

Mapping of Raw files to their short names Mapping source: file (user-defined) (automatic shortening of names was not sufficiently short – see 'best effort')

| original | short name | |
|------------------------|---------------|--------------|
| | | |
| Toni_20140521_GM_QC_01 | file 1 | 521_GM_QC_01 |
| Toni_20140521_GM_QC_02 | file 2 | 521_GM_QC_02 |
| Toni_20140522_GM_QC_01 | file 3 | 522_GM_QC_01 |
| Toni_20140531_FB_QC_02 | file 4 | 531_FB_QC_02 |
| Toni_20140608_FB_qc_01 | file 5 | 608_FB_qc_01 |

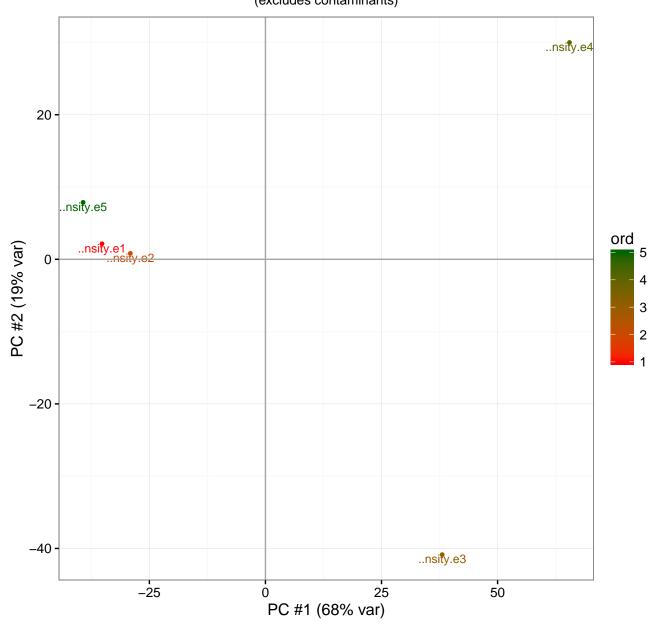
PAR: parameters

| parameter | value | parameter | value |
|------------------------------|---------------------------------------|------------------------------|----------------------------|
| Advanced ratios | False | MS/MS deisotoping (ITMS) | False |
| Alignment time window [min] | 100 | MS/MS deisotoping (TOF) | False |
| Cut peaks | True | MS/MS deisotoping (Unknown) | False |
| Decoy mode | revert | MS/MS recalibration | False |
| Discard unmodified counterpa | True | MS/MS tol. (FTMS) | 20 ppm |
| Find dependent peptides | False | MS/MS tol. (ITMS) | 0.5 Da |
| First pass AIF correlation | 0.8 | MS/MS tol. (TOF) | 0.1 Da |
| Fixed modifications | Carbamidomethyl (C) | MS/MS tol. (Unknown) | 0.5 Da |
| iBAQ | False | Peptides used for protein qu | Razor |
| iBAQ log fit | False | Protein FDR | 0.01 |
| Include contaminants | True | PSM FDR | 0.01 |
| Labeled amino acid filtering | True | Re-quantify | True |
| Match between runs | True | RT shift | False |
| Matching time window [min] | 1 | Site FDR | 0.01 |
| Min. delta score for modifie | 17 | Site quantification | Use least modified peptide |
| Min. delta score for unmodif | 0 | Site tables | Oxidation (M)Sites.txt |
| Min. peptide Length | 7 | Special AAs | KR |
| Min. peptides | 1 | Top MS/MS peaks per 100 Da | 12 |
| Min. ratio count | 2 | Top MS/MS peaks per 100 Da | 8 |
| Min. razor peptides | 1 | Top MS/MS peaks per 100 Da | 10 |
| Min. score for modified pept | 40 | Top MS/MS peaks per 100 Da | 10 |
| Min. score for unmodified pe | 0 | Use delta score | False |
| Min. unique peptides | 0 | Use Normalized Ratios For Oc | True |
| Modifications included in pr | Acetyl (Protein N-term) Oxidation (M) | Use only unmodified peptides | True |
| MS/MS deisotoping (FTMS) | True | Version | 1.4.1.2 |

uniprot_human_canonical_and_isoforms_20130513.fasta

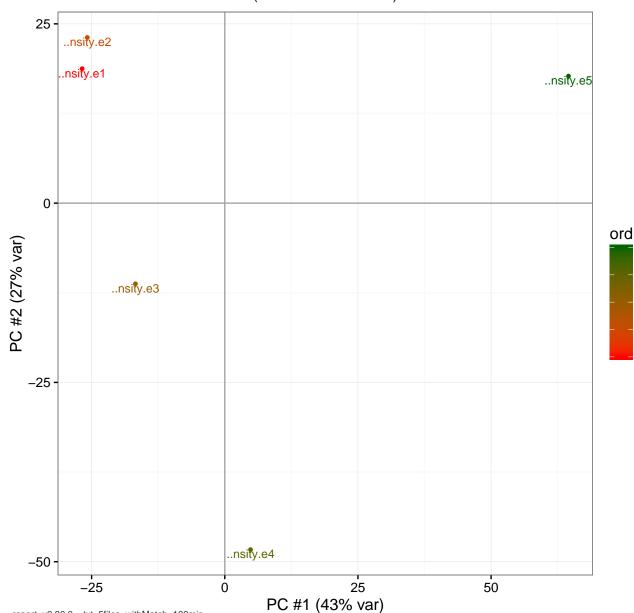
PG: PCA of 'raw intensity'





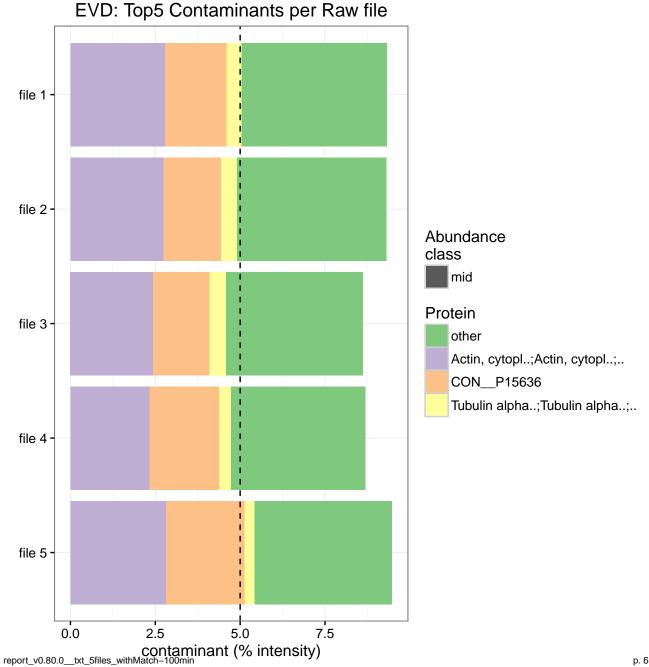
PG: PCA of 'Ifq intensity'

(excludes contaminants)



5

3



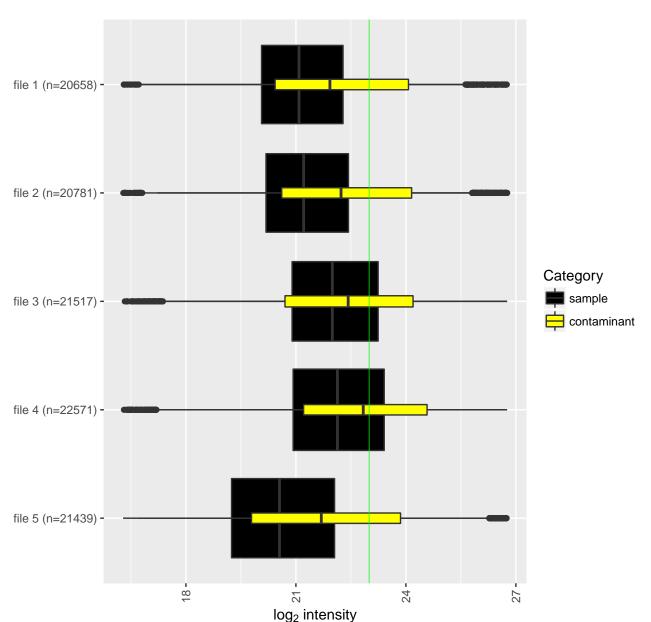
PG: Contaminants

Contaminant 'MYCOPLASMA' was not found in any sample.

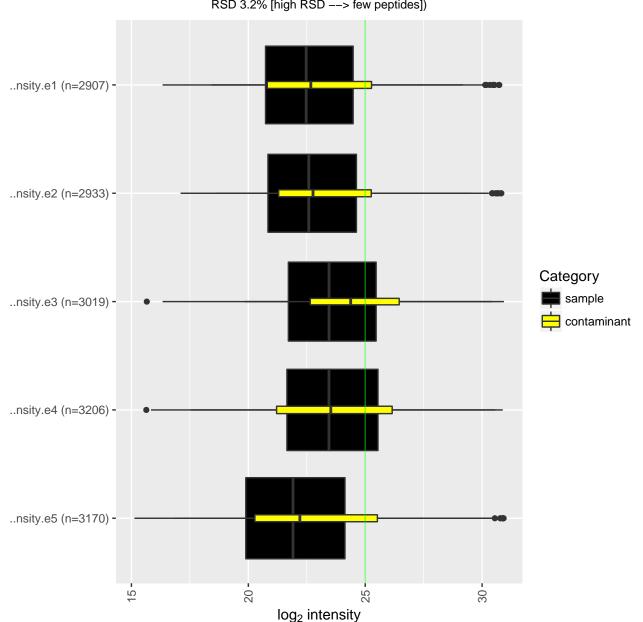
Did you use the correct database?

EVD: peptide intensity distribution

RSD 3.2% (expected < 5%)

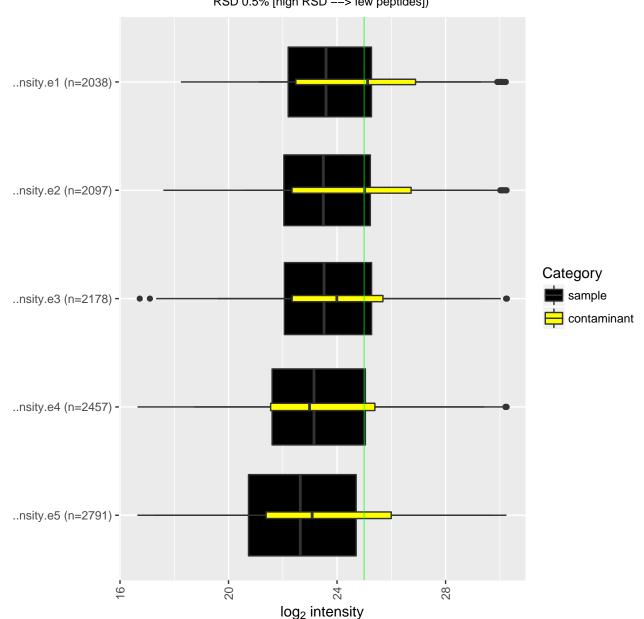


PG: intensity distribution
RSD 3% (w/o zero int.; expected < 5%)
RSD 3.2% [high RSD ---> few peptides])

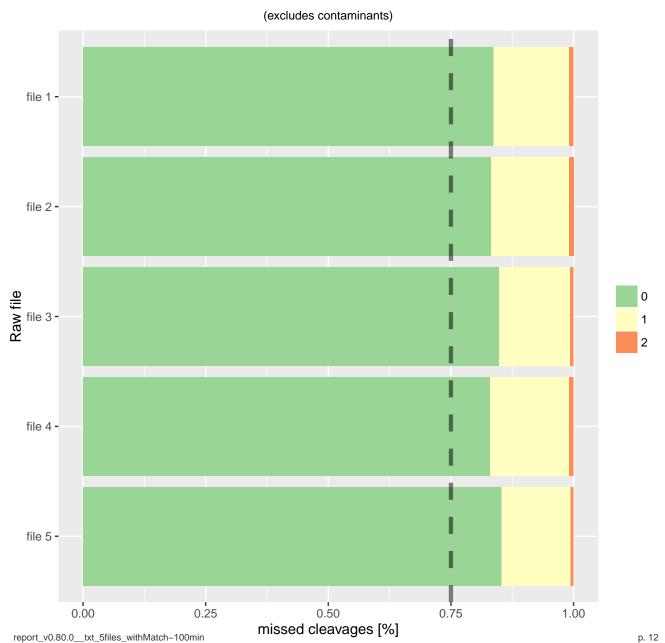


PG: LFQ intensity distribution RSD 1.7% (w/o zero int.; expected < 5%)

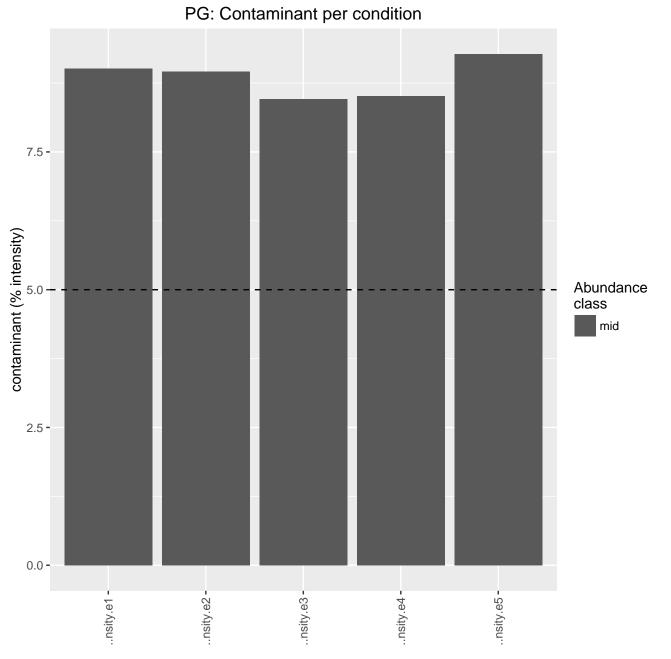
RSD 0.5% [high RSD --> few peptides])

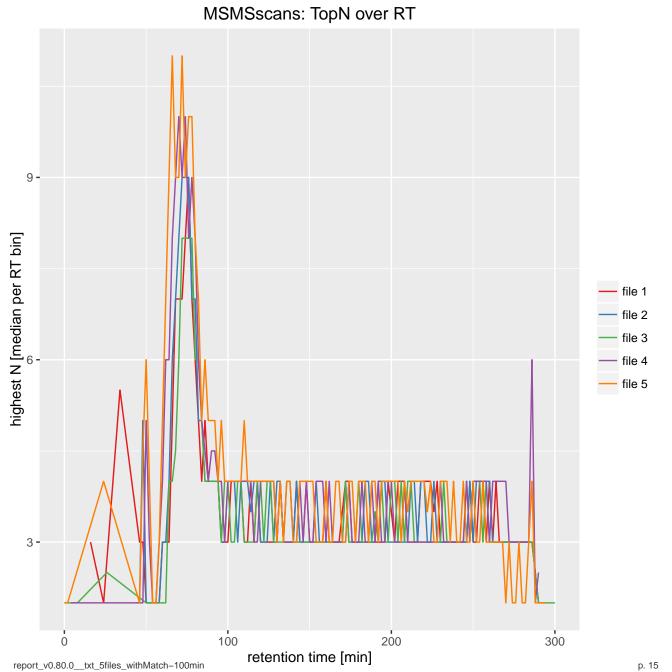


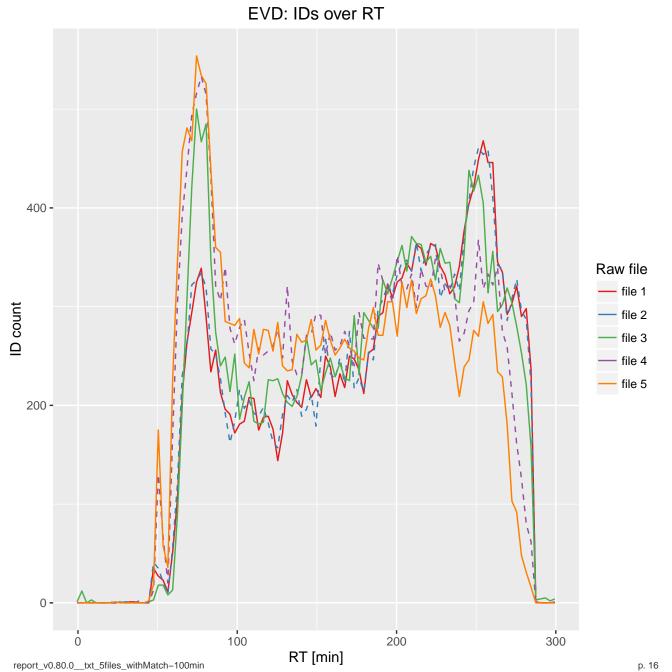
MSMS: Missed cleavages per Raw file

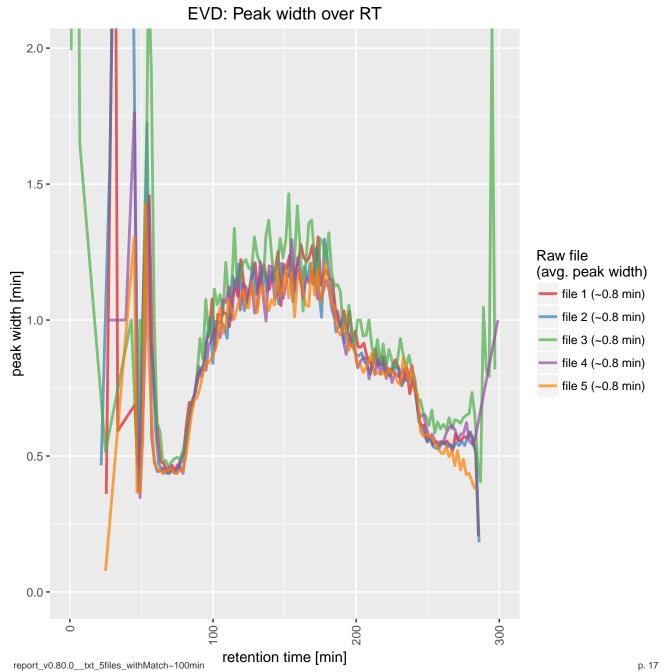


EVD: charge distribution file 1 file 2 charge 2 Raw file file 3 3 4 5 6 7 file 4 file 5 0.50 0.25 0.75 1.00 0.00 fraction [%] p. 13 report_v0.80.0__txt_5files_withMatch-100min





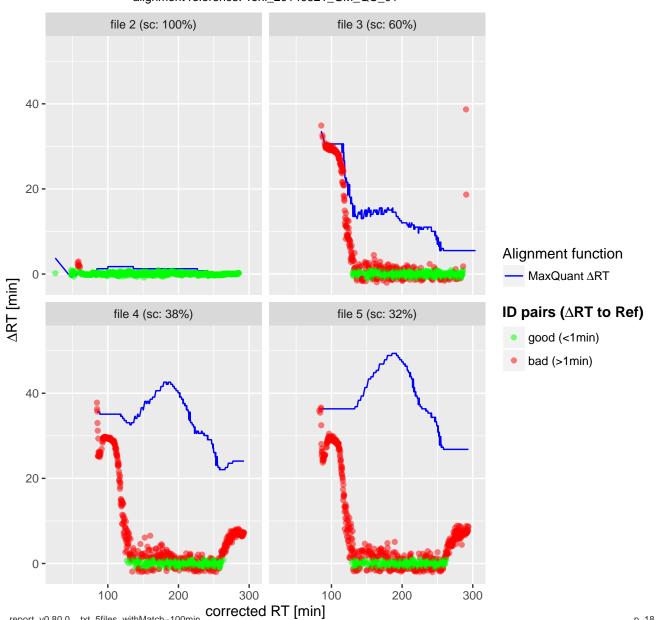




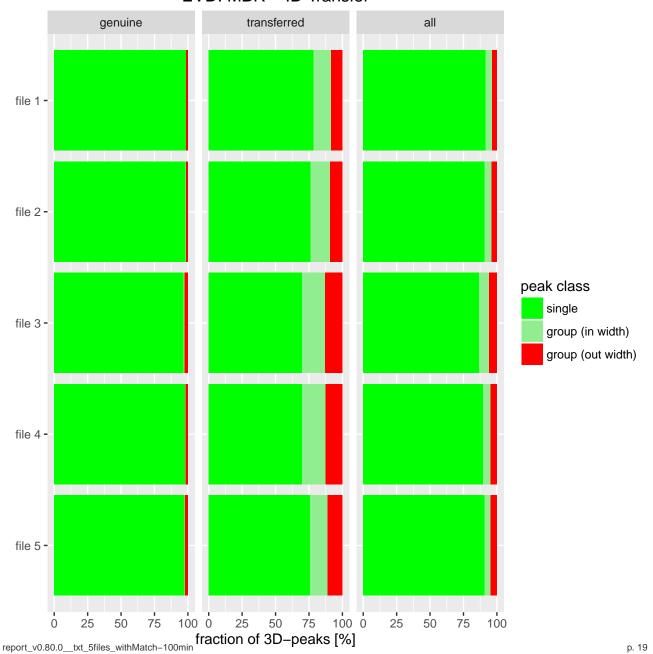
EVD: MBR – alignment

alignment reference: Toni_20140521_GM_QC_01

report_v0.80.0__txt_5files_withMatch=100min

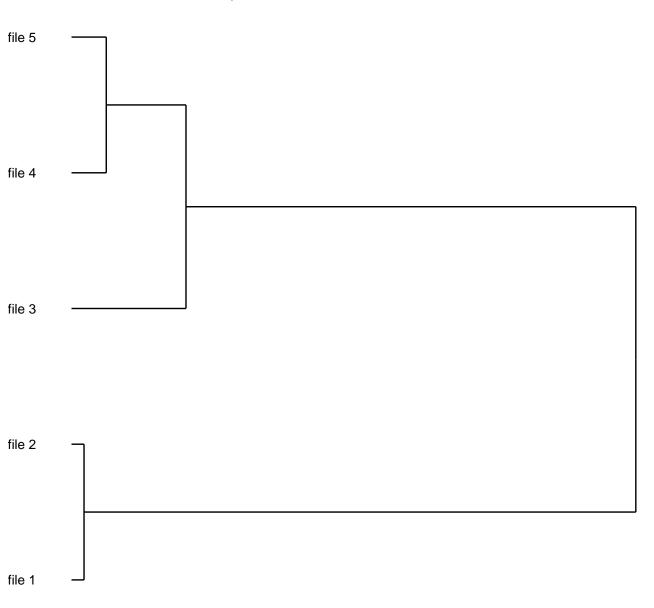


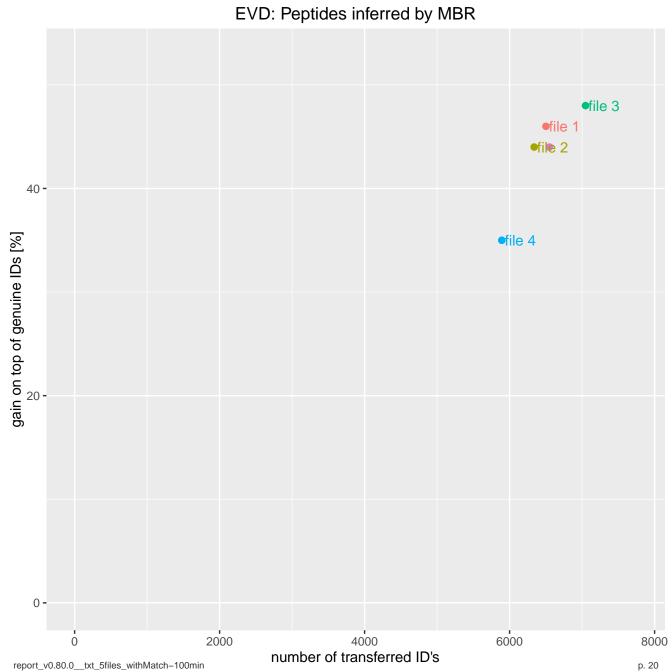
EVD: MBR - ID Transfer

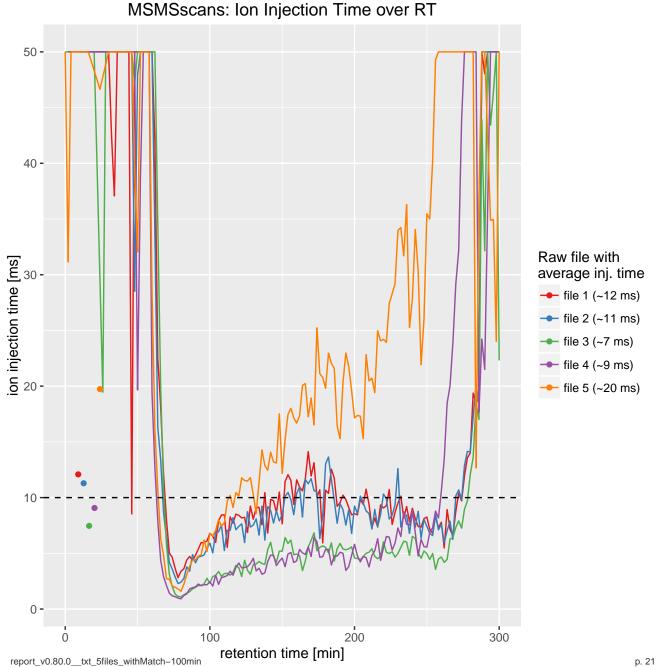


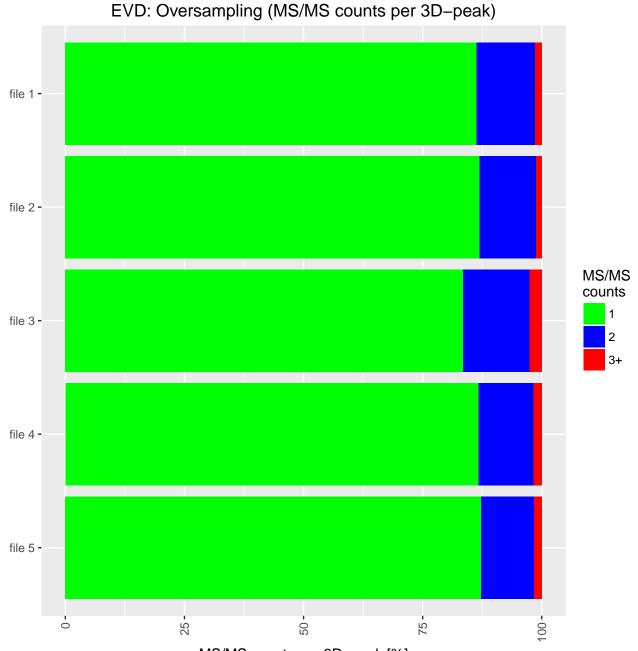
[experimental] EVD: Clustering Tree of Raw files

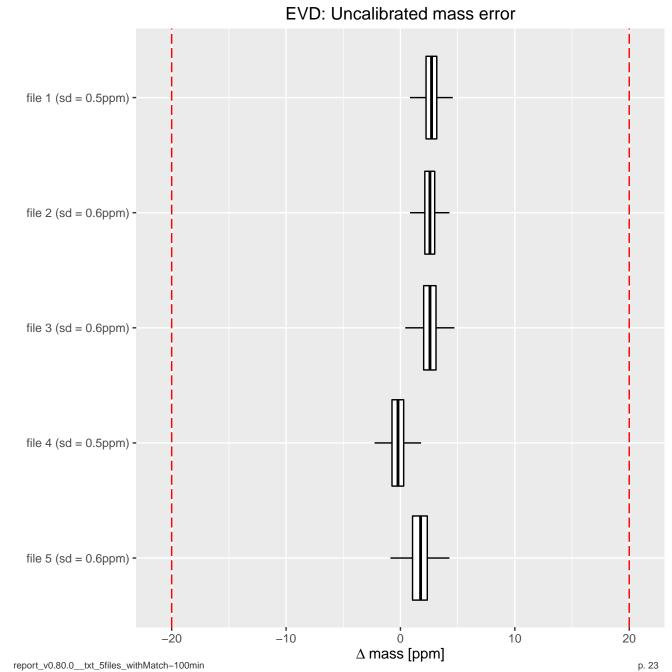
by Correlation of Corrected Retention Times

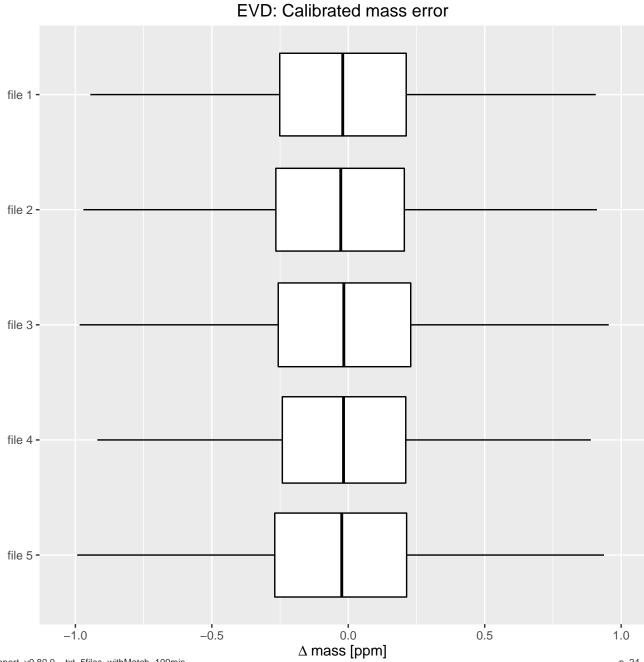




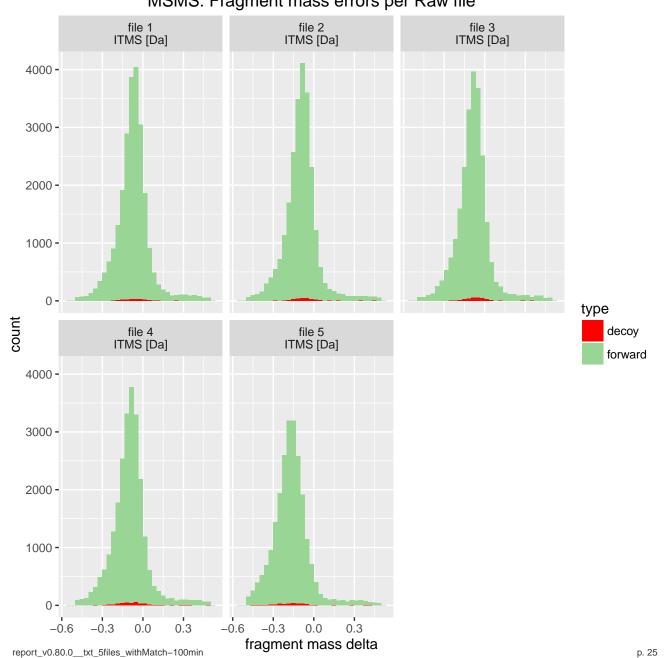


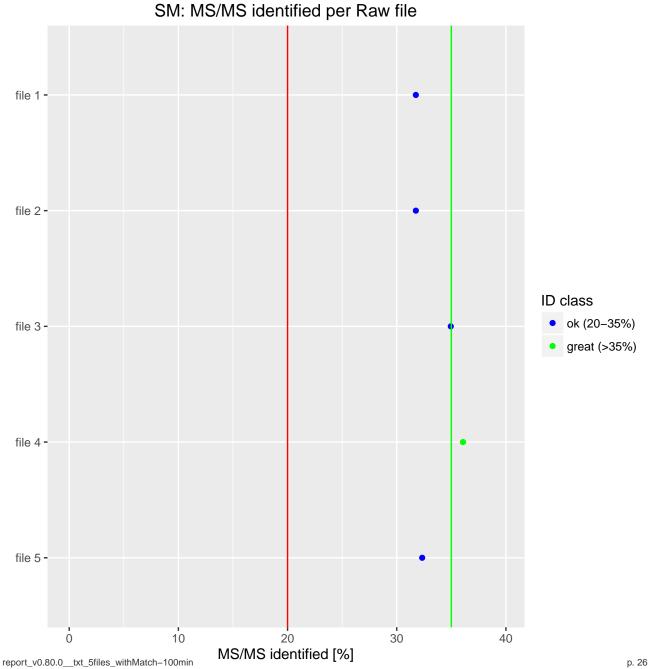


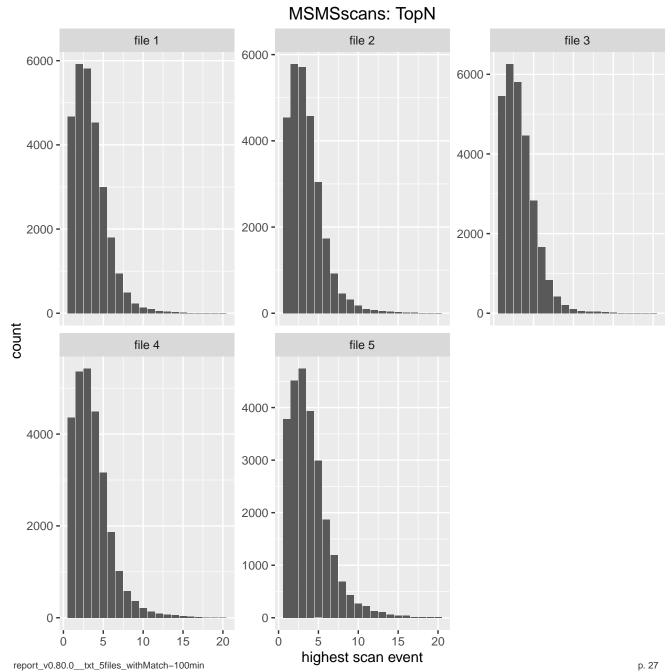




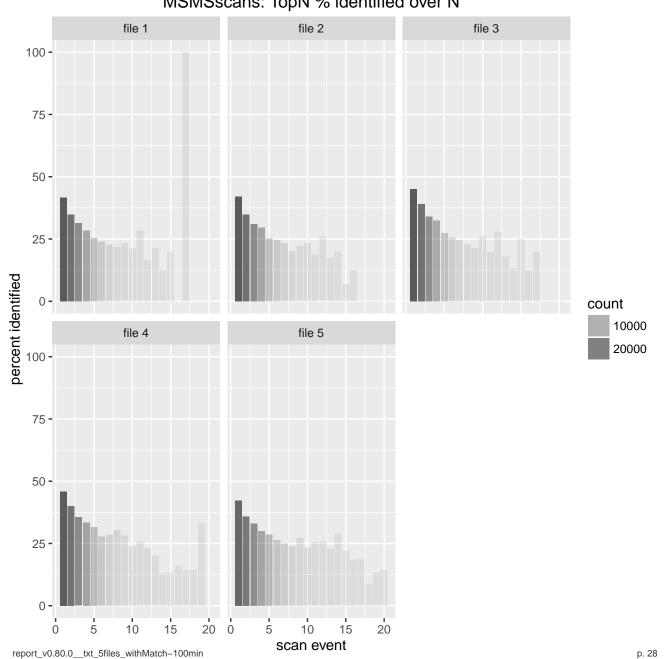
MSMS: Fragment mass errors per Raw file





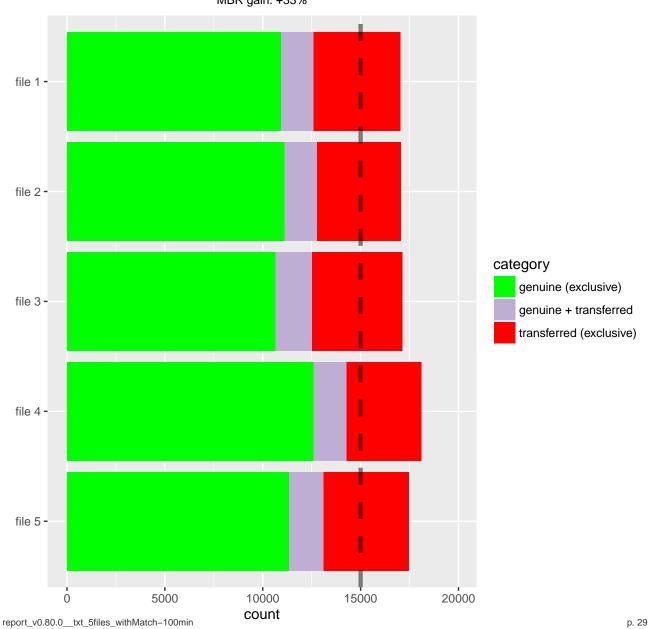


MSMSscans: TopN % identified over N



EVD: Peptide ID count

MBR gain: +33%



EVD: ProteinGroups count

