



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:05 AM UTC

PDB ID : 1QGK / pdb_00001qgk
Title : STRUCTURE OF IMPORTIN BETA BOUND TO THE IBB DOMAIN OF IMPORTIN ALPHA
Authors : Cingolani, G.; Petosa, C.; Weis, K.; Muller, C.W.
Deposited on : 1999-04-29
Resolution : 2.50 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

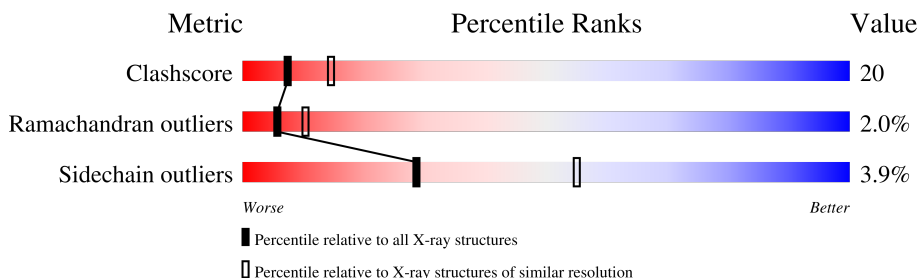
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	876	
2	B	44	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7229 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 PROTEIN (IMPORTIN BETA SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	876	Total	C	N	O	S	1	0	0
			6810	4288	1145	1330	47			

- Molecule 2 is a protein called PROTEIN (IMPORTIN ALPHA-2 SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	44	Total	C	N	O	S	0	0	0
			375	225	84	64	2			

- Molecule 3 is water.

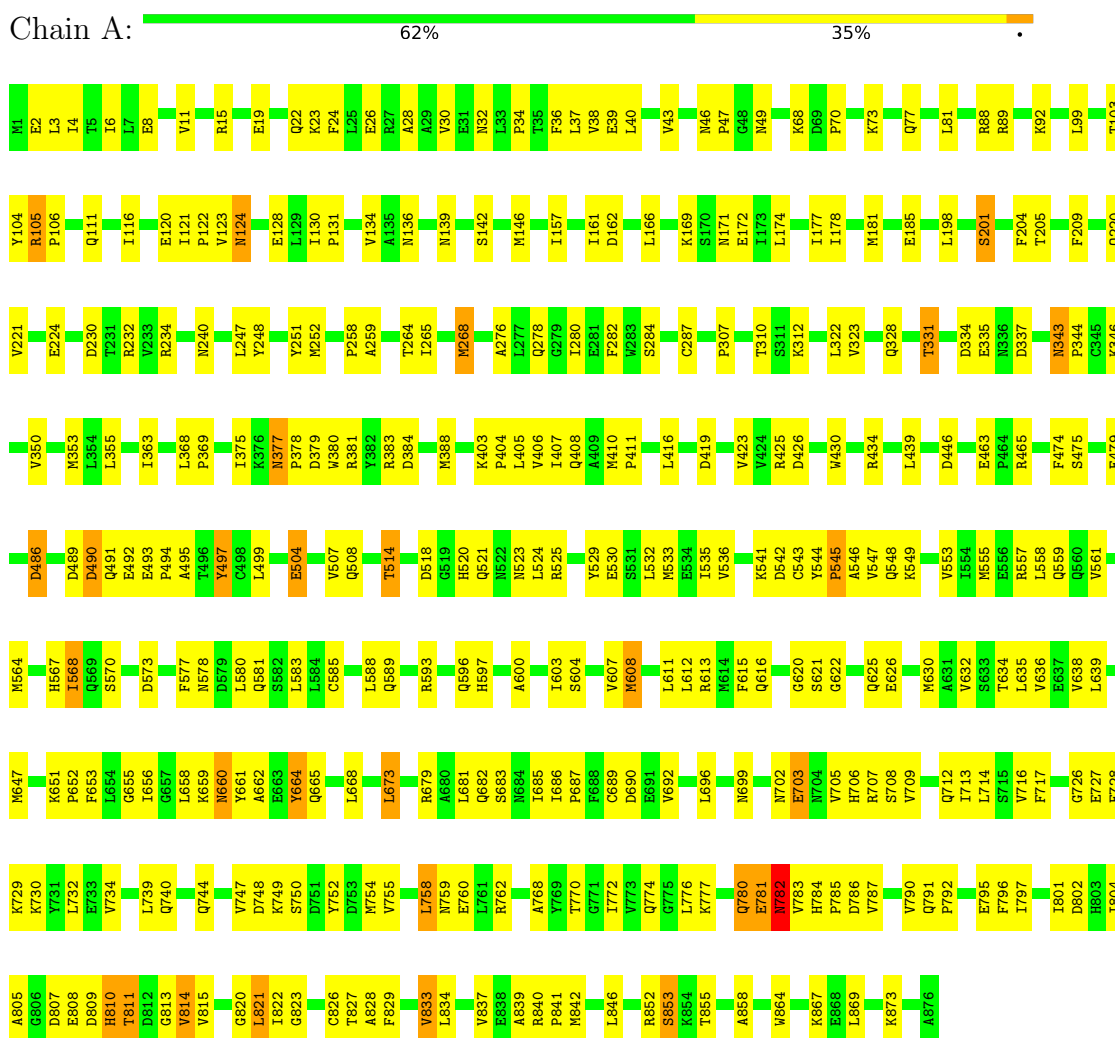
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	42	Total	O	0	0
			42	42		
3	B	2	Total	O	0	0
			2	2		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: PROTEIN (IMPORTIN BETA SUBUNIT)



- Molecule 2: PROTEIN (IMPORTIN ALPHA-2 SUBUNIT)





4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.58Å 97.61Å 83.19Å 90.00° 91.13° 90.00°	Depositor
Resolution (Å)	30.00 – 2.50	Depositor
% Data completeness (in resolution range)	97.5 (30.00-2.50)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.221 , 0.273	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7229	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

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	0.48	1/6922 (0.0%)	0.92	16/9397 (0.2%)
2	B	0.49	0/376	0.79	0/491
All	All	0.48	1/7298 (0.0%)	0.91	16/9888 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	268	MET	SD-CE	-5.29	1.66	1.79

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	810	HIS	CB-CA-C	-7.04	107.42	115.79
1	A	247	LEU	N-CA-C	6.73	118.61	111.28
1	A	252	MET	N-CA-C	6.36	120.64	113.01
1	A	363	ILE	N-CA-C	6.30	117.08	110.72
1	A	185	GLU	CA-C-N	6.28	125.50	118.97
1	A	185	GLU	C-N-CA	6.28	125.50	118.97
1	A	103	THR	N-CA-C	-5.98	106.15	113.50
1	A	258	PRO	N-CA-C	5.79	121.15	114.03
1	A	28	ALA	N-CA-C	-5.74	105.10	111.36
1	A	689	CYS	N-CA-C	5.66	117.53	111.36
1	A	337	ASP	N-CA-C	5.59	118.34	109.96
1	A	284	SER	N-CA-C	-5.42	105.37	111.28
1	A	760	GLU	N-CA-C	-5.37	107.28	113.88
1	A	323	VAL	CA-C-N	-5.22	113.25	119.05
1	A	323	VAL	C-N-CA	-5.22	113.25	119.05
1	A	201	SER	N-CA-C	5.17	119.58	113.28

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	6810	0	6805	285	0
2	B	375	0	403	21	0
3	A	42	0	0	6	0
3	B	2	0	0	0	0
All	All	7229	0	7208	294	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 20.

All (294) 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:181:MET:HE1	1:A:198:LEU:HD22	1.35	1.07
2:B:53:VAL:HG12	2:B:54:SER:H	1.34	0.91
1:A:377:ASN:ND2	1:A:379:ASP:H	1.71	0.88
1:A:608:MET:HE1	1:A:632:VAL:HG22	1.61	0.82
1:A:774:GLN:HG3	2:B:47:MET:HE1	1.62	0.80
1:A:328:GLN:O	1:A:331:THR:HB	1.81	0.80
1:A:375:ILE:HD11	1:A:416:LEU:HD11	1.63	0.80
1:A:543:CYS:C	1:A:545:PRO:HD2	2.06	0.80
1:A:377:ASN:HD22	1:A:378:PRO:N	1.81	0.79
1:A:668:LEU:HD22	1:A:712:GLN:HG3	1.65	0.78
1:A:613:ARG:HA	1:A:616:GLN:HE21	1.50	0.77
1:A:781:GLU:HG3	1:A:782:ASN:N	2.00	0.77
1:A:784:HIS:HB3	1:A:785:PRO:HD2	1.66	0.77
1:A:802:ASP:OD1	1:A:842:MET:HB2	1.85	0.77
1:A:377:ASN:HD22	1:A:379:ASP:H	1.33	0.76
1:A:369:PRO:HG2	3:A:917:HOH:O	1.87	0.74
1:A:532:LEU:O	1:A:535:ILE:HG22	1.85	0.74
1:A:811:THR:HG23	1:A:814:VAL:H	1.52	0.73
1:A:312:LYS:HE2	1:A:312:LYS:HA	1.71	0.73
1:A:544:TYR:N	1:A:545:PRO:HD2	2.04	0.72
1:A:172:GLU:H	1:A:172:GLU:CD	1.99	0.70
1:A:486:ASP:HB3	1:A:494:PRO:HD3	1.73	0.70
1:A:781:GLU:CG	1:A:782:ASN:H	2.03	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:VAL:HG12	1:A:11:VAL:O	1.91	0.70
1:A:608:MET:HE2	1:A:608:MET:HA	1.72	0.70
1:A:130:ILE:HG13	1:A:169:LYS:HD3	1.74	0.69
1:A:747:VAL:HG21	1:A:754:MET:HB3	1.75	0.69
1:A:781:GLU:CG	1:A:782:ASN:N	2.56	0.69
1:A:781:GLU:CD	1:A:782:ASN:H	2.01	0.69
1:A:826:CYS:SG	1:A:869:LEU:HD23	2.33	0.69
1:A:406:VAL:HG21	1:A:439:LEU:CD1	2.22	0.68
1:A:781:GLU:HG3	1:A:782:ASN:ND2	2.08	0.68
1:A:123:VAL:HG12	1:A:123:VAL:O	1.95	0.67
1:A:782:ASN:C	1:A:782:ASN:HD22	2.01	0.66
1:A:4:ILE:O	1:A:8:GLU:HG3	1.96	0.66
1:A:555:MET:SD	1:A:603:ILE:HG23	2.36	0.65
1:A:36:PHE:CZ	1:A:40:LEU:HD11	2.32	0.65
1:A:388:MET:HE1	2:B:18:LYS:CE	2.26	0.65
1:A:714:LEU:HD22	1:A:739:LEU:HD23	1.79	0.65
1:A:747:VAL:HG11	1:A:754:MET:HE3	1.79	0.65
1:A:105:ARG:HB3	1:A:106:PRO:CD	2.28	0.64
1:A:377:ASN:HD22	1:A:377:ASN:C	2.04	0.64
1:A:89:ARG:HG3	3:A:913:HOH:O	1.98	0.64
1:A:664:TYR:CE1	1:A:665:GLN:HG3	2.32	0.64
2:B:53:VAL:HG12	2:B:54:SER:N	2.10	0.64
1:A:388:MET:HE1	2:B:18:LYS:HE3	1.78	0.64
1:A:683:SER:O	1:A:686:ILE:HG13	1.98	0.63
1:A:81:LEU:HA	1:A:88:ARG:HH11	1.63	0.63
1:A:181:MET:CE	1:A:198:LEU:HD22	2.22	0.63
1:A:545:PRO:HG2	1:A:546:ALA:H	1.63	0.63
1:A:26:GLU:O	1:A:30:VAL:HG23	1.99	0.61
1:A:781:GLU:HG3	1:A:782:ASN:H	1.63	0.61
1:A:486:ASP:HA	1:A:489:ASP:OD1	2.00	0.61
1:A:81:LEU:HA	1:A:88:ARG:NH1	2.16	0.61
1:A:32:ASN:HD22	1:A:34:PRO:HD2	1.65	0.61
1:A:652:PRO:O	1:A:656:ILE:HG13	2.01	0.61
1:A:706:HIS:CD2	1:A:708:SER:HB2	2.35	0.61
1:A:514:THR:HG21	1:A:557:ARG:NH1	2.15	0.61
1:A:564:MET:HE3	1:A:564:MET:HA	1.84	0.60
1:A:406:VAL:HG21	1:A:439:LEU:HD12	1.82	0.60
1:A:426:ASP:OD2	1:A:465:ARG:HG2	2.01	0.60
1:A:157:ILE:HG22	1:A:166:LEU:HD11	1.82	0.60
1:A:823:GLY:HA3	1:A:864:TRP:CZ3	2.37	0.59
1:A:626:GLU:OE1	1:A:665:GLN:HB3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:651:LYS:N	1:A:652:PRO:HD2	2.18	0.59
1:A:616:GLN:HG2	1:A:653:PHE:CE1	2.38	0.59
1:A:647:MET:O	1:A:651:LYS:HG3	2.02	0.58
1:A:265:ILE:HA	1:A:268:MET:HE3	1.84	0.58
1:A:632:VAL:O	1:A:636:VAL:HG23	2.03	0.58
1:A:783:VAL:HG11	1:A:828:ALA:O	2.04	0.58
1:A:368:LEU:HD11	1:A:405:LEU:HD13	1.84	0.58
1:A:34:PRO:O	1:A:38:VAL:HG23	2.04	0.57
1:A:343:ASN:C	1:A:343:ASN:HD22	2.11	0.57
1:A:504:GLU:OE1	1:A:546:ALA:HB2	2.05	0.57
1:A:707:ARG:HG3	1:A:707:ARG:HH11	1.70	0.57
1:A:32:ASN:ND2	1:A:34:PRO:HD2	2.20	0.57
1:A:523:ASN:ND2	3:A:893:HOH:O	2.36	0.57
1:A:139:ASN:HB3	1:A:142:SER:OG	2.04	0.57
1:A:403:LYS:O	1:A:407:ILE:HG12	2.05	0.57
1:A:495:ALA:O	1:A:541:LYS:HG2	2.04	0.56
1:A:73:LYS:O	1:A:77:GLN:HG3	2.05	0.56
2:B:51:ARG:O	2:B:52:ASN:HB2	2.06	0.56
1:A:497:TYR:CE1	1:A:499:LEU:HB2	2.41	0.56
1:A:679:ARG:HH21	2:B:43:LYS:HE2	1.71	0.56
1:A:668:LEU:HD21	1:A:709:VAL:HA	1.88	0.55
1:A:807:ASP:HB3	1:A:809:ASP:OD2	2.06	0.55
1:A:544:TYR:N	1:A:545:PRO:CD	2.69	0.55
1:A:782:ASN:ND2	1:A:782:ASN:C	2.65	0.55
1:A:525:ARG:NH1	1:A:583:LEU:HD12	2.22	0.55
1:A:11:VAL:O	1:A:11:VAL:CG1	2.54	0.55
1:A:781:GLU:O	1:A:782:ASN:O	2.25	0.55
1:A:520:HIS:HB3	3:A:886:HOH:O	2.07	0.55
1:A:713:ILE:O	1:A:716:VAL:HG12	2.06	0.54
1:A:729:LYS:HD3	1:A:786:ASP:OD1	2.07	0.54
1:A:146:MET:HA	1:A:146:MET:HE3	1.90	0.54
1:A:541:LYS:HG3	1:A:542:ASP:N	2.22	0.54
1:A:178:ILE:CG2	1:A:221:VAL:HG21	2.37	0.54
1:A:833:VAL:O	1:A:837:VAL:HG23	2.07	0.54
1:A:604:SER:HB3	1:A:639:LEU:HD21	1.89	0.54
1:A:774:GLN:CG	2:B:47:MET:HE1	2.36	0.53
1:A:805:ALA:HB3	1:A:842:MET:HB3	1.90	0.53
1:A:530:GLU:HA	1:A:533:MET:HE3	1.90	0.53
1:A:780:GLN:O	1:A:781:GLU:O	2.26	0.53
1:A:525:ARG:HD3	3:A:896:HOH:O	2.08	0.53
1:A:659:LYS:O	1:A:660:ASN:C	2.52	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:660:ASN:C	1:A:662:ALA:H	2.16	0.53
1:A:593:ARG:NH2	2:B:35:ASN:OD1	2.42	0.53
1:A:171:ASN:HB3	1:A:172:GLU:OE2	2.09	0.53
1:A:726:GLY:C	1:A:728:PHE:H	2.16	0.53
1:A:105:ARG:HB3	1:A:106:PRO:HD3	1.90	0.53
1:A:752:TYR:HA	1:A:755:VAL:HB	1.91	0.53
1:A:613:ARG:HA	1:A:616:GLN:NE2	2.22	0.52
1:A:178:ILE:HG23	1:A:221:VAL:HG21	1.92	0.52
1:A:334:ASP:O	1:A:335:GLU:C	2.52	0.52
1:A:747:VAL:HG21	1:A:754:MET:CB	2.40	0.52
2:B:53:VAL:CG1	2:B:54:SER:H	2.16	0.52
1:A:834:LEU:HD21	1:A:873:LYS:HG2	1.92	0.52
1:A:15:ARG:O	1:A:19:GLU:HG2	2.10	0.52
1:A:368:LEU:N	1:A:369:PRO:HD2	2.25	0.52
1:A:201:SER:HB2	1:A:204:PHE:CZ	2.46	0.51
1:A:384:ASP:OD1	1:A:423:VAL:HG12	2.10	0.51
1:A:128:GLU:N	1:A:128:GLU:OE1	2.39	0.51
1:A:230:ASP:OD1	1:A:232:ARG:HB2	2.11	0.51
1:A:493:GLU:HG3	1:A:494:PRO:HD2	1.92	0.51
1:A:177:ILE:HG22	1:A:181:MET:HE2	1.91	0.51
1:A:690:ASP:CG	1:A:730:LYS:HZ3	2.19	0.51
1:A:791:GLN:N	1:A:792:PRO:HD2	2.25	0.51
1:A:852:ARG:O	1:A:853:SER:O	2.27	0.51
1:A:702:ASN:O	1:A:705:VAL:HG23	2.11	0.51
1:A:497:TYR:CZ	1:A:499:LEU:HB2	2.46	0.51
1:A:781:GLU:OE1	1:A:782:ASN:N	2.44	0.51
1:A:343:ASN:ND2	1:A:346:LYS:H	2.09	0.50
1:A:776:LEU:HD22	1:A:786:ASP:HB3	1.93	0.50
1:A:681:LEU:O	1:A:682:GLN:HB2	2.11	0.50
1:A:811:THR:HG23	1:A:814:VAL:HB	1.93	0.50
1:A:520:HIS:CB	3:A:886:HOH:O	2.60	0.50
1:A:613:ARG:HG3	1:A:613:ARG:HH11	1.76	0.50
1:A:514:THR:HG21	1:A:557:ARG:HH12	1.77	0.49
1:A:287:CYS:SG	1:A:355:LEU:HD23	2.52	0.49
1:A:6:ILE:HD13	1:A:24:PHE:CD1	2.46	0.49
1:A:626:GLU:O	1:A:630:MET:HG3	2.13	0.49
1:A:807:ASP:C	1:A:809:ASP:H	2.21	0.49
1:A:116:ILE:HG22	1:A:120:GLU:HG3	1.94	0.49
1:A:486:ASP:CB	1:A:494:PRO:HD3	2.42	0.49
1:A:514:THR:HG23	1:A:529:TYR:CE2	2.47	0.49
1:A:699:ASN:HD21	1:A:705:VAL:HG11	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:GLU:H	1:A:2:GLU:CD	2.21	0.48
1:A:748:ASP:C	1:A:750:SER:H	2.20	0.48
1:A:864:TRP:CD1	2:B:53:VAL:HG22	2.49	0.48
1:A:727:GLU:HG3	1:A:727:GLU:O	2.13	0.48
1:A:544:TYR:O	1:A:548:GLN:HG3	2.13	0.48
1:A:377:ASN:HD22	1:A:379:ASP:N	2.08	0.48
1:A:490:ASP:O	1:A:491:GLN:HG3	2.14	0.48
1:A:514:THR:HG21	1:A:529:TYR:CZ	2.49	0.48
1:A:810:HIS:HB3	1:A:811:THR:H	1.46	0.48
1:A:846:LEU:HD12	1:A:846:LEU:O	2.13	0.47
1:A:790:VAL:HG12	1:A:790:VAL:O	2.14	0.47
1:A:608:MET:O	1:A:612:LEU:HB2	2.15	0.47
1:A:810:HIS:C	1:A:811:THR:HG22	2.39	0.47
1:A:174:LEU:HD21	1:A:205:THR:HG21	1.96	0.47
1:A:425:ARG:NH2	1:A:463:GLU:OE1	2.43	0.47
1:A:855:THR:HB	1:A:858:ALA:HB3	1.96	0.47
1:A:864:TRP:CH2	2:B:51:ARG:HD3	2.49	0.47
1:A:608:MET:HG2	1:A:635:LEU:HD23	1.97	0.47
1:A:535:ILE:HG23	1:A:536:VAL:N	2.30	0.46
1:A:276:ALA:O	1:A:280:ILE:HG13	2.16	0.46
1:A:600:ALA:CB	1:A:638:VAL:HG11	2.46	0.46
1:A:632:VAL:CG1	1:A:673:LEU:HD11	2.46	0.46
1:A:248:TYR:HB3	1:A:251:TYR:HD2	1.80	0.46
1:A:796:PHE:O	1:A:797:ILE:C	2.58	0.46
1:A:377:ASN:ND2	1:A:378:PRO:N	2.60	0.46
1:A:570:SER:O	1:A:573:ASP:HB2	2.16	0.46
1:A:726:GLY:C	1:A:728:PHE:N	2.73	0.46
1:A:840:ARG:HH11	1:A:840:ARG:HG2	1.80	0.46
1:A:380:TRP:CZ2	1:A:381:ARG:NH1	2.84	0.46
1:A:804:ILE:O	1:A:810:HIS:NE2	2.47	0.46
1:A:815:VAL:HG13	1:A:846:LEU:HD11	1.96	0.46
2:B:48:LEU:HD23	2:B:48:LEU:O	2.16	0.46
1:A:603:ILE:HG22	1:A:607:VAL:CG2	2.46	0.46
1:A:770:THR:CG2	2:B:47:MET:HE2	2.45	0.46
1:A:783:VAL:HG23	1:A:783:VAL:O	2.15	0.46
1:A:220:GLN:O	1:A:224:GLU:HB2	2.15	0.45
1:A:783:VAL:HG12	1:A:828:ALA:HA	1.98	0.45
1:A:827:THR:O	1:A:827:THR:HG22	2.14	0.45
2:B:34:VAL:O	2:B:38:LEU:HB2	2.15	0.45
1:A:136:ASN:ND2	1:A:146:MET:CE	2.79	0.45
1:A:811:THR:CG2	1:A:814:VAL:HB	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:GLU:O	1:A:23:LYS:HG2	2.16	0.45
1:A:161:ILE:CG2	1:A:162:ASP:N	2.78	0.45
2:B:47:MET:SD	2:B:50:ARG:NH1	2.85	0.45
1:A:161:ILE:HG22	1:A:162:ASP:N	2.31	0.45
1:A:105:ARG:HD2	1:A:105:ARG:HA	1.52	0.45
1:A:493:GLU:CG	1:A:494:PRO:HD2	2.46	0.45
1:A:585:CYS:O	1:A:588:LEU:HB2	2.17	0.45
1:A:781:GLU:C	1:A:782:ASN:O	2.60	0.45
1:A:555:MET:O	1:A:558:LEU:HB3	2.17	0.45
1:A:616:GLN:HG2	1:A:653:PHE:CZ	2.52	0.45
1:A:758:LEU:HD22	1:A:762:ARG:HG3	1.98	0.45
1:A:410:MET:HB2	1:A:411:PRO:HD3	1.99	0.45
1:A:544:TYR:OH	1:A:596:GLN:HG3	2.17	0.45
1:A:728:PHE:C	1:A:730:LYS:N	2.72	0.45
1:A:784:HIS:O	1:A:787:VAL:HG22	2.17	0.45
1:A:864:TRP:CZ2	2:B:51:ARG:HB3	2.51	0.45
2:B:48:LEU:HD21	2:B:54:SER:OXT	2.17	0.45
1:A:647:MET:HE3	1:A:647:MET:HA	1.98	0.44
1:A:682:GLN:O	1:A:685:ILE:HG22	2.17	0.44
1:A:268:MET:HE1	1:A:322:LEU:CD2	2.47	0.44
1:A:759:ASN:HB3	1:A:811:THR:HG21	1.99	0.44
1:A:123:VAL:O	1:A:124:ASN:C	2.60	0.44
1:A:307:PRO:HG2	1:A:310:THR:HG22	1.99	0.44
1:A:377:ASN:ND2	1:A:378:PRO:HD2	2.32	0.44
1:A:839:ALA:O	1:A:841:PRO:HD3	2.17	0.44
1:A:2:GLU:CD	1:A:2:GLU:N	2.76	0.44
1:A:518:ASP:HB2	1:A:521:GLN:OE1	2.16	0.44
1:A:811:THR:C	1:A:813:GLY:N	2.74	0.44
1:A:39:GLU:O	1:A:43:VAL:HG23	2.17	0.44
1:A:626:GLU:CD	1:A:665:GLN:HB3	2.43	0.44
1:A:549:LYS:O	1:A:553:VAL:HG23	2.18	0.44
1:A:377:ASN:ND2	1:A:377:ASN:C	2.75	0.43
1:A:535:ILE:CG2	1:A:536:VAL:N	2.81	0.43
1:A:795:GLU:OE2	1:A:840:ARG:NH2	2.46	0.43
1:A:855:THR:O	1:A:858:ALA:N	2.51	0.43
1:A:198:LEU:HG	1:A:240:ASN:ND2	2.33	0.43
1:A:383:ARG:HH11	1:A:419:ASP:CG	2.26	0.43
1:A:426:ASP:OD2	1:A:465:ARG:CG	2.66	0.43
1:A:543:CYS:C	1:A:545:PRO:CD	2.86	0.43
1:A:92:LYS:NZ	1:A:120:GLU:OE1	2.41	0.43
1:A:130:ILE:N	1:A:131:PRO:HD2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:GLN:HA	1:A:634:THR:HG21	2.00	0.43
1:A:608:MET:HE3	1:A:635:LEU:HD23	2.01	0.43
1:A:104:TYR:CE2	1:A:106:PRO:O	2.71	0.43
1:A:46:ASN:HA	1:A:47:PRO:HD3	1.89	0.43
1:A:564:MET:HB3	1:A:577:PHE:CE1	2.53	0.43
1:A:777:LYS:HG2	1:A:777:LYS:O	2.18	0.43
1:A:350:VAL:HA	1:A:353:MET:HE2	2.01	0.43
1:A:406:VAL:HG21	1:A:439:LEU:HD13	2.00	0.43
1:A:634:THR:O	1:A:638:VAL:HG23	2.19	0.43
1:A:660:ASN:O	1:A:662:ALA:N	2.52	0.43
1:A:714:LEU:HD22	1:A:739:LEU:CD2	2.48	0.43
1:A:268:MET:HE1	1:A:322:LEU:HD23	2.00	0.43
1:A:559:GLN:C	1:A:561:VAL:N	2.75	0.43
1:A:564:MET:HB3	1:A:577:PHE:HE1	1.83	0.42
1:A:525:ARG:HG2	1:A:525:ARG:HH11	1.84	0.42
1:A:597:HIS:O	1:A:597:HIS:ND1	2.52	0.42
1:A:377:ASN:ND2	1:A:378:PRO:CD	2.82	0.42
1:A:820:GLY:HA3	2:B:51:ARG:HH12	1.83	0.42
1:A:564:MET:HE2	1:A:567:HIS:HD2	1.84	0.42
1:A:620:GLY:O	1:A:621:SER:HB3	2.18	0.42
1:A:692:VAL:O	1:A:696:LEU:HG	2.20	0.42
1:A:578:ASN:O	1:A:581:GLN:HB2	2.20	0.42
1:A:758:LEU:HD22	1:A:762:ARG:CD	2.49	0.42
1:A:46:ASN:HB3	1:A:49:ASN:ND2	2.35	0.42
1:A:740:GLN:NE2	1:A:796:PHE:CE2	2.88	0.42
1:A:121:ILE:N	1:A:122:PRO:HD2	2.35	0.42
1:A:430:TRP:CZ2	1:A:434:ARG:CD	3.03	0.42
1:A:430:TRP:NE1	2:B:18:LYS:HA	2.34	0.42
1:A:504:GLU:HA	1:A:507:VAL:HG22	2.02	0.42
1:A:136:ASN:HD22	1:A:146:MET:HE2	1.85	0.41
1:A:99:LEU:HB3	1:A:146:MET:HE1	2.02	0.41
1:A:714:LEU:O	1:A:717:PHE:HB2	2.20	0.41
1:A:768:ALA:O	1:A:772:ILE:HG13	2.19	0.41
1:A:791:GLN:HG2	1:A:829:PHE:CD1	2.56	0.41
1:A:707:ARG:HG3	1:A:707:ARG:NH1	2.33	0.41
1:A:68:LYS:O	1:A:70:PRO:HD3	2.20	0.41
1:A:123:VAL:O	1:A:123:VAL:CG1	2.65	0.41
1:A:343:ASN:HB2	1:A:344:PRO:CD	2.50	0.41
1:A:377:ASN:HA	1:A:378:PRO:HD3	1.87	0.41
1:A:541:LYS:C	1:A:543:CYS:H	2.28	0.41
1:A:37:LEU:HD23	1:A:37:LEU:HA	1.77	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:703:GLU:O	1:A:703:GLU:HG2	2.20	0.41
1:A:264:THR:HG21	1:A:282:PHE:CD2	2.55	0.41
1:A:655:GLY:O	1:A:658:LEU:N	2.54	0.41
1:A:685:ILE:O	1:A:685:ILE:HG12	2.21	0.41
1:A:686:ILE:HG23	1:A:727:GLU:HG2	2.03	0.41
1:A:740:GLN:O	1:A:744:GLN:HG3	2.20	0.41
1:A:780:GLN:HB2	1:A:781:GLU:H	1.69	0.41
1:A:811:THR:C	1:A:813:GLY:H	2.28	0.41
1:A:24:PHE:CD1	1:A:24:PHE:C	2.99	0.41
1:A:172:GLU:CD	1:A:172:GLU:N	2.75	0.41
1:A:475:SER:O	1:A:479:GLU:HG3	2.21	0.41
1:A:568:ILE:HG13	1:A:568:ILE:O	2.21	0.41
1:A:622:GLY:O	1:A:626:GLU:HG3	2.20	0.41
1:A:686:ILE:N	1:A:687:PRO:HD2	2.36	0.41
1:A:36:PHE:CE2	1:A:40:LEU:HD11	2.56	0.40
1:A:403:LYS:N	1:A:404:PRO:HD2	2.35	0.40
1:A:130:ILE:O	1:A:134:VAL:HG23	2.20	0.40
1:A:368:LEU:HD23	1:A:368:LEU:HA	1.94	0.40
1:A:543:CYS:O	1:A:547:VAL:HG23	2.22	0.40
1:A:608:MET:HE2	1:A:611:LEU:HD12	2.03	0.40
1:A:801:ILE:HG21	1:A:822:ILE:HD11	2.04	0.40
2:B:51:ARG:O	2:B:52:ASN:CB	2.70	0.40
1:A:615:PHE:CE1	1:A:625:GLN:NE2	2.90	0.40
1:A:699:ASN:ND2	1:A:705:VAL:HG11	2.36	0.40
1:A:821:LEU:O	1:A:822:ILE:C	2.65	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	874/876 (100%)	795 (91%)	62 (7%)	17 (2%)	6 11

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	42/44 (96%)	39 (93%)	2 (5%)	1 (2%)	4 8
All	All	916/920 (100%)	834 (91%)	64 (7%)	18 (2%)	6 10

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	492	GLU
1	A	781	GLU
1	A	782	ASN
1	A	811	THR
1	A	853	SER
1	A	568	ILE
1	A	661	TYR
1	A	660	ASN
1	A	664	TYR
1	A	703	GLU
1	A	749	LYS
1	A	808	GLU
2	B	52	ASN
1	A	124	ASN
1	A	259	ALA
1	A	833	VAL
1	A	105	ARG
1	A	545	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	751/751 (100%)	721 (96%)	30 (4%)	28 54
2	B	40/40 (100%)	39 (98%)	1 (2%)	42 69
All	All	791/791 (100%)	760 (96%)	31 (4%)	28 55

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	22	GLN
1	A	111	GLN
1	A	209	PHE
1	A	234	ARG
1	A	278	GLN
1	A	331	THR
1	A	343	ASN
1	A	377	ASN
1	A	408	GLN
1	A	446	ASP
1	A	474	PHE
1	A	486	ASP
1	A	490	ASP
1	A	497	TYR
1	A	504	GLU
1	A	508	GLN
1	A	514	THR
1	A	524	LEU
1	A	580	LEU
1	A	608	MET
1	A	673	LEU
1	A	732	LEU
1	A	734	VAL
1	A	758	LEU
1	A	780	GLN
1	A	782	ASN
1	A	814	VAL
1	A	821	LEU
1	A	867	LYS
2	B	25	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	32	ASN
1	A	60	GLN
1	A	132	GLN
1	A	136	ASN
1	A	141	ASN
1	A	179	GLN
1	A	216	HIS

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Mol	Chain	Res	Type
1	A	239	GLN
1	A	240	ASN
1	A	285	ASN
1	A	303	GLN
1	A	343	ASN
1	A	366	HIS
1	A	377	ASN
1	A	401	GLN
1	A	454	GLN
1	A	520	HIS
1	A	548	GLN
1	A	567	HIS
1	A	576	GLN
1	A	581	GLN
1	A	589	GLN
1	A	596	GLN
1	A	616	GLN
1	A	699	ASN
1	A	702	ASN
1	A	759	ASN
1	A	782	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 oligosaccharides 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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.