



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 02:57 pm BST

PDB ID : 2Q6D
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

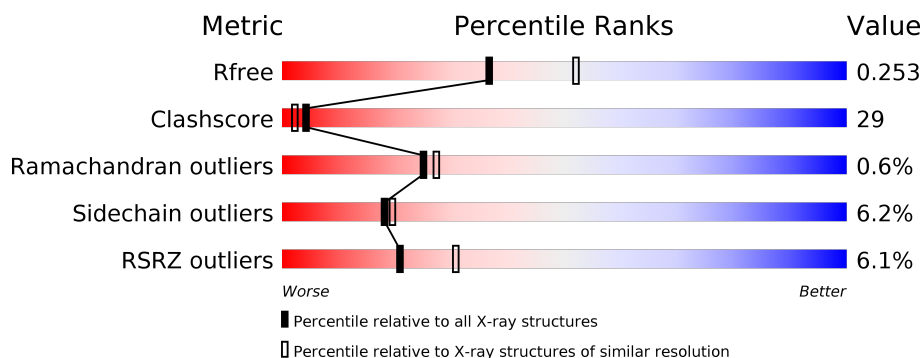
1 Overall quality at a glance

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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (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 on 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

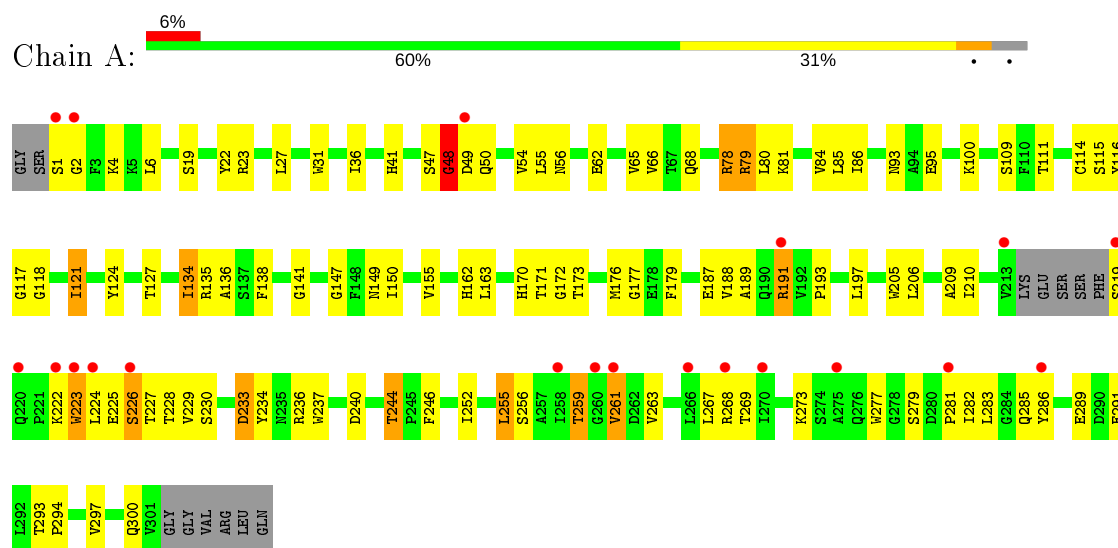
- Molecule 2 is water.

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		

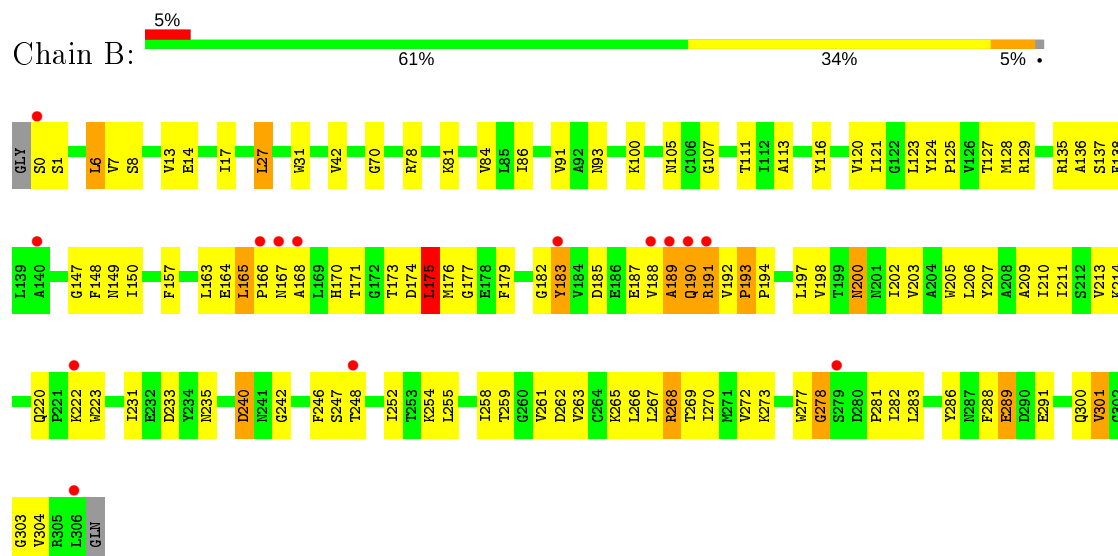
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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

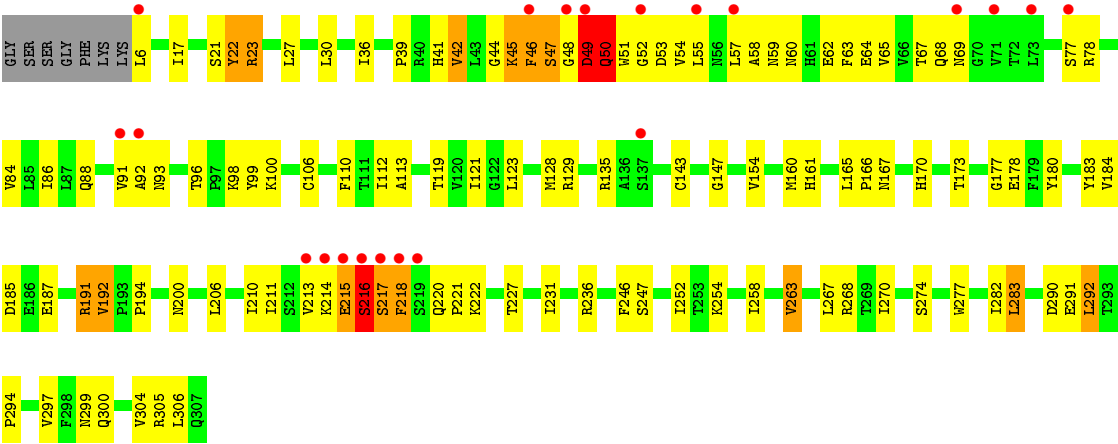


- Molecule 1: Infectious bronchitis virus (IBV) main protease



- 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 40.04 – 2.35	Depositor EDS
% Data completeness (in resolution range)	95.6 (50.00-2.35) 95.7 (40.04-2.35)	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 , 56.2	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	0.50	1/2325 (0.0%)	0.72	2/3165 (0.1%)
1	B	0.43	1/2408 (0.0%)	0.80	6/3276 (0.2%)
1	C	0.51	1/2372 (0.0%)	0.79	5/3229 (0.2%)
All	All	0.48	3/7105 (0.0%)	0.77	13/9670 (0.1%)

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 (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	49	ASP	C-N	-9.53	1.12	1.34
1	A	244	THR	C-N	-6.36	1.22	1.34
1	B	183	TYR	C-N	5.79	1.47	1.34

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	189	ALA	CB-CA-C	-17.41	83.98	110.10
1	C	49	ASP	O-C-N	-9.41	107.64	122.70
1	B	190	GLN	N-CA-CB	-7.58	96.96	110.60
1	C	6	LEU	CA-C-N	-6.97	101.86	117.20
1	C	50	GLN	N-CA-CB	6.74	122.73	110.60
1	C	49	ASP	CB-CA-C	-6.28	97.83	110.40
1	B	175	LEU	CA-CB-CG	-6.28	100.86	115.30
1	A	223	TRP	N-CA-CB	-6.12	99.59	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	259	THR	C-N-CA	-5.99	109.72	122.30
1	B	191	ARG	N-CA-C	-5.61	95.85	111.00
1	C	45	LYS	CB-CA-C	5.39	121.19	110.40
1	B	165	LEU	N-CA-C	-5.38	96.47	111.00
1	B	174	ASP	N-CA-C	-5.14	97.11	111.00

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	115	0
1	B	2355	0	2317	145	0
1	C	2320	0	2273	153	0
2	A	167	0	0	21	0
2	B	194	0	0	34	0
2	C	205	0	0	36	0
All	All	7515	0	6824	405	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 29.

All (405) 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:C:48:GLY:CA	1:C:49:ASP:HB2	1.65	1.26
1:B:127:THR:HB	2:B:483:HOH:O	1.10	1.25

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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:166:PRO:HD2	1:B:190:GLN:CD	1.71	1.09
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: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:B:166:PRO:CD	1:B:190:GLN:CD	2.34	0.95
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:HH11	1:C:23:ARG:HG3	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: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:47:SER:C	1:C:50:GLN:HE21	1.79	0.85
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:262:ASP:HB3	2:B:345:HOH:O	1.78	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:166:PRO:HG2	1:B:190:GLN:CD	2.03	0.79
1:C:84:VAL:HG23	1:C:177:GLY:HA2	1.63	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: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: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:166:PRO:CG	1:B:190:GLN:CD	2.56	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:A:237:TRP:C	2:A:470:HOH:O	2.27	0.71
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:C:47:SER:C	1:C:50:GLN:HG3	2.10	0.71
1:B:166:PRO:CD	1:B:190:GLN:OE1	2.38	0.71
1:C:23:ARG:N	1:C:23:ARG:HD3	2.05	0.71
1:C:64:GLU:HG3	2:C:505:HOH:O	1.90	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:231:ILE:HD13	1:C:267:LEU:CD1	2.23	0.68
1:C:46:PHE:CE1	1:C:54:VAL:CG1	2.73	0.68
1:B:0:SER:CB	2:B:488:HOH:O	2.40	0.68
1:C:48:GLY:CA	1:C:49:ASP:CB	2.46	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: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:A:47:SER:O	1:A:48:GLY:C	2.34	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:197:LEU:O	1:A:244:THR:HG23	1.97	0.64
1:B:183:TYR:CD2	1:B:183:TYR:N	2.64	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:47:SER:C	1:C:50:GLN:NE2	2.52	0.63
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
1:C:165:LEU:CD1	1:C:192:VAL:HG13	2.29	0.63
1:C:55:LEU:CD1	2:C:502:HOH:O	2.10	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:22:TYR:C	1:C:22:TYR:CD2	2.72	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:1:SER:N	2:A:425:HOH:O	2.21	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:A:138:PHE:O	1:B:303:GLY:HA3	2.01	0.60
1:B:187:GLU:O	1:B:187:GLU:HG2	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:281:PRO:C	1:A:282:ILE:HD12	2.22	0.60
1:A:19:SER:OG	1:A:66:VAL:HB	2.01	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: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: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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:ARG:NH1	1:C:23:ARG:HG3	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:214:LYS:C	1:C:217:SER:H	2.07	0.56
1:C:50:GLN:O	1:C:54:VAL:HG23	2.07	0.55
1:C:213:VAL:O	1:C:213:VAL:HG12	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:B:107:GLY:C	2:B:497:HOH:O	2.45	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:44:GLY:O	1:C:45:LYS:C	2.43	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: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:78:ARG:C	1:A:79:ARG:HG3	2.27	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:55:LEU:HD11	1:A:78:ARG:HG3	1.90	0.53
1:A:36:ILE:CD1	1:A:65:VAL:HG11	2.38	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:NH1	1:C:23:ARG:CG	2.59	0.53
1:B:214:LYS:NZ	1:B:259:THR:O	2.42	0.53
1:A:163:LEU:O	1:A:163:LEU:HD12	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:175:LEU:HD12	2:B:493:HOH:O	1.94	0.52
1:B:166:PRO:CG	1:B:190:GLN:OE1	2.56	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:106:CYS:SG	1:C:128:MET:HE2	2.51	0.51
1:C:59:ASN:HB3	1:C:62:GLU:HG3	1.92	0.51
1:B:17:ILE:HD13	1:B:120:VAL:HG21	1.93	0.51
1:B:7:VAL:CG2	1:B:125:PRO:HD3	2.40	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:A:226:SER:O	1:A:226:SER:OG	2.29	0.51
1:B:261:VAL:HG12	1:B:262:ASP:N	2.26	0.51
1:A:86:ILE:HD12	1:A:86:ILE:N	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:135:ARG:HD2	2:C:483:HOH:O	2.11	0.50
1:C:47:SER:CA	1:C:50:GLN:NE2	2.71	0.50
1:A:149:ASN:C	1:A:149:ASN:OD1	2.50	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
1:A:100:LYS:HE3	2:A:464:HOH:O	2.12	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:214:LYS:HE2	1:B:220:GLN:NE2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:161:HIS:CE1	1:C:170:HIS:CD2	3.00	0.49
1:C:86:ILE:HD12	1:C:86:ILE:N	2.27	0.49
1:C:268:ARG:HG3	1:C:268:ARG:HH11	1.76	0.49
1:A:49:ASP:O	1:A:49:ASP:CG	2.50	0.49
1:C:23:ARG:CD	1:C:23:ARG:N	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:234:TYR:CD2	1:A:267:LEU:HB3	2.48	0.48
1:A:223:TRP:CE3	1:A:269:THR:HG21	2.48	0.48
1:A:41:HIS:CD2	2:A:453:HOH:O	2.64	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:HD2	1:B:183:TYR:N	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:143:CYS:HA	1:C:161:HIS:CD2	2.46	0.48
1:C:88:GLN:HB2	2:C:496:HOH:O	2.14	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:22:TYR:CG	1:A:23:ARG:N	2.82	0.47
1:A:281:PRO:HB3	1:A:286:TYR:CE2	2.49	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:HD12	1:B:86:ILE:N	2.29	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:C:270:ILE:O	1:C:274:SER:HB2	2.16	0.46
1:C:99:TYR:N	2:C:495:HOH:O	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:ASN:HD22	1:B:200:ASN:N	2.13	0.46
1:C:129:ARG:HH22	1:C:290:ASP:CG	2.18	0.46
1:C:112:ILE:HD13	1:C:147:GLY:HA2	1.97	0.46
1:B:223:TRP:CE2	1:B:282:ILE:HG13	2.51	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:C:55:LEU:CD2	2:C:487:HOH:O	2.60	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: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:A:85:LEU:C	1:A:86:ILE:HD12	2.37	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
1:B:210:ILE:HD13	1:B:261:VAL:HG21	1.99	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: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:N	1:B:185:ASP:OD2	2.47	0.44
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:B:149:ASN:C	1:B:150:ILE:HD12	2.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:281:PRO:HB3	1:A:286:TYR:CZ	2.54	0.43
1:B:138:PHE:HD2	1:B:170:HIS:CD2	2.36	0.43
1:C:231:ILE:HD12	1:C:246:PHE:HD2	1.76	0.43
1:B:165:LEU:O	1:B:166:PRO:C	2.57	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:301:VAL:O	1:B:304:VAL:HG13	2.18	0.43
1:B:70:GLY:O	2:B:464:HOH:O	2.21	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:HA	1:C:283:LEU:HD12	1.88	0.42
1:A:162:HIS:HB2	1:A:172:GLY:HA2	2.00	0.42
1:A:49:ASP:CA	2:A:462:HOH:O	2.66	0.42
1:A:47:SER:N	1:A:50:GLN:NE2	2.68	0.42
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:B:31:TRP:CE2	1:B:93:ASN:HB2	2.55	0.42
1:C:218:PHE:HA	1:C:218:PHE:HD2	1.63	0.42
1:C:277:TRP:CZ2	1:C:282:ILE:HD12	2.54	0.42
1:A:124:TYR:OH	1:A:136:ALA:HB1	2.20	0.42
1:A:31:TRP:CE2	1:A:93:ASN:HB2	2.55	0.42
1:B:121:ILE:HD12	1:B:121:ILE:C	2.40	0.42
1:B:149:ASN:C	1:B:149:ASN:OD1	2.58	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:211:ILE:O	1:C:215:GLU:CG	2.67	0.42
1:C:23:ARG:H	1:C:23:ARG:HD3	1.81	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: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:C:78:ARG:HG3	1:C:78:ARG:NH1	2.35	0.42
1:A:163:LEU:C	1:A:163:LEU:HD12	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:128:MET:HE3	1:C:180:TYR:HB3	2.01	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:A:223:TRP:O	1:A:224:LEU:C	2.56	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:112:ILE:HD13	1:C:147:GLY:HA3	2.02	0.41
1:C:53:ASP:O	1:C:57:LEU:HG	2.21	0.41
1:A:191:ARG:O	1:A:193:PRO:HD3	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:B:81:LYS:HE2	1:B:176:MET:O	2.21	0.41
1:C:154:VAL:CG2	2:C:499:HOH:O	2.68	0.41
1:C:46:PHE:HB2	1:C:51:TRP:CH2	2.56	0.41
1:B:166:PRO:O	1:B:167:ASN:HB2	2.20	0.41
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:B:6:LEU:HD21	1:B:300:GLN:CD	2.41	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%)	22	23
1	B	305/309 (99%)	282 (92%)	20 (7%)	3 (1%)	15	15
1	C	300/309 (97%)	285 (95%)	15 (5%)	0	100	100
All	All	897/927 (97%)	845 (94%)	47 (5%)	5 (1%)	25	27

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 [i](#)

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%)	236 (94%)	14 (6%)	21	23
1	B	259/260 (100%)	246 (95%)	13 (5%)	24	28
1	C	255/260 (98%)	235 (92%)	20 (8%)	12	12
All	All	764/780 (98%)	717 (94%)	47 (6%)	18	19

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

Mol	Chain	Res	Type
1	A	27	LEU
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	255	LEU
1	A	285	GLN
1	A	289	GLU
1	B	6	LEU
1	B	27	LEU
1	B	78	ARG
1	B	171	THR
1	B	175	LEU
1	B	192	VAL
1	B	200	ASN
1	B	233	ASP
1	B	240	ASP
1	B	254	LYS
1	B	268	ARG
1	B	289	GLU
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	191	ARG
1	C	192	VAL
1	C	215	GLU
1	C	216	SER
1	C	218	PHE
1	C	227	THR

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Mol	Chain	Res	Type
1	C	236	ARG
1	C	263	VAL
1	C	283	LEU
1	C	292	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	A	50	GLN
1	A	88	GLN
1	A	105	ASN
1	A	196	ASN
1	A	299	ASN
1	A	300	GLN
1	B	200	ASN
1	B	220	GLN
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	307	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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](#)

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%)	0.20	20 (6%)	17 25	21, 46, 100, 129	0
1	B	307/309 (99%)	0.26	14 (4%)	32 45	23, 49, 87, 115	0
1	C	302/309 (97%)	0.10	21 (6%)	16 24	21, 42, 90, 112	0
All	All	905/927 (97%)	0.19	55 (6%)	21 31	21, 45, 93, 129	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	218	PHE	7.1
1	A	220	GLN	5.5
1	B	191	ARG	5.3
1	B	166	PRO	5.3
1	A	261	VAL	5.2
1	B	189	ALA	4.7
1	A	258	ILE	4.6
1	B	190	GLN	4.4
1	C	6	LEU	4.1
1	C	217	SER	4.0
1	B	140	ALA	3.9
1	A	266	LEU	3.8
1	C	55	LEU	3.6
1	C	215	GLU	3.6
1	A	275	ALA	3.5
1	C	46	PHE	3.5
1	B	306	LEU	3.5
1	B	279	SER	3.5
1	C	219	SER	3.4
1	A	281	PRO	3.4
1	B	167	ASN	3.3
1	A	213	VAL	3.2
1	B	188	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	216	SER	3.2
1	A	260	GLY	3.1
1	C	69	ASN	3.1
1	C	91	VAL	2.9
1	A	223	TRP	2.8
1	C	49	ASP	2.8
1	A	224	LEU	2.8
1	A	222	LYS	2.8
1	A	219	SER	2.8
1	A	191	ARG	2.6
1	A	268	ARG	2.5
1	C	48	GLY	2.5
1	C	213	VAL	2.5
1	C	73	LEU	2.5
1	B	168	ALA	2.4
1	C	52	GLY	2.4
1	A	286	TYR	2.4
1	C	92	ALA	2.3
1	B	248	THR	2.3
1	A	270	ILE	2.3
1	A	226	SER	2.3
1	A	2	GLY	2.3
1	A	1	SER	2.2
1	B	183	TYR	2.2
1	B	222	LYS	2.1
1	C	71	VAL	2.1
1	C	77	SER	2.1
1	A	49	ASP	2.1
1	C	137	SER	2.1
1	C	57	LEU	2.0
1	B	0	SER	2.0
1	C	214	LYS	2.0

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.