



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:46 PM GMT

PDB ID : 1M10
Title : Crystal structure of the complex of Glycoprotein Ib alpha and the von Willebrand Factor A1 Domain
Authors : Huizinga, E.G.; Tsuji, S.; Romijn, R.A.P.; Schiphorst, M.E.; de Groot, P.G.; Sixma, J.J.; Gros, P.
Deposited on : 2002-06-16
Resolution : 3.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

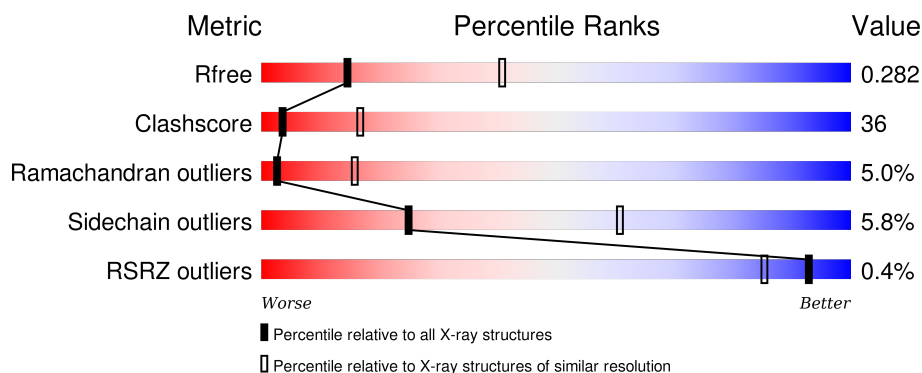
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

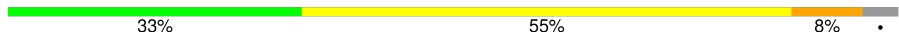
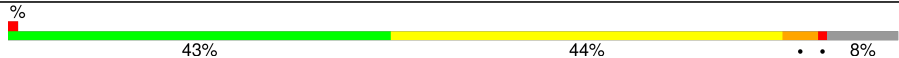
The reported resolution of this entry is 3.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1114 (3.14-3.06)
Clashscore	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	208	 33% 55% 8% .
2	B	290	 % 43% 44% . . 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3688 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called von Willebrand Factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	199	Total	C	N	O	S	0	0	0
			1601	1022	282	291	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	543	GLN	ARG	ENGINEERED	UNP P04275

- Molecule 2 is a protein called Glycoprotein Ib alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	267	Total	C	N	O	S	0	0	0
			2087	1344	346	389	8			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	21	GLN	ASN	ENGINEERED	UNP P07359
B	159	GLN	ASN	ENGINEERED	UNP P07359
B	239	VAL	MET	ENGINEERED	UNP P07359

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	89.84Å 89.84Å 124.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.50 – 3.10 30.50 – 3.09	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.50-3.10) 100.0 (30.50-3.09)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 3.11Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.296 0.230 , 0.282	Depositor DCC
R_{free} test set	529 reflections (5.09%)	DCC
Wilson B-factor (Å ²)	64.0	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.4	EDS
Estimated twinning fraction	0.057 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 10454 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3688	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.45	0/1630	0.70	0/2197
2	B	0.46	0/2134	0.75	0/2916
All	All	0.45	0/3764	0.73	0/5113

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1601	0	1645	146	0
2	B	2087	0	2130	128	0
All	All	3688	0	3775	269	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 36.

All (269) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:559:HIS:HA	1:A:592:ALA:HA	1.24	1.16
1:A:636:ARG:HH11	1:A:636:ARG:HA	1.33	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:545:ARG:HH22	1:A:703:PRO:HG3	1.29	0.93
1:A:559:HIS:CD2	1:A:560:ASP:H	1.89	0.90
1:A:512:LEU:HD22	1:A:611:ARG:HH21	1.40	0.84
1:A:636:ARG:NH1	1:A:636:ARG:HA	1.93	0.83
2:B:23:THR:HA	2:B:42:LEU:O	1.77	0.82
2:B:158:ASN:HB2	2:B:182:ASN:HD21	1.43	0.81
2:B:203:HIS:HE1	2:B:230:TRP:H	1.29	0.80
2:B:137:LYS:O	2:B:160:LEU:HA	1.84	0.77
2:B:232:GLN:NE2	2:B:237:LYS:HE2	2.00	0.77
2:B:92:LEU:HD22	2:B:115:LEU:HD22	1.65	0.77
1:A:682:GLU:HA	1:A:685:GLN:NE2	2.01	0.76
2:B:187:ILE:HG13	2:B:216:PHE:CD1	2.20	0.76
2:B:49:ALA:HB2	2:B:73:ASP:HB2	1.68	0.76
1:A:519:LEU:HD12	1:A:519:LEU:N	2.00	0.76
1:A:545:ARG:HH22	1:A:703:PRO:CG	2.00	0.75
2:B:203:HIS:CE1	2:B:230:TRP:H	2.05	0.74
1:A:599:LYS:O	1:A:603:PHE:HD2	1.70	0.73
2:B:95:LEU:HD22	2:B:99:LEU:HD12	1.68	0.73
1:A:511:ARG:NH1	1:A:695:CYS:HB3	2.05	0.72
1:A:548:GLN:HA	1:A:552:ARG:HG2	1.70	0.72
1:A:576:GLU:OE1	1:A:576:GLU:HA	1.89	0.71
1:A:639:GLN:HG2	1:A:668:GLN:OE1	1.91	0.70
2:B:232:GLN:NE2	2:B:237:LYS:CE	2.55	0.70
1:A:545:ARG:NH2	1:A:703:PRO:HG3	2.04	0.70
1:A:559:HIS:CA	1:A:592:ALA:HA	2.14	0.69
1:A:550:TRP:CE3	1:A:550:TRP:HA	2.28	0.68
1:A:559:HIS:HA	1:A:592:ALA:CA	2.15	0.68
2:B:144:LEU:HD11	2:B:153:LEU:HD22	1.76	0.68
2:B:75:THR:HA	2:B:98:THR:HG23	1.74	0.68
1:A:543:GLN:O	1:A:544:LEU:HG	1.93	0.67
1:A:676:VAL:C	1:A:677:LEU:HD12	2.14	0.67
1:A:513:LEU:HB3	1:A:551:VAL:HG22	1.75	0.67
1:A:517:PHE:CE1	1:A:619:LEU:HD23	2.30	0.67
1:A:546:ILE:HG21	1:A:577:LEU:HD12	1.76	0.67
2:B:128:GLU:HB3	2:B:130:TYR:CE1	2.30	0.67
2:B:1:HIS:CG	2:B:2:PRO:HD2	2.30	0.66
1:A:559:HIS:HD2	1:A:560:ASP:H	1.43	0.66
2:B:22:LEU:HD12	2:B:41:ASN:ND2	2.12	0.65
1:A:538:VAL:O	1:A:542:GLU:HG3	1.96	0.65
1:A:523:SER:C	1:A:525:LEU:H	2.00	0.65
1:A:658:ASN:OD1	1:A:661:GLN:HG3	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:666:GLU:HA	1:A:672:ASN:O	1.96	0.64
1:A:511:ARG:HA	1:A:701:ALA:HB2	1.79	0.64
1:A:609:ILE:HG23	1:A:616:ARG:NH2	2.13	0.64
1:A:645:LYS:HD2	1:A:645:LYS:N	2.13	0.64
2:B:34:THR:HB	2:B:58:THR:HG22	1.80	0.64
1:A:508:TYR:C	1:A:508:TYR:CD1	2.71	0.63
2:B:248:CYS:HB2	2:B:251:SER:HB2	1.79	0.63
2:B:115:LEU:O	2:B:140:PRO:HG3	1.99	0.63
1:A:510:SER:O	1:A:511:ARG:HG3	1.99	0.63
2:B:259:TYR:CE2	2:B:261:GLY:HA2	2.33	0.63
2:B:111:ARG:HD3	2:B:135:GLU:OE2	2.00	0.62
1:A:545:ARG:NH1	1:A:703:PRO:HA	2.14	0.62
1:A:533:LEU:HD11	1:A:621:LEU:HD13	1.81	0.62
2:B:16:ASN:HD22	2:B:18:ASP:CG	2.02	0.62
1:A:539:ASP:HA	1:A:542:GLU:OE1	1.99	0.62
2:B:180:GLN:HA	2:B:203:HIS:O	2.00	0.61
1:A:609:ILE:HD13	1:A:616:ARG:HH22	1.65	0.61
1:A:550:TRP:HE3	1:A:550:TRP:HA	1.65	0.61
1:A:527:GLU:HA	1:A:586:TYR:CD1	2.36	0.61
2:B:257:TYR:CE2	2:B:258:LYS:HG3	2.36	0.60
1:A:610:ASP:O	1:A:612:PRO:HD3	2.01	0.60
1:A:603:PHE:O	2:B:152:LYS:NZ	2.32	0.60
2:B:132:LYS:HG2	2:B:133:GLY:N	2.16	0.60
1:A:546:ILE:HG22	1:A:574:PRO:HG3	1.84	0.60
1:A:650:PRO:HD3	1:A:665:ILE:HG21	1.83	0.60
1:A:617:ILE:HD13	1:A:694:LEU:HD22	1.84	0.59
2:B:22:LEU:HD12	2:B:41:ASN:HD21	1.65	0.59
2:B:17:CYS:HB2	2:B:38:LEU:HD23	1.83	0.59
1:A:539:ASP:OD2	1:A:687:ARG:NH1	2.36	0.59
1:A:609:ILE:CD1	1:A:616:ARG:HH22	2.16	0.59
2:B:34:THR:O	2:B:57:LEU:HD12	2.02	0.58
2:B:234:VAL:O	2:B:234:VAL:HG12	2.02	0.58
2:B:43:LEU:C	2:B:45:THR:H	2.07	0.58
2:B:266:THR:O	2:B:267:LEU:HB2	2.02	0.58
1:A:691:VAL:HA	1:A:694:LEU:HD12	1.84	0.58
2:B:76:LEU:HD12	2:B:99:LEU:HD21	1.85	0.58
1:A:517:PHE:HE1	1:A:619:LEU:HD23	1.69	0.58
2:B:89:LEU:H	2:B:110:ASN:HD22	1.51	0.57
1:A:677:LEU:HD22	1:A:683:LEU:HD23	1.87	0.57
2:B:95:LEU:HD22	2:B:99:LEU:CD1	2.34	0.57
1:A:652:GLY:C	1:A:653:ILE:HG13	2.25	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:546:ILE:CG2	1:A:577:LEU:HD12	2.34	0.57
2:B:103:THR:HA	2:B:126:LEU:HA	1.86	0.57
1:A:545:ARG:HH12	1:A:703:PRO:HA	1.69	0.57
1:A:526:SER:OG	1:A:529:GLU:HG3	2.05	0.56
1:A:510:SER:OG	1:A:545:ARG:HG3	2.05	0.56
2:B:180:GLN:N	2:B:180:GLN:OE1	2.38	0.56
2:B:187:ILE:HG13	2:B:216:PHE:CG	2.41	0.56
2:B:187:ILE:HB	2:B:216:PHE:HB2	1.87	0.56
2:B:52:MET:N	2:B:53:PRO:CD	2.69	0.55
2:B:126:LEU:HD23	2:B:147:THR:HG21	1.88	0.55
2:B:233:GLY:O	2:B:235:ASP:N	2.40	0.55
2:B:107:VAL:HG12	2:B:107:VAL:O	2.06	0.55
1:A:519:LEU:CD1	1:A:519:LEU:N	2.68	0.55
2:B:17:CYS:HB3	2:B:22:LEU:HD11	1.88	0.55
1:A:559:HIS:HB2	1:A:591:VAL:O	2.07	0.54
1:A:676:VAL:O	1:A:677:LEU:HD12	2.07	0.54
1:A:579:ARG:O	1:A:582:SER:N	2.40	0.54
2:B:254:PHE:HE2	2:B:259:TYR:HA	1.73	0.54
1:A:547:SER:C	1:A:549:LYS:H	2.11	0.53
2:B:1:HIS:CE1	2:B:2:PRO:HG2	2.44	0.53
2:B:180:GLN:HB2	2:B:203:HIS:CD2	2.44	0.53
1:A:559:HIS:CG	1:A:560:ASP:H	2.25	0.53
2:B:1:HIS:CE1	2:B:3:ILE:H	2.27	0.53
1:A:543:GLN:NE2	1:A:691:VAL:HG11	2.24	0.52
2:B:1:HIS:CD2	2:B:29:LEU:HD23	2.44	0.52
2:B:186:THR:HA	2:B:213:ILE:CD1	2.38	0.52
2:B:96:GLY:HA3	2:B:121:ARG:O	2.09	0.52
2:B:116:PRO:O	2:B:119:ALA:HB2	2.08	0.52
1:A:693:TYR:CE2	1:A:697:LEU:HD11	2.44	0.52
2:B:128:GLU:HG2	2:B:152:LYS:HG3	1.91	0.52
1:A:511:ARG:HG2	1:A:700:GLU:HA	1.92	0.52
1:A:509:CYS:O	1:A:544:LEU:HD23	2.10	0.52
1:A:511:ARG:HH11	1:A:695:CYS:C	2.13	0.52
2:B:156:ALA:HB2	2:B:180:GLN:OE1	2.10	0.51
1:A:511:ARG:HA	1:A:701:ALA:CB	2.40	0.51
1:A:513:LEU:HD11	1:A:617:ILE:HD12	1.91	0.51
1:A:609:ILE:CG1	1:A:616:ARG:HH22	2.23	0.51
1:A:663:ARG:NH1	1:A:666:GLU:OE1	2.38	0.51
2:B:6:VAL:HG12	2:B:7:SER:N	2.24	0.51
2:B:250:ASN:O	2:B:252:ASP:N	2.44	0.51
2:B:17:CYS:HB3	2:B:22:LEU:CD1	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:593:SER:HB3	1:A:596:GLU:HB2	1.92	0.51
1:A:609:ILE:CG2	1:A:616:ARG:NH2	2.73	0.51
2:B:89:LEU:HB2	2:B:110:ASN:HD22	1.76	0.51
1:A:684:GLU:O	1:A:684:GLU:HG3	2.11	0.50
2:B:128:GLU:HB3	2:B:130:TYR:HE1	1.77	0.50
2:B:89:LEU:HB2	2:B:110:ASN:ND2	2.26	0.50
2:B:145:THR:HB	2:B:146:PRO:HD3	1.94	0.50
1:A:559:HIS:CD2	1:A:560:ASP:N	2.70	0.50
1:A:650:PRO:CD	1:A:665:ILE:HG21	2.41	0.50
2:B:242:ASN:HD21	2:B:244:ALA:HB3	1.77	0.50
2:B:68:THR:O	2:B:69:LYS:HG3	2.11	0.50
1:A:543:GLN:HE22	1:A:691:VAL:HG11	1.76	0.50
1:A:506:ASP:OD2	1:A:688:ASP:OD2	2.28	0.50
2:B:151:GLU:O	2:B:175:ASP:OD1	2.30	0.50
1:A:548:GLN:HG2	1:A:552:ARG:HD3	1.94	0.49
1:A:586:TYR:CE2	1:A:588:GLY:HA3	2.47	0.49
1:A:524:ARG:HD2	1:A:656:HIS:CD2	2.47	0.49
2:B:155:LEU:HB2	2:B:179:LEU:HD23	1.95	0.49
2:B:16:ASN:HB3	2:B:18:ASP:OD2	2.13	0.49
2:B:251:SER:O	2:B:253:LYS:N	2.46	0.49
2:B:180:GLN:HG2	2:B:181:GLU:HG2	1.95	0.48
2:B:224:ALA:HA	2:B:243:VAL:HB	1.95	0.48
2:B:206:PRO:O	2:B:249:ASP:HB2	2.12	0.48
2:B:52:MET:CE	2:B:74:GLY:HA3	2.43	0.48
1:A:511:ARG:HG2	1:A:699:PRO:O	2.13	0.48
1:A:690:ILE:HG22	1:A:691:VAL:N	2.29	0.48
1:A:539:ASP:HB3	1:A:687:ARG:HD2	1.95	0.48
2:B:76:LEU:HB2	2:B:99:LEU:HD23	1.95	0.48
1:A:550:TRP:CH2	1:A:702:PRO:HD2	2.49	0.48
1:A:670:PRO:HG2	1:A:671:GLU:OE1	2.14	0.47
1:A:686:GLN:O	1:A:687:ARG:C	2.52	0.47
1:A:636:ARG:CA	1:A:636:ARG:NH1	2.71	0.47
1:A:523:SER:C	1:A:525:LEU:N	2.64	0.47
2:B:4:CYS:HB3	2:B:16:ASN:O	2.14	0.47
1:A:660:LYS:NZ	1:A:664:LEU:HG	2.29	0.47
2:B:19:LYS:HG3	2:B:19:LYS:O	2.13	0.47
1:A:533:LEU:O	1:A:536:PHE:HB3	2.15	0.47
1:A:683:LEU:C	1:A:685:GLN:H	2.18	0.47
1:A:599:LYS:O	1:A:603:PHE:CD2	2.61	0.47
2:B:64:ARG:NH2	2:B:86:HIS:CD2	2.83	0.47
2:B:5:GLU:HG2	2:B:6:VAL:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:89:LEU:H	2:B:110:ASN:ND2	2.13	0.47
1:A:545:ARG:HH22	1:A:703:PRO:CB	2.28	0.46
1:A:523:SER:HB3	1:A:588:GLY:O	2.14	0.46
2:B:75:THR:O	2:B:77:PRO:HD3	2.15	0.46
2:B:249:ASP:C	2:B:249:ASP:OD1	2.53	0.46
1:A:542:GLU:C	1:A:544:LEU:H	2.17	0.46
1:A:674:ALA:C	1:A:675:PHE:CD1	2.89	0.46
2:B:109:PHE:HA	2:B:133:GLY:O	2.16	0.46
1:A:627:PRO:O	1:A:630:MET:HG3	2.16	0.46
1:A:524:ARG:HG3	1:A:624:SER:HB3	1.98	0.46
1:A:665:ILE:HG22	1:A:666:GLU:N	2.31	0.46
1:A:683:LEU:C	1:A:685:GLN:N	2.68	0.46
1:A:523:SER:O	1:A:525:LEU:N	2.48	0.46
2:B:172:GLU:OE2	2:B:173:ASN:ND2	2.49	0.46
2:B:6:VAL:O	2:B:7:SER:HB3	2.15	0.46
1:A:527:GLU:HA	1:A:586:TYR:CE1	2.50	0.46
1:A:625:GLN:HG3	1:A:656:HIS:O	2.15	0.46
1:A:601:THR:HG23	1:A:605:ILE:HD12	1.97	0.45
2:B:248:CYS:O	2:B:249:ASP:C	2.52	0.45
2:B:127:GLN:C	2:B:150:LEU:HD12	2.36	0.45
2:B:232:GLN:HE21	2:B:237:LYS:CE	2.29	0.45
2:B:233:GLY:C	2:B:235:ASP:H	2.19	0.45
2:B:228:TYR:CE2	2:B:241:SER:HB2	2.51	0.45
2:B:218:ARG:HA	2:B:221:GLN:OE1	2.17	0.45
1:A:577:LEU:O	1:A:578:ARG:C	2.55	0.45
2:B:89:LEU:N	2:B:110:ASN:HD22	2.15	0.45
1:A:549:LYS:O	1:A:550:TRP:CE3	2.70	0.45
1:A:610:ASP:O	1:A:612:PRO:CD	2.65	0.45
2:B:174:LEU:HD23	2:B:197:LEU:HD13	1.98	0.45
1:A:613:GLU:OE1	1:A:613:GLU:N	2.29	0.45
1:A:690:ILE:C	1:A:692:SER:N	2.71	0.44
1:A:577:LEU:O	1:A:579:ARG:N	2.51	0.44
1:A:663:ARG:O	1:A:667:LYS:HG2	2.18	0.44
1:A:632:ARG:NH2	2:B:225:GLU:CD	2.71	0.44
1:A:616:ARG:HD2	1:A:646:VAL:HG22	1.99	0.44
1:A:692:SER:O	1:A:695:CYS:N	2.51	0.44
2:B:143:LEU:HG	2:B:144:LEU:HD23	2.00	0.44
1:A:556:VAL:HG11	1:A:562:SER:HB3	1.99	0.44
2:B:34:THR:HB	2:B:58:THR:CG2	2.45	0.44
2:B:43:LEU:O	2:B:45:THR:N	2.49	0.44
2:B:232:GLN:HE21	2:B:237:LYS:HE3	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:674:ALA:O	1:A:675:PHE:CD1	2.71	0.44
2:B:218:ARG:O	2:B:221:GLN:HB2	2.18	0.43
2:B:32:ASP:O	2:B:33:THR:C	2.56	0.43
1:A:556:VAL:CG1	1:A:562:SER:HB3	2.48	0.43
2:B:85:SER:HA	2:B:108:SER:O	2.17	0.43
1:A:508:TYR:C	1:A:508:TYR:HD1	2.21	0.43
2:B:13:LEU:CD2	2:B:32:ASP:HB2	2.48	0.43
1:A:609:ILE:HG12	1:A:616:ARG:HH22	1.82	0.43
2:B:43:LEU:O	2:B:44:TYR:HB2	2.19	0.43
2:B:127:GLN:NE2	2:B:151:GLU:OE2	2.51	0.43
1:A:622:MET:SD	1:A:662:ILE:HG12	2.59	0.43
1:A:649:ILE:O	1:A:649:ILE:HG22	2.19	0.43
1:A:682:GLU:O	1:A:682:GLU:HG2	2.19	0.42
1:A:547:SER:C	1:A:549:LYS:N	2.72	0.42
1:A:617:ILE:HA	1:A:647:ILE:O	2.18	0.42
1:A:633:ASN:O	1:A:634:PHE:C	2.57	0.42
1:A:543:GLN:O	1:A:544:LEU:CG	2.66	0.42
1:A:659:LEU:O	1:A:663:ARG:HG2	2.18	0.42
2:B:43:LEU:C	2:B:45:THR:N	2.73	0.42
2:B:51:LEU:C	2:B:53:PRO:HD2	2.39	0.42
2:B:6:VAL:O	2:B:7:SER:CB	2.68	0.42
2:B:66:GLU:O	2:B:67:LEU:C	2.58	0.42
1:A:505:HIS:O	1:A:506:ASP:O	2.36	0.42
2:B:214:LEU:HD23	2:B:214:LEU:HA	1.87	0.42
2:B:259:TYR:CZ	2:B:261:GLY:HA2	2.54	0.42
2:B:13:LEU:HD22	2:B:32:ASP:HB2	2.02	0.42
2:B:144:LEU:HD11	2:B:153:LEU:CD2	2.46	0.42
1:A:533:LEU:O	1:A:537:VAL:HG23	2.20	0.42
1:A:671:GLU:HA	1:A:693:TYR:OH	2.19	0.41
1:A:546:ILE:O	1:A:574:PRO:HG3	2.19	0.41
2:B:29:LEU:HA	2:B:29:LEU:HD23	1.88	0.41
1:A:651:VAL:HG12	1:A:651:VAL:O	2.18	0.41
1:A:677:LEU:CD2	1:A:683:LEU:HD23	2.50	0.41
1:A:579:ARG:O	1:A:580:ILE:C	2.58	0.41
2:B:9:VAL:HG21	2:B:14:GLU:OE2	2.20	0.41
1:A:561:GLY:HA3	2:B:240:THR:HA	2.01	0.41
1:A:690:ILE:O	1:A:692:SER:N	2.53	0.41
2:B:242:ASN:ND2	2:B:244:ALA:HB3	2.35	0.41
2:B:66:GLU:O	2:B:68:THR:HG23	2.20	0.41
1:A:647:ILE:HG21	1:A:671:GLU:HB3	2.02	0.41
2:B:191:PHE:H	2:B:215:TYR:HE2	1.66	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:160:LEU:H	2:B:182:ASN:HD22	1.69	0.41
2:B:39:SER:OG	2:B:63:ASP:OD1	2.35	0.41
1:A:560:ASP:O	2:B:240:THR:HG23	2.21	0.41
1:A:507:PHE:HZ	1:A:692:SER:HG	1.67	0.41
1:A:545:ARG:CZ	1:A:703:PRO:HA	2.50	0.41
2:B:228:TYR:CD2	2:B:241:SER:HA	2.56	0.41
1:A:644:LYS:O	1:A:645:LYS:HB2	2.20	0.41
2:B:127:GLN:HA	2:B:150:LEU:HA	2.02	0.41
2:B:49:ALA:HB1	2:B:52:MET:CE	2.51	0.41
2:B:19:LYS:CG	2:B:19:LYS:O	2.69	0.41
1:A:595:SER:OG	1:A:626:GLU:OE2	2.39	0.41
1:A:625:GLN:HB3	1:A:658:ASN:HD22	1.86	0.41
2:B:110:ASN:HB2	2:B:134:ASN:HD21	1.86	0.41
2:B:110:ASN:HB3	2:B:111:ARG:H	1.73	0.41
2:B:267:LEU:HD12	2:B:267:LEU:HA	1.79	0.41
1:A:515:LEU:O	1:A:553:VAL:HA	2.21	0.40
1:A:598:LEU:HD23	1:A:598:LEU:HA	1.91	0.40
2:B:167:LEU:HB3	2:B:191:PHE:HE1	1.87	0.40
1:A:670:PRO:HG2	1:A:671:GLU:CD	2.41	0.40
1:A:632:ARG:NH2	2:B:225:GLU:OE1	2.55	0.40
2:B:185:TYR:HB2	2:B:208:LEU:HB2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	197/208 (95%)	159 (81%)	28 (14%)	10 (5%)	2	15
2	B	265/290 (91%)	221 (83%)	31 (12%)	13 (5%)	3	16
All	All	462/498 (93%)	380 (82%)	59 (13%)	23 (5%)	3	16

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	506	ASP
1	A	559	HIS
2	B	28	ASP
2	B	234	VAL
2	B	249	ASP
2	B	252	ASP
1	A	628	GLN
2	B	7	SER
2	B	96	GLY
2	B	251	SER
1	A	524	ARG
1	A	543	GLN
1	A	545	ARG
1	A	577	LEU
1	A	578	ARG
1	A	687	ARG
2	B	33	THR
2	B	141	PRO
2	B	57	LEU
1	A	693	TYR
2	B	27	PRO
2	B	206	PRO
2	B	265	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	177/186 (95%)	165 (93%)	12 (7%)	20	55
2	B	240/260 (92%)	228 (95%)	12 (5%)	30	67
All	All	417/446 (94%)	393 (94%)	24 (6%)	25	61

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	508	TYR
1	A	512	LEU
1	A	519	LEU
1	A	550	TRP
1	A	560	ASP
1	A	573	ARG
1	A	576	GLU
1	A	589	SER
1	A	597	VAL
1	A	613	GLU
1	A	629	ARG
1	A	668	GLN
2	B	23	THR
2	B	71	GLN
2	B	91	SER
2	B	108	SER
2	B	113	THR
2	B	162	GLU
2	B	172	GLU
2	B	235	ASP
2	B	251	SER
2	B	252	ASP
2	B	257	TYR
2	B	266	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	543	GLN
1	A	559	HIS
1	A	563	HIS
1	A	625	GLN
1	A	639	GLN
1	A	656	HIS
2	B	16	ASN
2	B	59	GLN
2	B	86	HIS
2	B	110	ASN
2	B	134	ASN
2	B	159	GLN
2	B	182	ASN
2	B	203	HIS
2	B	232	GLN

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Mol	Chain	Res	Type
2	B	242	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	199/208 (95%)	-0.24	0	100 100	25, 60, 86, 92	0
2	B	267/290 (92%)	-0.40	2 (0%)	89 78	24, 46, 61, 72	0
All	All	466/498 (93%)	-0.33	2 (0%)	93 85	24, 50, 80, 92	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	234	VAL	2.6
2	B	267	LEU	2.3

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.