## class 10: strucutral bioinformatics p1

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### The PDB Database

The main repository of biomolecular structure data is called the Protein Data Bank (PDB for short). It is the second oldest database (after GenBank).

What is currently in the PDB?

```
stats <- read.csv("Data Export Summary.csv", row.names = 1)
stats</pre>
```

	V 53.4	EM	NMD	Multiple methods	Noutron	0+bor
5	X.ray			Multiple.methods		
Protein (only)	171,959	18,083	12,622	210	84	32
Protein/Oligosaccharide	10,018	2,968	34	10	2	0
Protein/NA	8,847	5,376	286	7	0	0
Nucleic acid (only)	2,947	185	1,535	14	3	1
Other	170	10	33	Θ	0	0
Oligosaccharide (only)	11	0	6	1	0	4
	Total					
Protein (only)	202,990					
Protein/Oligosaccharide	13,032					
Protein/NA	14,516					
Nucleic acid (only)	4,685					
Other	213					
Oligosaccharide (only)	22					

Q1: What percentage of structures in the PDB are solved by X-Ray and Electron Microscopy.

```
x <- stats$X.ray
y <- gsub(",", "", stats$X.ray)
sum(as.numeric(y))</pre>
```

#### [1] 193952

```
comma.sum <- function(x) {
  y <- gsub(",", "", x)

  return( sum(as.numeric(y)))
}</pre>
```

```
xray.sum <- comma.sum(stats$X.ray)
xray.sum</pre>
```

#### [1] 193952

```
xray.sum <- comma.sum(stats$X.ray)
em.sum <- comma.sum(stats$EM)
total.sum <- comma.sum(stats$Total)</pre>
```

```
xray.sum/total.sum * 100
```

```
[1] 82.37223
```

```
em.sum/total.sum * 100
```

```
[1] 11.30648
```

Q2: What proportion of structures in the PDB are protein?

```
total.num <- as.numeric(gsub(",", "", stats$Total))
sum(total.num[1:3]/total.sum)</pre>
```

```
[1] 0.9791046
```

Q3: Type HIV in the PDB website search box on the home page and determine how many HIV-1 protease structures are in the current PDB?

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## 2. Visualizing with Mol-star

Explore the hIV-1 protease structure with PDB code: 1HSG Mol-star homepage at https:://molstar.org/viewer/.



Figure 1: Figure 1. A first view of HIV-Pr

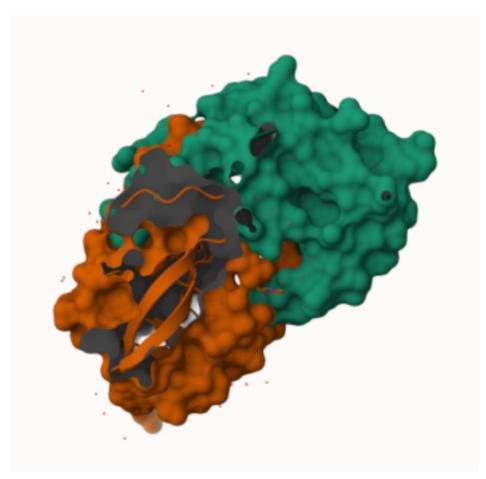


Figure 2: Figure 2. Space-fill model for whole protein



Figure 3: Figure 3: Aspartic Acid and Water Spacefill

# 3. Visualization in R

library(bio3d)

pdb <- read.pdb("1HSG")</pre>

Note: Accessing on-line PDB file

pdb

```
Call: read.pdb(file = "1HSG")

Total Models#: 1
   Total Atoms#: 1686, XYZs#: 5058 Chains#: 2 (values: A B)

Protein Atoms#: 1514 (residues/Calpha atoms#: 198)
Nucleic acid Atoms#: 0 (residues/phosphate atoms#: 0)

Non-protein/nucleic Atoms#: 172 (residues: 128)
Non-protein/nucleic resid values: [ HOH (127), MK1 (1) ]

Protein sequence:
   PQITLWQRPLVTIKIGGQLKEALLDTGADDTVLEEMSLPGRWKPKMIGGIGGFIKVRQYD
   QILIEICGHKAIGTVLVGPTPVNIIGRNLLTQIGCTLNFPQITLWQRPLVTIKIGGQLKE
   ALLDTGADDTVLEEMSLPGRWKPKMIGGIGGFIKVRQYDQILIEICGHKAIGTVLVGPTP
   VNIIGRNLLTQIGCTLNF
+ attr: atom, xyz, seqres, helix, sheet,
   calpha, remark, call
```

### attributes(pdb\$atom)

\$names [1] "type" [9] "x"		] "type" "eleno"		,		"alt" "o"	"resid" "b"			"chain" "segid"		"resno" "elesy"		ert" rge"
\$row.na	\$row.names													
[1]	1	2	3	4	5	6	7	8	9	10	11	12	13	14
[15]	15	16	17	18	19	20	21	22	23	24	25	26	27	28
[29]	29	30	31	32	33	34	35	36	37	38	39	40	41	42
[43]	43	44	45	46	47	48	49	50	51	52	53	54	55	56
[57]	57	58	59	60	61	62	63	64	65	66	67	68	69	70
[71]	71	72	73	74	75	76	77	78	79	80	81	82	83	84
[85]	85	86	87	88	89	90	91	92	93	94	95	96	97	98
[99]	99	100	101	102	103	104	105	106	107	108	109	110	111	112
[113]	113	114	115	116	117	118	119	120	121	122	123	124	125	126
[127]	127	128	129	130	131	132	133	134	135	136	137	138	139	140
[141]	141	142	143	144	145	146	147	148	149	150	151	152	153	154
[155]	155	156	157	158	159	160	161	162	163	164	165	166	167	168

[169]	169	170	171	172	173	174	175	176	177	178	179	180	181	182
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[197]	197	198	199	200	201	202	203	204	205	206	207	208	209	210
[211]	211	212	213	214	215	216	217	218	219	220	221	222	223	224
[225]	225	226	227	228	229	230	231	232	233	234	235	236	237	238
[239]	239	240	241	242	243	244	245	246	247	248	249	250	251	252
[253]	253	254	255	256	257	258	259	260	261	262	263	264	265	266
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[295]	295	296	297	298	299	300	301	302	303	304	305	306	307	308
[309]	309	310	311	312	313	314	315	316	317	318	319	320	321	322
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[561]	561	562	563	564	565	566	567	568	569	570	571	572	573	574
[575]	575	576	577	578	579	580	581	582	583	584	585	586	587	588
[589]	589	590	591	592	593	594	595	596	597	598	599	600	601	602
[603]	603	604	605	606	607	608	609	610	611	612	613	614	615	616
[617]	617	618	619	620	621	622	623	624	625	626	627	628	629	630
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[645]	645	646	647	648	649	650	651	652	653	654	655	656	657	658
[659]	659	660	661	662	663	664	665	666	667	668	669	670	671	672
[673]	673	674	675	676	677	678	679	680	681	682	683	684	685	686
[687]	687	688	689	690	691	692	693	694	695	696	697	698	699	700
[701]	701	702	703	704	705	706	707	708	709	710	711	712	713	714
[715]	715	716	717	718	719	720	721	722	723	724	725	726	727	728
[729]	729	730	731	732	733	734	735	736	737	738	739	740	741	742
[743]	743	744	745	746	747	748	749	750	751	752	753	754	755	756
[757]	757	758	759	760	761	762	763	764	765	766	767	768	769	770
[771]	771	772	773	774	775	776	777	778	779	780	781	782	783	784
[785]	785	786	787	788	789	790	791	792	793	794	795	796	797	798
[799]	799	800	801	802	803	804	805	806	807	808	809	810	811	812
[813]	813	814	815	816	817	818	819	820	821	822	823	824	825	826
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[1121] 1121 1122 1123 1124 1125 1126 1127 1128 1129 1130 1131 1132 1133 1134
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[1205] 1205 1206 1207 1208 1209 1210 1211 1212 1213 1214 1215 1216 1217 1218
[1219] 1219 1220 1221 1222 1223 1224 1225 1226 1227 1228 1229 1230 1231 1232
[1233] 1233 1234 1235 1236 1237 1238 1239 1240 1241 1242 1243 1244 1245 1246
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[1261] 1261 1262 1263 1264 1265 1266 1267 1268 1269 1270 1271 1272 1273 1274
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[1471] 1471 1472 1473 1474 1475 1476 1477 1478 1479 1480 1481 1482 1483 1484
[1485] 1485 1486 1487 1488 1489 1490 1491 1492 1493 1494 1495 1496 1497 1498
[1499] 1499 1500 1501 1502 1503 1504 1505 1506 1507 1508 1509 1510 1511 1512
```

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[1513] 1513 1514 1515 1516 1517 1518 1519 1520 1521 1522 1523 1524 1525 1526 [1527] 1527 1528 1529 1530 1531 1532 1533 1534 1535 1536 1537 1538 1539 1540 [1541] 1541 1542 1543 1544 1545 1546 1547 1548 1549 1550 1551 1552 1553 1554 [1555] 1555 1556 1557 1558 1559 1560 1561 1562 1563 1564 1565 1566 1567 1568 [1569] 1569 1570 1571 1572 1573 1574 1575 1576 1577 1578 1579 1580 1581 1582 [1583] 1583 1584 1585 1586 1587 1588 1589 1590 1591 1592 1593 1594 1595 1596 [1597] 1597 1598 1599 1600 1601 1602 1603 1604 1605 1606 1607 1608 1609 1610 [1611] 1611 1612 1613 1614 1615 1616 1617 1618 1619 1620 1621 1622 1623 1624 [1625] 1625 1626 1627 1628 1629 1630 1631 1632 1633 1634 1635 1636 1637 1638 [1639] 1639 1640 1641 1642 1643 1644 1645 1646 1647 1648 1649 1650 1651 1652 [1653] 1653 1654 1655 1656 1657 1658 1659 1660 1661 1662 1663 1664 1665 1666 [1667] 1667 1668 1669 1670 1671 1672 1673 1674 1675 1676 1677 1678 1679 1680 [1681] 1681 1682 1683 1684 1685 1686
```

We can see atom data with pdb\$atom:

```
head (pdb$atom)
```

```
type eleno elety alt resid chain resno insert
                                                       Х
                                                                     Z 0
                                                              У
                          PR0
1 ATOM
           1
                 N <NA>
                                  Α
                                         1
                                             <NA> 29.361 39.686 5.862 1 38.10
2 ATOM
           2
                CA <NA>
                          PR0
                                  Α
                                         1
                                             <NA> 30.307 38.663 5.319 1 40.62
                C <NA>
3 ATOM
           3
                          PR0
                                        1
                                             <NA> 29.760 38.071 4.022 1 42.64
                                  Α
                 0 <NA>
                          PR0
4 ATOM
           4
                                        1
                                             <NA> 28.600 38.302 3.676 1 43.40
5 ATOM
           5
                CB <NA>
                          PR0
                                  Α
                                        1
                                             <NA> 30.508 37.541 6.342 1 37.87
6 ATOM
           6
                CG <NA>
                          PR0
                                             <NA> 29.296 37.591 7.162 1 38.40
  segid elesy charge
1 <NA>
                <NA>
            N
2 <NA>
            C
                <NA>
3
  <NA>
            C
                <NA>
4
  <NA>
            0
                <NA>
5
  <NA>
            C
                <NA>
6 <NA>
                <NA>
```

```
head(pdbseq(pdb))
```

```
1 2 3 4 5 6
"P" "Q" "I" "T" "L" "W"
```

We can make quick 3d viz.

```
library(bio3dview)
```

### library(NGLVieweR)

```
#view.pdb(pdb) |>
    #setSpin()
```

```
#view.pdb(pdb, backgroundColor = "pink", colorScheme = "sse")
```

```
library(bio3dview)
library(NGLVieweR)
```

## Predicting functional motions of a single structure

We can finish off today with a bioinformatics prediction of the functional motions of a protein.

We will run a Normal Mode Analysis. (NMA)

```
#adk <- read.pdb("6s36")
#adk</pre>
```

```
#m <- nma(adk)
#plot(m)</pre>
```

```
#view.nma(m)
```

We can write out a trajectory of the predicted dynamics and view this in Mol-star.

```
#mktrj(m, file="nma.pdb")
```