

Reading material for Stora Tentan at DBB

Everything marked in yellow is mandatory but a general understanding of all topics mentioned below is needed.

1. Resources from Intro course on the web

a. Biological Databases

- i. http://en.wikipedia.org/wiki/Biological_databases
- ii. <http://www.ploscompbiol.org/article/info%3Adoi%2F10.1371%2Fjournal.pcbi.0010034>
- iii. <http://www.uniprot.org/help/about>
- iv. http://en.wikipedia.org/wiki/List_of_biological_databases
- v. EMBD <http://www.emdatabank.org>
- vi. <http://en.wikipedia.org/wiki/Entrez>
- vii. http://metadatabase.org/wiki/Main_Page

b. Genes, sequencing and genomes

- i. <http://www.yourgenome.org/facts/what-is-a-genome>
- ii. <http://en.wikipedia.org/wiki/Bioinformatics>
- iii. <http://en.wikipedia.org/wiki/Genome>
- iv. https://en.wikipedia.org/wiki/Gene_prediction
- v. <https://en.wikipedia.org/wiki/Synteny>
- vi. http://en.wikipedia.org/wiki/Introduction_to_genetics
- vii. http://en.wikipedia.org/wiki/Human_genome
- viii. http://en.wikipedia.org/wiki/Genome_evolution
- ix. <http://en.wikipedia.org/wiki/Sequencing>
- x. [Are there laws of genome evolution ?](#)
- xi. <https://www.ebi.ac.uk/training/online/course/ebi-next-generation-sequencing-practical-course/what-you-will-learn/what-next-generation-dna->

c. Sequence alignment and searches

- i. https://en.wikipedia.org/wiki/Sequence_homology
- ii. http://en.wikipedia.org/wiki/Sequence_alignment
- iii. http://en.wikipedia.org/wiki/Needleman-Wunsch_algorithm
- iv. http://en.wikipedia.org/wiki/Substitution_matrix
- v. <http://en.wikipedia.org/wiki/BLOSUM>
- vi. http://en.wikipedia.org/wiki/Point_accepted_mutation
- vii. <http://bioinfo.se/courses/introduction-to-bioinformatics/lessons/sequence-alignments/>
- viii. http://en.wikipedia.org/wiki/Smith%E2%80%93Waterman_algorithm
- ix. <http://www.ncbi.nlm.nih.gov/blast/tutorial/>
- x. <http://en.wikipedia.org/wiki/BLAST>
- xi. <http://www.youtube.com/watch?v=HXEpBnUbAMo>
- xii. <http://en.wikipedia.org/wiki/FASTA>
- xiii. http://en.wikipedia.org/wiki/Sequence_alignment_software
- xiv. <http://www.ncbi.nlm.nih.gov/books/NBK1734/>
- xv. http://openwetware.org/wiki/Wikiomics:BLAST_tutorial

d. Multiple sequence alignments

- i. https://en.wikipedia.org/wiki/Multiple_sequence_alignment
- ii. https://en.wikipedia.org/wiki/Hidden_Markov_model
- e. Phylogeny and evolution
 - i. http://evolution.berkeley.edu/evolibrary/article/phylogenetics_01
 - ii. https://en.wikipedia.org/wiki/Phylogenetic_tree
 - 1. <https://en.wikipedia.org/wiki/UPGMA>
 - 2. https://en.wikipedia.org/wiki/Neighbor_joining
 - iii. <https://en.wikipedia.org/wiki/Phylogenomics>
 - iv. https://en.wikipedia.org/wiki/Phylogenetic_network
 - v. https://en.wikipedia.org/wiki/List_of_phylogenetics_software
 - vi. https://en.wikipedia.org/wiki/Phylogenetic_tree_viewers
 - vii. <https://en.wikipedia.org/wiki/Phylogenetics>
 - viii. https://www.ted.com/talks/svante_paaebo_dna_clues_to_our_inner_neanderthal?language=en
- f. Machine learning
 - i. https://en.wikipedia.org/wiki/Machine_learning
 - ii. https://en.wikipedia.org/wiki/Deep_learning
 - iii. <http://journals.plos.org/ploscompbiol/article?id=10.1371/journal.pcbi.0030116>
- g. Protein structure and folding
 - i. https://en.wikipedia.org/wiki/Protein_folding
 - ii. https://en.wikipedia.org/wiki/Homology_modeling
 - iii. https://en.wikipedia.org/wiki/Protein_structure_prediction
 - iv. https://en.wikipedia.org/wiki/Membrane_topology
- h. Systems biology
 - i. https://en.wikipedia.org/wiki/Biological_network
 - ii. <https://en.wikipedia.org/wiki/Interactome>
 - iii. https://en.wikipedia.org/wiki/Systems_biology
 - iv. https://en.wikipedia.org/wiki/Flux_balance_analysis
 - v. https://en.wikipedia.org/wiki/Metabolic_network_modelling

2. Alignments and sequence searches

- a. Needleman SB, Wunsch CD. **A general method applicable to the search for similarities in the amino acid sequence of two proteins.** J Mol Biol. 1970 Mar;48(3):443-53. [PubMed](#)
- b. Dayhoff et al, jA model of evolutionary change in proteins. Atlas of Protein Sequence and Structure 5 (3): 345–352 (1978) [PDF](#)
- c. Henikoff & Henikoff: Amino acid substitution matrices from protein blocks. PNAS 89 (22): 10915–9
- d. Smith TF, Waterman MS. **Identification of common molecular subsequences.** J Mol Biol. 1981 Mar 25;147(1):195-7. [PubMed](#)
- e. Altschul SF, Gish W, Miller W, Myers EW, Lipman DJ. **Basic local alignment search tool.** J Mol Biol. 1990 Oct 5;215(3):403-10.
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- g. Kent WJ. **BLAT--the BLAST-like alignment tool.** Genome Res. 2002 Apr;12(4):656-64. PMID: [PubMed](#)

- h. Hidden Markov models in computational biology. Applications to protein modeling. Krogh A, Brown M, Mian IS, Sjölander K, Haussler D. *J Mol Biol.* 1994 Feb 4;235(5):1501-31. [pdf](#)
 - i. A probabilistic model of local sequence alignment that simplifies statistical significance estimation. Eddy SR. *PLoS Comput Biol.* 2008 May 30;4(5):e1000069. [PDF](#)
 - j. A tutorial on hidden Markov models and selected applications in speech recognition. Lawrence R Rabiner, [PDF](#)
 - k. Söding J. **Protein homology detection by HMM-HMM comparison.** *Bioinformatics.* 2005 Apr 1;21(7):951-60. Epub 2004 Nov 5. Erratum in: *Bioinformatics.* 2005 May 1;21(9):2144. PMID: [PubMed](#)
 - l. **Review of Common Sequence Alignment Methods: Clues to Enhance Reliability** Christophe Lambert*, Jean-Marc Van Campenhout, Xavier DeBolle and Eric Depiereux *Current Genomics*, 2003, 4, 131-146
3. Fold Recognition
- a. Bowie JU, Luthy R, Eisenberg D. **A method to identify protein sequences that fold into a known three-dimensional structure.** *Science.* 1991 Jul 12;253(5016):164-70, [PubMed](#)
 - b. Jones DT, Taylor WR, Thornton JM. **A new approach to protein fold recognition.** *Nature.* 1992 Jul 2;358(6381):86-9, [PubMed](#)
 - c. Park J, Karplus K, Barret C, Hughey R, Haussler D, Hubbard T, Chothia C. **Sequence Comparisons Using Multiple Sequences Detect Three Times as Many Remote Homologues as Pairwise Methods.** *J. Mol. Biol.*(1998) 284, 1201-1210. [PubMed](#)
 - d. Simons, KT, Ruczinski, I, & Kooperberg, C. (1999). Improved recognition of native like protein structures using a combination of sequence-dependent and sequence independent features of proteins. *Proteins: Structure.*
 - e. Fischer D, Eisenberg D. **Protein fold recognition using sequence-derived predictions.** *Protein Sci.* 1996 May;5(5):947-55. [PubMed](#)
4. Protein Domains
- a. Holm L, Sander C. **Mapping the protein universe.** *Science.* 1996 Aug 2;273(5275):595-603. Review. [PubMed](#)
 - b. Murzin AG, Brenner SE, Hubbard T, Chothia C. **Abstract SCOP: a structural classification of proteins database for the investigation of sequences and structures.** *J Mol Biol.* 1995 Apr 7;247(4):536-40. [PubMed](#) [PDF](#)
 - c. Sonnhammer EL, Eddy SR, Durbin R. **Abstract Pfam: a comprehensive database of protein domain families based on seed alignments.** *Proteins.* 1997 Jul;28(3):405-20. [PubMed](#)
 - d. Gordana Apic, Julian Gough, Sarah A Teichmann, **Domain combinations in archaeal, eubacterial and eukaryotic proteomes,** *Journal of Molecular Biology*, Volume 310, Issue 2, 6 July 2001, Pages 311-325, ISSN 0022-2836 [Link](#)
 - e. Gabrielle A. Reeves, Timothy J. Dallman, Oliver C. Redfern, Adrian Akpor, Christine A. Orengo, **Structural Diversity of Domain Superfamilies in the CATH Database** *Journal of Molecular Biology*, Volume 360, Issue 3, 14 July 2006, Pages 725-741 [Link](#)
 - f. Levy ED, Boeri Erba E, Robinson CV, Teichmann SA. **Assembly reflects evolution of protein complexes.** *Nature.* 2008 Jun 26;453(7199):1262-5. Epub 2008 Jun 18. [DOI](#)

- g. **Domain Rearrangements in Protein Evolution** Asa K. Bjorklund, Diana Ekman, Sara Light, Johannes Frey-Skott and Arne Elofsson *J. Mol. Biol.* (2005) 353, 911–923
5. Secondary Structure Predictions
 - a. Chou PY, Fasman GD. **Conformational parameters for amino acids in helical, beta-sheet, and random coil regions calculated from proteins.** *Biochemistry.* 1974 Jan 15;13(2):211-22. [PDF](#)
 - b. Chou PY, Fasman GD. **Prediction of protein conformation.** *Biochemistry.* 1974 Jan 15;13(2):222-45. [PubMed](#) [PDF](#)
 - c. **Analysis of the accuracy and implications of simple methods for predicting the secondary structure of globular proteins.** Garnier J, Osguthorpe DJ, Robson B. *J Mol Biol.* 1978 Mar 25;120(1):97-120. PMID: [PubMed](#)
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 6. Protein structure modeling
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 7. Protein Structure Prediction
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 - d. Weigt, Martin, White, Robert A, Szurmant, Hendrik, Hoch, James A, & Hwa, Terence. (2009). **Identification of direct residue contacts in protein-protein interaction by message passing.** *Proc. Natl. Acad. Sci* **106**(1), 67–72.
 - e. Simons, KT., Kooperberg, C., Huang, E., & Baker, D. (1997). Assembly of protein tertiary structures from fragments with similar local sequences using simulated annealing and Bayesian scoring functions. *J. Mol. Biol.* **268**(1), 209–225.
 - f. Simons, KT, Ruczinski, I, & Kooperberg, C. (1999). Improved recognition of native like protein structures using a combination of sequence-dependent and sequence independent features of proteins. *Proteins: Structure.*
 - g. Marks DS, Colwell LJ, Sheridan R, Hopf TA, Pagnani A, Zecchina R, Sander C. **Protein 3D structure computed from evolutionary sequence variation.** *PLoS One.* 2011;6(12):e28766.
 - h. Marks DS, Hopf TA, Sander C. **Protein structure prediction from sequence variation.** *Nat Biotechnol.* 2012 Nov;30(11):1072-80
 8. Motif finding
 - a. Lawrence, C. E., Altschul, S. F., Boguski, M. S., Liu, J. S., Neuwald, A. F., & Wootton, J. C. **Detecting subtle sequence signals: a Gibbs sampling strategy for multiple alignment.** *Science,* 1993, 262(5131), 208-14. [PubMed](#)
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 9. Subcellular sorting

- a. **Identification of prokaryotic and eukaryotic signal peptides and prediction of their cleavage sites.** Nielsen H, Engelbrecht J, Brunak S, von Heijne G. *Protein Eng.* 1997 Jan;10(1):1-6. PMID: [PubMed](#) [PDF](#)
 - b. **A new method for predicting signal sequence cleavage sites.** von Heijne G. *Nucleic Acids Res.* 1986 Jun 11;14(11):4683-90. PMID: [PubMed](#) [PDF](#)
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- a. **Predicting transmembrane protein topology with a hidden Markov model: application to complete genomes.** Krogh A, Larsson B, von Heijne G, Sonnhammer EL. *J Mol Biol.* 2001 Jan 19;305(3):567-80. PMID: [PubMed](#) [PDF](#)
 - b. **Membrane protein structure prediction. Hydrophobicity analysis and the positive-inside rule.** von Heijne G. *J Mol Biol.* 1992 May 20;225(2):487-94. PMID: [PubMed](#) [PDF](#)
 - c. **Membrane protein structure: prediction versus reality.** Elofsson A, von Heijne G. *Annu Rev Biochem.* 2007;76:125-40. Review. PMID: 17579561
11. Molecular Modelling and computational chemistry
- a. Chothia C, Levitt M, Richardson D. **Structure of proteins: Packing of alpha-helices and pleated sheets** *Proc Natl Acad Sci U S A.* 1977 Oct;74(10):4130-4. [PDF](#)
 - b. McCammon JA, Gelin BR, Karplus M. **Dynamics of folded proteins.** *Nature.* 1977 Jun 16;267(5612):585-90. [PDF](#)
 - c. Levinthal C. **Molecular model-building by computer.** *Sci Am.* 1966 Jun;214(6):42-52. [PDF](#)
 - d. Bradley P, Misura KM, Baker D. **Toward high-resolution de novo structure prediction for small proteins.** *Science.* 2005 Sep 16;309(5742):1868-71. [PDF](#) [Supplement](#)
 - e. Kohlhoff, K. et al, Cloud-based simulations on Google Exacycle reveal ligand modulation of GPCR activation pathways, *Nature Chem.* 6 (2014), 15-21
12. Molecular Docking
- a. Kuntz, ID; Blaney, JM; OATley, SJ; et al. **A geometric approach to macromolecule-ligand interactions** *Journal of Molecular BIOLOGY* Volume: 161 Issue: 2 Pages: 269-288
 - b. Meng, EC; SHOichet, BK; Kuntz, ID **Automated docking with grid-based energy evaluation** *JOURNAL OF COMPUTATIONAL CHEMISTRY* Volume: 13 Issue: 4 Pages: 505- 524 DOI: 10.1002/jcc.540130412 Published: MAY 1992
 - c. BOHM, HJ **The development of a simple empirical scoring function to estimate the binding constant for a protein ligand complex of known 3- dimensional structure** *Journal of Computer-Aided Molecular Design* Volume: 8 Issue: 3 Pages: 243-256 DOI: 10.1007/BF00126743 Published: JUN 1994
 - d. Muegge, I; Martin, YC **A general and fast scoring function for protein-ligand interactions: A simplified potential approach** *Journal of Medicinal Chemistry* Volume: 42 Issue: 5 Pages: 791-804 DOI: 10.1021/jm980536j Published: MAR 11 1999
 - e. Warren, Gregory L.; Andrews, C. Webster; Capelli, Anna-Maria; et al. **A critical assessment of docking programs and scoring functions** Source: *JOURNAL OF MEDICINAL CHEMISTRY* Volume: 49 Issue: 20 Pages: 5912- 5931 DOI: 10.1021/jm050362n Published: OCT 5 2006
13. Protein Design

- a. Dahiyat BI, Mayo SL. **De novo protein design: fully automated sequence selection.** *Science*. 1997 Oct 3;278(5335):82-7. [PubMed](#)
- b. Dalal S, Balasubramanian S, Regan L. **Protein alchemy: changing beta-sheet into alpha-helix.** *Nat Struct Biol*. 1997 Jul;4(7):548-52. [PubMed](#) [PDF](#)
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- f. Procko E, Berguig GY, Shen BW et al. **A computationally designed inhibitor of an Epstein-Barr viral Bcl-2 protein induces apoptosis in infected cells.** *Cell*. 2014 Jun 19;157(7):1644-56, [PubMed](#)
- g. Huang PS, Boyken SE, Baker D. **The coming of age of de novo protein design.** *Nature*. 2016 Sep 15;537(7620):320-7. [PubMed](#)

14. Protein protein interactions

- a. Tsoka S, Ouzounis CA. **Prediction of protein interactions: metabolic enzymes are frequently involved in gene fusion.** *Nat Genet*. 2000 Oct;26(2):141-2, [PubMed](#)
- b. Marcotte EM, Pellegrini M, Thompson MJ, Yeates TO, Eisenberg D. **A combined algorithm for genome-wide prediction of protein function.** *Nature*. 1999 Nov 4;402(6757):83-6 [PubMed](#)
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- d. Doolittle RF. **Do you dig my groove?** *Nature Genetics*, 1999, 23, 6-8 [PDF](#)
- e. Enright AJ, Ouzounis CA. **Functional associations of proteins in entire genomes by means of exhaustive detection of gene fusions** *Genome Biology*, 2001, 2, 0034.1-0034.7 [PDF](#)
- f. Snel B, Bork P, Huynen M. **Genome evolution: Gene fusion versus gene fission** *TIG*, 2000, 16, 9-11 [PDF](#)
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- h. **What properties characterize the hub proteins of the protein-protein interaction network of *Saccharomyces cerevisiae*?** Diana Ekman, Sara Light, Åsa K Björklund and Arne Elofsson *Genome Biology* 2006, 7:R45 (doi:10.1186/gb-2006-7-6-r45)
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15. Protein Networks

- a. Ideker TE, Thorsson V, Karp RM. **Discovery of regulatory interactions through perturbation: inference and experimental design.** *Pac Symp Biocomput* 2000;:305-16. [PDF](#), [PubMed](#)

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16. Sequencing and genomes

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- c. Margulies M et al. **Genome sequencing in microfabricated high-density picolitre reactors.** Nature. 2005 Sep 15;437(7057):376-80 [PubMed](#)
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17. Gene assignment

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18. Genome assembly

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19. Gene mapping and genotyping

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20. Molecular Evolution

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