

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
Data File : BF100876.D
Acq On : 23 Nov 2017 15:31
Operator : SJ/JU
Sample : I6567-08
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
112017-JETT-F-U-S

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-BF110817.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	5.069	504	507	513	rBV	150002	146926	5.77%	1.097%
2	5.469	571	575	579	rBV	650054	562690	22.09%	4.201%
3	6.475	743	746	756	rBB	406322	359160	14.10%	2.682%
4	6.622	767	771	775	rBV	1177399	1122636	44.07%	8.382%
5	6.857	807	811	814	rBV	530165	418649	16.43%	3.126%
6	7.016	833	838	841	rBV	1861783	1659999	65.16%	12.394%
7	7.422	901	907	910	rBV	1223536	1368153	53.70%	10.215%
8	8.139	1025	1029	1032	rBV	684818	549491	21.57%	4.103%
9	9.216	1206	1212	1215	rBV	2631356	2547555	100.00%	19.021%
10	9.892	1321	1327	1331	rBV2	686205	620360	24.35%	4.632%
11	10.680	1457	1461	1465	rBV	970721	897587	35.23%	6.702%
12	11.380	1576	1580	1583	rBV	674377	552875	21.70%	4.128%
13	12.968	1845	1850	1853	rBV	1919580	1828730	71.78%	13.654%
14	14.015	2025	2028	2032	rBV	391473	358405	14.07%	2.676%
15	15.486	2271	2278	2287	rBV	309907	400229	15.71%	2.988%

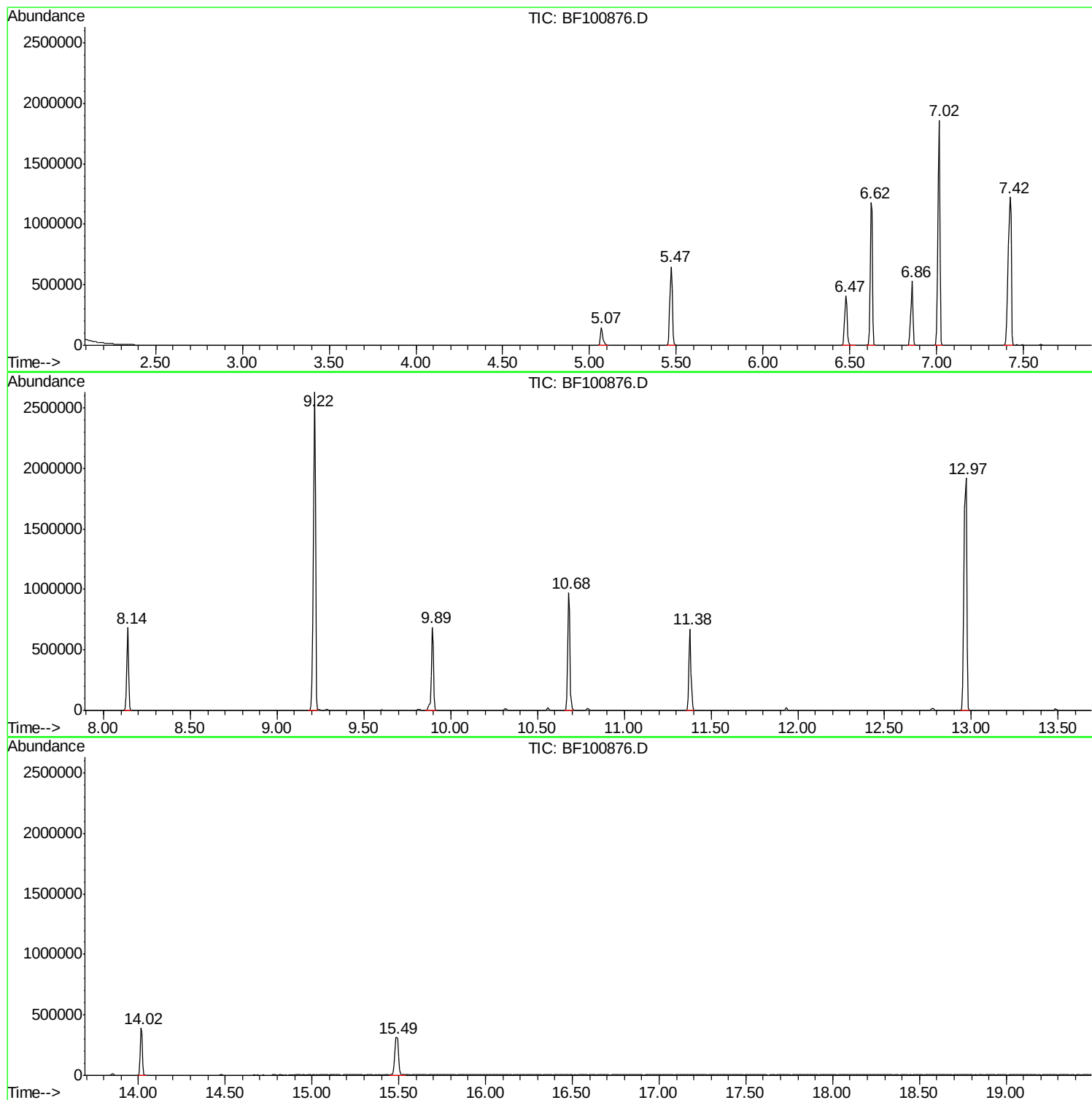
Sum of corrected areas: 13393445

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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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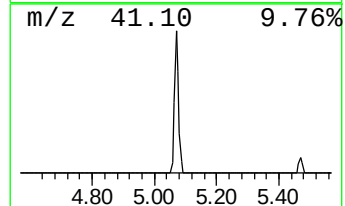
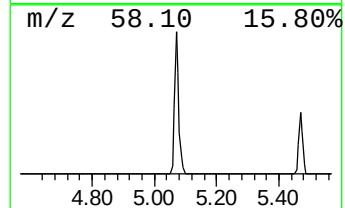
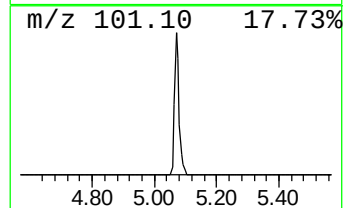
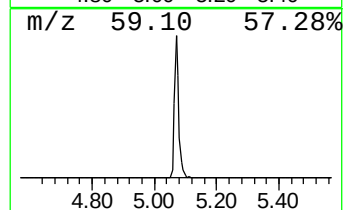
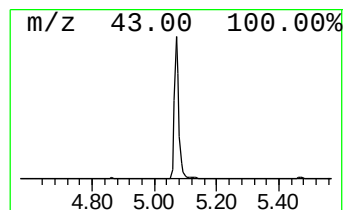
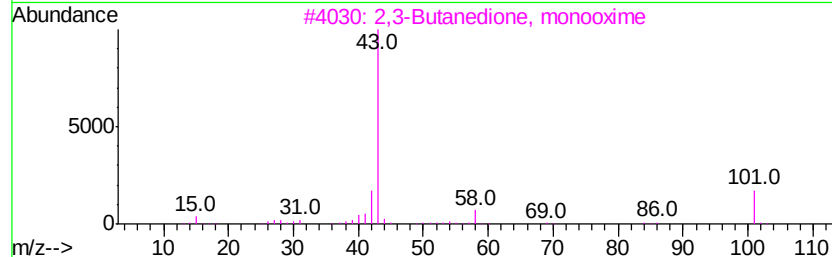
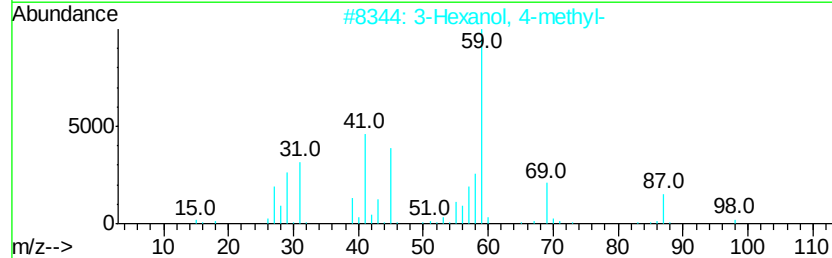
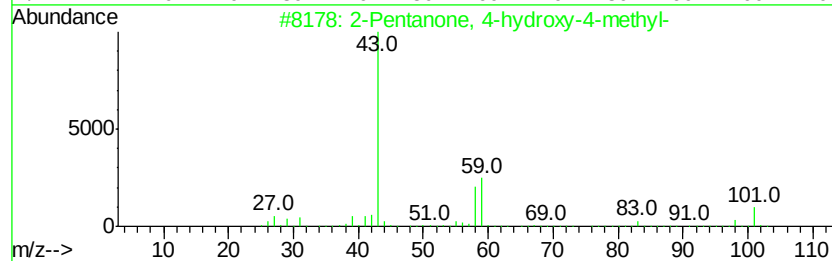
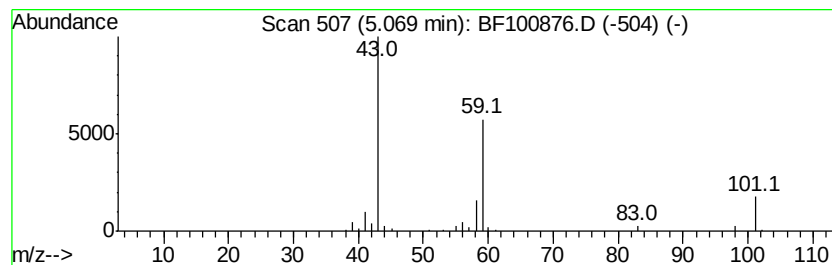
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	7.02 ng	146926	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	9
5		Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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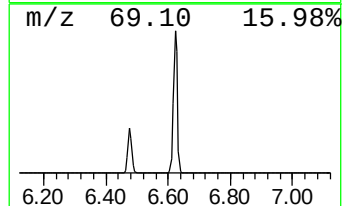
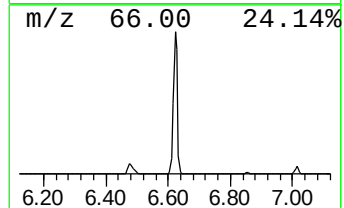
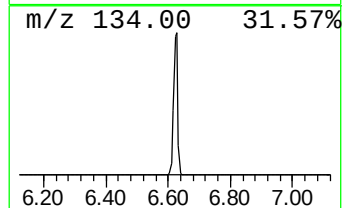
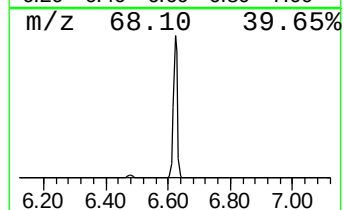
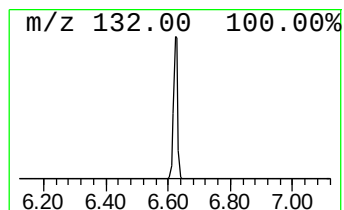
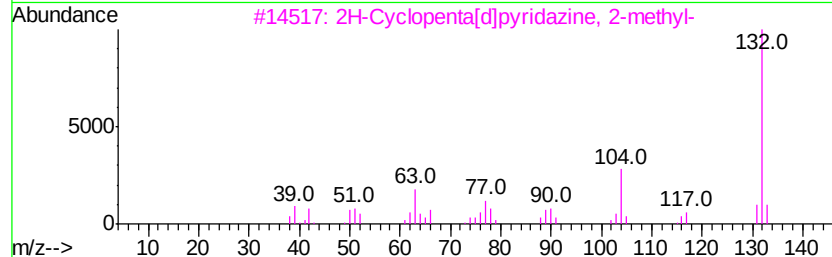
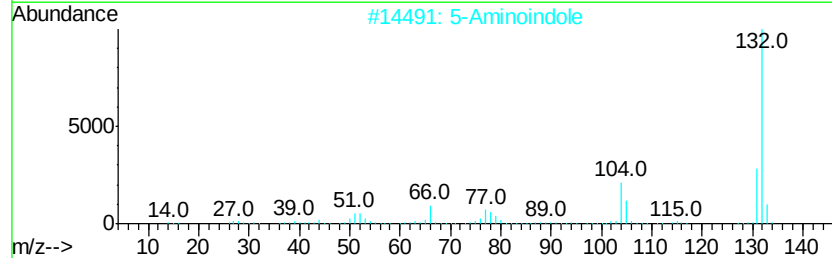
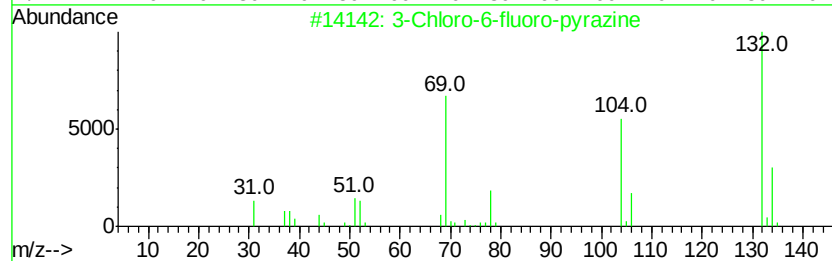
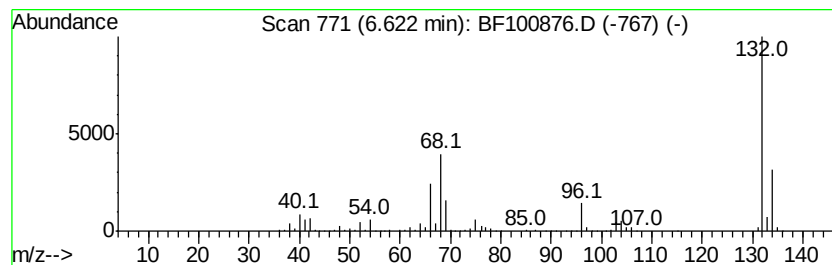
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Peak Number 2 unknown6.62 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	53.63 ng	1122640	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4			4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	7.0	ng	146926	1	6.86	418649	20.0
unknown6.62	6.62	53.6	ng	1122640	1	6.86	418649	20.0