

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120616\
 Data File : BF091452.D
 Acq On : 6 Dec 2016 12:01
 Operator : UM/SJ
 Sample : H5802-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FIELDBLANK

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-BF112916.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	5.019	236	239	242	rBV	475769	564306	9.81%	1.692%
2	5.441	273	276	278	rBV	2193335	2097074	36.47%	6.287%
3	6.470	363	366	368	rVB	1372383	1360434	23.66%	4.079%
4	6.607	375	378	380	rBV	3552401	3311834	57.60%	9.929%
5	6.836	396	398	400	rVB	890605	764885	13.30%	2.293%
6	6.996	409	412	414	rBV	3287606	2981666	51.86%	8.939%
7	7.407	444	448	450	rBV	2509213	2924362	50.86%	8.767%
8	8.127	508	511	513	rBV	902721	1030871	17.93%	3.090%
9	9.202	602	605	608	rBV	4803624	4683338	81.46%	14.040%
10	9.876	661	664	666	rBV	1527297	1322315	23.00%	3.964%
11	10.665	730	733	735	rBV	3434976	3489909	60.70%	10.463%
12	11.351	791	793	796	rVB	970312	1219169	21.20%	3.655%
13	12.951	929	933	935	rBV	4764572	5749552	100.00%	17.237%
14	13.991	1021	1024	1026	rBV	828998	1135285	19.75%	3.404%
15	15.408	1145	1148	1151	rBV	631945	721207	12.54%	2.162%

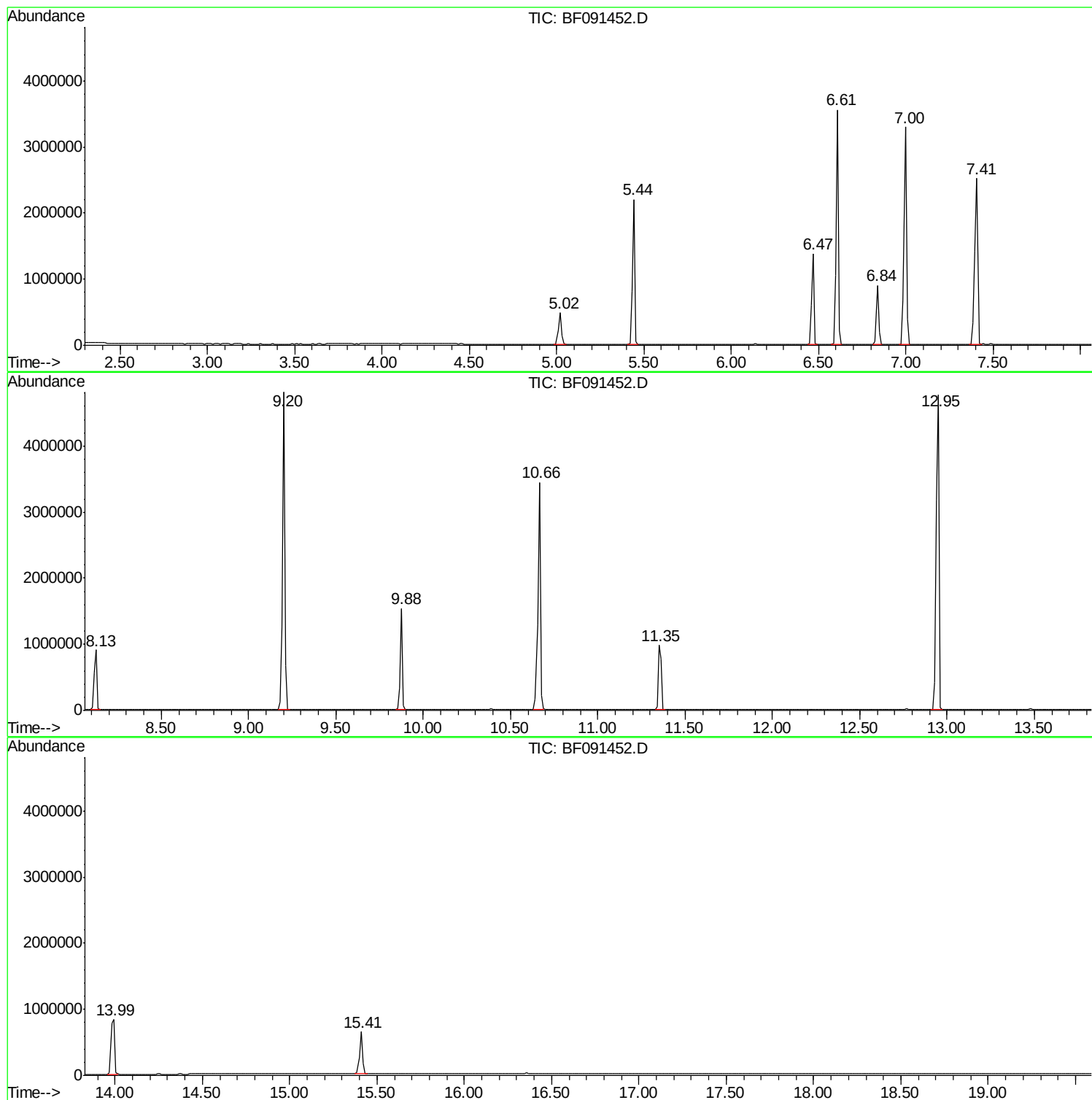
Sum of corrected areas: 33356207

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.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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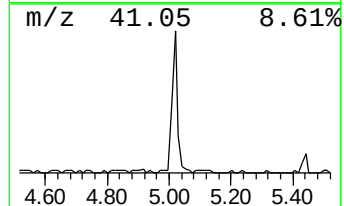
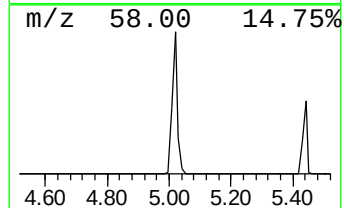
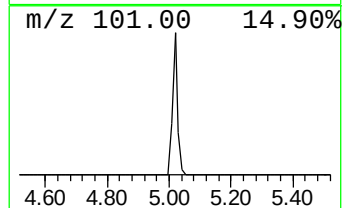
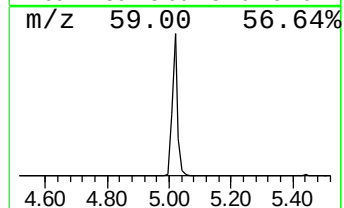
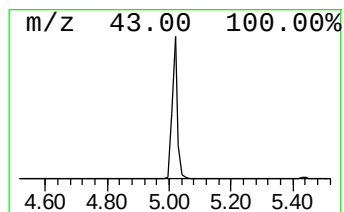
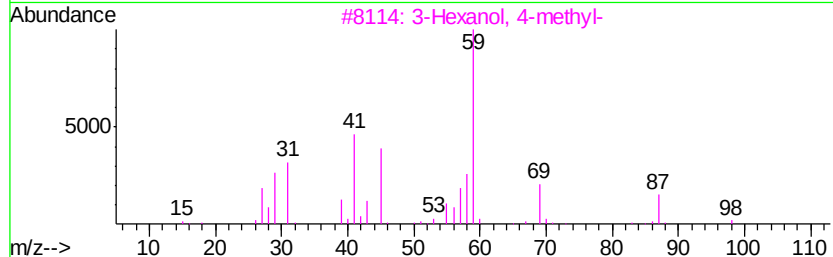
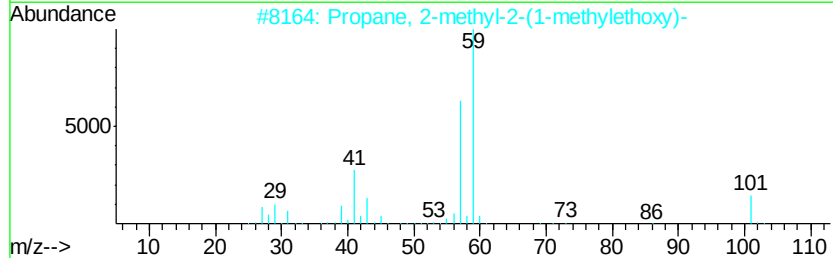
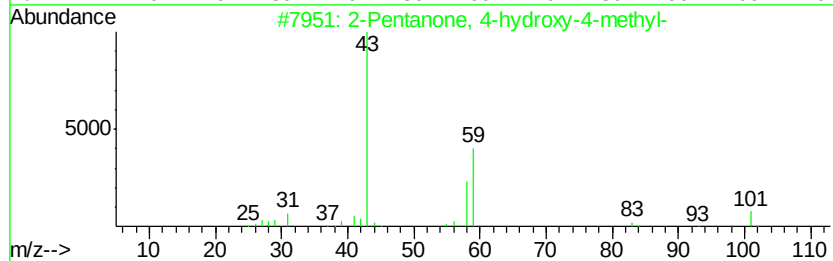
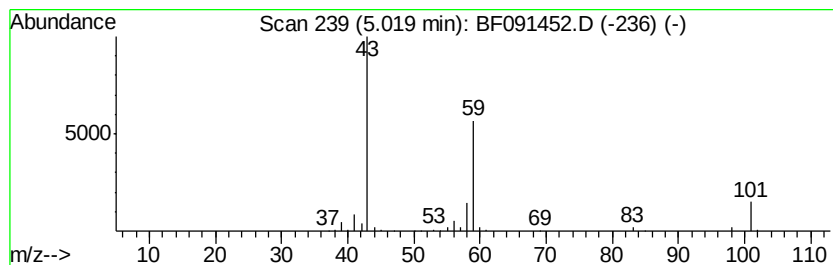
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.02	14.76 ng	564306	1,4-Dichlorobenzene-d4	6.84	
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	28
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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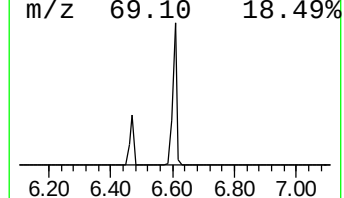
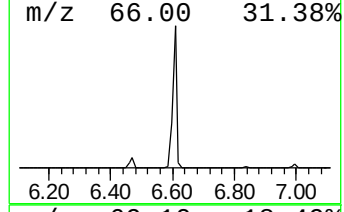
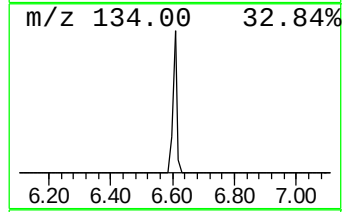
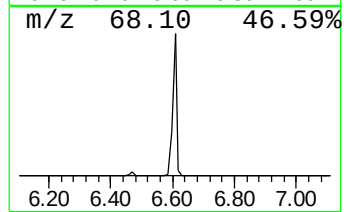
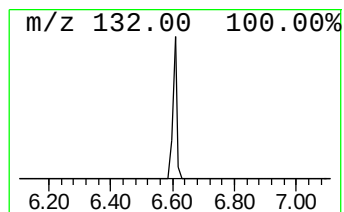
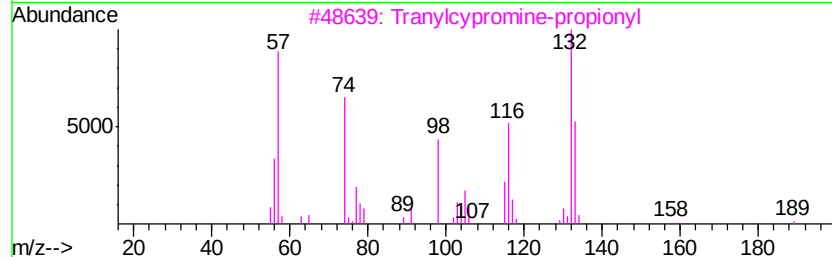
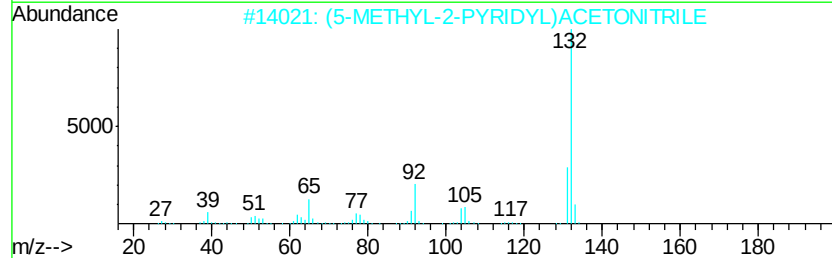
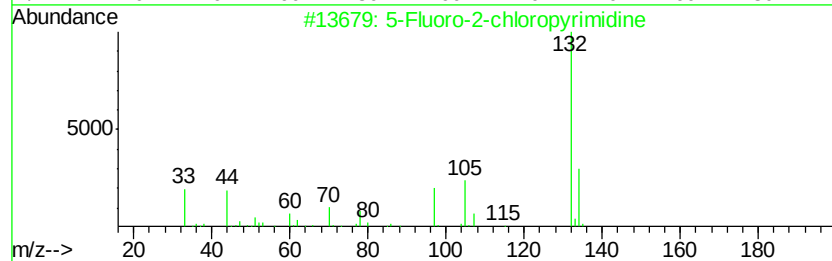
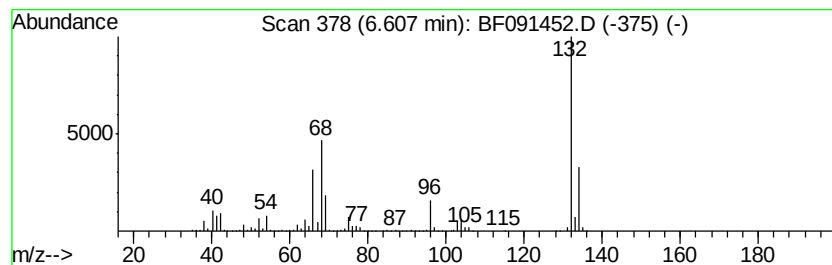
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Peak Number 2 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	86.60 ng	3311830	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2			(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5			Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.02	14.8	ng	564306	1	6.84	764885	20.0
unknown6.61	6.61	86.6	ng	3311830	1	6.84	764885	20.0