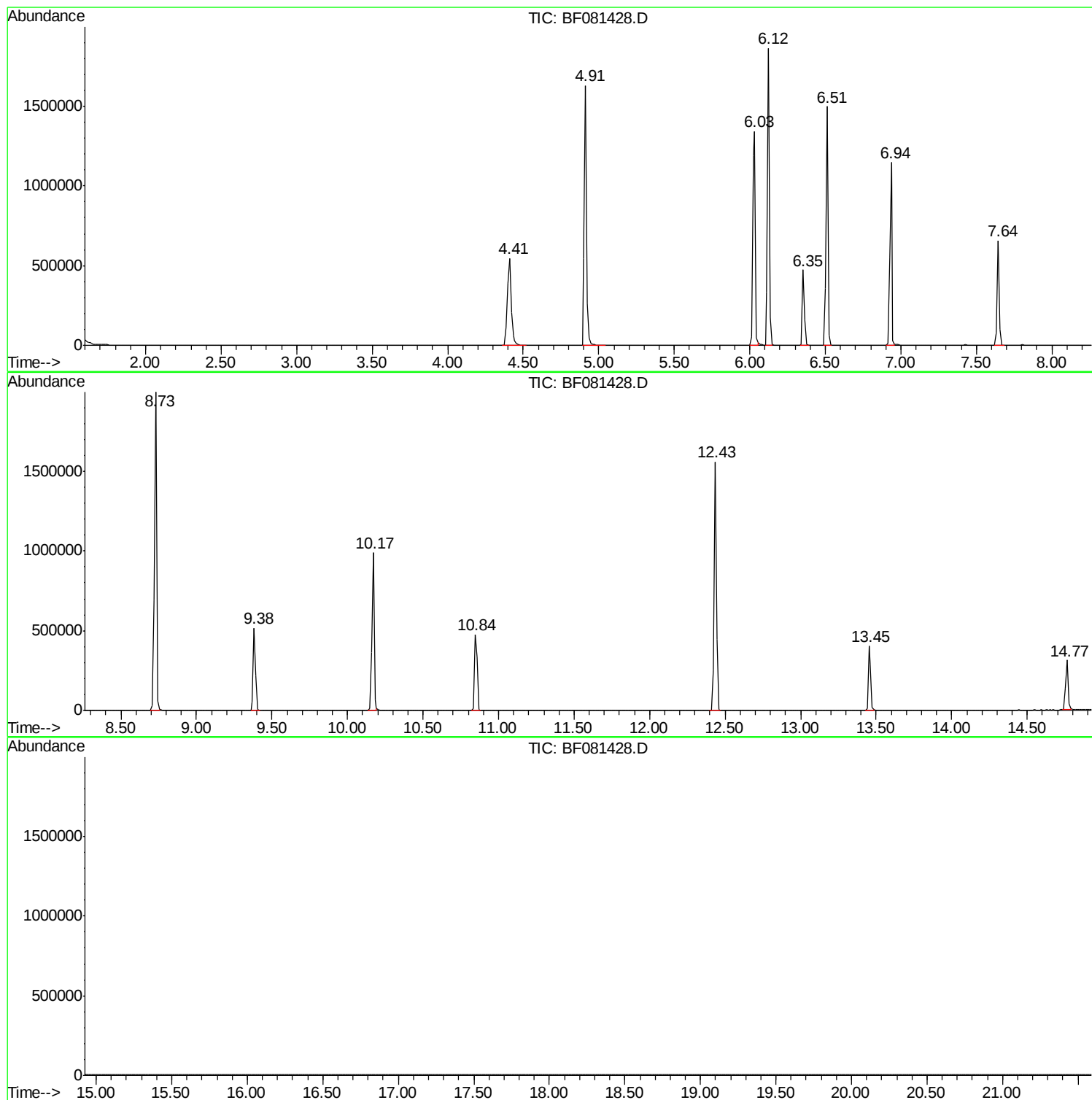
Sum of corrected areas: 16042353

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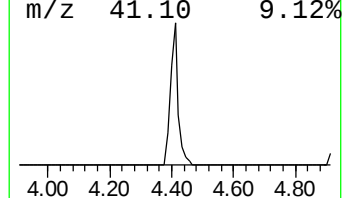
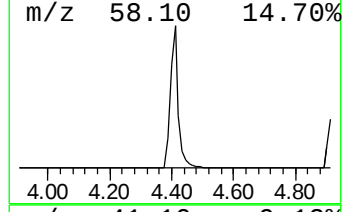
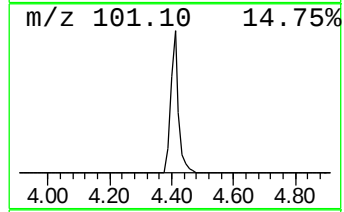
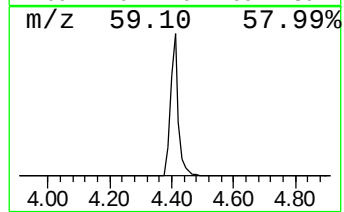
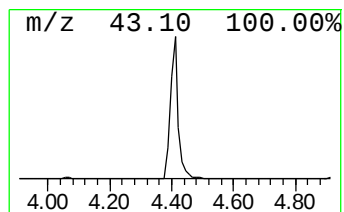
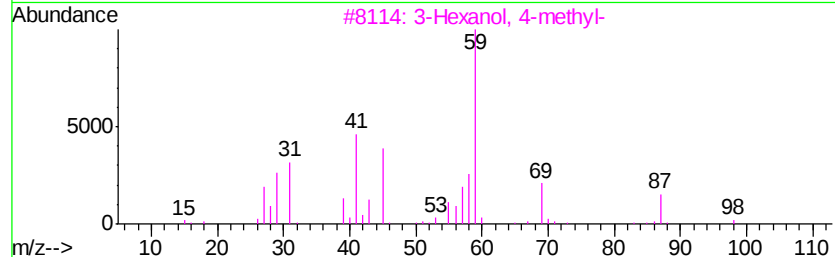
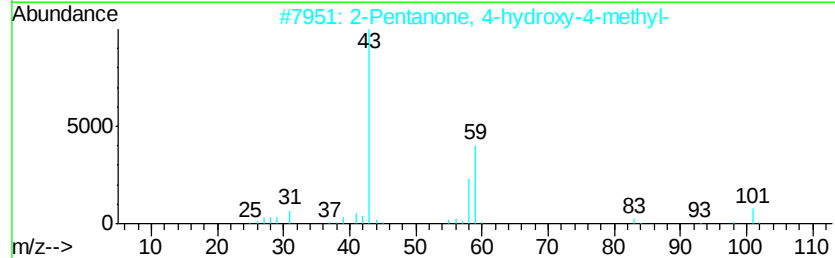
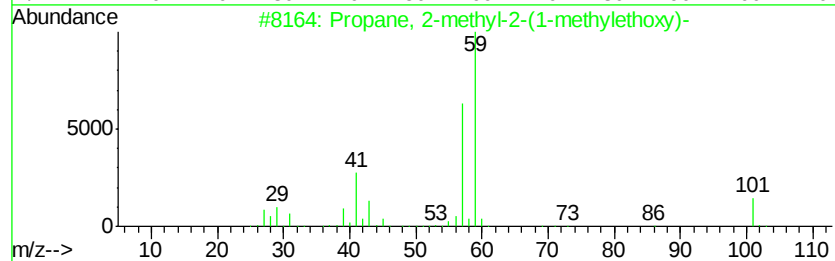
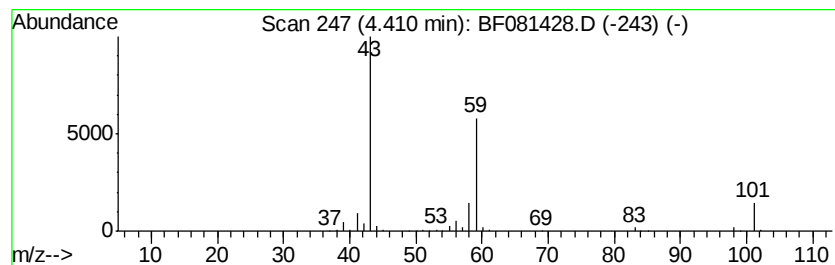
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 2-methyl-2-(1-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	42.64 ng	942702	1,4-Dichlorobenzene-d4	6.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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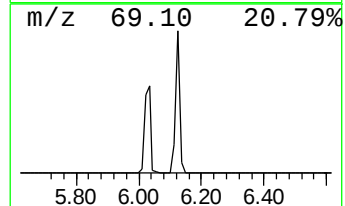
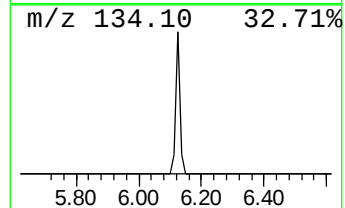
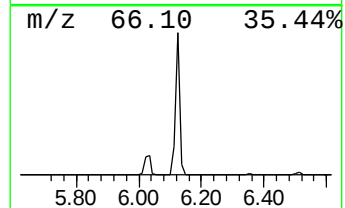
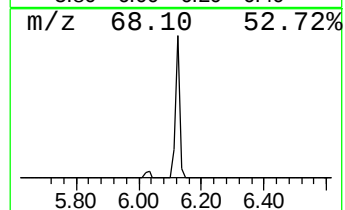
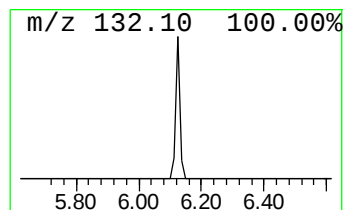
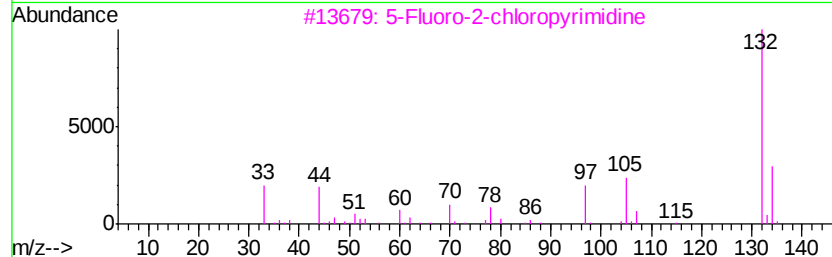
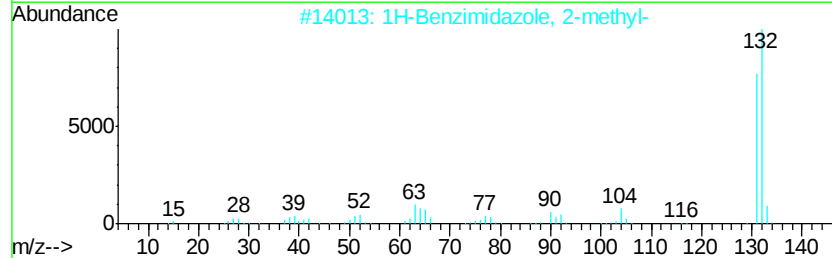
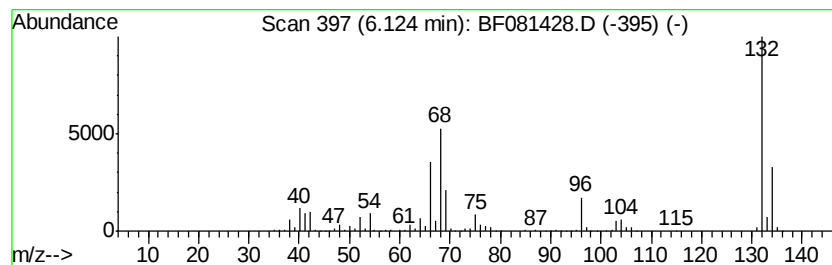
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Peak Number 2 unknown6.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	73.61 ng	1627580	1,4-Dichlorobenzene-d4	6.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	1H-Benzimidazole, 2-methyl-			132	C8H8N2	000615-15-6	25
3	5-Fluoro-2-chloropyrimidine			132	C4H2ClFN2	062802-42-0	16
4	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	10
5	(5-METHYL-2-PYRIDYL)ACETONITRILE			132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2-methyl...	4.41	42.6	ng	942702	1	6.35	442212	20.0
unknown6.12	6.12	73.6	ng	1627580	1	6.35	442212	20.0