

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092217\
 Data File : BF098687.D
 Acq On : 22 Sep 2017 9:20
 Operator : SJ/JU
 Sample : PB102454BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102454BL

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-BF092117.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	5.169	519	524	534	rBV	514932	660430	14.75%	1.807%
2	5.551	583	589	592	rBV	3742226	3917646	87.50%	10.717%
3	6.545	751	758	761	rBV	3544757	3771901	84.25%	10.319%
4	6.692	778	783	787	rBV	3738384	3955564	88.35%	10.821%
5	6.922	819	822	826	rVB	1069898	950266	21.22%	2.600%
6	7.081	845	849	853	rBV	3330750	3119087	69.67%	8.533%
7	7.486	912	918	921	rBV	2340092	2464452	55.05%	6.742%
8	8.204	1035	1040	1044	rBV	1478670	1316363	29.40%	3.601%
9	9.280	1217	1223	1226	rBV	4418935	4477149	100.00%	12.248%
10	9.963	1333	1339	1342	rBV	1441104	1424178	31.81%	3.896%
11	10.751	1467	1473	1476	rBV	2493044	2610689	58.31%	7.142%
12	11.445	1587	1591	1595	rVB	1683179	1418108	31.67%	3.879%
13	13.033	1856	1861	1864	rBV	4259039	4176559	93.29%	11.426%
14	14.080	2036	2039	2043	rVB	1365828	1267113	28.30%	3.466%
15	15.574	2287	2293	2302	rBV	820126	1024508	22.88%	2.803%

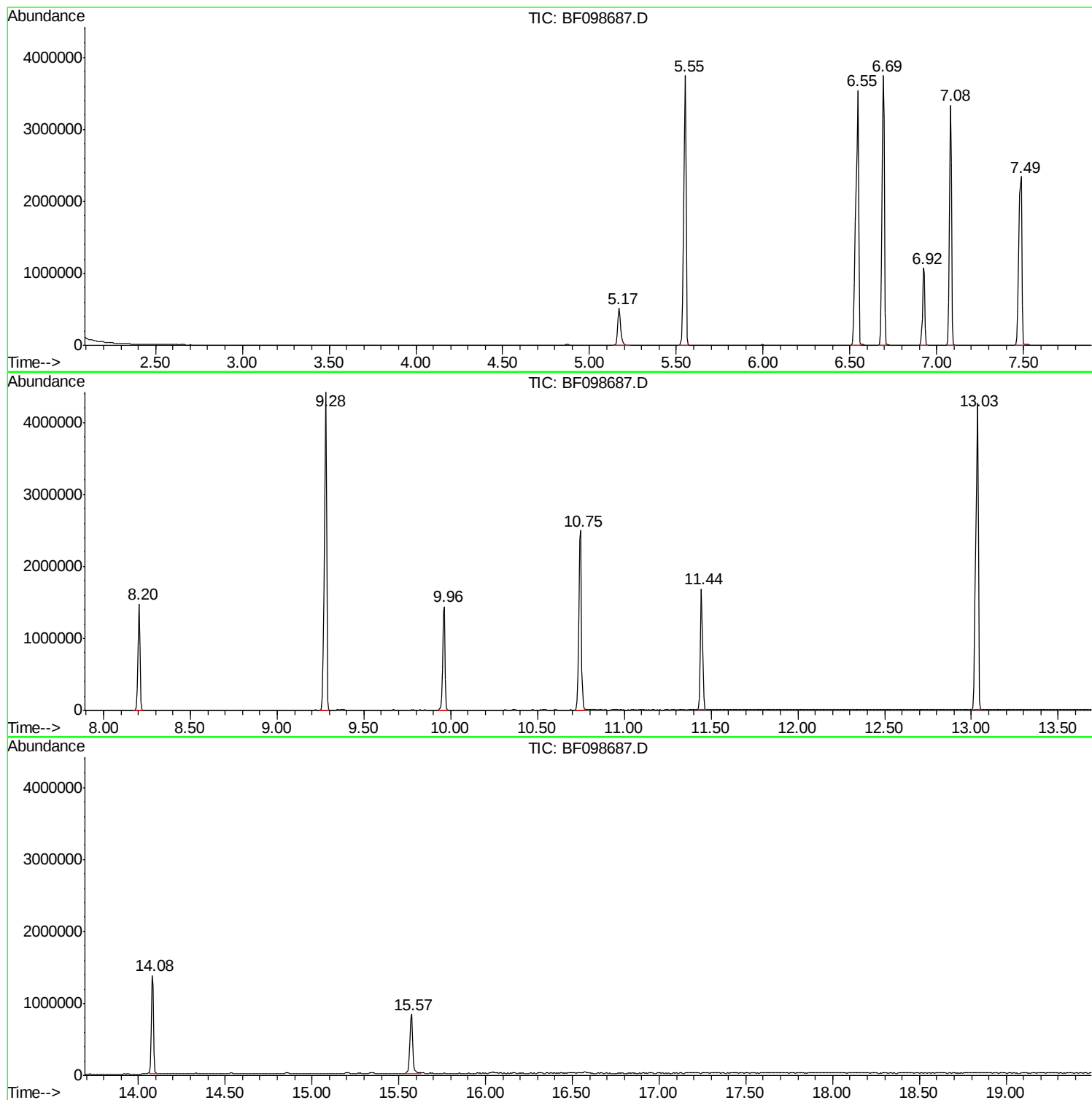
Sum of corrected areas: 36554013

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092117.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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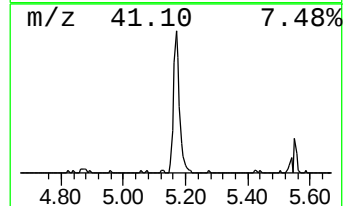
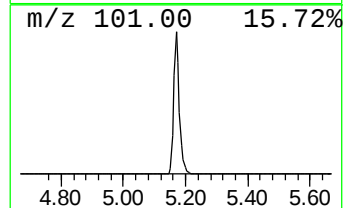
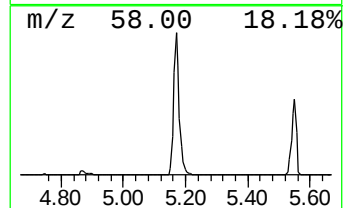
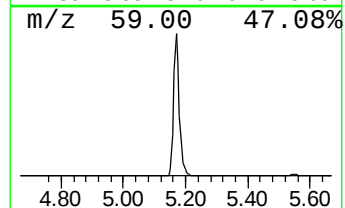
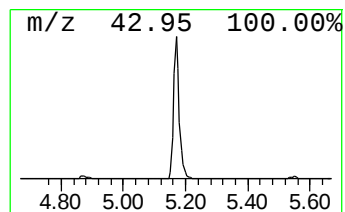
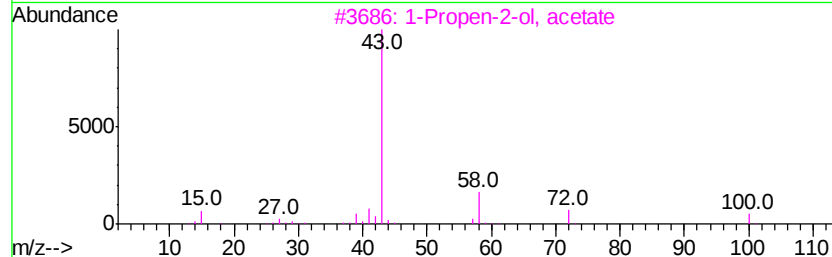
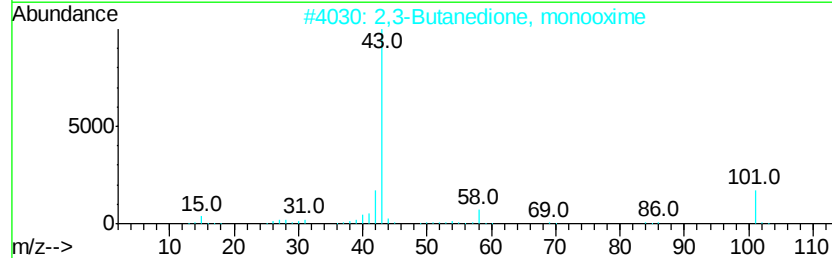
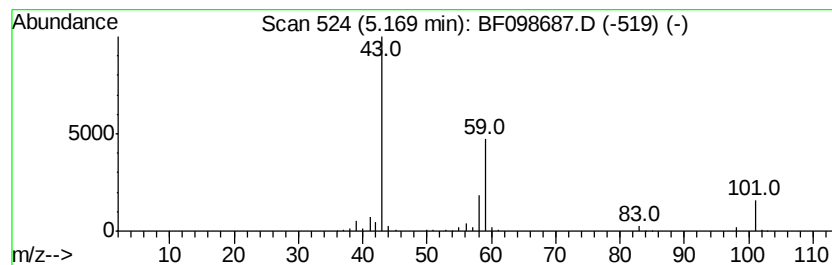
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	13.90 ng	660430	1,4-Dichlorobenzene-d4	6.92

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		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	7
5		Acetone	58	C3H6O	000067-64-1	7



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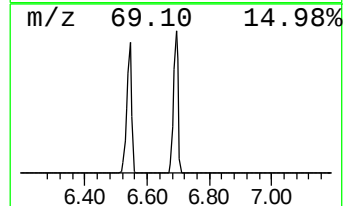
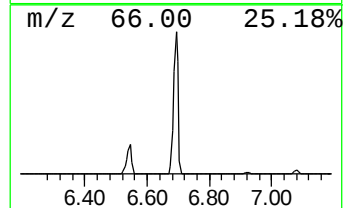
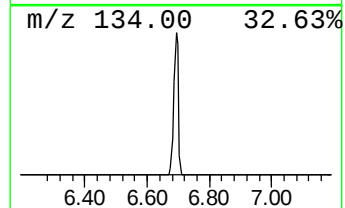
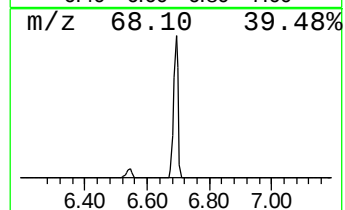
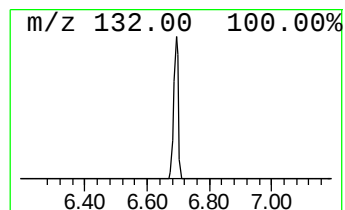
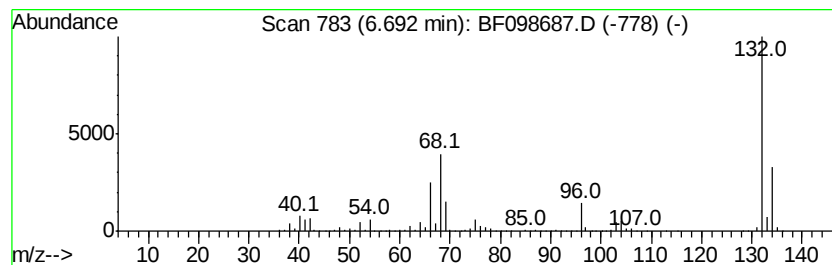
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Peak Number 2 unknown6.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	83.25 ng	3955560	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	13.9	ng	660430	1	6.92	950266	20.0
unknown6.69	6.69	83.3	ng	3955560	1	6.92	950266	20.0