



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:31 AM UTC

PDB ID : 2Q6D / pdb_00002q6d
Title : Crystal structure of infectious bronchitis virus (IBV) main protease
Authors : Xue, X.Y.; Yang, H.T.; Xue, F.; Bartlam, M.; Rao, Z.H.
Deposited on : 2007-06-04
Resolution : 2.35 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)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

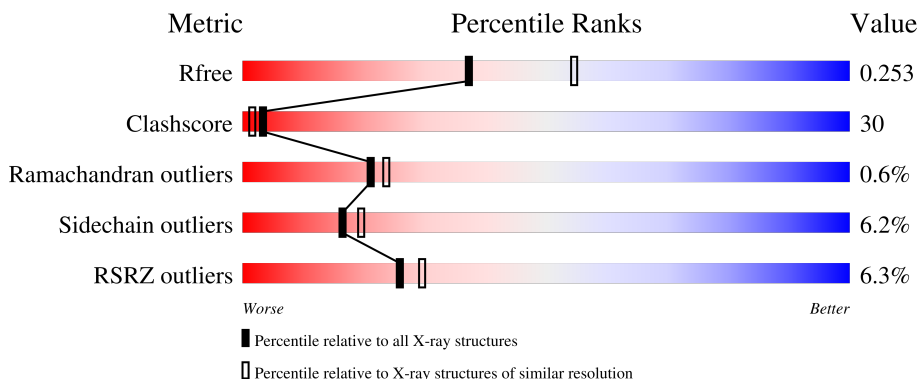
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	180053	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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 $\leq 5\%$. 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	309	
1	B	309	
1	C	309	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7515 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 Infectious bronchitis virus (IBV) main protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2274	1450	381	433	10			
1	B	307	Total	C	N	O	S	0	0	0
			2355	1500	396	449	10			
1	C	302	Total	C	N	O	S	0	0	0
			2320	1476	390	444	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q3Y5H1
A	0	SER	-	expression tag	UNP Q3Y5H1
B	-1	GLY	-	expression tag	UNP Q3Y5H1
B	0	SER	-	expression tag	UNP Q3Y5H1
C	-1	GLY	-	expression tag	UNP Q3Y5H1
C	0	SER	-	expression tag	UNP Q3Y5H1

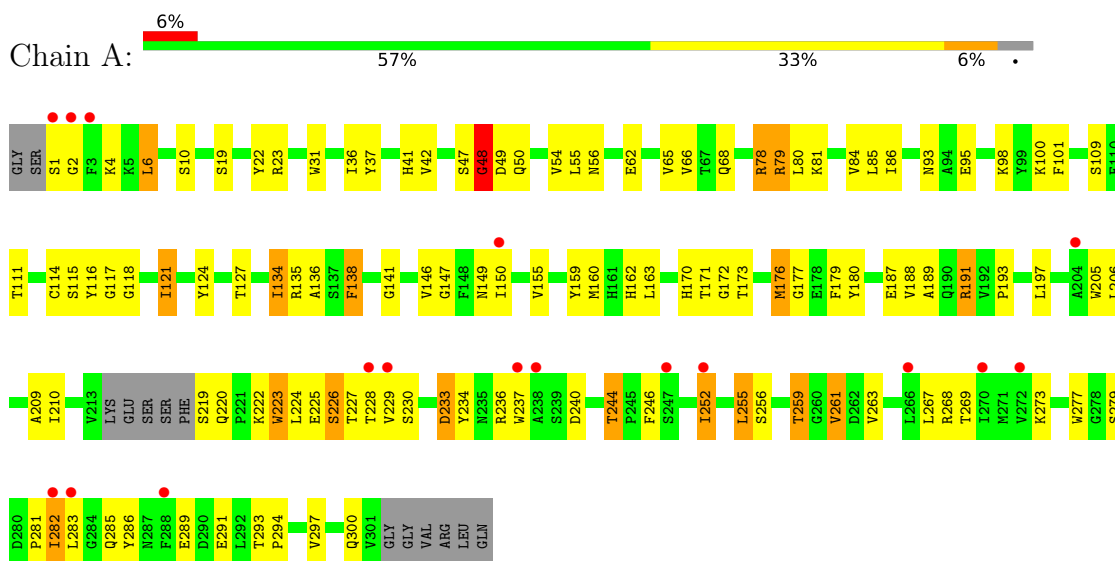
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	167	Total	O	0	0
			167	167		
2	B	194	Total	O	0	0
			194	194		
2	C	205	Total	O	0	0
			205	205		

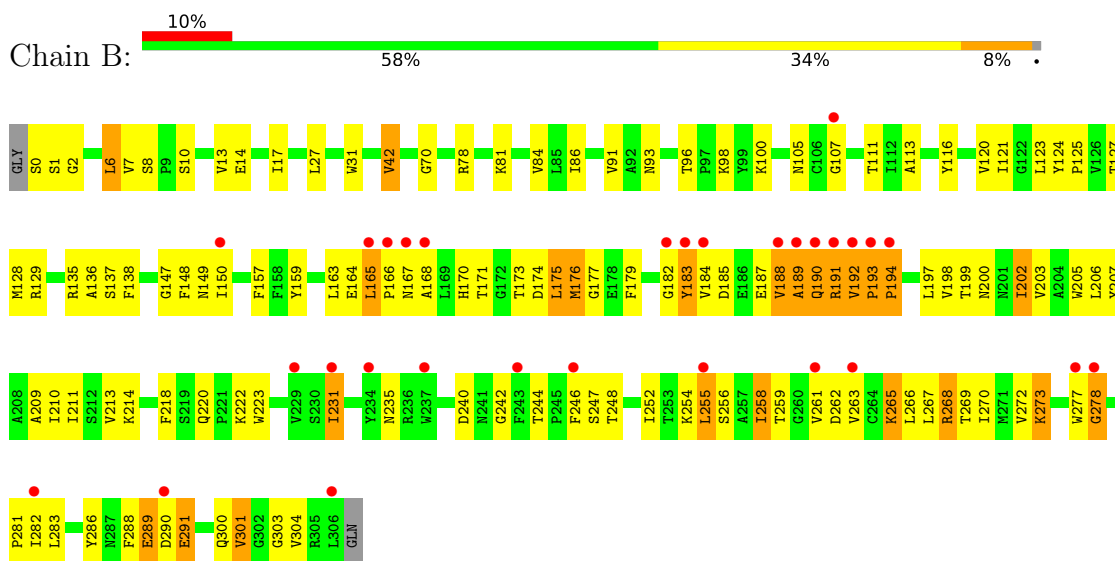
3 Residue-property plots [i](#)

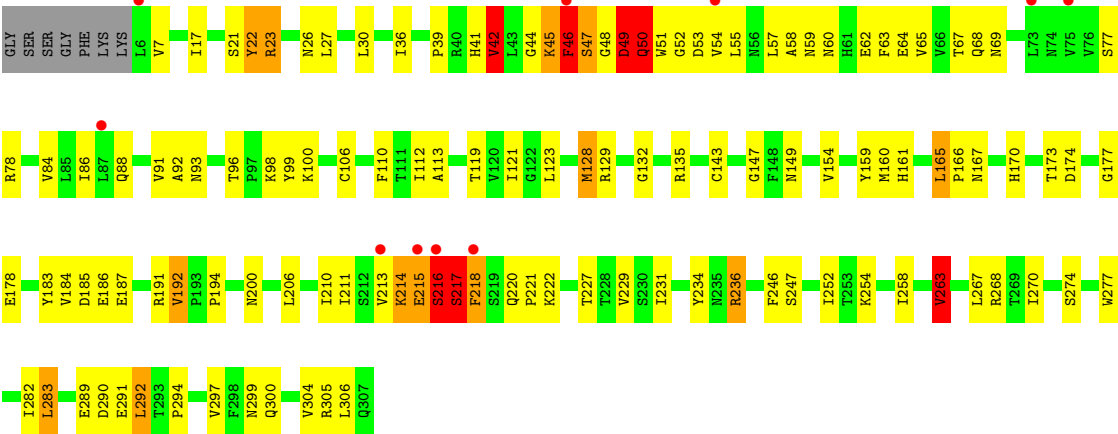
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

- Molecule 1: Infectious bronchitis virus (IBV) main protease



- Molecule 1: Infectious bronchitis virus (IBV) main protease





4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	118.90Å 118.90Å 270.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.35 50.00 – 2.35	Depositor EDS
% Data completeness (in resolution range)	95.6 (50.00-2.35) 95.7 (50.00-2.35)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.76 (at 2.34Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.226 , 0.259 0.219 , 0.253	Depositor DCC
R_{free} test set	4788 reflections (10.08%)	wwPDB-VP
Wilson B-factor (Å ²)	41.1	Xtriage
Anisotropy	0.316	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7515	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.61% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.61	1/2325 (0.0%)	1.06	15/3165 (0.5%)
1	B	0.54	2/2408 (0.1%)	1.13	23/3276 (0.7%)
1	C	0.62	3/2372 (0.1%)	1.12	21/3229 (0.7%)
All	All	0.59	6/7105 (0.1%)	1.11	59/9670 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	3
All	All	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	49	ASP	C-N	-12.12	1.12	1.33
1	A	244	THR	C-N	-10.51	1.22	1.33
1	B	183	TYR	C-N	8.87	1.47	1.33
1	B	194	PRO	C-O	8.13	1.32	1.23
1	C	7	VAL	N-CA	-6.64	1.38	1.46
1	C	128	MET	SD-CE	-5.49	1.65	1.79

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	189	ALA	CB-CA-C	-15.65	83.98	109.80
1	B	165	LEU	N-CA-C	-10.60	96.47	109.72
1	C	49	ASP	O-C-N	-9.60	107.64	123.00
1	B	190	GLN	N-CA-CB	-9.33	96.96	111.23
1	C	217	SER	CB-CA-C	-9.23	100.70	109.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	256	SER	N-CA-C	-8.92	100.05	112.45
1	A	223	TRP	N-CA-CB	-8.50	99.59	110.90
1	C	186	GLU	N-CA-C	-8.40	102.07	111.82
1	A	259	THR	CA-C-N	-8.27	109.72	120.22
1	A	259	THR	C-N-CA	-8.27	109.72	120.22
1	B	174	ASP	N-CA-C	-8.22	97.11	109.86
1	C	45	LYS	N-CA-C	-8.20	98.18	110.24
1	A	220	GLN	CA-C-N	8.11	128.16	119.89
1	A	220	GLN	C-N-CA	8.11	128.16	119.89
1	B	192	VAL	CA-C-N	8.06	128.68	120.38
1	B	192	VAL	C-N-CA	8.06	128.68	120.38
1	C	7	VAL	N-CA-C	7.91	119.24	108.17
1	C	45	LYS	CB-CA-C	7.56	121.19	109.84
1	C	50	GLN	N-CA-CB	7.51	122.73	111.01
1	B	191	ARG	N-CA-C	-7.28	95.85	108.52
1	C	159	TYR	N-CA-C	6.99	119.79	109.24
1	B	218	PHE	N-CA-C	-6.74	105.06	113.55
1	B	96	THR	N-CA-C	-6.53	101.84	109.93
1	C	49	ASP	CB-CA-C	-6.46	97.83	110.10
1	B	291	GLU	N-CA-C	6.40	120.40	112.59
1	C	46	PHE	N-CA-C	-6.37	100.42	108.45
1	C	42	VAL	N-CA-C	-6.36	103.14	111.09
1	C	50	GLN	N-CA-C	-6.19	104.49	114.09
1	B	98	LYS	N-CA-C	-6.11	100.79	109.96
1	B	10	SER	N-CA-C	6.10	120.43	112.92
1	B	176	MET	N-CA-C	-6.10	105.84	113.28
1	C	214	LYS	N-CA-C	-6.09	104.56	112.23
1	C	26	ASN	N-CA-C	6.06	119.78	110.20
1	B	231	ILE	N-CA-C	-5.85	104.81	110.42
1	A	146	VAL	N-CA-C	5.81	117.98	108.97
1	A	223	TRP	N-CA-C	5.80	121.79	114.31
1	C	289	GLU	N-CA-C	-5.80	99.28	108.73
1	B	265	LYS	N-CA-C	-5.72	105.19	111.82
1	A	159	TYR	N-CA-C	5.71	117.86	109.24
1	C	7	VAL	N-CA-CB	-5.70	103.50	111.25
1	A	6	LEU	N-CA-C	5.68	118.68	110.28
1	B	202	ILE	N-CA-C	-5.67	104.99	110.72
1	B	2	GLY	N-CA-C	-5.67	105.93	115.62
1	A	138	PHE	N-CA-C	-5.51	105.37	111.71
1	B	184	VAL	O-C-N	-5.51	117.25	122.98
1	B	42	VAL	N-CA-C	-5.45	104.40	112.04
1	C	229	VAL	N-CA-C	-5.39	100.08	108.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	188	VAL	N-CA-C	5.35	120.47	109.34
1	A	98	LYS	N-CA-C	-5.30	102.21	110.10
1	A	10	SER	N-CA-C	5.20	121.41	113.61
1	C	174	ASP	N-CA-C	-5.16	102.36	110.36
1	B	290	ASP	N-CA-C	5.15	119.84	113.50
1	C	263	VAL	CB-CA-C	-5.14	104.54	112.05
1	A	180	TYR	N-CA-C	-5.13	102.16	110.32
1	B	159	TYR	N-CA-C	5.09	117.13	109.23
1	C	46	PHE	CB-CA-C	-5.08	101.39	113.02
1	A	42	VAL	N-CA-C	-5.07	104.58	111.89
1	C	236	ARG	N-CA-C	-5.07	105.83	111.36
1	A	176	MET	N-CA-C	-5.02	107.21	113.38

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	48	GLY	Peptide
1	C	216	SER	Peptide
1	C	217	SER	Peptide
1	C	49	ASP	Mainchain

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	2274	0	2234	121	0
1	B	2355	0	2317	147	0
1	C	2320	0	2273	157	0
2	A	167	0	0	22	0
2	B	194	0	0	34	0
2	C	205	0	0	36	0
All	All	7515	0	6824	417	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 30.

All (417) 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:B:127:THR:HB	2:B:483:HOH:O	1.10	1.25
1:B:166:PRO:HG2	1:B:190:GLN:NE2	1.49	1.25
1:C:143:CYS:HB3	2:C:501:HOH:O	1.32	1.24
1:C:55:LEU:HD12	2:C:502:HOH:O	1.31	1.23
1:C:252:ILE:HG13	2:C:478:HOH:O	1.39	1.21
1:C:48:GLY:CA	1:C:49:ASP:HB2	1.65	1.21
1:C:98:LYS:HA	2:C:495:HOH:O	1.43	1.17
1:B:163:LEU:HD12	2:B:478:HOH:O	1.44	1.15
1:C:47:SER:O	1:C:50:GLN:HG3	1.47	1.15
1:C:291:GLU:HG2	2:C:511:HOH:O	1.43	1.15
1:B:166:PRO:HD2	1:B:190:GLN:CD	1.71	1.14
1:A:49:ASP:HB3	2:A:459:HOH:O	1.47	1.13
1:B:166:PRO:HG2	1:B:190:GLN:HE22	0.94	1.11
1:C:46:PHE:CE1	1:C:54:VAL:HG11	1.86	1.11
1:B:258:ILE:HG13	2:B:494:HOH:O	1.52	1.08
1:A:230:SER:HB3	1:A:233:ASP:HB2	1.38	1.05
1:A:95:GLU:HG3	2:A:466:HOH:O	1.57	1.04
1:B:166:PRO:CG	1:B:190:GLN:NE2	2.21	1.04
1:C:194:PRO:HG3	2:C:484:HOH:O	1.56	1.04
1:C:46:PHE:HE1	1:C:54:VAL:HG11	1.22	1.01
1:B:166:PRO:CD	1:B:190:GLN:CD	2.34	1.00
1:C:48:GLY:HA2	1:C:49:ASP:HB2	1.40	1.00
1:C:48:GLY:HA3	1:C:49:ASP:HB2	1.40	1.00
1:B:262:ASP:HB2	2:B:500:HOH:O	1.62	0.99
1:A:4:LYS:H	1:A:300:GLN:HE22	1.11	0.98
1:A:55:LEU:HB3	2:A:458:HOH:O	1.67	0.94
1:B:175:LEU:CD1	2:B:493:HOH:O	2.14	0.93
1:B:166:PRO:HD2	1:B:190:GLN:OE1	1.66	0.93
1:B:166:PRO:CG	1:B:190:GLN:HE22	1.81	0.92
1:C:48:GLY:HA3	1:C:49:ASP:CB	1.97	0.91
1:B:164:GLU:HB2	2:B:491:HOH:O	1.70	0.91
1:C:23:ARG:HG3	1:C:23:ARG:HH11	1.35	0.91
1:C:160:MET:SD	2:C:504:HOH:O	2.28	0.91
1:A:49:ASP:HB2	2:A:462:HOH:O	1.70	0.91
1:B:191:ARG:O	1:B:193:PRO:HD3	1.69	0.90
1:C:47:SER:C	1:C:50:GLN:HE21	1.79	0.90
1:A:163:LEU:HD11	1:A:171:THR:HG22	1.52	0.89
1:B:13:VAL:HG12	1:B:17:ILE:HD11	1.55	0.88
1:A:277:TRP:CZ2	1:A:282:ILE:HD13	2.09	0.88
1:B:188:VAL:HG13	1:B:189:ALA:H	1.39	0.87
1:C:106:CYS:HA	1:C:128:MET:HE2	1.58	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:LEU:HD11	1:C:192:VAL:HG13	1.58	0.85
1:A:150:ILE:HD12	1:A:155:VAL:HG22	1.59	0.84
1:B:166:PRO:HG2	1:B:190:GLN:CD	2.03	0.84
1:B:262:ASP:HB3	2:B:345:HOH:O	1.78	0.84
1:C:23:ARG:HH11	1:C:23:ARG:CG	1.92	0.83
1:C:210:ILE:HG22	1:C:214:LYS:HE2	1.59	0.82
1:B:8:SER:HB3	1:B:150:ILE:HD13	1.61	0.80
1:A:277:TRP:HZ2	1:A:282:ILE:HD13	1.47	0.80
1:C:47:SER:N	1:C:50:GLN:HE21	1.80	0.80
1:A:114:CYS:O	1:A:121:ILE:HD13	1.82	0.79
1:C:84:VAL:HG23	1:C:177:GLY:HA2	1.63	0.78
1:B:166:PRO:CG	1:B:190:GLN:CD	2.56	0.78
1:C:214:LYS:O	1:C:217:SER:N	2.17	0.78
1:C:77:SER:HB3	1:C:88:GLN:HE21	1.49	0.77
1:C:47:SER:H	1:C:50:GLN:HE21	1.31	0.77
1:A:134:ILE:HD13	1:A:170:HIS:O	1.84	0.77
1:C:222:LYS:HG3	2:C:503:HOH:O	1.85	0.76
1:C:47:SER:C	1:C:50:GLN:HG3	2.10	0.76
1:A:150:ILE:CD1	1:A:155:VAL:HG22	2.15	0.76
1:B:107:GLY:HA3	2:B:497:HOH:O	1.84	0.76
1:A:237:TRP:C	2:A:470:HOH:O	2.27	0.76
1:A:252:ILE:HA	1:A:255:LEU:HD12	1.68	0.76
1:C:47:SER:H	1:C:50:GLN:NE2	1.83	0.75
1:A:114:CYS:SG	1:A:121:ILE:HD11	2.26	0.75
1:B:170:HIS:CE1	2:B:474:HOH:O	2.41	0.74
1:C:214:LYS:O	1:C:217:SER:HA	1.87	0.74
1:C:218:PHE:CE1	2:C:507:HOH:O	2.41	0.74
1:A:163:LEU:CD1	1:A:171:THR:HG22	2.17	0.73
1:A:4:LYS:H	1:A:300:GLN:NE2	1.84	0.73
1:B:17:ILE:HD13	1:B:120:VAL:CG2	2.19	0.73
1:B:202:ILE:HG21	1:B:270:ILE:HD12	1.71	0.73
1:C:47:SER:CA	1:C:50:GLN:HE21	2.01	0.72
1:B:157:PHE:HB3	1:B:175:LEU:HD13	1.70	0.72
1:C:129:ARG:NH2	1:C:290:ASP:OD1	2.22	0.72
1:C:231:ILE:HD12	1:C:246:PHE:CD2	2.25	0.72
1:C:47:SER:N	1:C:50:GLN:NE2	2.37	0.71
1:C:41:HIS:CE1	2:C:501:HOH:O	2.42	0.71
1:B:265:LYS:HD2	2:B:486:HOH:O	1.89	0.71
1:A:47:SER:O	1:A:48:GLY:C	2.34	0.71
1:C:23:ARG:HD3	1:C:23:ARG:N	2.05	0.71
1:A:1:SER:N	2:A:425:HOH:O	2.21	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:64:GLU:HG3	2:C:505:HOH:O	1.90	0.70
1:C:128:MET:HE3	1:C:132:GLY:HA2	1.73	0.70
1:A:68:GLN:HG3	2:A:455:HOH:O	1.91	0.70
1:C:252:ILE:CG1	2:C:478:HOH:O	2.11	0.70
1:C:213:VAL:HG12	2:C:508:HOH:O	1.91	0.70
1:A:109:SER:HB3	1:A:127:THR:HG22	1.74	0.70
1:B:8:SER:HB3	1:B:150:ILE:CD1	2.22	0.70
1:C:22:TYR:O	1:C:22:TYR:CD2	2.45	0.70
1:B:127:THR:OG1	1:B:291:GLU:HG2	1.91	0.69
1:A:210:ILE:HD13	1:A:261:VAL:HG21	1.74	0.69
1:A:273:LYS:HE2	2:A:473:HOH:O	1.91	0.69
1:A:279:SER:O	2:A:446:HOH:O	2.10	0.69
1:C:46:PHE:CE1	1:C:54:VAL:CG1	2.73	0.68
1:C:231:ILE:HD13	1:C:267:LEU:CD1	2.23	0.68
1:C:47:SER:C	1:C:50:GLN:NE2	2.52	0.68
1:B:0:SER:CB	2:B:488:HOH:O	2.40	0.68
1:C:27:LEU:HD21	1:C:42:VAL:HG22	1.76	0.67
1:B:231:ILE:HD12	1:B:246:PHE:CD2	2.28	0.67
1:B:277:TRP:CZ2	1:B:282:ILE:HD13	2.30	0.67
1:C:22:TYR:CD2	1:C:22:TYR:C	2.72	0.67
1:A:101:PHE:HD1	1:A:176:MET:HE3	1.59	0.67
1:C:27:LEU:HG	1:C:42:VAL:HG13	1.77	0.67
1:A:222:LYS:NZ	2:A:448:HOH:O	2.28	0.67
1:C:165:LEU:HD11	1:C:192:VAL:CG1	2.23	0.67
1:A:23:ARG:HD3	1:A:62:GLU:OE2	1.96	0.66
1:C:47:SER:O	1:C:50:GLN:CG	2.35	0.65
1:B:170:HIS:HE1	2:B:474:HOH:O	1.79	0.65
1:C:143:CYS:CB	2:C:501:HOH:O	2.09	0.65
1:B:222:LYS:HG3	2:B:481:HOH:O	1.96	0.65
1:C:210:ILE:CG2	1:C:214:LYS:HE2	2.27	0.65
1:B:248:THR:CG2	2:B:475:HOH:O	2.43	0.65
1:A:281:PRO:C	1:A:282:ILE:HD12	2.22	0.64
1:A:197:LEU:O	1:A:244:THR:HG23	1.97	0.64
1:B:183:TYR:N	1:B:183:TYR:CD2	2.64	0.64
1:B:166:PRO:CD	1:B:190:GLN:OE1	2.38	0.64
1:B:175:LEU:HG	2:B:493:HOH:O	1.97	0.64
1:A:55:LEU:HD22	1:A:80:LEU:HG	1.79	0.64
1:C:187:GLU:O	1:C:191:ARG:HD2	1.98	0.63
1:A:230:SER:CB	1:A:233:ASP:HB2	2.23	0.63
1:C:63:PHE:HE2	1:C:78:ARG:HD3	1.63	0.63
1:C:22:TYR:O	1:C:22:TYR:HD2	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:LEU:CD1	2:C:502:HOH:O	2.10	0.63
1:C:165:LEU:CD1	1:C:192:VAL:HG13	2.29	0.63
1:C:106:CYS:SG	1:C:128:MET:CE	2.87	0.63
1:C:143:CYS:HA	1:C:161:HIS:HD2	1.62	0.63
1:C:214:LYS:O	1:C:217:SER:CA	2.47	0.63
1:C:112:ILE:HD13	1:C:147:GLY:CA	2.29	0.62
1:A:149:ASN:O	1:A:150:ILE:HD13	1.98	0.62
1:B:166:PRO:CD	1:B:190:GLN:NE2	2.58	0.62
1:C:184:VAL:HG23	1:C:185:ASP:H	1.64	0.62
1:C:98:LYS:HG3	2:C:500:HOH:O	2.00	0.62
1:B:13:VAL:O	1:B:17:ILE:HD12	1.99	0.62
1:A:121:ILE:HD13	1:A:121:ILE:H	1.65	0.61
1:A:47:SER:O	1:A:49:ASP:N	2.33	0.61
1:B:129:ARG:NH2	1:B:135:ARG:HH21	1.98	0.60
1:B:202:ILE:HD12	1:B:288:PHE:HB3	1.82	0.60
1:C:58:ALA:O	1:C:78:ARG:NH2	2.34	0.60
1:A:50:GLN:O	1:A:54:VAL:HG23	2.00	0.60
1:C:214:LYS:C	1:C:217:SER:H	2.07	0.60
1:A:138:PHE:O	1:B:303:GLY:HA3	2.01	0.60
1:B:187:GLU:HG2	1:B:187:GLU:O	2.01	0.60
1:B:188:VAL:HG22	1:B:189:ALA:N	2.16	0.60
1:B:198:VAL:O	1:B:202:ILE:HG12	2.01	0.60
1:A:41:HIS:HD2	2:A:453:HOH:O	1.85	0.60
1:C:299:ASN:CB	2:C:486:HOH:O	2.49	0.60
1:B:301:VAL:O	1:B:304:VAL:HG22	2.02	0.60
1:C:67:THR:HG22	1:C:68:GLN:N	2.15	0.60
1:A:19:SER:OG	1:A:66:VAL:HB	2.01	0.59
1:C:44:GLY:O	1:C:45:LYS:C	2.43	0.59
1:C:210:ILE:HD13	1:C:221:PRO:HG2	1.85	0.59
1:A:19:SER:HB2	2:A:457:HOH:O	2.01	0.59
1:A:176:MET:HA	1:A:176:MET:HE2	1.85	0.59
1:B:107:GLY:C	2:B:497:HOH:O	2.45	0.59
1:C:213:VAL:CG1	2:C:508:HOH:O	2.48	0.59
1:A:163:LEU:HD11	1:A:171:THR:CG2	2.30	0.59
1:B:193:PRO:HB3	1:B:194:PRO:HD2	1.84	0.59
1:B:206:LEU:HD11	1:B:270:ILE:HD11	1.85	0.59
1:A:78:ARG:C	1:A:79:ARG:HG3	2.27	0.58
1:C:78:ARG:HG3	1:C:78:ARG:HH11	1.69	0.58
1:A:109:SER:CB	1:A:127:THR:HG22	2.34	0.58
1:C:299:ASN:HB2	2:C:486:HOH:O	2.03	0.58
1:B:107:GLY:CA	2:B:497:HOH:O	2.46	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:299:ASN:HA	2:C:486:HOH:O	2.03	0.58
1:B:211:ILE:HD13	1:B:301:VAL:CG2	2.34	0.57
1:A:37:TYR:O	1:A:160:MET:HE1	2.04	0.57
1:A:116:TYR:HD2	1:A:121:ILE:HD12	1.69	0.57
1:C:52:GLY:O	1:C:55:LEU:HB3	2.04	0.57
1:C:55:LEU:HD21	2:C:487:HOH:O	2.02	0.57
1:C:46:PHE:CE1	1:C:54:VAL:HG21	2.39	0.57
1:B:209:ALA:O	1:B:213:VAL:HG23	2.05	0.57
1:B:188:VAL:HG13	1:B:189:ALA:N	2.15	0.57
1:A:19:SER:CB	2:A:457:HOH:O	2.53	0.56
1:C:23:ARG:HG3	1:C:23:ARG:NH1	2.14	0.56
1:A:127:THR:OG1	1:A:291:GLU:OE2	2.19	0.56
1:B:14:GLU:HA	1:B:17:ILE:HD12	1.87	0.56
1:B:206:LEU:HD11	1:B:270:ILE:CD1	2.35	0.56
1:A:286:TYR:HB2	1:B:286:TYR:HB2	1.88	0.56
1:A:282:ILE:HD12	1:A:282:ILE:N	2.21	0.56
1:C:161:HIS:CE1	1:C:170:HIS:HD2	2.24	0.56
1:A:236:ARG:NH1	2:A:447:HOH:O	2.38	0.56
1:B:148:PHE:HE1	1:B:150:ILE:HD11	1.70	0.56
1:C:48:GLY:CA	1:C:49:ASP:CB	2.46	0.55
1:C:50:GLN:O	1:C:54:VAL:HG23	2.07	0.55
1:C:213:VAL:HG12	1:C:213:VAL:O	2.05	0.55
1:B:175:LEU:HD11	2:B:493:HOH:O	1.92	0.55
1:B:268:ARG:O	1:B:272:VAL:HG23	2.06	0.55
1:A:149:ASN:C	1:A:149:ASN:OD1	2.50	0.55
1:B:200:ASN:HD21	1:B:247:SER:H	1.53	0.55
1:C:17:ILE:HD12	2:C:485:HOH:O	2.06	0.55
1:C:47:SER:HB2	2:C:313:HOH:O	2.06	0.55
1:C:297:VAL:O	1:C:300:GLN:HB2	2.06	0.54
1:B:168:ALA:HB3	2:B:469:HOH:O	2.05	0.54
1:A:81:LYS:HB3	1:A:86:ILE:HD11	1.90	0.54
1:B:129:ARG:CZ	1:B:135:ARG:HH21	2.20	0.54
1:C:106:CYS:CA	1:C:128:MET:HE2	2.34	0.54
1:A:224:LEU:HD12	1:A:225:GLU:N	2.23	0.54
1:C:77:SER:HB3	1:C:88:GLN:HB3	1.90	0.54
1:A:36:ILE:HD13	1:A:65:VAL:HG11	1.89	0.54
1:B:17:ILE:HD13	1:B:120:VAL:HG23	1.88	0.54
1:C:63:PHE:CE2	1:C:78:ARG:HD3	2.41	0.54
1:A:116:TYR:CZ	1:A:141:GLY:HA2	2.43	0.54
1:A:36:ILE:CD1	1:A:65:VAL:HG11	2.38	0.53
1:A:49:ASP:CG	1:A:49:ASP:O	2.50	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LEU:HD11	1:A:78:ARG:HG3	1.90	0.53
1:B:206:LEU:CD1	1:B:270:ILE:HD11	2.38	0.53
1:A:173:THR:HG22	1:A:179:PHE:HA	1.90	0.53
1:C:23:ARG:CG	1:C:23:ARG:NH1	2.59	0.53
1:B:214:LYS:NZ	1:B:259:THR:O	2.42	0.53
1:A:163:LEU:HD12	1:A:163:LEU:O	2.09	0.53
1:B:193:PRO:CB	1:B:194:PRO:HD2	2.38	0.53
1:C:113:ALA:HA	1:C:123:LEU:HD23	1.90	0.53
1:B:14:GLU:HA	1:B:17:ILE:CD1	2.39	0.53
1:A:224:LEU:HD12	1:A:225:GLU:H	1.74	0.53
1:C:292:LEU:HD22	1:C:297:VAL:HG23	1.91	0.53
1:A:205:TRP:NE1	1:A:283:LEU:HD12	2.23	0.53
1:B:0:SER:HB3	2:B:501:HOH:O	2.08	0.52
1:B:270:ILE:HD12	1:B:288:PHE:HE2	1.74	0.52
1:C:161:HIS:HE1	1:C:170:HIS:HD2	1.57	0.52
1:B:166:PRO:CG	1:B:190:GLN:OE1	2.56	0.52
1:B:175:LEU:HD12	2:B:493:HOH:O	1.94	0.52
1:B:150:ILE:N	1:B:150:ILE:HD12	2.25	0.52
1:A:19:SER:O	1:A:65:VAL:HA	2.10	0.52
1:C:161:HIS:HE1	1:C:170:HIS:CD2	2.28	0.52
1:A:229:VAL:CG2	1:A:268:ARG:HE	2.22	0.52
1:A:256:SER:O	1:A:259:THR:O	2.27	0.52
1:C:254:LYS:O	1:C:258:ILE:HG13	2.10	0.52
1:C:218:PHE:HE1	2:C:507:HOH:O	1.88	0.51
1:C:231:ILE:HD13	1:C:267:LEU:HD12	1.91	0.51
1:B:197:LEU:HD12	2:B:470:HOH:O	2.11	0.51
1:C:59:ASN:HB3	1:C:62:GLU:HG3	1.92	0.51
1:B:7:VAL:CG2	1:B:125:PRO:HD3	2.40	0.51
1:B:17:ILE:HD13	1:B:120:VAL:HG21	1.93	0.51
1:B:211:ILE:HD13	1:B:301:VAL:HG21	1.92	0.51
1:B:202:ILE:HD13	1:B:289:GLU:O	2.11	0.51
1:C:129:ARG:HH22	1:C:290:ASP:CG	2.18	0.51
1:B:261:VAL:HG12	1:B:262:ASP:N	2.26	0.51
1:A:86:ILE:N	1:A:86:ILE:HD12	2.25	0.50
1:B:203:VAL:CG1	1:B:252:ILE:HD12	2.41	0.50
1:A:111:THR:O	1:A:147:GLY:HA2	2.11	0.50
1:A:229:VAL:HG22	1:A:268:ARG:HE	1.75	0.50
1:C:47:SER:CA	1:C:50:GLN:NE2	2.71	0.50
1:C:135:ARG:HD2	2:C:483:HOH:O	2.11	0.50
1:C:93:ASN:HB3	1:C:96:THR:OG1	2.11	0.50
1:C:67:THR:CG2	1:C:68:GLN:N	2.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:LYS:HE3	2:A:464:HOH:O	2.12	0.50
1:A:85:LEU:C	1:A:86:ILE:HD12	2.37	0.50
1:B:7:VAL:HG13	1:B:111:THR:CG2	2.42	0.50
1:C:143:CYS:SG	2:C:501:HOH:O	2.57	0.50
1:A:233:ASP:O	1:A:236:ARG:HB3	2.12	0.49
1:A:117:GLY:HA2	2:A:452:HOH:O	2.11	0.49
1:B:273:LYS:NZ	1:B:278:GLY:H	2.09	0.49
1:A:47:SER:H	1:A:50:GLN:NE2	2.09	0.49
1:B:27:LEU:HD21	1:B:42:VAL:HB	1.94	0.49
1:B:84:VAL:HG23	1:B:177:GLY:HA2	1.94	0.49
1:B:214:LYS:HE2	1:B:220:GLN:NE2	2.27	0.49
1:C:86:ILE:HD12	1:C:86:ILE:N	2.27	0.49
1:C:161:HIS:CE1	1:C:170:HIS:CD2	3.00	0.49
1:B:149:ASN:C	1:B:150:ILE:HD12	2.38	0.49
1:C:268:ARG:HG3	1:C:268:ARG:HH11	1.76	0.49
1:C:23:ARG:N	1:C:23:ARG:CD	2.75	0.49
1:C:299:ASN:CA	2:C:486:HOH:O	2.61	0.49
1:A:79:ARG:HD3	2:A:396:HOH:O	2.13	0.48
1:B:265:LYS:CD	2:B:486:HOH:O	2.56	0.48
1:B:0:SER:HB2	2:B:488:HOH:O	2.11	0.48
1:A:41:HIS:CD2	2:A:453:HOH:O	2.64	0.48
1:A:176:MET:HE2	2:A:400:HOH:O	2.13	0.48
1:A:223:TRP:CE3	1:A:269:THR:HG21	2.48	0.48
1:A:234:TYR:CD2	1:A:267:LEU:HB3	2.48	0.48
1:B:113:ALA:HA	1:B:123:LEU:HD23	1.95	0.48
1:A:223:TRP:CE3	1:A:224:LEU:N	2.81	0.48
1:B:183:TYR:N	1:B:183:TYR:HD2	2.10	0.48
1:C:210:ILE:O	1:C:214:LYS:HG2	2.14	0.48
1:A:4:LYS:N	1:A:300:GLN:HE22	1.94	0.48
1:C:88:GLN:HB2	2:C:496:HOH:O	2.14	0.48
1:C:143:CYS:HA	1:C:161:HIS:CD2	2.46	0.48
1:A:227:THR:HB	1:A:268:ARG:HD3	1.95	0.48
1:A:163:LEU:HD22	1:C:304:VAL:HG11	1.95	0.47
1:A:281:PRO:HB3	1:A:286:TYR:CE2	2.49	0.47
1:B:165:LEU:O	1:B:166:PRO:C	2.57	0.47
1:C:67:THR:HG22	1:C:69:ASN:H	1.79	0.47
1:B:281:PRO:HD3	1:B:286:TYR:CE2	2.49	0.47
1:B:127:THR:CG2	1:B:291:GLU:HG2	2.44	0.47
1:B:240:ASP:HA	2:B:476:HOH:O	2.13	0.47
1:B:173:THR:HG22	1:B:179:PHE:HA	1.97	0.47
1:B:86:ILE:N	1:B:86:ILE:HD12	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:ILE:C	1:B:121:ILE:HD12	2.40	0.47
1:B:165:LEU:O	1:B:166:PRO:O	2.33	0.47
1:B:202:ILE:CG2	1:B:270:ILE:HD12	2.43	0.47
1:A:188:VAL:HG12	1:A:189:ALA:N	2.28	0.47
1:C:60:ASN:OD1	1:C:78:ARG:NE	2.47	0.47
1:C:121:ILE:HD12	1:C:121:ILE:N	2.30	0.47
1:C:216:SER:O	1:C:217:SER:HB2	2.15	0.47
1:C:39:PRO:O	1:C:42:VAL:CG2	2.63	0.46
1:A:49:ASP:CB	2:A:462:HOH:O	2.46	0.46
1:B:149:ASN:C	1:B:149:ASN:OD1	2.58	0.46
1:A:101:PHE:CD1	1:A:176:MET:HE3	2.46	0.46
1:C:99:TYR:N	2:C:495:HOH:O	2.49	0.46
1:C:270:ILE:O	1:C:274:SER:HB2	2.16	0.46
1:B:199:THR:H	1:B:244:THR:HG1	1.62	0.46
1:B:200:ASN:HD22	1:B:200:ASN:N	2.13	0.46
1:C:112:ILE:HD13	1:C:147:GLY:HA2	1.97	0.46
1:A:223:TRP:O	1:A:224:LEU:C	2.56	0.46
1:B:223:TRP:CE2	1:B:282:ILE:HG13	2.51	0.46
1:C:55:LEU:CD2	2:C:487:HOH:O	2.60	0.46
1:C:213:VAL:CG1	1:C:213:VAL:O	2.63	0.46
1:C:220:GLN:NE2	2:C:357:HOH:O	2.47	0.46
1:A:163:LEU:HD12	1:A:163:LEU:C	2.40	0.46
1:A:188:VAL:CG1	1:A:189:ALA:N	2.79	0.46
1:B:242:GLY:O	2:B:470:HOH:O	2.21	0.46
1:A:31:TRP:CD2	1:A:93:ASN:HB2	2.50	0.46
1:C:78:ARG:HD2	2:C:502:HOH:O	2.15	0.46
1:A:135:ARG:HD2	1:B:0:SER:OG	2.15	0.45
1:A:47:SER:H	1:A:50:GLN:HE21	1.64	0.45
1:A:187:GLU:HG3	1:C:306:LEU:HG	1.98	0.45
1:A:227:THR:HG22	1:A:228:THR:N	2.32	0.45
1:B:6:LEU:HD21	1:B:300:GLN:CD	2.41	0.45
1:C:263:VAL:HG22	2:C:322:HOH:O	2.16	0.45
1:A:116:TYR:CD2	1:A:121:ILE:HD12	2.49	0.45
1:C:252:ILE:HD11	1:C:294:PRO:CG	2.46	0.45
1:A:84:VAL:HG23	1:A:177:GLY:HA2	1.99	0.45
1:B:148:PHE:CE1	1:B:150:ILE:HD11	2.51	0.45
1:B:265:LYS:HD2	2:B:500:HOH:O	2.16	0.45
1:C:166:PRO:O	1:C:167:ASN:HB2	2.16	0.45
1:C:184:VAL:HG23	1:C:185:ASP:N	2.29	0.45
1:A:206:LEU:O	1:A:209:ALA:HB3	2.16	0.44
1:A:116:TYR:CE2	1:A:141:GLY:HA2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:ILE:HD13	1:B:261:VAL:HG21	1.99	0.44
1:C:234:TYR:CD1	1:C:234:TYR:C	2.96	0.44
1:B:31:TRP:CD2	1:B:93:ASN:HB2	2.52	0.44
1:A:6:LEU:CD2	1:B:124:TYR:HD2	2.30	0.44
1:B:166:PRO:HG2	1:B:190:GLN:OE1	2.16	0.44
1:B:282:ILE:N	1:B:282:ILE:HD12	2.32	0.44
1:C:206:LEU:O	1:C:210:ILE:HG12	2.17	0.44
1:C:268:ARG:HG3	1:C:268:ARG:NH1	2.33	0.44
1:C:91:VAL:HG22	1:C:92:ALA:N	2.32	0.44
1:A:226:SER:O	1:A:226:SER:OG	2.29	0.44
1:B:248:THR:HG22	2:B:475:HOH:O	2.12	0.44
1:B:263:VAL:HA	1:B:266:LEU:HD13	1.99	0.44
1:A:237:TRP:CA	2:A:470:HOH:O	2.65	0.44
1:B:185:ASP:OD2	1:B:185:ASP:N	2.47	0.44
1:A:22:TYR:CG	1:A:23:ARG:N	2.82	0.43
1:C:194:PRO:CG	2:C:484:HOH:O	2.38	0.43
1:A:47:SER:N	1:A:50:GLN:HE21	2.16	0.43
1:C:110:PHE:HE1	1:C:112:ILE:HD11	1.84	0.43
1:C:200:ASN:HD21	1:C:247:SER:H	1.66	0.43
1:B:100:LYS:HD3	2:B:334:HOH:O	2.18	0.43
1:B:222:LYS:CG	2:B:481:HOH:O	2.61	0.43
1:B:255:LEU:HA	1:B:258:ILE:HB	2.01	0.43
1:C:106:CYS:SG	1:C:128:MET:HE2	2.59	0.43
1:A:281:PRO:HB3	1:A:286:TYR:CZ	2.54	0.43
1:A:234:TYR:CD1	1:A:234:TYR:C	2.97	0.43
1:B:138:PHE:HD2	1:B:170:HIS:CD2	2.36	0.43
1:B:273:LYS:HZ1	1:B:278:GLY:H	1.65	0.43
1:C:231:ILE:HD12	1:C:246:PHE:HD2	1.76	0.43
1:C:252:ILE:HD11	1:C:294:PRO:HG3	2.00	0.43
1:A:56:ASN:HA	1:A:78:ARG:HH21	1.83	0.43
1:B:70:GLY:O	2:B:464:HOH:O	2.21	0.43
1:B:301:VAL:O	1:B:304:VAL:HG13	2.18	0.43
1:B:116:TYR:CD2	1:B:121:ILE:HG12	2.54	0.43
1:B:203:VAL:HG12	1:B:252:ILE:HD12	2.00	0.43
1:B:1:SER:HB3	1:B:283:LEU:HD22	2.00	0.42
1:B:222:LYS:CB	2:B:481:HOH:O	2.67	0.42
1:C:283:LEU:HD12	1:C:283:LEU:HA	1.88	0.42
1:A:162:HIS:HB2	1:A:172:GLY:HA2	2.00	0.42
1:A:47:SER:N	1:A:50:GLN:NE2	2.68	0.42
1:A:49:ASP:CA	2:A:462:HOH:O	2.66	0.42
1:B:31:TRP:CE2	1:B:93:ASN:HB2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:TYR:HD2	1:B:121:ILE:HG12	1.84	0.42
1:B:124:TYR:OH	1:B:136:ALA:HB1	2.20	0.42
1:C:218:PHE:HD2	1:C:218:PHE:HA	1.63	0.42
1:C:277:TRP:CZ2	1:C:282:ILE:HD12	2.54	0.42
1:A:31:TRP:CE2	1:A:93:ASN:HB2	2.55	0.42
1:A:124:TYR:OH	1:A:136:ALA:HB1	2.20	0.42
1:B:202:ILE:HG21	1:B:270:ILE:CD1	2.44	0.42
1:B:165:LEU:HD22	1:B:190:GLN:HG2	2.01	0.42
1:C:23:ARG:HD3	1:C:23:ARG:H	1.81	0.42
1:C:211:ILE:O	1:C:215:GLU:CG	2.67	0.42
1:A:282:ILE:N	1:A:282:ILE:CD1	2.82	0.42
1:A:293:THR:O	1:A:297:VAL:HG23	2.19	0.42
1:B:111:THR:O	1:B:147:GLY:HA2	2.20	0.42
1:C:78:ARG:HG3	1:C:78:ARG:NH1	2.35	0.42
1:C:149:ASN:OD1	1:C:149:ASN:C	2.62	0.42
1:C:183:TYR:HA	2:C:432:HOH:O	2.19	0.42
1:C:231:ILE:CD1	1:C:267:LEU:HD11	2.50	0.42
1:B:91:VAL:HB	2:B:460:HOH:O	2.20	0.42
1:C:99:TYR:O	1:C:100:LYS:HG3	2.20	0.42
1:B:202:ILE:CD1	1:B:288:PHE:HB3	2.50	0.41
1:B:207:TYR:CG	1:B:255:LEU:HD13	2.55	0.41
1:A:293:THR:HB	1:A:294:PRO:HD2	2.02	0.41
1:A:81:LYS:HG3	1:A:176:MET:O	2.20	0.41
1:B:266:LEU:O	1:B:269:THR:HB	2.20	0.41
1:C:231:ILE:HD13	1:C:267:LEU:HD11	2.01	0.41
1:A:246:PHE:CZ	1:A:263:VAL:HG11	2.55	0.41
1:B:105:ASN:O	1:B:128:MET:HB3	2.20	0.41
1:A:2:GLY:O	1:B:137:SER:HB2	2.20	0.41
1:A:187:GLU:HG2	1:C:305:ARG:HA	2.03	0.41
1:B:246:PHE:CZ	1:B:263:VAL:HG11	2.56	0.41
1:A:68:GLN:HG3	1:A:118:GLY:O	2.20	0.41
1:B:202:ILE:CG2	1:B:270:ILE:CD1	2.98	0.41
1:C:53:ASP:O	1:C:57:LEU:HG	2.21	0.41
1:C:112:ILE:HD13	1:C:147:GLY:HA3	2.02	0.41
1:A:191:ARG:O	1:A:193:PRO:HD3	2.21	0.41
1:B:81:LYS:HE2	1:B:176:MET:O	2.21	0.41
1:B:183:TYR:OH	1:B:193:PRO:HD2	2.20	0.41
1:B:235:ASN:HB2	2:B:482:HOH:O	2.20	0.41
1:C:46:PHE:HB2	1:C:51:TRP:CH2	2.56	0.41
1:C:154:VAL:CG2	2:C:499:HOH:O	2.68	0.41
1:B:166:PRO:O	1:B:167:ASN:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:PRO:O	1:C:42:VAL:HG22	2.21	0.40
1:C:47:SER:O	1:C:50:GLN:NE2	2.49	0.40
1:B:166:PRO:HD2	1:B:190:GLN:CG	2.46	0.40
1:B:263:VAL:O	1:B:267:LEU:HG	2.20	0.40
1:C:36:ILE:HD13	1:C:65:VAL:HG11	2.03	0.40
1:C:173:THR:HA	1:C:178:GLU:O	2.21	0.40
1:B:205:TRP:CE2	1:B:289:GLU:HB2	2.56	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	292/309 (94%)	278 (95%)	12 (4%)	2 (1%)	18	20
1	B	305/309 (99%)	282 (92%)	20 (7%)	3 (1%)	12	12
1	C	300/309 (97%)	285 (95%)	15 (5%)	0	100	100
All	All	897/927 (97%)	845 (94%)	47 (5%)	5 (1%)	21	24

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	48	GLY
1	B	182	GLY
1	B	193	PRO
1	A	261	VAL
1	B	278	GLY

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	250/260 (96%)	235 (94%)	15 (6%)	17	21
1	B	259/260 (100%)	247 (95%)	12 (5%)	24	31
1	C	255/260 (98%)	235 (92%)	20 (8%)	11	12
All	All	764/780 (98%)	717 (94%)	47 (6%)	16	19

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	ARG
1	A	79	ARG
1	A	115	SER
1	A	121	ILE
1	A	134	ILE
1	A	191	ARG
1	A	219	SER
1	A	226	SER
1	A	233	ASP
1	A	240	ASP
1	A	252	ILE
1	A	255	LEU
1	A	282	ILE
1	A	285	GLN
1	A	289	GLU
1	B	6	LEU
1	B	78	ARG
1	B	171	THR
1	B	175	LEU
1	B	192	VAL
1	B	254	LYS
1	B	255	LEU
1	B	258	ILE
1	B	268	ARG
1	B	273	LYS
1	B	289	GLU

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Mol	Chain	Res	Type
1	B	301	VAL
1	C	21	SER
1	C	22	TYR
1	C	23	ARG
1	C	30	LEU
1	C	42	VAL
1	C	46	PHE
1	C	47	SER
1	C	49	ASP
1	C	50	GLN
1	C	119	THR
1	C	165	LEU
1	C	192	VAL
1	C	215	GLU
1	C	216	SER
1	C	218	PHE
1	C	227	THR
1	C	236	ARG
1	C	263	VAL
1	C	283	LEU
1	C	292	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	A	50	GLN
1	A	61	HIS
1	A	74	ASN
1	A	105	ASN
1	A	167	ASN
1	A	196	ASN
1	A	285	GLN
1	A	299	ASN
1	A	300	GLN
1	B	68	GLN
1	B	170	HIS
1	B	196	ASN
1	B	200	ASN
1	B	220	GLN
1	B	235	ASN
1	B	241	ASN

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Mol	Chain	Res	Type
1	C	41	HIS
1	C	50	GLN
1	C	88	GLN
1	C	161	HIS
1	C	170	HIS
1	C	190	GLN
1	C	196	ASN
1	C	200	ASN
1	C	285	GLN
1	C	307	GLN

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	49:ASP	C	50:GLN	N	1.12

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	296/309 (95%)	0.41	17 (5%) 29 34	21, 46, 100, 129	0
1	B	307/309 (99%)	0.54	30 (9%) 13 15	23, 49, 87, 115	0
1	C	302/309 (97%)	0.22	10 (3%) 49 56	21, 42, 90, 112	0
All	All	905/927 (97%)	0.39	57 (6%) 26 29	21, 45, 93, 129	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	189	ALA	4.1
1	C	46	PHE	4.1
1	B	192	VAL	3.9
1	B	191	ARG	3.5
1	B	263	VAL	3.4
1	C	213	VAL	3.3
1	B	166	PRO	3.3
1	B	229	VAL	3.2
1	C	6	LEU	3.1
1	A	266	LEU	3.0
1	B	231	ILE	3.0
1	C	54	VAL	2.9
1	B	282	ILE	2.8
1	B	183	TYR	2.8
1	B	306	LEU	2.8
1	A	283	LEU	2.7
1	A	272	VAL	2.7
1	A	2	GLY	2.7
1	A	270	ILE	2.6
1	B	290	ASP	2.6
1	B	184	VAL	2.6
1	B	255	LEU	2.6
1	A	238	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	288	PHE	2.5
1	B	168	ALA	2.5
1	B	277	TRP	2.4
1	B	190	GLN	2.4
1	B	182	GLY	2.4
1	B	246	PHE	2.4
1	B	237	TRP	2.4
1	B	188	VAL	2.3
1	A	252	ILE	2.3
1	B	278	GLY	2.3
1	A	3	PHE	2.3
1	C	218	PHE	2.3
1	B	165	LEU	2.3
1	B	107	GLY	2.3
1	B	234	TYR	2.3
1	A	237	TRP	2.3
1	A	229	VAL	2.3
1	B	261	VAL	2.3
1	A	1	SER	2.2
1	C	75	VAL	2.2
1	A	204	ALA	2.2
1	B	167	ASN	2.2
1	A	228	THR	2.2
1	C	216	SER	2.2
1	A	150	ILE	2.2
1	A	282	ILE	2.1
1	C	215	GLU	2.1
1	B	243	PHE	2.1
1	A	247	SER	2.1
1	C	73	LEU	2.1
1	B	193	PRO	2.1
1	B	194	PRO	2.1
1	C	87	LEU	2.0
1	B	150	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.