

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030216\
 Data File : BF085111.D
 Acq On : 2 Mar 2016 11:23
 Operator : SJ/IZ
 Sample : PB88627BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88627BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030116.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	5.201	252	255	264	rBV	709230	892485	14.65%	1.936%
2	5.602	286	290	292	rBV	4268138	5016193	82.37%	10.883%
3	6.607	375	378	380	rBV	4398953	4657393	76.47%	10.104%
4	6.756	388	391	393	rBV	4749620	4939727	81.11%	10.717%
5	6.985	409	411	414	rBV	1235975	1085752	17.83%	2.356%
6	7.145	422	425	427	rBV	4252427	3779699	62.06%	8.200%
7	7.545	457	460	463	rBV	2747444	2900997	47.63%	6.294%
8	8.276	521	524	526	rBV	1062615	1457414	23.93%	3.162%
9	9.350	615	618	620	rBV	5625054	5812650	95.44%	12.611%
10	10.025	675	677	680	rVB	1731272	1631430	26.79%	3.539%
11	10.813	743	746	749	rBV	3414996	3432131	56.36%	7.446%
12	11.511	805	807	810	rVB	2021647	1705511	28.00%	3.700%
13	13.099	943	946	948	rBV	6208543	6090184	100.00%	13.213%
14	14.151	1035	1038	1040	rBV	1206838	1596558	26.22%	3.464%
15	15.625	1163	1167	1176	rBV	693292	1094626	17.97%	2.375%

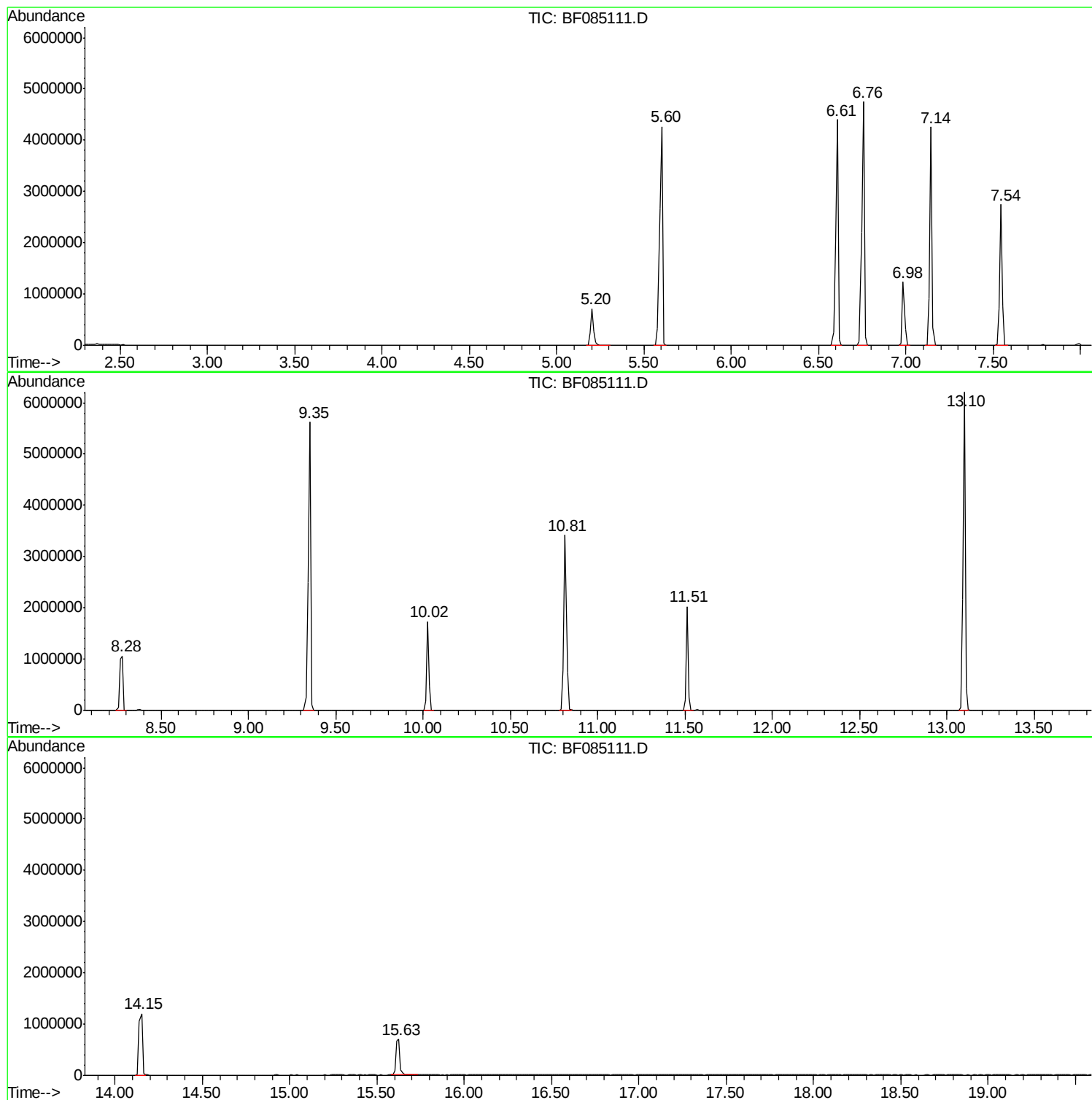
Sum of corrected areas: 46092750

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030116.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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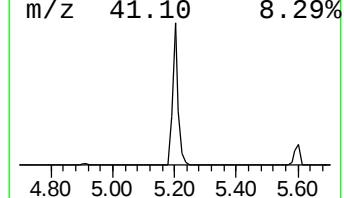
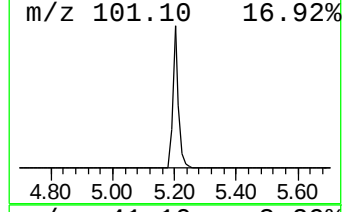
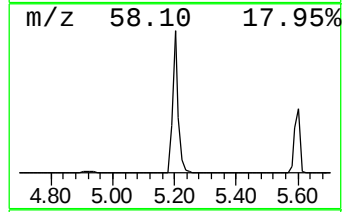
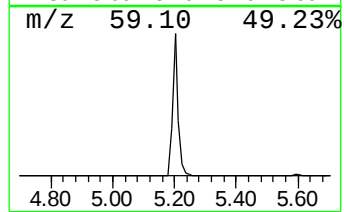
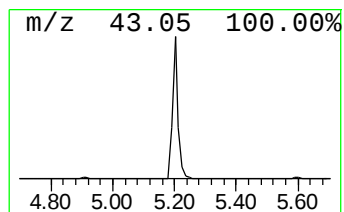
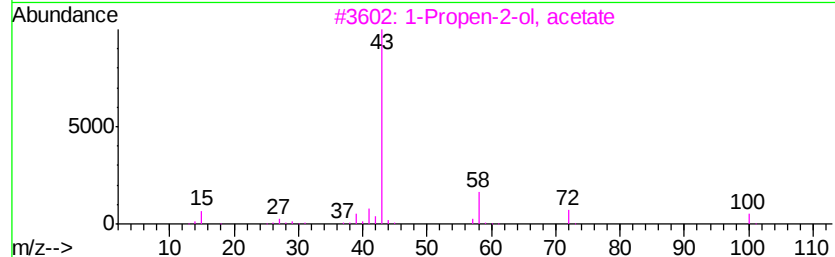
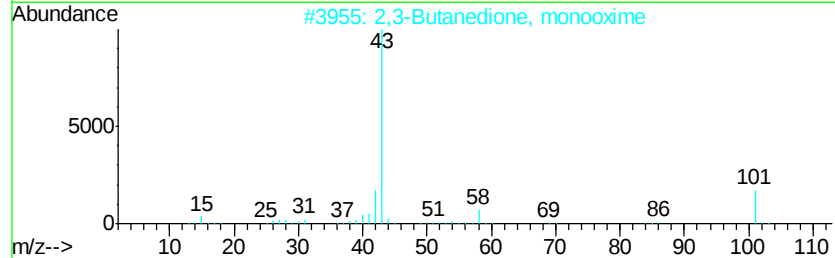
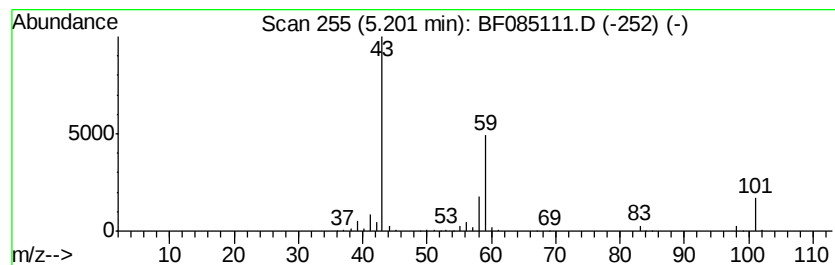
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	16.44 ng	892485	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		Butane	58	C4H10	000106-97-8	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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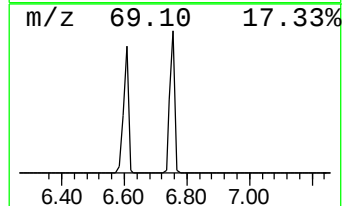
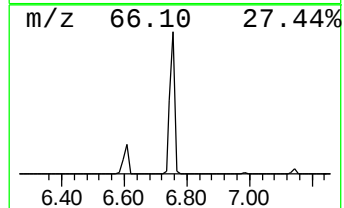
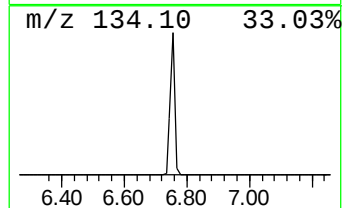
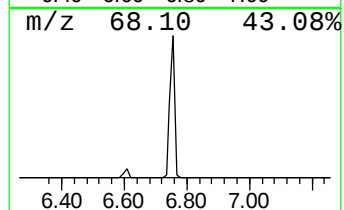
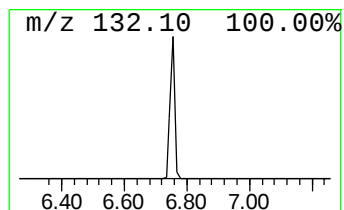
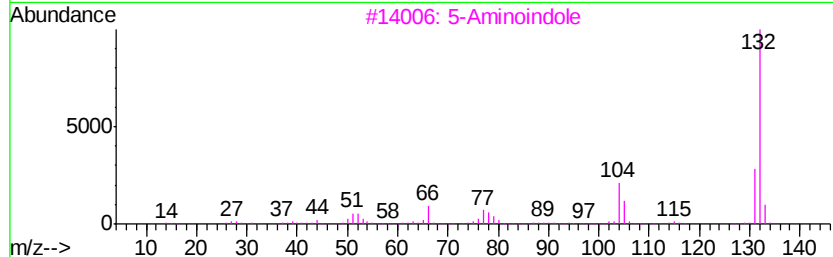
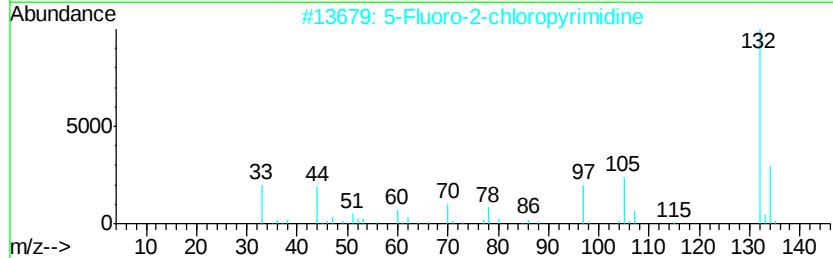
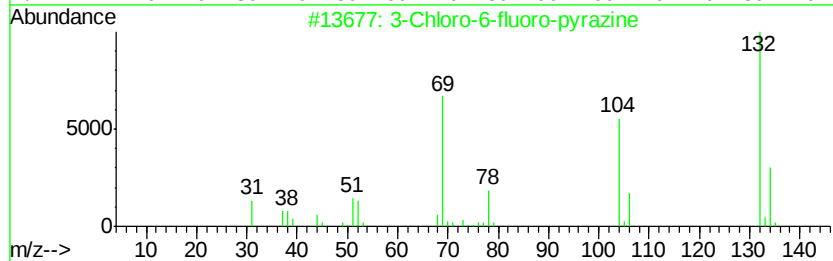
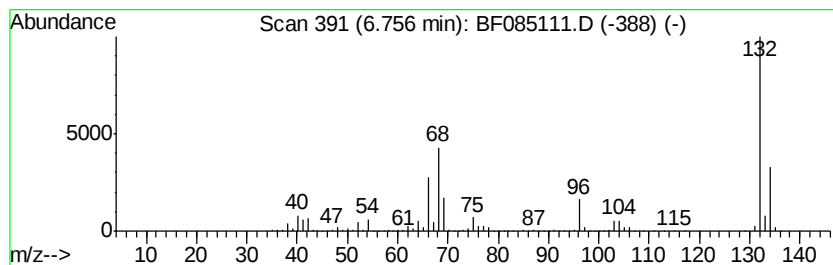
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Peak Number 2 Benzene, 1,2,4-trifluoro-6.76 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	90.99 ng	4939730	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	16.4	ng	892485	1	6.98	1085750	20.0
Benzene, 1,2,4-tr...	6.76	91.0	ng	4939730	1	6.98	1085750	20.0