

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101917\
Data File : BF099663.D
Acq On : 19 Oct 2017 18:06
Operator : SJ/JU
Sample : I5869-02RE
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-5S-10-11-17RE

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	5.122	497	516	519	rBV	2962241	12248157	100.00%	42.803%
2	6.457	741	743	755	rBV	279520	320252	2.61%	1.119%
3	6.592	763	766	776	rVB	306935	253590	2.07%	0.886%
4	6.816	801	804	811	rBV	765582	689945	5.63%	2.411%
5	6.975	827	831	835	rBV	2368771	2304084	18.81%	8.052%
6	7.386	896	901	904	rBV	1630144	1858817	15.18%	6.496%
7	8.098	1018	1022	1026	rBV	1112549	955086	7.80%	3.338%
8	9.175	1199	1205	1208	rBV	3272063	3471570	28.34%	12.132%
9	9.851	1316	1320	1324	rVB	1204919	1075183	8.78%	3.757%
10	11.339	1569	1573	1577	rBV	1223646	1069476	8.73%	3.737%
11	12.927	1837	1843	1846	rBV	2772921	2959300	24.16%	10.342%
12	13.980	2018	2022	2026	rBV	886741	772239	6.30%	2.699%
13	15.439	2263	2270	2277	rBV	492026	637309	5.20%	2.227%

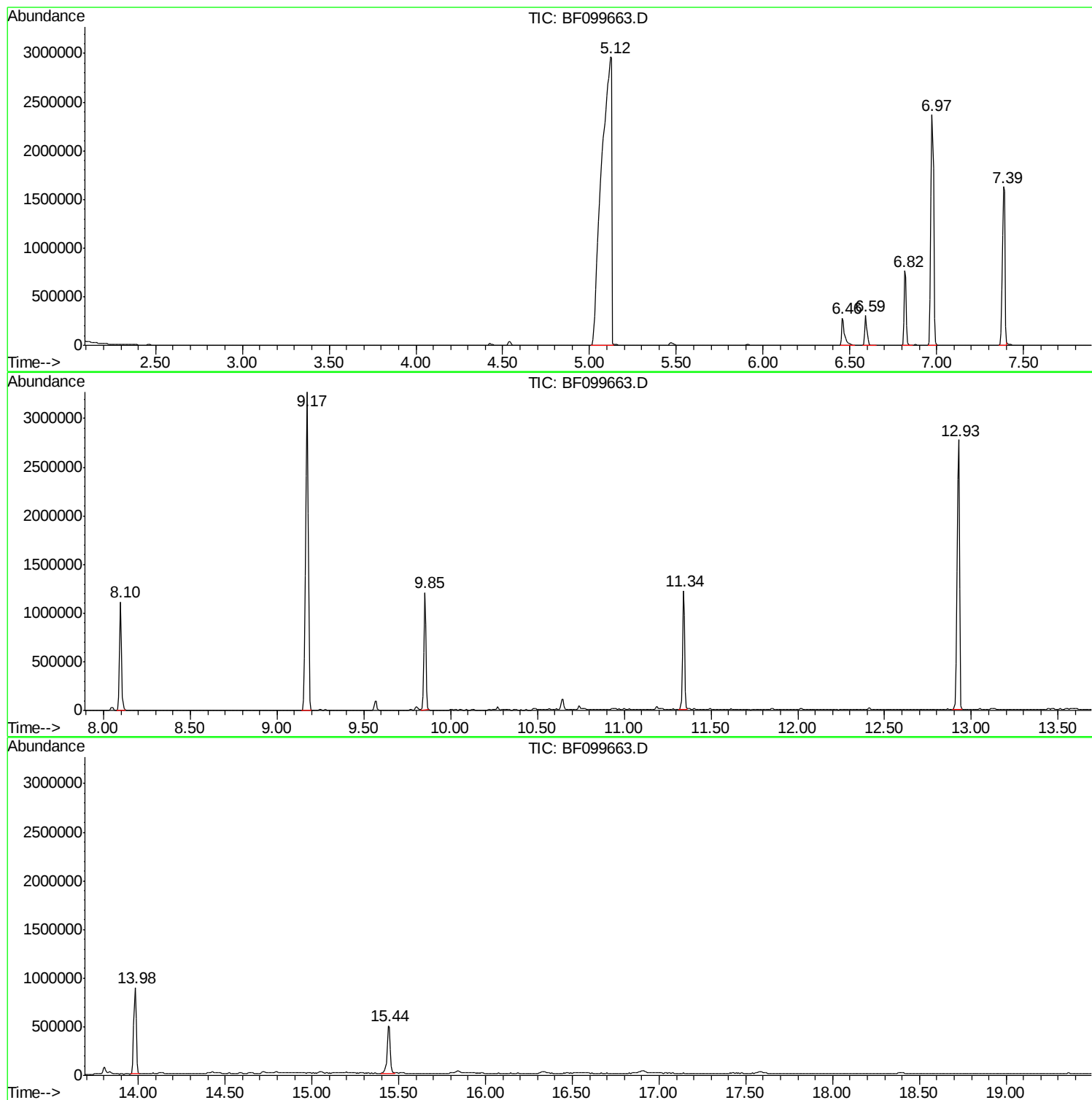
Sum of corrected areas: 28615008

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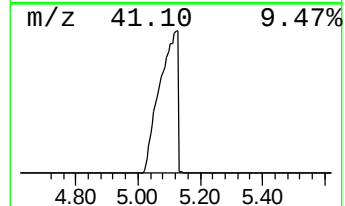
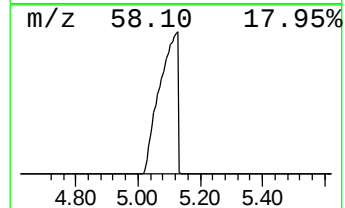
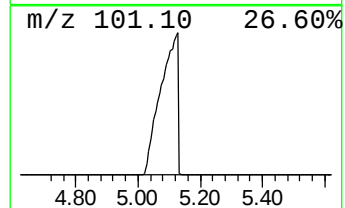
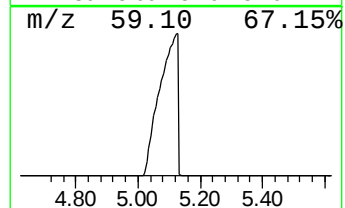
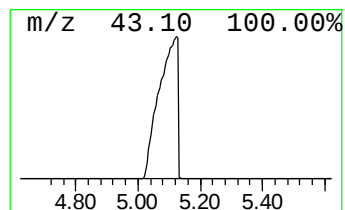
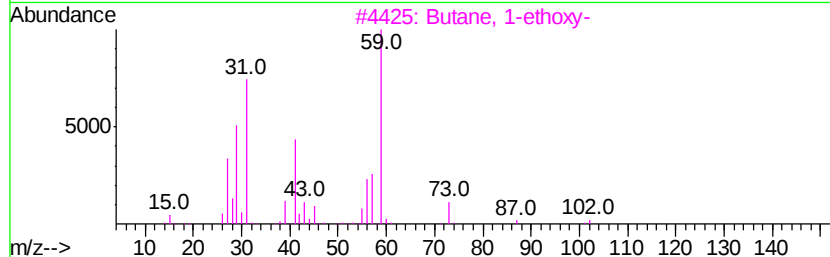
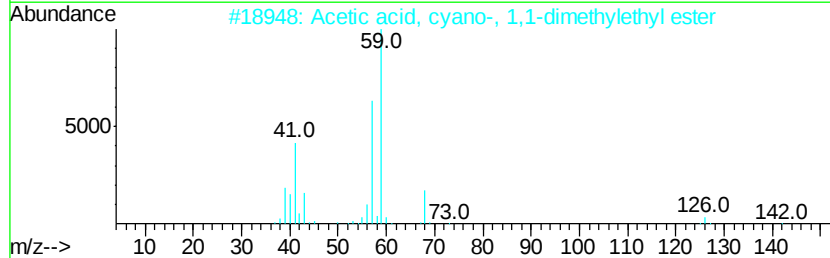
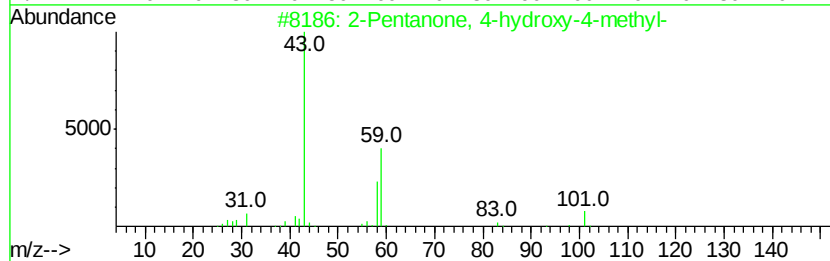
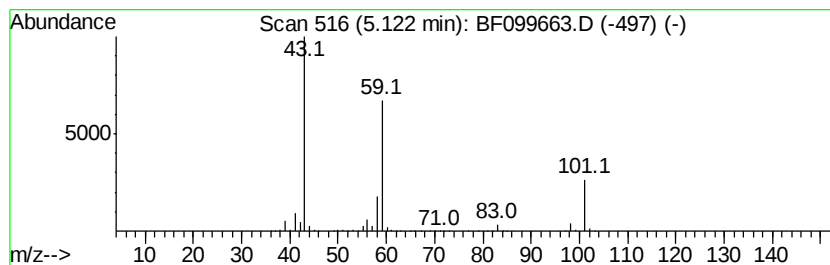
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	355.05 ng	12248200	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Acetone	58	C3H6O	000067-64-1	9



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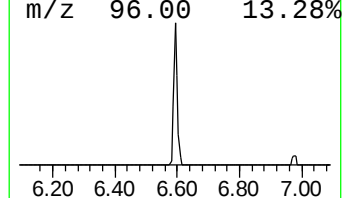
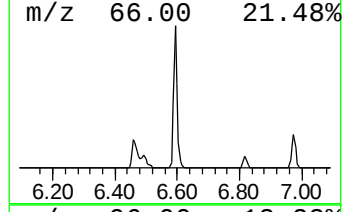
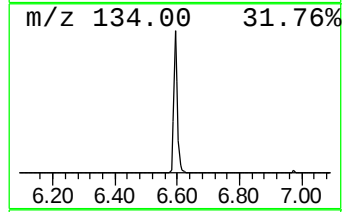
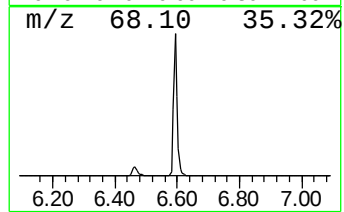
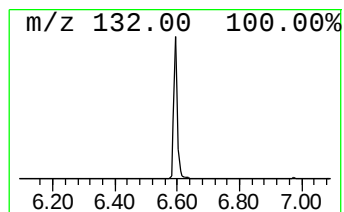
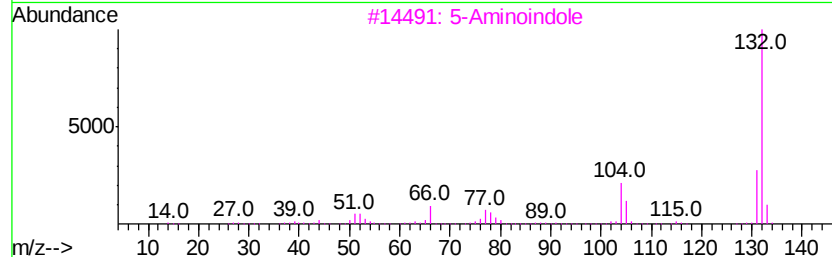
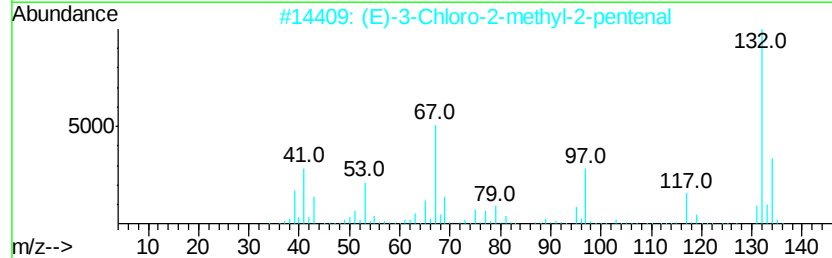
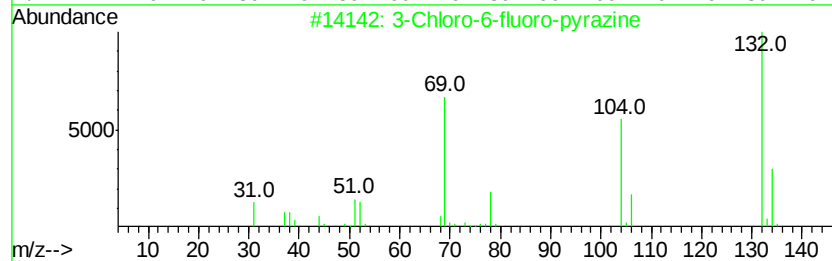
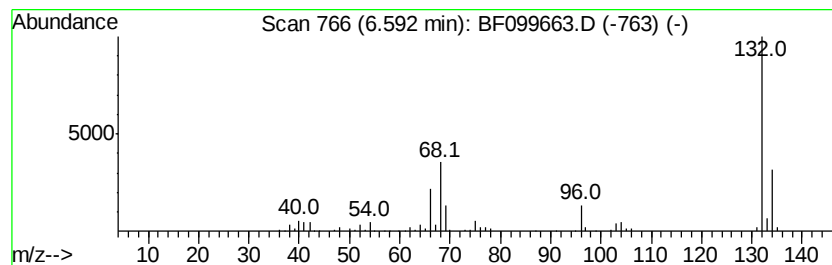
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Peak Number 2 unknown6.59 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	7.35 ng	253590	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3			5-Aminoindole	132	C8H8N2	005192-03-0	12
4			4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5			1,3,5-Trifluorobenzene	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...	5.12	355.1	ng	12248200	1	6.82	689945	20.0
unknown6.59	6.59	7.3	ng	253590	1	6.82	689945	20.0