

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030816\
 Data File : BF085259.D
 Acq On : 8 Mar 2016 16:53
 Operator : UM/SJ
 Sample : H1713-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-FB-05

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-BF030316.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	5.156	248	251	255	rVB	226587	301050	5.84%	0.938%
2	5.556	284	286	289	rBV	1789942	1711880	33.19%	5.335%
3	6.573	373	375	377	rBV	1346559	1098019	21.29%	3.422%
4	6.722	386	388	391	rBV	2954678	3246315	62.94%	10.116%
5	6.962	407	409	411	rBV	920371	861748	16.71%	2.685%
6	7.122	420	423	425	rVB	3043916	3585409	69.52%	11.173%
7	7.522	455	458	460	rBV	2620092	2630935	51.01%	8.199%
8	8.242	519	521	524	rBV	1212334	1089272	21.12%	3.394%
9	9.328	613	616	618	rBV	3869379	5088563	98.67%	15.857%
10	10.002	672	675	677	rBV	1428493	1250514	24.25%	3.897%
11	10.791	741	744	746	rBV	2886937	2959551	57.38%	9.223%
12	11.488	802	805	807	rBV	1256546	1284558	24.91%	4.003%
13	13.076	941	944	946	rBV	4410687	5157402	100.00%	16.072%
14	14.117	1033	1035	1038	rBV	1134429	1040392	20.17%	3.242%
15	15.580	1160	1163	1166	rBV	557033	784089	15.20%	2.443%

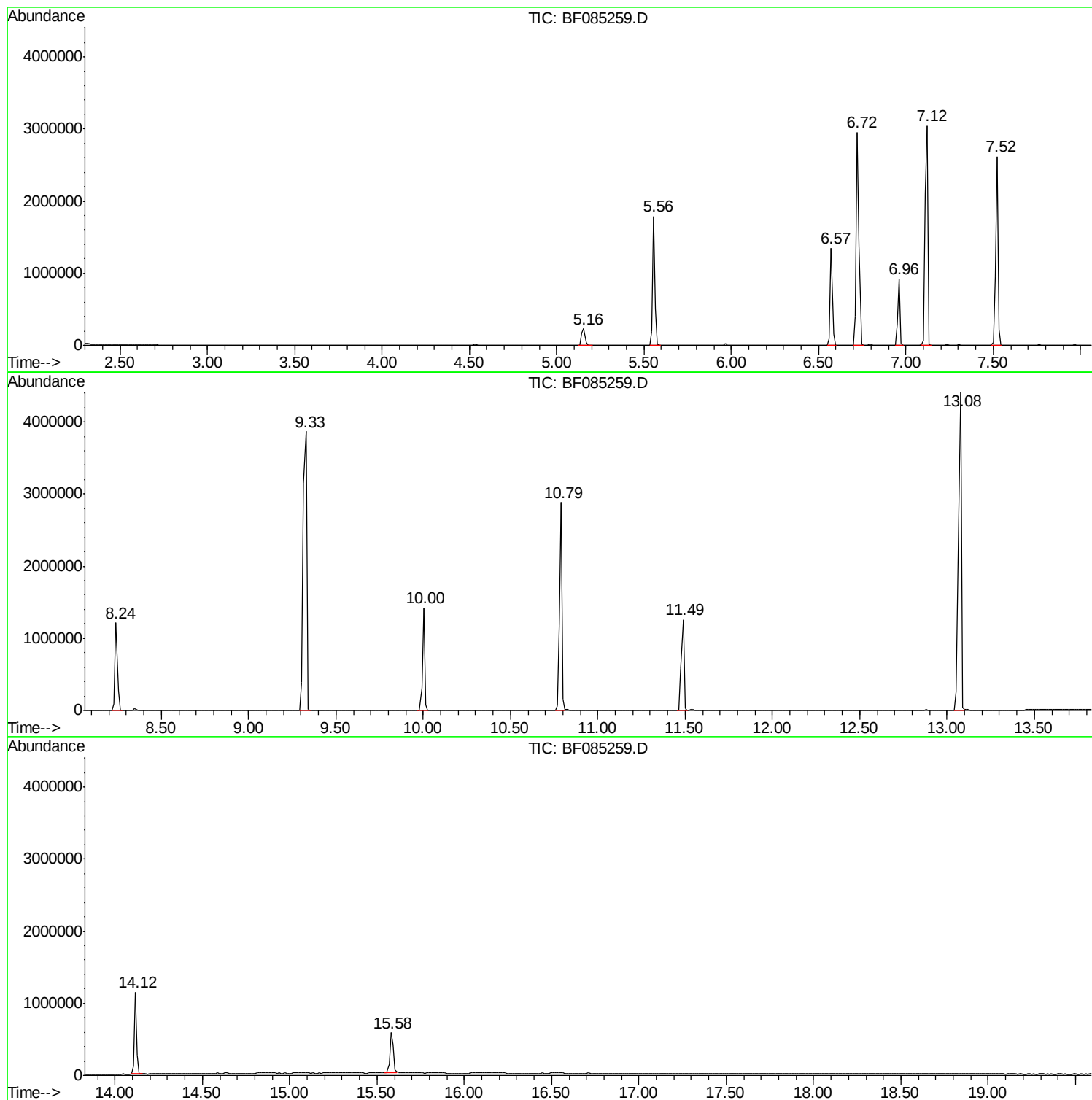
Sum of corrected areas: 32089697

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.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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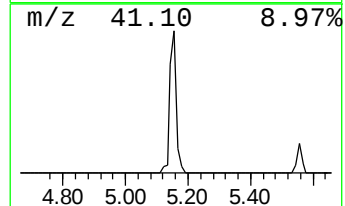
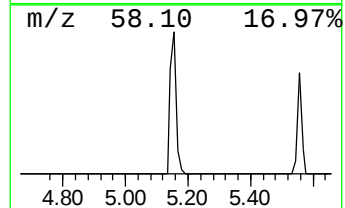
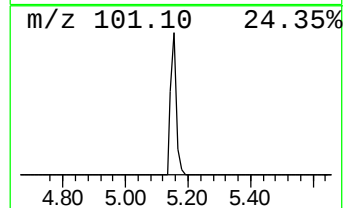
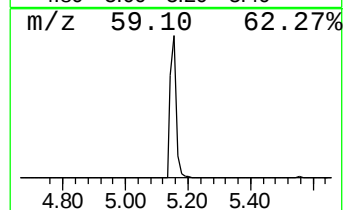
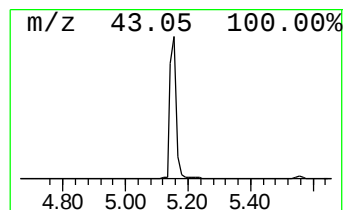
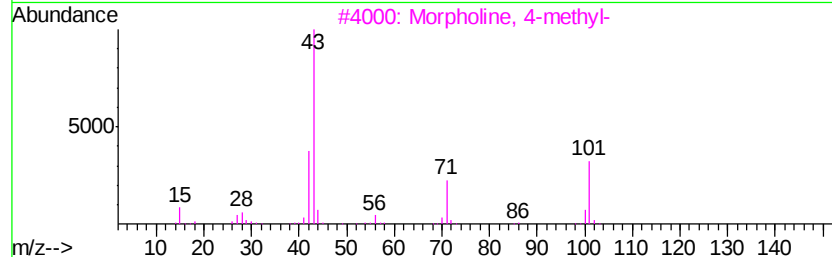
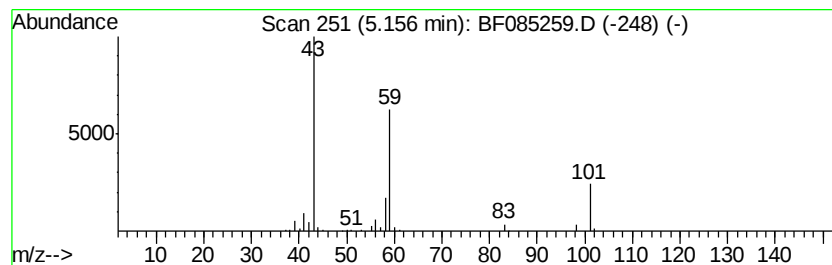
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	6.99 ng	301050	1,4-Dichlorobenzene-d4	6.96

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-dimethy...	141	C7H11NO2	001116-98-9	17
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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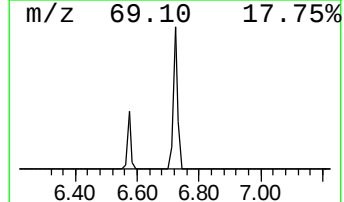
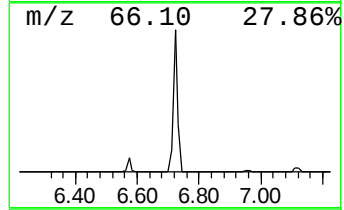
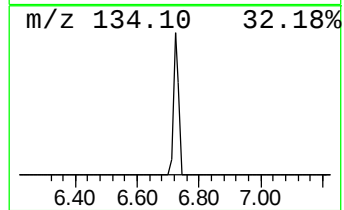
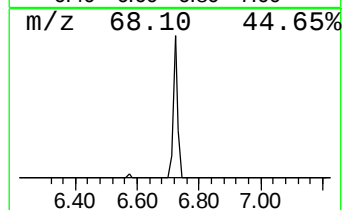
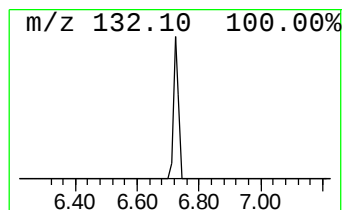
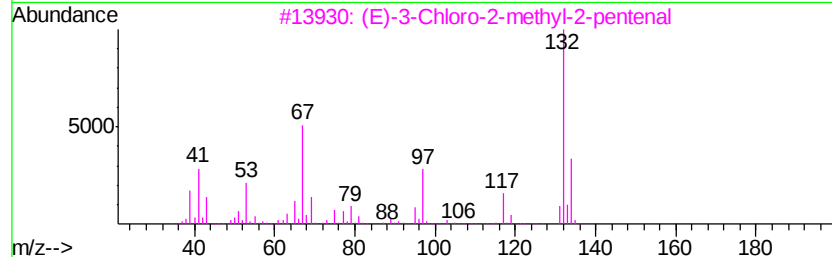
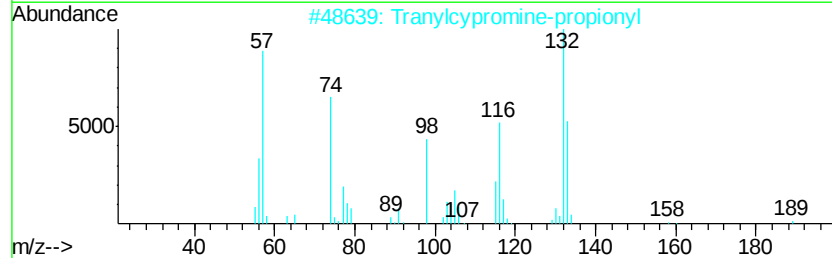
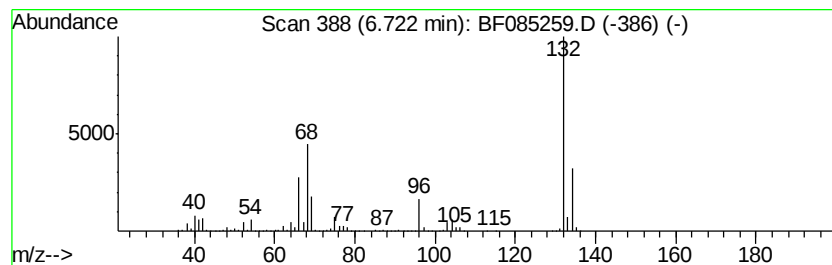
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Peak Number 2 unknown6.72 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	75.34 ng	3246320	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
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...	5.16	7.0	ng	301050	1	6.96	861748	20.0
unknown6.72	6.72	75.3	ng	3246320	1	6.96	861748	20.0