

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
 Data File : BF103869.D
 Acq On : 23 Mar 2018 10:22
 Operator : JU/SJ
 Sample : PB107555BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB107555BL

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-BF031518.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	5.222	527	533	541	rBV	121261	161904	3.71%	0.485%
2	5.604	592	598	601	rBV	3119329	3339511	76.51%	9.996%
3	6.598	761	767	770	rBV	3183062	3270180	74.92%	9.789%
4	6.745	787	792	795	rBV	3529012	3404597	78.00%	10.191%
5	6.975	827	831	834	rBV	957934	832999	19.08%	2.493%
6	7.128	853	857	861	rBV	2825076	2746782	62.93%	8.222%
7	7.534	921	926	929	rBV	2072697	2172698	49.78%	6.504%
8	8.257	1044	1049	1052	rBV	1281535	1135162	26.01%	3.398%
9	9.333	1226	1232	1235	rBV	3842803	4034556	92.44%	12.077%
10	10.016	1344	1348	1356	rVB	1491004	1303875	29.87%	3.903%
11	10.810	1478	1483	1486	rBV	2519287	2497288	57.22%	7.475%
12	11.510	1598	1602	1605	rBV	1484431	1338138	30.66%	4.005%
13	13.098	1867	1872	1875	rBV	4079926	4364715	100.00%	13.065%
14	14.162	2048	2053	2056	rBV	1568406	1430865	32.78%	4.283%
15	15.704	2309	2315	2326	rBV	1001611	1374638	31.49%	4.115%

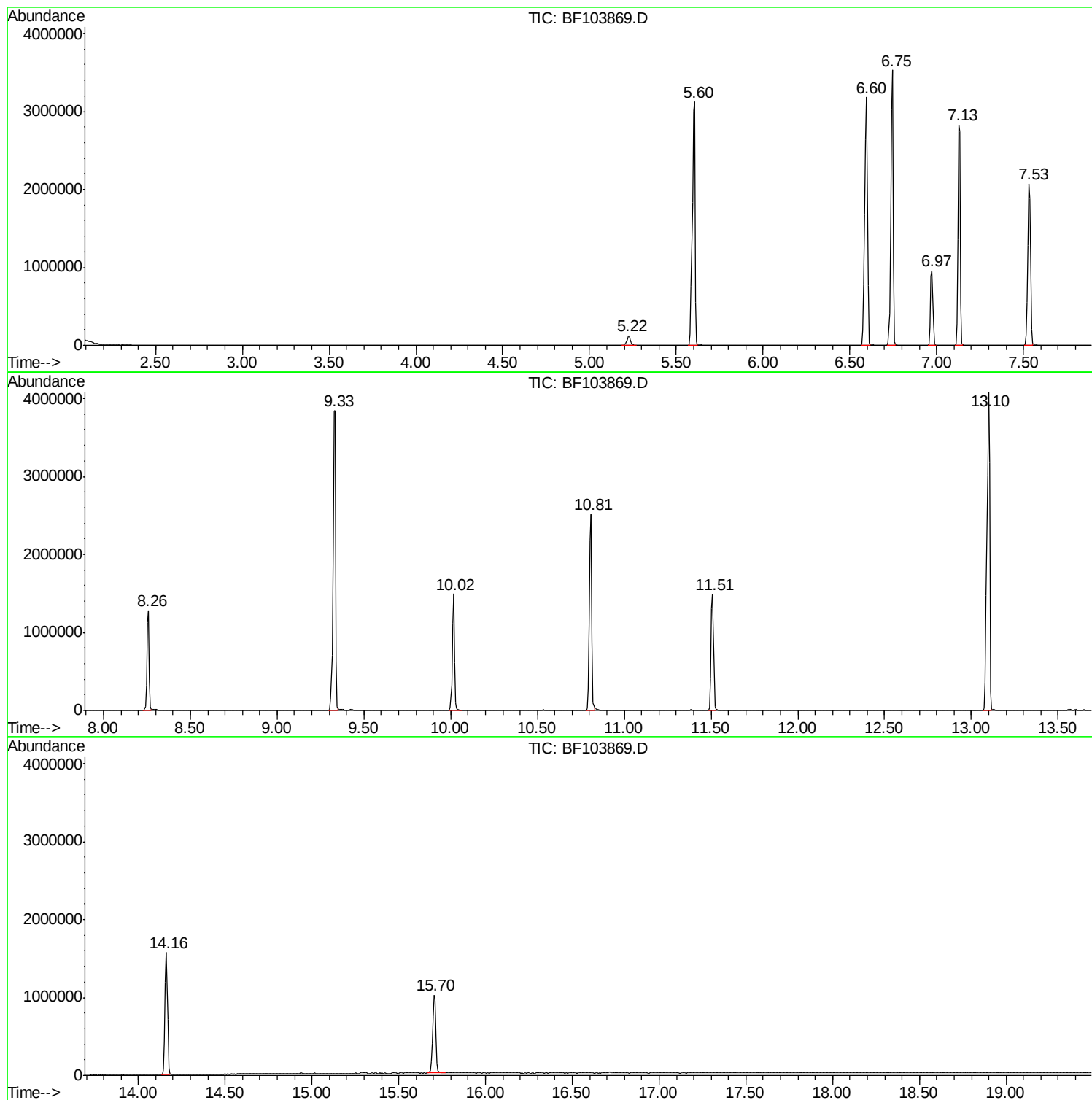
Sum of corrected areas: 33407908

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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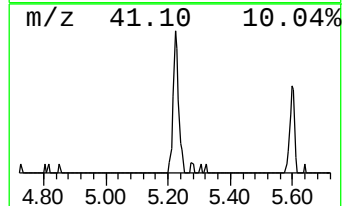
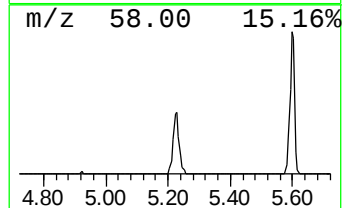
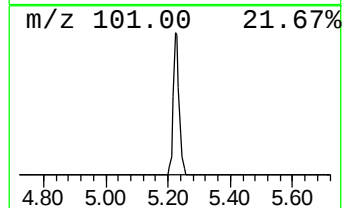
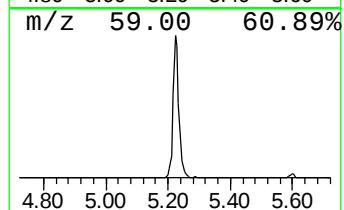
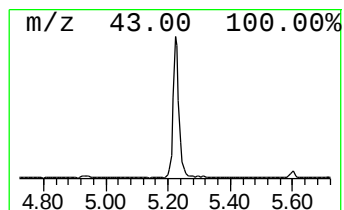
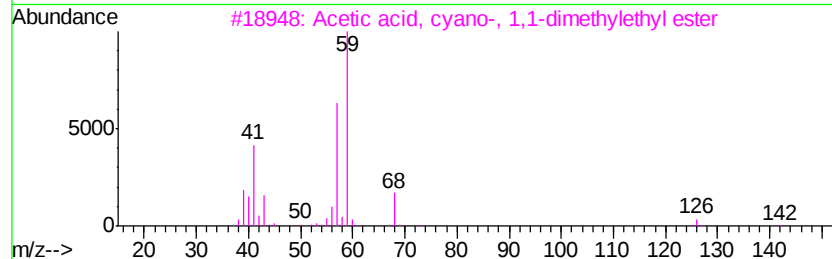
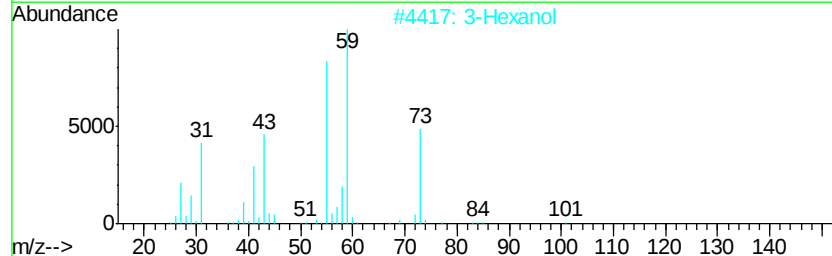
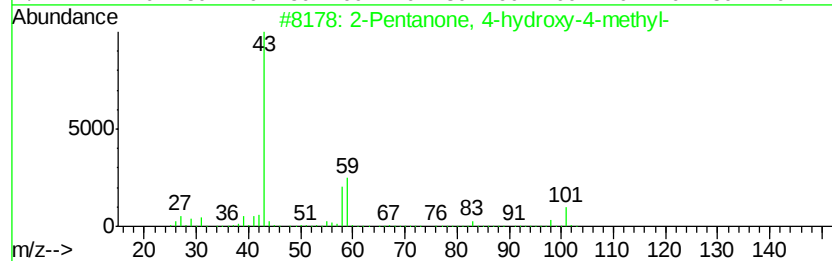
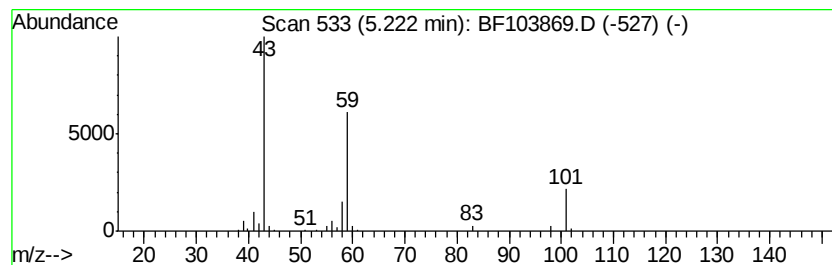
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	3.89 ng	161904	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			3-Hexanol	102	C6H14O	000623-37-0	28
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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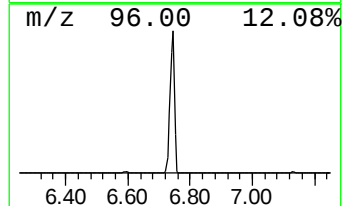
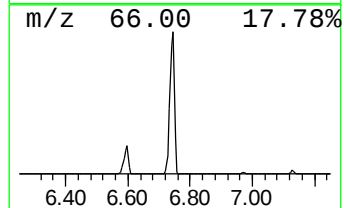
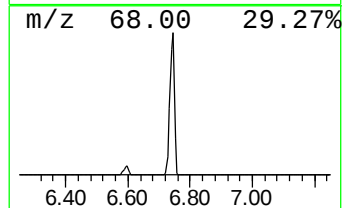
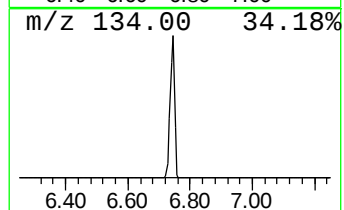
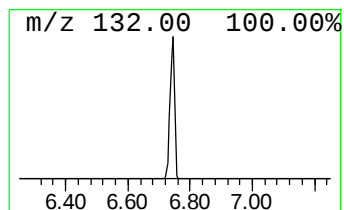
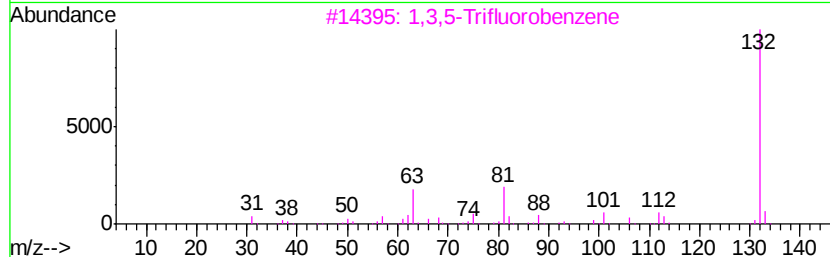
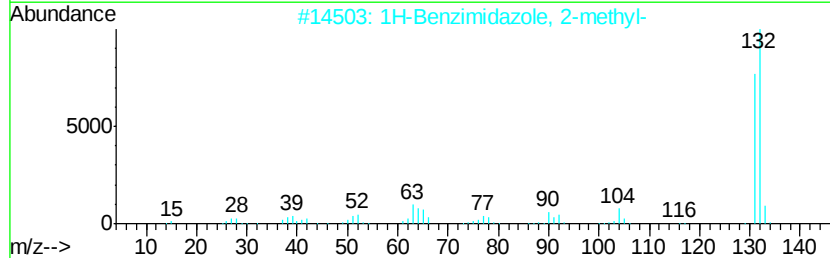
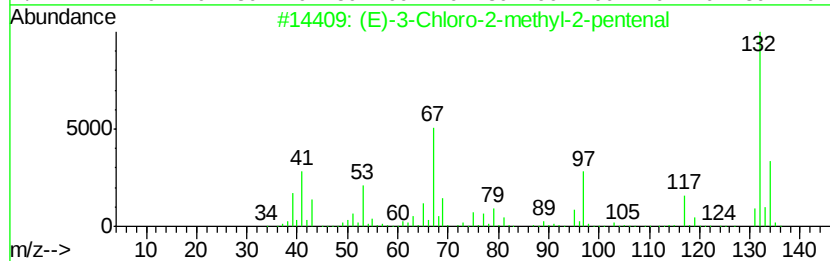
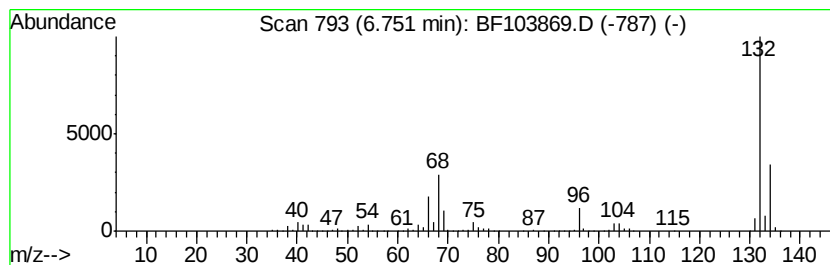
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Peak Number 2 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	81.74 ng	3404600	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	3.9	ng	161904	1	6.97	832999	20.0
unknown6.75	6.75	81.7	ng	3404600	1	6.97	832999	20.0