

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101717\
 Data File : BF099601.D
 Acq On : 18 Oct 2017 00:31
 Operator : SJ/JU
 Sample : I5869-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-09-10-11-17

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-BF092317.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	2.463	52	64	73	rVB	812649	1246206	43.55%	6.848%
2	5.045	500	503	517	rBV	108755	129024	4.51%	0.709%
3	5.457	569	573	593	rBV	1078626	1031712	36.06%	5.669%
4	6.469	742	745	758	rBV	684533	629267	21.99%	3.458%
5	6.598	763	767	771	rBV	1992839	1918559	67.05%	10.542%
6	6.828	802	806	813	rBB	695268	585469	20.46%	3.217%
7	6.981	828	832	836	rBV	2054628	1941273	67.85%	10.667%
8	7.392	897	902	905	rBV	1496128	1512943	52.88%	8.313%
9	8.104	1020	1023	1027	rBV	870638	799347	27.94%	4.392%
10	9.180	1201	1206	1209	rBV	2937951	2861240	100.00%	15.722%
11	9.863	1318	1322	1326	rVB	958415	849335	29.68%	4.667%
12	10.651	1452	1456	1460	rBV	1186164	1117412	39.05%	6.140%
13	11.351	1571	1575	1578	rBV	836114	726138	25.38%	3.990%
14	12.927	1839	1843	1847	rBV	1846411	1812609	63.35%	9.960%
15	13.986	2020	2023	2027	rBV	542443	493655	17.25%	2.713%
16	15.451	2267	2272	2284	rVB	440854	544608	19.03%	2.993%

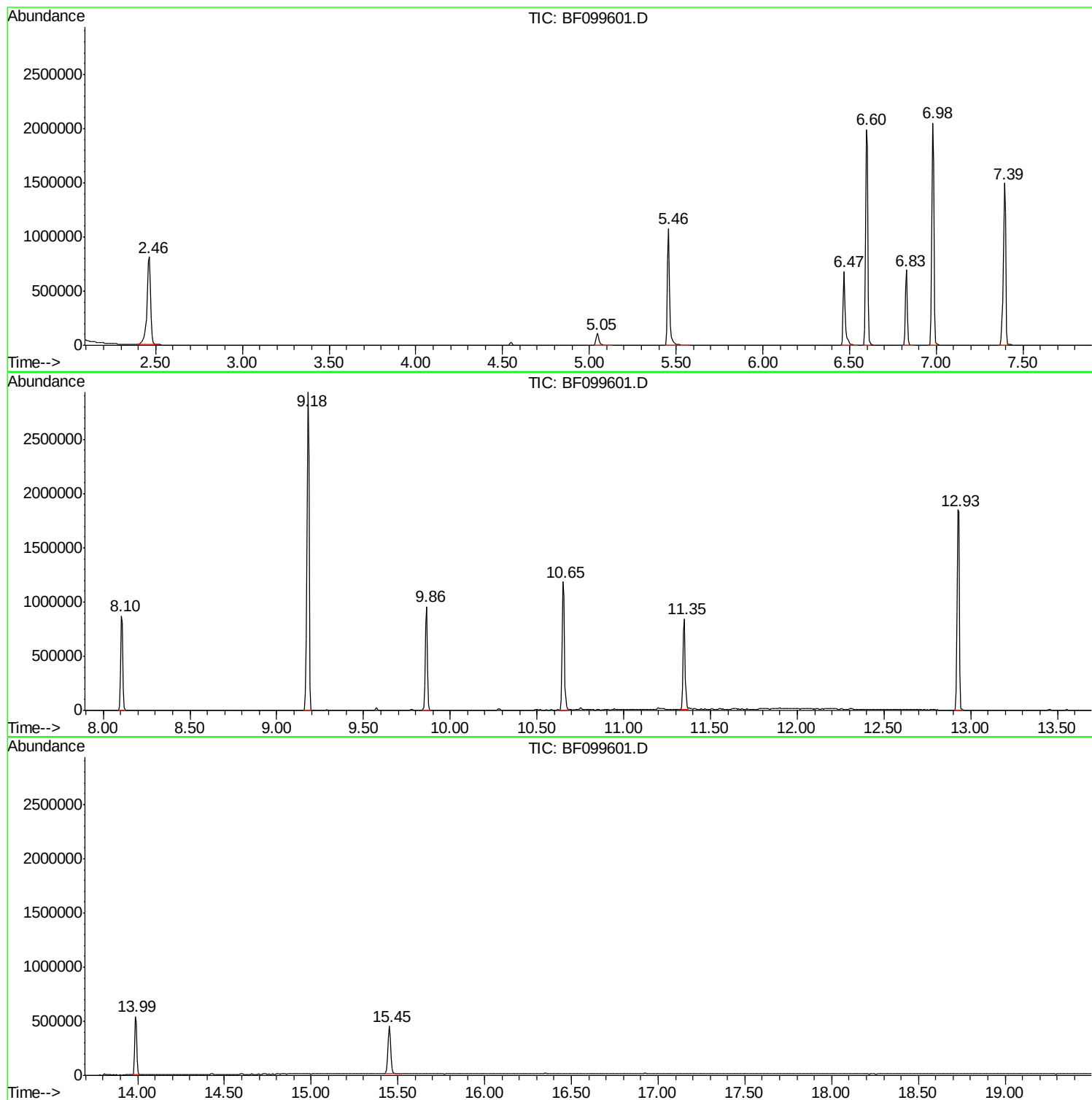
Sum of corrected areas: 18198797

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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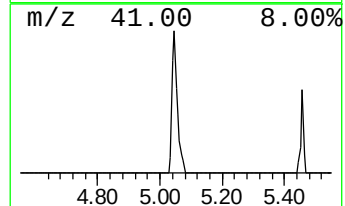
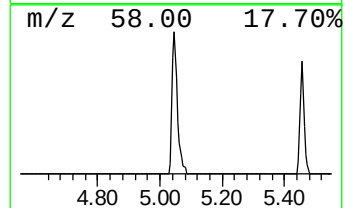
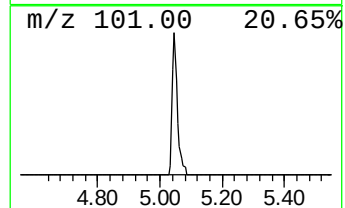
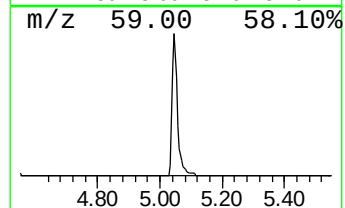
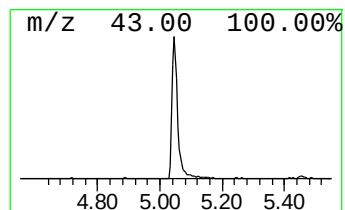
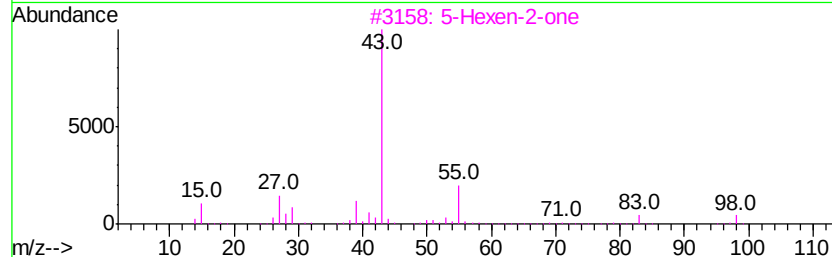
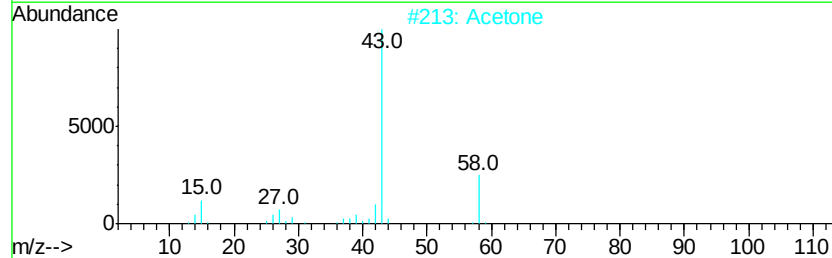
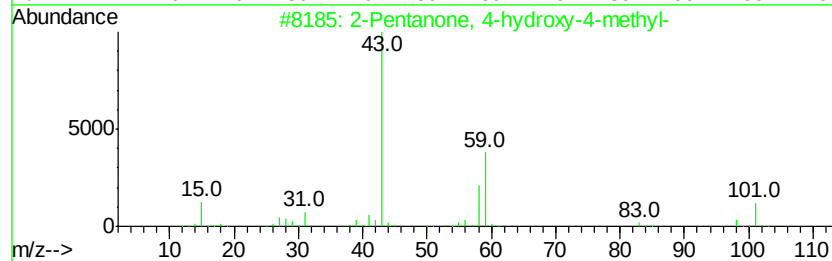
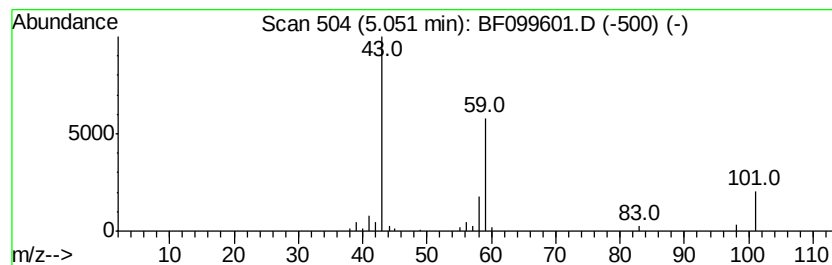
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	4.41 ng	129024	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetone	58	C3H6O	000067-64-1	9
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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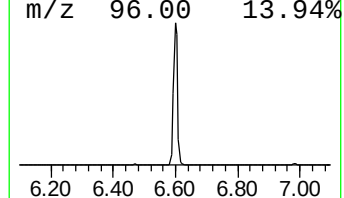
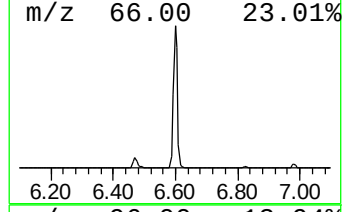
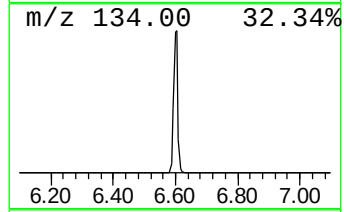
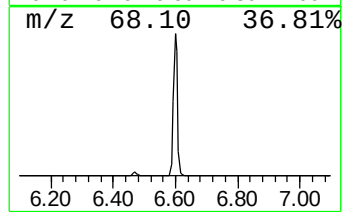
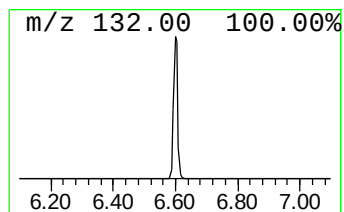
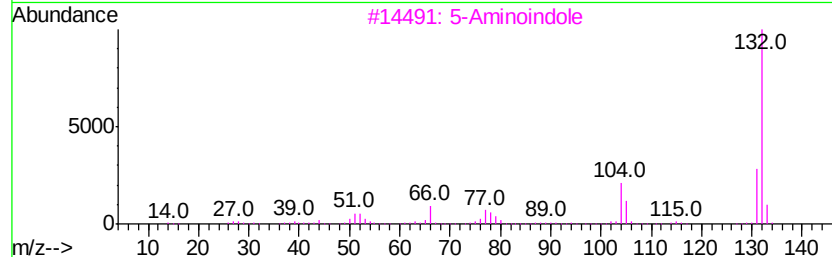
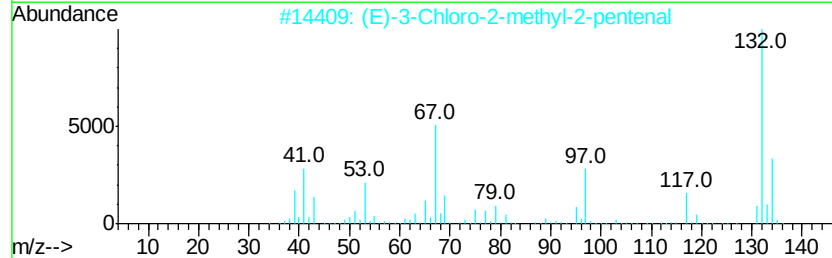
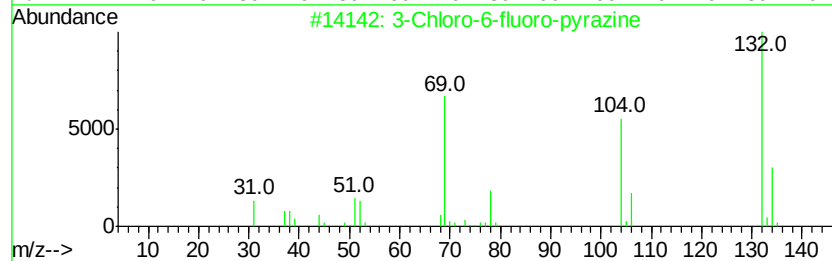
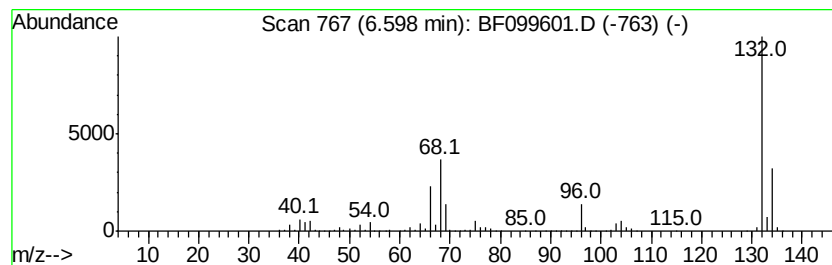
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Peak Number 3 unknown6.60 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	65.54 ng	1918560	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.05	4.4	ng	129024	1	6.83	585469	20.0
unknown6.60	6.60	65.5	ng	1918560	1	6.83	585469	20.0