

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021017\
 Data File : BF092913.D
 Acq On : 11 Feb 2017 8:21
 Operator : SJ/MA
 Sample : I1767-02
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 020817-PRRT-F-U-M

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-BF013017.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	5.059	258	260	273	rVB	452662	469605	8.61%	1.550%
2	5.493	295	298	300	rBV	987219	1365400	25.05%	4.506%
3	6.522	386	388	394	rBV	1050653	921399	16.90%	3.041%
4	6.659	397	400	402	rBV	2855029	2490779	45.69%	8.220%
5	6.887	418	420	423	rBV	802563	896169	16.44%	2.958%
6	7.047	432	434	437	rBV	3002667	3315006	60.81%	10.940%
7	7.459	467	470	473	rBV	2317387	2780709	51.01%	9.177%
8	8.179	531	533	536	rBV	1301114	1203965	22.09%	3.973%
9	9.265	624	628	630	rBV	4111996	5451454	100.00%	17.991%
10	9.939	684	687	689	rBV	1257319	1421567	26.08%	4.691%
11	10.728	753	756	758	rBV	1790402	1973008	36.19%	6.511%
12	10.842	764	766	770	rVB	50938	83469	1.53%	0.275%
13	11.425	814	817	819	rBV	1228665	1422748	26.10%	4.695%
14	13.014	953	956	958	rBV	3558022	4560478	83.66%	15.050%
15	13.905	1031	1034	1041	rBV	37729	86184	1.58%	0.284%
16	14.054	1045	1047	1050	rVB	1076412	1014092	18.60%	3.347%
17	15.505	1171	1174	1182	rBV	604970	845265	15.51%	2.790%

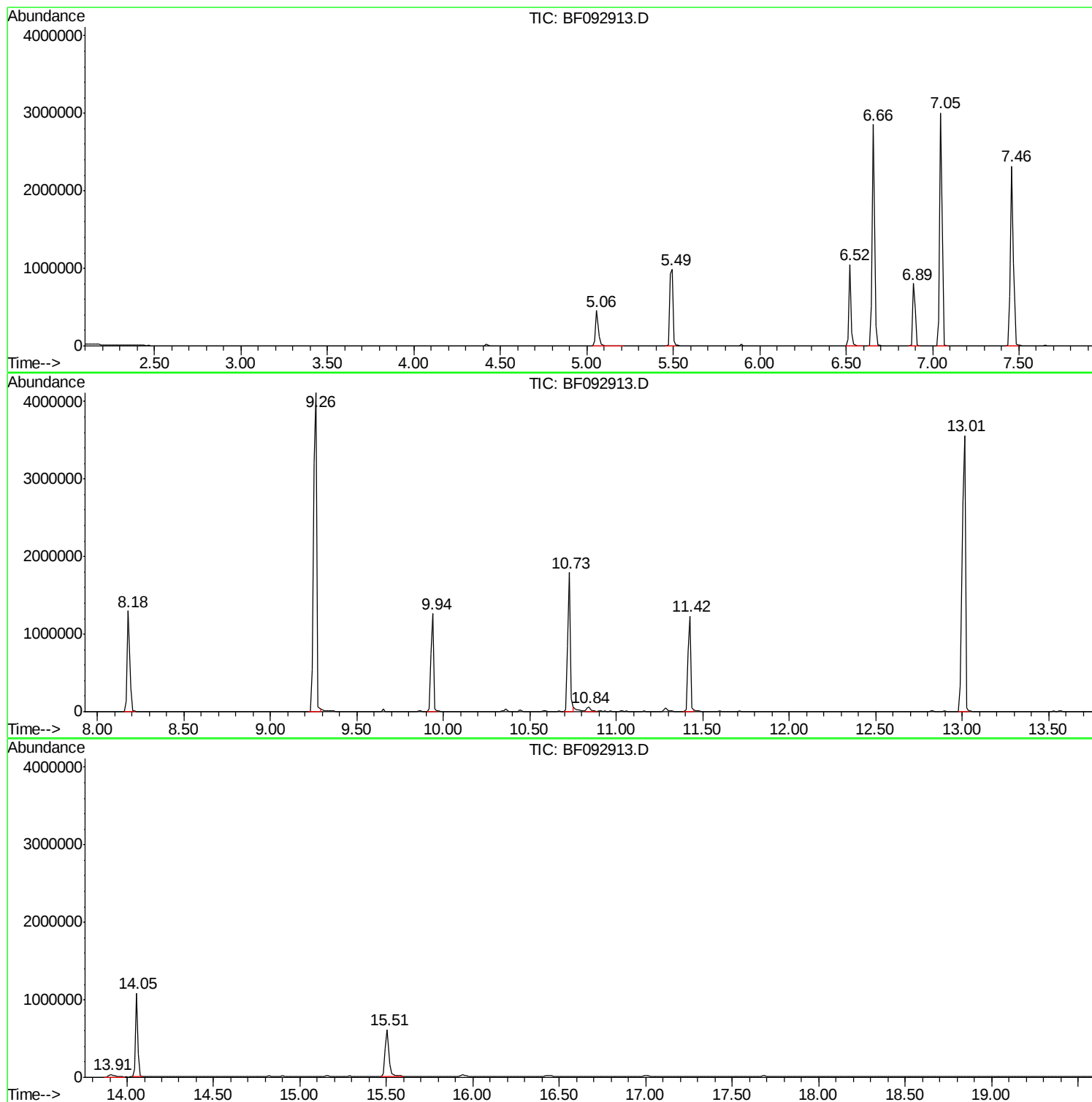
Sum of corrected areas: 30301297

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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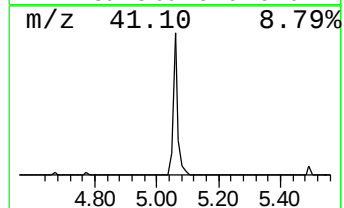
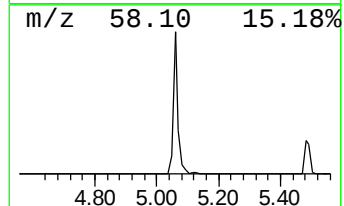
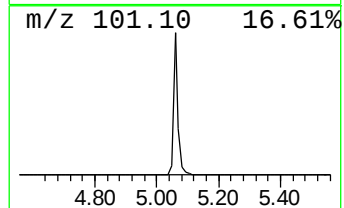
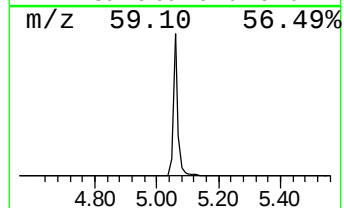
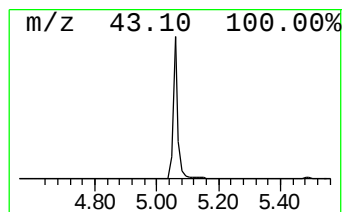
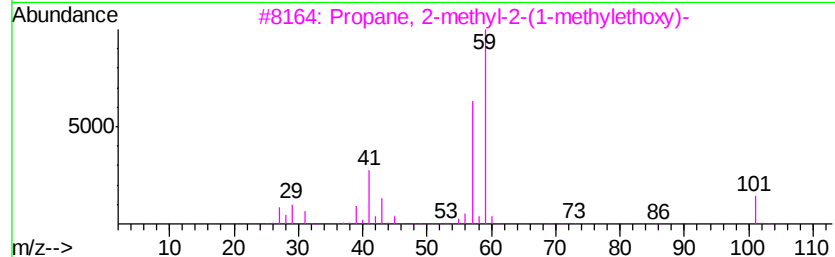
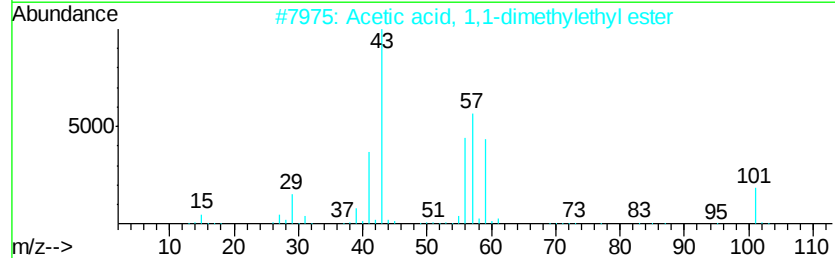
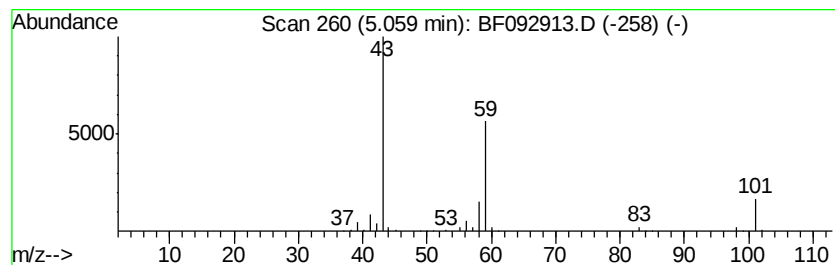
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	10.48 ng	469605	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5		3-Octanol	130	C8H18O	000589-98-0	33



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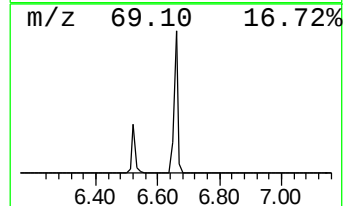
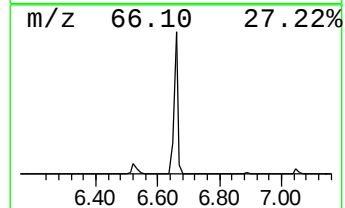
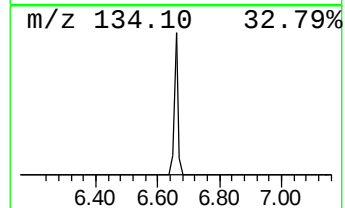
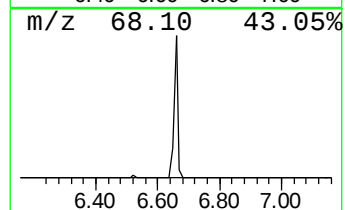
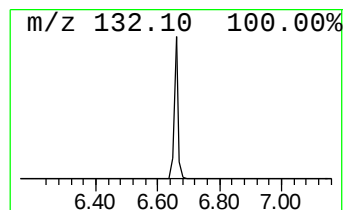
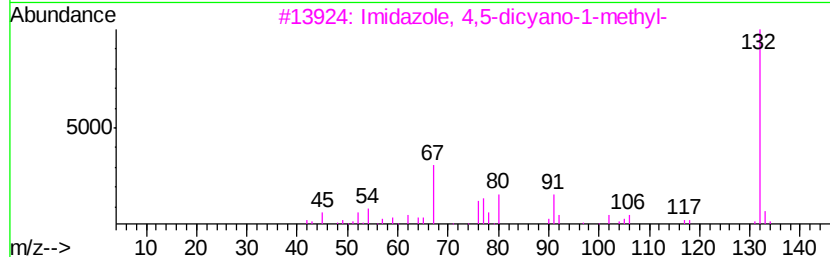
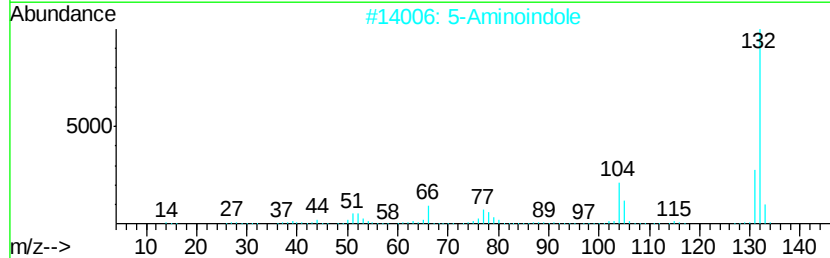
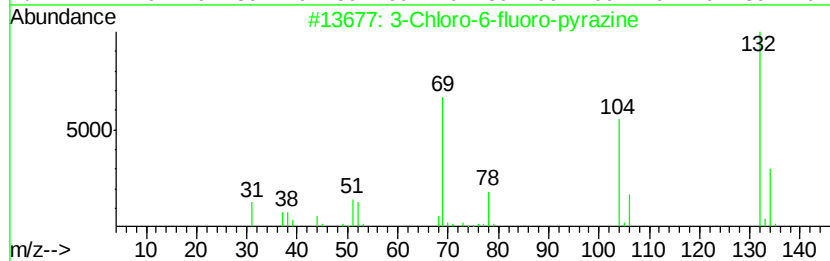
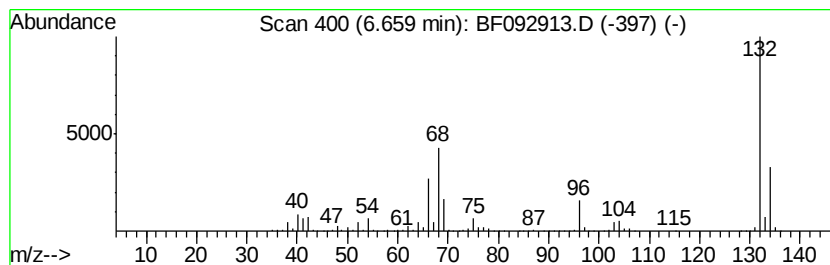
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Peak Number 2 unknown6.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	55.59 ng	2490780	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	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...	5.06	10.5	ng	469605	1	6.89	896169	20.0
unknown6.66	6.66	55.6	ng	2490780	1	6.89	896169	20.0