



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 04:45 PM GMT

PDB ID : 1U4C
Title : Structure of spindle checkpoint protein Bub3
Authors : Larsen, N.A.; Harrison, S.C.
Deposited on : 2004-07-23
Resolution : 2.35 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

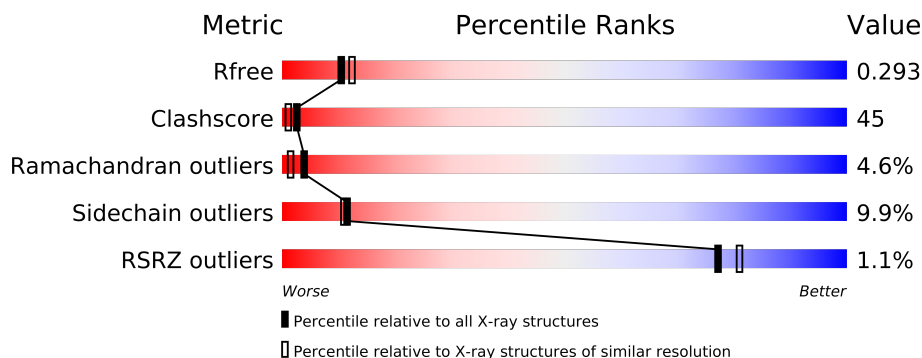
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	3327 (2.40-2.32)
Clashscore	79885	1064 (2.38-2.34)
Ramachandran outliers	78287	1048 (2.38-2.34)
Sidechain outliers	78261	1049 (2.38-2.34)
RSRZ outliers	66119	3330 (2.40-2.32)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	349	
1	B	349	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5511 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Cell cycle arrest protein BUB3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	Se	0	0	0
			2547	1619	428	487	10	3			
1	B	331	Total	C	N	O	S	Se	0	0	0
			2620	1662	439	506	10	3			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP P26449
A	156	MSE	MET	MODIFIED RESIDUE	UNP P26449
A	167	MSE	MET	MODIFIED RESIDUE	UNP P26449
A	342	LEU	-	EXPRESSION TAG	UNP P26449
A	343	GLU	-	EXPRESSION TAG	UNP P26449
A	344	HIS	-	EXPRESSION TAG	UNP P26449
A	345	HIS	-	EXPRESSION TAG	UNP P26449
A	346	HIS	-	EXPRESSION TAG	UNP P26449
A	347	HIS	-	EXPRESSION TAG	UNP P26449
A	348	HIS	-	EXPRESSION TAG	UNP P26449
A	349	HIS	-	EXPRESSION TAG	UNP P26449
B	1	MSE	MET	MODIFIED RESIDUE	UNP P26449
B	156	MSE	MET	MODIFIED RESIDUE	UNP P26449
B	167	MSE	MET	MODIFIED RESIDUE	UNP P26449
B	342	LEU	-	EXPRESSION TAG	UNP P26449
B	343	GLU	-	EXPRESSION TAG	UNP P26449
B	344	HIS	-	EXPRESSION TAG	UNP P26449
B	345	HIS	-	EXPRESSION TAG	UNP P26449
B	346	HIS	-	EXPRESSION TAG	UNP P26449
B	347	HIS	-	EXPRESSION TAG	UNP P26449
B	348	HIS	-	EXPRESSION TAG	UNP P26449
B	349	HIS	-	EXPRESSION TAG	UNP P26449

- Molecule 2 is water.

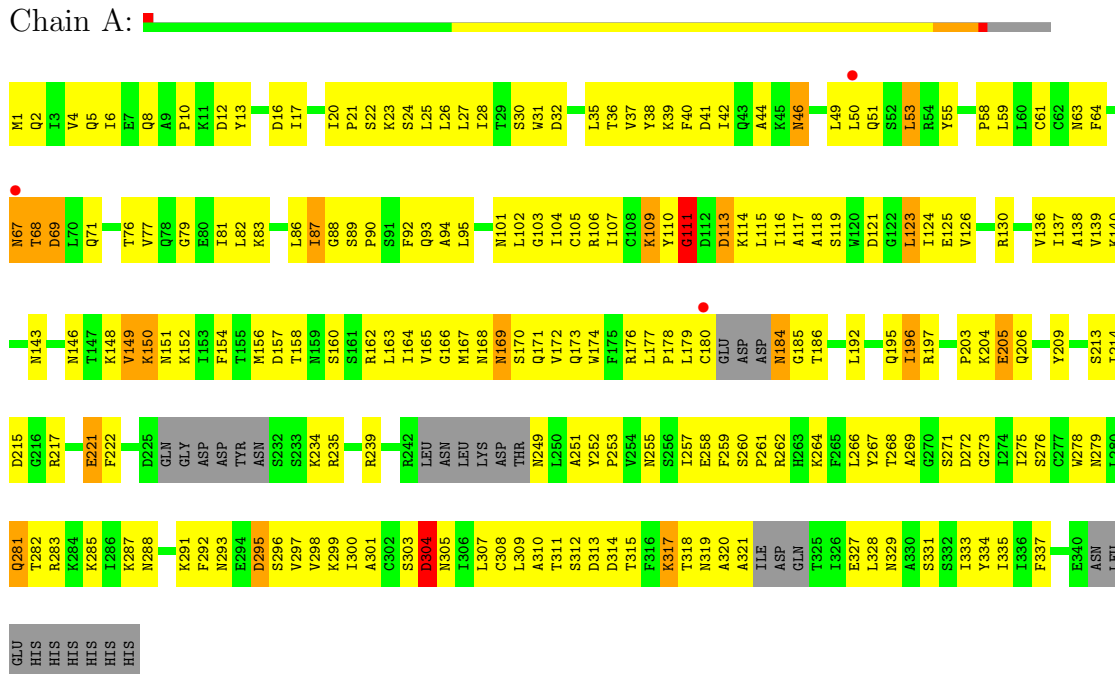
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	167	Total 167	O 167	0	0
2	B	177	Total 177	O 177	0	0

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 errors 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.

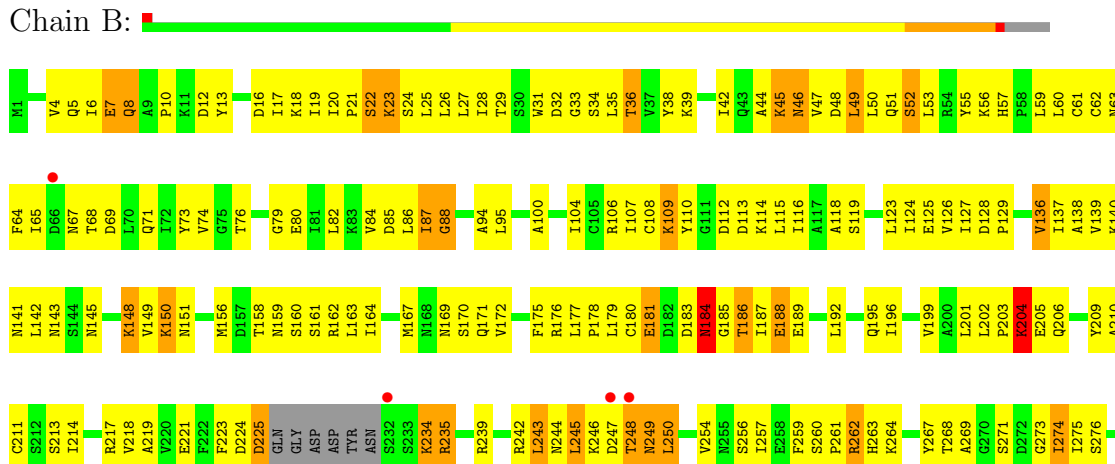
• Molecule 1: Cell cycle arrest protein BUB3

Chain A:



• Molecule 1: Cell cycle arrest protein BUB3

Chain B:



L280	Q281	T282	R283	K284	N288	K291	F292	N293	E294	D295	S296	V297	V298	K299	D304	N305	I306	L307	C308	L309	A310	T311	S312	D313	D314	T315	F316	K317	T318	A321	ILE	ASP	GLN	T325	I326	E327	L328	S331	S332	I333	Y334	E340	ASN	LEU	GLU	HIS	HIS	HIS	HIS	HIS
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	73.27Å 73.27Å 110.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 2.35 41.65 – 2.35	Depositor EDS
% Data completeness (in resolution range)	89.2 (40.00-2.35) 94.5 (41.65-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 2.34Å)	Xtriage
Refinement program	SHELXL-97, CNS	Depositor
R, R_{free}	0.196 , 0.288 0.206 , 0.293	Depositor DCC
R_{free} test set	1074 reflections (4.31%)	DCC
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 77.8	EDS
Estimated twinning fraction	0.016 for -h,-k,l 0.487 for h,-h-k,-l 0.014 for -k,-h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 51781 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5511	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.44	0/2587	0.70	1/3492 (0.0%)
1	B	0.44	0/2662	0.71	1/3597 (0.0%)
All	All	0.44	0/5249	0.70	2/7089 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	111	GLY	N-CA-C	-5.44	99.50	113.10
1	B	254	VAL	N-CA-C	-5.24	96.86	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2547	0	2546	220	0
1	B	2620	0	2614	242	0
2	A	167	0	0	12	0
2	B	177	0	0	14	0
All	All	5511	0	5160	461	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 45.

All (461) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:148:LYS:HG2	1:B:149:VAL:H	1.12	1.13
1:A:271:SER:HA	1:A:296:SER:HB2	1.38	1.06
1:A:148:LYS:HG2	1:A:149:VAL:H	1.14	1.04
1:B:28:ILE:HD13	1:B:333:ILE:HD13	1.46	0.98
1:B:18:LYS:HE2	1:B:18:LYS:HA	1.47	0.95
1:B:149:VAL:HG23	1:B:150:LYS:H	1.32	0.94
1:A:149:VAL:HG13	1:A:150:LYS:H	1.32	0.94
1:B:142:LEU:HD13	1:B:167:MSE:HE3	1.48	0.93
1:B:326:ILE:HD13	1:B:326:ILE:H	1.32	0.93
1:B:164:ILE:HD11	1:B:172:VAL:HG11	1.50	0.93
1:B:243:LEU:H	1:B:243:LEU:HD23	1.34	0.92
1:A:217:ARG:HG2	1:A:217:ARG:HH11	1.35	0.92
1:A:109:LYS:H	1:A:109:LYS:HE3	1.36	0.89
1:B:51:GLN:HE21	1:B:86:LEU:HB3	1.37	0.88
1:A:2:GLN:HA	2:A:375:HOH:O	1.74	0.86
1:B:260:SER:HB3	1:B:267:TYR:HE1	1.40	0.85
1:B:325:THR:HB	1:B:326:ILE:HD13	1.59	0.84
1:A:12:ASP:OD1	1:A:32:ASP:HB3	1.78	0.84
1:B:148:LYS:HG2	1:B:149:VAL:N	1.93	0.84
1:A:315:THR:HB	1:A:328:LEU:HD22	1.57	0.83
1:B:87:ILE:HG12	1:B:88:GLY:H	1.43	0.83
1:B:67:ASN:HD22	1:B:113:ASP:HB2	1.43	0.83
1:A:46:ASN:HD22	1:A:46:ASN:C	1.82	0.82
1:A:164:ILE:HD11	1:A:172:VAL:HG11	1.61	0.82
1:A:148:LYS:HG2	1:A:149:VAL:N	1.95	0.81
1:A:152:LYS:HD2	1:A:154:PHE:HE2	1.47	0.80
1:B:139:VAL:HG12	1:B:140:LYS:HG2	1.63	0.79
1:B:184:ASN:ND2	1:B:185:GLY:H	1.81	0.79
1:A:110:TYR:HB2	1:A:116:ILE:HG13	1.64	0.79
1:A:63:ASN:ND2	1:A:115:LEU:HD22	1.97	0.78
1:B:87:ILE:HD13	1:B:87:ILE:H	1.49	0.78
1:B:20:ILE:HD12	1:B:25:LEU:HD12	1.66	0.78
1:B:10:PRO:HG3	1:B:36:THR:HG21	1.65	0.77
1:A:184:ASN:ND2	1:A:185:GLY:H	1.82	0.77
1:A:214:ILE:HA	1:A:253:PRO:HB3	1.65	0.76
1:B:12:ASP:OD1	1:B:32:ASP:HB3	1.84	0.76
1:B:61:CYS:SG	1:B:107:ILE:HG13	2.25	0.76
1:A:1:MSE:HE1	1:A:287:LYS:NZ	2.01	0.76
1:B:125:GLU:HG2	1:B:138:ALA:HA	1.66	0.75
1:B:187:ILE:HG22	1:B:188:GLU:N	2.02	0.75
1:A:295:ASP:HB3	1:A:312:SER:O	1.86	0.75
1:A:87:ILE:HG12	1:A:88:GLY:H	1.52	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:276:SER:HB3	1:A:288:ASN:ND2	2.03	0.74
1:B:260:SER:HB3	1:B:267:TYR:CE1	2.23	0.73
1:B:23:LYS:N	1:B:23:LYS:HD3	2.03	0.73
1:A:152:LYS:HD2	1:A:154:PHE:CE2	2.22	0.73
1:B:195:GLN:HB3	1:B:214:ILE:HD13	1.70	0.73
1:A:170:SER:HA	2:A:383:HOH:O	1.91	0.71
1:B:28:ILE:CD1	1:B:333:ILE:HD13	2.20	0.71
1:A:87:ILE:HD13	1:A:87:ILE:H	1.55	0.71
1:A:313:ASP:HB2	1:A:329:ASN:HB2	1.72	0.70
1:B:250:LEU:H	1:B:250:LEU:HD22	1.56	0.70
1:A:221:GLU:HG3	1:A:235:ARG:HD2	1.72	0.70
1:B:5:GLN:HB2	1:B:334:TYR:CE2	2.26	0.70
1:B:39:LYS:HB2	1:B:50:LEU:HD21	1.73	0.69
1:A:109:LYS:H	1:A:109:LYS:CE	2.04	0.69
1:B:167:MSE:HE2	2:B:392:HOH:O	1.91	0.69
1:B:246:LYS:HD3	1:B:247:ASP:H	1.57	0.68
1:B:295:ASP:HB3	1:B:312:SER:O	1.94	0.68
1:A:192:LEU:HD22	1:A:213:SER:HB3	1.74	0.68
1:B:217:ARG:HH11	1:B:217:ARG:HG2	1.59	0.68
1:B:28:ILE:HD13	1:B:333:ILE:CD1	2.20	0.68
1:B:210:ALA:HA	1:B:219:ALA:O	1.94	0.68
1:A:81:ILE:HD12	1:A:136:VAL:HG21	1.75	0.68
1:B:148:LYS:CG	1:B:149:VAL:H	1.96	0.68
1:B:148:LYS:HE3	1:B:150:LYS:HG2	1.76	0.68
1:B:53:LEU:HD12	1:B:53:LEU:N	2.08	0.67
1:A:166:GLY:HA3	1:A:196:ILE:HD11	1.76	0.67
1:A:298:VAL:HG21	1:A:312:SER:HB2	1.75	0.67
1:A:101:ASN:O	1:A:102:LEU:HD23	1.95	0.67
1:A:281:GLN:H	1:A:281:GLN:NE2	1.93	0.67
1:B:150:LYS:NZ	1:B:150:LYS:HA	2.10	0.67
1:A:26:LEU:HD21	1:A:335:ILE:HD11	1.77	0.67
1:B:51:GLN:HG3	1:B:52:SER:N	2.10	0.67
1:A:1:MSE:HE1	1:A:287:LYS:HZ1	1.57	0.67
1:A:192:LEU:HD12	1:A:196:ILE:HG22	1.75	0.67
1:B:315:THR:HA	1:B:318:THR:HB	1.77	0.66
1:A:119:SER:OG	1:A:123:LEU:HB2	1.95	0.66
1:B:67:ASN:HD22	1:B:113:ASP:CB	2.09	0.66
1:A:53:LEU:HD12	1:A:53:LEU:H	1.61	0.66
1:A:17:ILE:HD11	1:A:310:ALA:CB	2.26	0.66
1:B:298:VAL:HG21	1:B:312:SER:HB2	1.76	0.66
1:B:51:GLN:HG3	1:B:52:SER:H	1.61	0.65
1:B:18:LYS:HB2	1:B:27:LEU:HB2	1.77	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:260:SER:HB3	1:A:267:TYR:HE1	1.62	0.65
1:B:87:ILE:HD13	1:B:87:ILE:N	2.10	0.65
1:B:326:ILE:CD1	1:B:326:ILE:H	2.00	0.65
1:B:292:PHE:CE2	1:B:309:LEU:HD21	2.31	0.65
1:A:257:ILE:HG22	1:A:268:THR:HG22	1.79	0.65
1:B:149:VAL:O	1:B:150:LYS:HB2	1.98	0.64
1:A:149:VAL:HG13	1:A:150:LYS:N	2.09	0.64
1:A:217:ARG:HG2	1:A:217:ARG:NH1	2.09	0.64
1:B:184:ASN:ND2	1:B:185:GLY:N	2.46	0.64
1:A:125:GLU:HB3	1:A:136:VAL:HG13	1.78	0.64
1:B:7:GLU:HA	1:B:7:GLU:OE1	1.97	0.64
1:A:160:SER:OG	1:A:204:LYS:HD3	1.98	0.64
1:B:326:ILE:N	1:B:326:ILE:HD13	2.09	0.64
1:B:184:ASN:CG	1:B:185:GLY:H	2.01	0.63
1:A:39:LYS:HB2	1:A:50:LEU:HD11	1.81	0.63
1:A:46:ASN:ND2	1:A:46:ASN:C	2.51	0.62
1:A:105:CYS:HB2	2:A:507:HOH:O	1.98	0.62
1:A:148:LYS:CG	1:A:149:VAL:H	1.95	0.62
1:A:81:ILE:CD1	1:A:136:VAL:HG21	2.30	0.62
1:B:143:ASN:OD1	1:B:151:ASN:HB2	1.99	0.61
1:A:16:ASP:OD1	1:A:299:LYS:HE3	1.99	0.61
1:A:20:ILE:HG12	1:A:64:PHE:CE2	2.35	0.61
1:B:21:PRO:HG2	1:B:22:SER:H	1.66	0.61
1:A:28:ILE:HD13	1:A:333:ILE:HD13	1.82	0.61
1:A:49:LEU:HD23	1:A:50:LEU:N	2.15	0.61
1:A:163:LEU:HD23	1:A:164:ILE:N	2.16	0.61
1:A:125:GLU:HB3	1:A:136:VAL:CG1	2.31	0.61
1:B:256:SER:H	1:B:269:ALA:HB3	1.66	0.61
1:B:39:LYS:HB2	1:B:50:LEU:CD2	2.30	0.61
1:A:126:VAL:O	1:A:137:ILE:HG22	2.01	0.61
1:B:180:CYS:SG	1:B:181:GLU:N	2.73	0.60
1:B:42:ILE:HG12	2:B:374:HOH:O	2.00	0.60
1:A:260:SER:O	1:A:264:LYS:HA	2.01	0.60
1:B:192:LEU:HD22	1:B:213:SER:HB3	1.83	0.60
1:A:139:VAL:HG11	1:A:179:LEU:HD13	1.84	0.59
1:B:142:LEU:CD1	1:B:167:MSE:HE3	2.26	0.59
1:A:53:LEU:HD12	1:A:53:LEU:N	2.17	0.59
1:B:184:ASN:N	1:B:184:ASN:HD22	1.99	0.59
1:B:32:ASP:OD2	1:B:34:SER:N	2.34	0.59
1:B:100:ALA:HA	1:B:123:LEU:HD12	1.84	0.59
1:A:184:ASN:N	1:A:184:ASN:HD22	1.99	0.59
1:B:149:VAL:HG23	1:B:150:LYS:N	2.11	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:261:PRO:HA	1:B:264:LYS:NZ	2.17	0.59
1:A:279:ASN:HB3	1:A:282:THR:HG22	1.82	0.59
1:A:79:GLY:H	1:A:103:GLY:HA2	1.67	0.59
1:A:271:SER:HA	1:A:296:SER:CB	2.22	0.59
1:B:65:ILE:HB	1:B:71:GLN:HG3	1.85	0.59
1:A:55:TYR:CE2	1:A:82:LEU:HD11	2.38	0.59
1:A:217:ARG:NH1	1:A:239:ARG:CD	2.66	0.58
1:A:315:THR:HA	1:A:318:THR:HB	1.85	0.58
1:A:156:MSE:HG2	1:A:157:ASP:N	2.17	0.58
1:B:163:LEU:HB2	1:B:177:LEU:HD11	1.86	0.58
1:A:67:ASN:O	1:B:284:LYS:HE2	2.03	0.58
1:A:176:ARG:HH22	1:A:180:CYS:C	2.07	0.58
1:B:118:ALA:HA	1:B:123:LEU:O	2.04	0.58
1:B:187:ILE:CG2	1:B:188:GLU:N	2.67	0.58
1:B:307:LEU:C	1:B:307:LEU:HD23	2.23	0.58
1:B:246:LYS:CD	1:B:247:ASP:H	2.17	0.58
1:B:187:ILE:HG22	1:B:188:GLU:H	1.67	0.58
1:A:160:SER:HA	2:A:456:HOH:O	2.03	0.58
1:B:164:ILE:HD11	1:B:172:VAL:CG1	2.30	0.57
1:A:156:MSE:CG	1:A:157:ASP:N	2.67	0.57
1:B:26:LEU:HB3	1:B:38:TYR:HB2	1.86	0.57
1:B:267:TYR:HA	1:B:276:SER:O	2.03	0.57
1:A:276:SER:HB3	1:A:288:ASN:HD22	1.69	0.57
1:B:281:GLN:HE21	1:B:281:GLN:H	1.51	0.57
1:B:6:ILE:HD12	1:B:333:ILE:HG21	1.85	0.57
1:B:6:ILE:HG22	1:B:8:GLN:H	1.70	0.57
1:A:21:PRO:HG2	1:A:22:SER:H	1.69	0.57
1:A:217:ARG:HH12	1:A:239:ARG:CD	2.18	0.57
1:A:307:LEU:C	1:A:307:LEU:HD23	2.24	0.57
1:A:64:PHE:O	1:A:115:LEU:HD21	2.05	0.56
1:A:17:ILE:HD11	1:A:310:ALA:HB2	1.86	0.56
1:B:17:ILE:HD12	1:B:299:LYS:HB3	1.86	0.56
1:A:217:ARG:HH11	1:A:217:ARG:CG	2.06	0.56
1:A:63:ASN:HD22	1:A:115:LEU:HD13	1.71	0.56
1:B:257:ILE:HG22	1:B:268:THR:HG22	1.87	0.56
1:A:41:ASP:OD1	1:A:44:ALA:HB3	2.04	0.56
1:A:292:PHE:CE2	1:A:309:LEU:HD21	2.41	0.56
1:B:140:LYS:HE3	1:B:175:PHE:CD2	2.41	0.56
1:A:143:ASN:HB3	1:A:146:ASN:OD1	2.05	0.56
1:A:272:ASP:O	1:A:291:LYS:HE2	2.06	0.56
1:B:217:ARG:NH1	1:B:217:ARG:HG2	2.22	0.55
1:B:73:TYR:CE2	1:B:127:ILE:HG21	2.42	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:152:LYS:O	1:A:167:MSE:HA	2.06	0.55
1:B:142:LEU:HD13	1:B:167:MSE:CE	2.29	0.55
1:A:6:ILE:HD12	1:A:333:ILE:CG2	2.37	0.55
1:B:114:LYS:HD3	1:B:128:ASP:OD2	2.07	0.55
1:A:81:ILE:O	1:A:94:ALA:HA	2.06	0.55
1:A:279:ASN:CG	1:A:282:THR:HG22	2.27	0.55
1:A:67:ASN:OD1	1:A:68:THR:N	2.39	0.55
1:B:127:ILE:O	1:B:129:PRO:HD3	2.07	0.55
1:B:315:THR:HB	1:B:328:LEU:HD22	1.89	0.55
1:B:250:LEU:HD13	1:B:250:LEU:N	2.23	0.54
1:A:26:LEU:HD12	1:A:26:LEU:O	2.07	0.54
1:B:79:GLY:HA2	1:B:104:ILE:HG12	1.87	0.54
1:A:87:ILE:HD13	1:A:87:ILE:N	2.22	0.54
1:B:116:ILE:HD13	1:B:163:LEU:CD1	2.37	0.54
1:A:87:ILE:HG12	1:A:88:GLY:N	2.21	0.54
1:A:177:LEU:HA	1:A:178:PRO:C	2.27	0.54
1:B:51:GLN:HG2	1:B:86:LEU:HD22	1.89	0.54
1:B:244:ASN:HB2	2:B:516:HOH:O	2.08	0.54
1:B:17:ILE:HD11	1:B:310:ALA:HB2	1.90	0.54
1:B:6:ILE:HG12	1:B:47:VAL:HG11	1.90	0.54
1:B:195:GLN:O	1:B:214:ILE:HG23	2.08	0.54
1:A:63:ASN:HD22	1:A:115:LEU:HD22	1.73	0.54
1:A:261:PRO:HG2	1:A:304:ASP:N	2.23	0.54
1:A:164:ILE:CD1	1:A:172:VAL:HG11	2.35	0.53
1:A:143:ASN:OD1	1:A:151:ASN:HB2	2.08	0.53
1:B:6:ILE:HD12	1:B:333:ILE:CG2	2.39	0.53
1:A:184:ASN:ND2	1:A:185:GLY:N	2.54	0.53
1:B:244:ASN:C	1:B:246:LYS:H	2.12	0.53
1:A:30:SER:HB3	1:A:32:ASP:OD2	2.09	0.53
1:A:28:ILE:CD1	1:A:333:ILE:HD13	2.39	0.53
1:A:279:ASN:CB	1:A:282:THR:HG22	2.38	0.53
1:A:20:ILE:HD11	1:A:27:LEU:HD11	1.89	0.53
1:B:293:ASN:OD1	1:B:295:ASP:HB2	2.09	0.53
1:A:39:LYS:CB	1:A:50:LEU:HD11	2.38	0.53
1:A:217:ARG:NH1	1:A:217:ARG:CG	2.67	0.53
1:B:49:LEU:HD21	1:B:51:GLN:O	2.09	0.53
1:B:87:ILE:HG12	1:B:88:GLY:N	2.20	0.53
1:A:195:GLN:O	1:A:214:ILE:HG23	2.09	0.53
1:B:8:GLN:HA	1:B:8:GLN:OE1	2.09	0.52
1:B:187:ILE:CG2	1:B:188:GLU:H	2.21	0.52
1:B:139:VAL:CG1	1:B:140:LYS:HE2	2.40	0.52
1:A:149:VAL:CG1	1:A:150:LYS:H	2.15	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:161:SER:O	1:B:162:ARG:HG3	2.10	0.52
1:A:167:MSE:HE3	1:A:171:GLN:HE21	1.74	0.52
1:A:156:MSE:CG	1:A:157:ASP:H	2.22	0.52
1:A:269:ALA:HB1	1:A:297:VAL:HB	1.92	0.52
1:B:114:LYS:NZ	1:B:114:LYS:HB2	2.24	0.52
1:B:261:PRO:HA	1:B:264:LYS:HZ2	1.73	0.52
1:B:35:LEU:CD2	1:B:84:VAL:HG11	2.40	0.52
1:B:273:GLY:O	1:B:291:LYS:HG3	2.09	0.52
1:B:124:ILE:O	1:B:139:VAL:HB	2.10	0.52
1:B:248:THR:HG23	1:B:249:ASN:OD1	2.10	0.52
1:B:192:LEU:HD13	1:B:213:SER:HB3	1.91	0.52
1:A:63:ASN:HD21	1:A:115:LEU:HD22	1.71	0.52
1:B:145:ASN:HB2	2:B:377:HOH:O	2.09	0.51
1:A:320:ALA:O	1:A:321:ALA:HB2	2.10	0.51
1:B:48:ASP:O	1:B:50:LEU:HD12	2.10	0.51
1:B:5:GLN:HB2	1:B:334:TYR:CD2	2.44	0.51
1:A:285:LYS:NZ	1:A:288:ASN:HD21	2.08	0.51
1:B:186:THR:HA	2:B:367:HOH:O	2.11	0.51
1:A:114:LYS:O	1:A:115:LEU:HD23	2.11	0.51
1:A:203:PRO:HB2	1:A:205:GLU:OE2	2.11	0.51
1:B:68:THR:HG22	1:B:69:ASP:OD1	2.10	0.51
1:B:202:LEU:HD13	1:B:259:PHE:CZ	2.46	0.51
1:B:65:ILE:HB	1:B:71:GLN:CG	2.40	0.51
1:A:311:THR:O	1:A:331:SER:HB2	2.10	0.51
1:A:12:ASP:CG	1:A:32:ASP:HB3	2.29	0.51
1:B:239:ARG:NH1	1:B:243:LEU:HD11	2.25	0.50
1:A:140:LYS:NZ	1:A:184:ASN:HB3	2.27	0.50
1:B:20:ILE:HG12	1:B:64:PHE:CE2	2.47	0.50
1:A:95:LEU:HD12	1:A:136:VAL:HG23	1.94	0.50
1:A:149:VAL:HG23	1:A:152:LYS:HG2	1.92	0.50
1:B:82:LEU:CD2	1:B:94:ALA:HB2	2.42	0.50
1:A:273:GLY:O	1:A:291:LYS:HG3	2.11	0.50
1:A:169:ASN:HB2	1:A:171:GLN:HG2	1.93	0.50
1:A:168:ASN:O	1:A:170:SER:N	2.36	0.50
1:B:248:THR:O	1:B:249:ASN:O	2.29	0.50
1:B:189:GLU:HB2	2:B:423:HOH:O	2.12	0.50
1:B:184:ASN:ND2	1:B:184:ASN:N	2.59	0.50
1:B:281:GLN:H	1:B:281:GLN:NE2	2.10	0.49
1:A:61:CYS:SG	1:A:107:ILE:HG13	2.52	0.49
1:B:59:LEU:HD23	1:B:76:THR:HG22	1.94	0.49
1:B:50:LEU:N	1:B:50:LEU:HD12	2.27	0.49
1:B:262:ARG:HH21	1:B:340:GLU:HG2	1.78	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:31:TRP:O	1:A:58:PRO:HB3	2.13	0.49
1:B:52:SER:C	1:B:53:LEU:HD12	2.31	0.49
1:B:262:ARG:HB2	1:B:340:GLU:OE1	2.13	0.49
1:A:184:ASN:ND2	1:A:184:ASN:N	2.60	0.49
1:A:148:LYS:HG2	1:A:149:VAL:HG12	1.94	0.49
1:A:109:LYS:CE	1:A:109:LYS:N	2.74	0.49
1:B:309:LEU:HD12	1:B:311:THR:CG2	2.43	0.49
1:A:164:ILE:O	1:A:164:ILE:HG23	2.13	0.49
1:B:139:VAL:HG12	1:B:140:LYS:CG	2.40	0.49
1:B:23:LYS:H	1:B:23:LYS:HD3	1.78	0.49
1:B:309:LEU:HD12	1:B:311:THR:HG22	1.94	0.49
1:A:10:PRO:HD2	1:A:333:ILE:HD11	1.95	0.49
1:B:160:SER:OG	1:B:204:LYS:HE2	2.13	0.49
1:B:211:CYS:O	1:B:218:VAL:HA	2.14	0.48
1:B:239:ARG:HH11	1:B:243:LEU:HD21	1.78	0.48
1:A:278:TRP:CH2	1:A:285:LYS:HE3	2.48	0.48
1:A:259:PHE:CZ	1:A:266:LEU:HD13	2.48	0.48
1:A:197:ARG:HH11	1:A:197:ARG:HG3	1.78	0.48
1:A:116:ILE:HG12	1:A:126:VAL:HG22	1.95	0.48
1:B:20:ILE:HD12	1:B:25:LEU:CD1	2.41	0.48
1:A:205:GLU:CD	1:A:205:GLU:H	2.16	0.48
1:B:48:ASP:O	1:B:50:LEU:CD1	2.61	0.48
1:B:22:SER:HB2	1:B:23:LYS:HD3	1.94	0.48
1:B:110:TYR:O	1:B:114:LYS:HB2	2.14	0.48
1:B:39:LYS:CB	1:B:50:LEU:HD11	2.43	0.48
1:B:31:TRP:CZ2	1:B:60:LEU:HD21	2.49	0.48
1:A:124:ILE:HD11	1:A:165:VAL:HG11	1.96	0.48
1:B:177:LEU:HA	1:B:178:PRO:C	2.35	0.48
1:B:149:VAL:CG2	1:B:150:LYS:H	2.11	0.47
1:B:244:ASN:O	1:B:246:LYS:N	2.45	0.47
1:B:13:TYR:CE1	1:B:317:LYS:HD2	2.49	0.47
1:A:221:GLU:CG	1:A:235:ARG:HD2	2.42	0.47
1:A:69:ASP:HB2	2:A:424:HOH:O	2.13	0.47
1:B:112:ASP:CG	1:B:112:ASP:O	2.52	0.47
1:A:13:TYR:CZ	1:A:317:LYS:HD2	2.49	0.47
1:B:95:LEU:HA	2:B:420:HOH:O	2.14	0.47
1:B:16:ASP:HB2	1:B:61:CYS:HA	1.96	0.47
1:B:104:ILE:HD13	1:B:119:SER:HB3	1.96	0.47
1:A:158:THR:HA	1:A:162:ARG:O	2.15	0.47
1:A:36:THR:HG23	2:A:469:HOH:O	2.15	0.47
1:B:315:THR:HB	1:B:328:LEU:CD2	2.43	0.47
1:B:263:HIS:O	1:B:264:LYS:HB2	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:252:TYR:HA	1:A:253:PRO:HD2	1.81	0.47
1:A:176:ARG:HG2	1:A:176:ARG:HH11	1.79	0.47
1:B:203:PRO:O	1:B:205:GLU:N	2.48	0.47
1:A:24:SER:HA	1:A:40:PHE:CZ	2.50	0.47
1:A:217:ARG:HH12	1:A:239:ARG:HD2	1.79	0.47
1:B:139:VAL:CG1	1:B:179:LEU:HD13	2.44	0.47
1:A:303:SER:O	1:A:305:ASN:N	2.48	0.47
1:B:234:LYS:HA	2:B:457:HOH:O	2.13	0.47
1:A:49:LEU:C	1:A:50:LEU:HD12	2.36	0.46
1:A:67:ASN:HD22	1:A:113:ASP:HA	1.80	0.46
1:B:139:VAL:HG11	1:B:179:LEU:HD13	1.96	0.46
1:B:61:CYS:HB2	2:B:387:HOH:O	2.15	0.46
1:B:159:ASN:ND2	1:B:201:LEU:HD13	2.29	0.46
1:B:167:MSE:SE	1:B:171:GLN:HE21	2.48	0.46
1:B:275:ILE:CD1	1:B:309:LEU:HD22	2.45	0.46
1:A:307:LEU:C	1:A:307:LEU:CD2	2.82	0.46
1:B:169:ASN:C	1:B:171:GLN:N	2.69	0.46
1:A:51:GLN:HG3	1:A:53:LEU:CD1	2.45	0.46
1:A:260:SER:HB3	1:A:267:TYR:CE1	2.45	0.46
1:A:217:ARG:NH1	1:A:239:ARG:HG3	2.30	0.46
1:A:156:MSE:HA	1:A:164:ILE:O	2.16	0.46
1:B:162:ARG:NH2	1:B:225:ASP:OD2	2.49	0.46
1:B:142:LEU:HB3	1:B:151:ASN:HB3	1.98	0.46
1:B:271:SER:C	1:B:273:GLY:H	2.19	0.46
1:B:39:LYS:CB	1:B:50:LEU:HD21	2.45	0.46
1:B:202:LEU:HD23	1:B:209:TYR:N	2.30	0.46
1:B:62:CYS:HB3	1:B:74:VAL:HG12	1.98	0.45
1:A:215:ASP:HA	1:A:251:ALA:HB1	1.97	0.45
1:B:176:ARG:NH1	2:B:469:HOH:O	2.48	0.45
1:A:285:LYS:HD3	1:A:288:ASN:OD1	2.16	0.45
1:B:42:ILE:O	1:B:45:LYS:HE3	2.15	0.45
1:A:59:LEU:HD23	1:A:76:THR:HG22	1.97	0.45
1:B:260:SER:O	1:B:264:LYS:HA	2.16	0.45
1:B:35:LEU:HD23	1:B:84:VAL:HG11	1.99	0.45
1:B:202:LEU:HD23	1:B:209:TYR:CA	2.46	0.45
1:A:222:PHE:HE2	1:A:234:LYS:O	1.99	0.45
1:B:164:ILE:HD13	1:B:199:VAL:HG21	1.99	0.45
1:B:87:ILE:N	1:B:87:ILE:CD1	2.79	0.45
1:A:1:MSE:HE1	1:A:287:LYS:HZ2	1.78	0.45
1:A:279:ASN:HB3	1:A:282:THR:CG2	2.45	0.45
1:B:282:THR:OG1	1:B:284:LYS:HD3	2.17	0.45
1:B:282:THR:OG1	1:B:284:LYS:CD	2.64	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:146:ASN:CG	1:A:150:LYS:HD2	2.37	0.45
1:B:53:LEU:CD1	1:B:53:LEU:N	2.79	0.45
1:B:62:CYS:O	1:B:63:ASN:HB2	2.15	0.45
1:B:84:VAL:HG12	1:B:85:ASP:N	2.32	0.45
1:A:4:VAL:HG11	1:A:337:PHE:HE2	1.81	0.45
1:A:124:ILE:HB	1:A:140:LYS:HB2	1.99	0.45
1:A:6:ILE:HD12	1:A:333:ILE:HG22	1.98	0.45
1:A:140:LYS:NZ	1:A:184:ASN:CB	2.80	0.45
1:B:109:LYS:HA	1:B:115:LEU:HD23	1.98	0.45
1:B:65:ILE:HG23	1:B:113:ASP:HA	1.97	0.45
1:B:100:ALA:N	1:B:123:LEU:HD13	2.32	0.45
1:A:192:LEU:HD23	2:A:448:HOH:O	2.15	0.44
1:A:106:ARG:HH11	1:A:106:ARG:HG3	1.82	0.44
1:B:175:PHE:CD1	1:B:175:PHE:N	2.84	0.44
1:B:119:SER:OG	1:B:123:LEU:HB2	2.17	0.44
1:A:58:PRO:HB2	1:A:77:VAL:CG2	2.47	0.44
1:B:276:SER:OG	1:B:288:ASN:ND2	2.50	0.44
1:A:110:TYR:C	1:A:111:GLY:O	2.56	0.44
1:B:32:ASP:OD2	1:B:33:GLY:N	2.50	0.44
1:B:116:ILE:O	1:B:156:MSE:HE1	2.16	0.44
1:B:126:VAL:CG1	1:B:127:ILE:N	2.81	0.44
1:A:138:ALA:HB3	2:A:467:HOH:O	2.17	0.44
1:B:49:LEU:C	1:B:50:LEU:HD12	2.38	0.44
1:A:110:TYR:O	1:A:111:GLY:O	2.35	0.44
1:B:184:ASN:CG	1:B:185:GLY:N	2.70	0.44
1:A:184:ASN:CG	1:A:185:GLY:H	2.21	0.44
1:B:292:PHE:CD1	1:B:292:PHE:N	2.85	0.44
1:A:160:SER:OG	1:A:204:LYS:CD	2.64	0.44
1:A:110:TYR:CE2	1:A:177:LEU:HD22	2.52	0.44
1:B:6:ILE:HG21	1:B:38:TYR:CE2	2.53	0.44
1:B:125:GLU:HB3	1:B:136:VAL:CG1	2.47	0.44
1:A:16:ASP:HA	1:A:299:LYS:HG3	1.98	0.44
1:B:19:ILE:HD13	1:B:306:ILE:HD11	1.98	0.44
1:A:214:ILE:CA	1:A:253:PRO:HB3	2.40	0.44
1:A:5:GLN:HB2	1:A:334:TYR:CE2	2.52	0.44
1:A:149:VAL:O	1:A:151:ASN:N	2.51	0.44
1:A:172:VAL:HG12	1:A:173:GLN:N	2.33	0.44
1:B:76:THR:OG1	1:B:80:GLU:HB2	2.18	0.44
1:B:282:THR:O	1:B:283:ARG:HB2	2.18	0.43
1:B:109:LYS:HB2	1:B:109:LYS:HE2	1.90	0.43
1:B:263:HIS:CD2	1:B:340:GLU:HB2	2.53	0.43
1:A:55:TYR:CE2	1:A:76:THR:HG21	2.54	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:162:ARG:NH1	1:B:223:PHE:CD1	2.86	0.43
1:B:205:GLU:HG3	1:B:206:GLN:HG3	2.00	0.43
1:A:249:ASN:HA	2:A:475:HOH:O	2.17	0.43
1:B:33:GLY:O	1:B:55:TYR:N	2.50	0.43
1:B:247:ASP:O	1:B:248:THR:HG22	2.18	0.43
1:A:113:ASP:HB2	1:A:130:ARG:HD3	1.99	0.43
1:B:16:ASP:HB3	1:B:29:THR:OG1	2.19	0.43
1:A:37:VAL:HG12	1:A:38:TYR:N	2.33	0.43
1:A:20:ILE:HA	1:A:21:PRO:HD2	1.66	0.43
1:B:23:LYS:CD	1:B:23:LYS:N	2.79	0.43
1:A:203:PRO:HD2	1:A:206:GLN:OE1	2.18	0.43
1:B:202:LEU:HA	1:B:259:PHE:CE1	2.54	0.43
1:B:51:GLN:NE2	1:B:86:LEU:HB3	2.19	0.43
1:A:192:LEU:HD12	1:A:196:ILE:CG2	2.48	0.43
1:B:17:ILE:HD11	1:B:310:ALA:CB	2.48	0.43
1:B:202:LEU:HA	1:B:259:PHE:CZ	2.54	0.43
1:A:117:ALA:O	1:A:124:ILE:HA	2.19	0.43
1:B:82:LEU:HA	1:B:94:ALA:HA	2.01	0.43
1:A:83:LYS:O	1:A:92:PHE:HA	2.18	0.43
1:A:164:ILE:HG13	1:A:174:TRP:HB3	2.01	0.42
1:A:257:ILE:HG13	1:A:257:ILE:O	2.19	0.42
1:A:28:ILE:O	1:A:35:LEU:HA	2.19	0.42
1:A:307:LEU:HD23	1:A:308:CYS:N	2.34	0.42
1:A:217:ARG:HH12	1:A:239:ARG:HG3	1.84	0.42
1:A:315:THR:HG23	1:A:319:ASN:HB2	2.00	0.42
1:A:205:GLU:OE2	1:A:205:GLU:N	2.52	0.42
1:B:260:SER:O	1:B:264:LYS:HD3	2.18	0.42
1:B:242:ARG:HD3	2:B:516:HOH:O	2.18	0.42
1:A:102:LEU:HG	1:A:121:ASP:HB3	2.01	0.42
1:B:126:VAL:HB	1:B:137:ILE:HG23	2.01	0.42
1:B:239:ARG:HH12	1:B:243:LEU:HD11	1.84	0.42
1:B:32:ASP:OD2	1:B:32:ASP:C	2.58	0.42
1:A:25:LEU:HD23	1:A:39:LYS:HD2	2.02	0.42
1:A:118:ALA:HA	1:A:123:LEU:O	2.19	0.42
1:A:268:THR:O	1:A:275:ILE:HA	2.20	0.42
1:A:192:LEU:CD1	1:A:196:ILE:HG22	2.45	0.42
1:B:108:CYS:HB2	1:B:156:MSE:HE3	2.01	0.42
1:B:235:ARG:HG3	1:B:235:ARG:H	1.58	0.42
1:A:167:MSE:CE	1:A:171:GLN:HE21	2.31	0.42
1:B:18:LYS:HG3	1:B:63:ASN:HA	2.01	0.42
1:B:176:ARG:HD3	2:B:480:HOH:O	2.20	0.42
1:B:44:ALA:O	1:B:45:LYS:HB2	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:301:ALA:O	1:A:307:LEU:HA	2.20	0.42
1:B:202:LEU:HD11	1:B:280:LEU:HD12	2.02	0.42
1:A:239:ARG:HG3	2:A:404:HOH:O	2.18	0.41
1:B:139:VAL:O	1:B:140:LYS:HD3	2.20	0.41
1:B:20:ILE:CD1	1:B:25:LEU:HD12	2.45	0.41
1:B:312:SER:HA	1:B:331:SER:CB	2.49	0.41
1:A:137:ILE:HG23	1:A:137:ILE:O	2.20	0.41
1:B:203:PRO:HD2	1:B:206:GLN:OE1	2.20	0.41
1:A:39:LYS:HB2	1:A:50:LEU:CD1	2.49	0.41
1:A:164:ILE:HG12	1:A:172:VAL:CG1	2.51	0.41
1:B:158:THR:O	1:B:159:ASN:HB3	2.19	0.41
1:A:269:ALA:HB2	1:A:300:ILE:HG23	2.02	0.41
1:A:13:TYR:O	1:A:31:TRP:HD1	2.02	0.41
1:A:249:ASN:N	2:A:391:HOH:O	2.54	0.41
1:B:150:LYS:HZ1	1:B:150:LYS:HA	1.83	0.41
1:A:209:TYR:CZ	1:A:221:GLU:HB3	2.56	0.41
1:A:39:LYS:N	1:A:50:LEU:HD13	2.36	0.41
1:A:303:SER:O	1:A:304:ASP:C	2.57	0.41
1:A:169:ASN:O	1:A:170:SER:C	2.59	0.41
1:B:169:ASN:O	1:B:170:SER:C	2.59	0.41
1:A:89:SER:HA	1:A:90:PRO:C	2.40	0.41
1:B:46:ASN:HD22	1:B:47:VAL:N	2.18	0.41
1:B:108:CYS:SG	1:B:156:MSE:HE3	2.61	0.41
1:A:217:ARG:NH1	1:A:239:ARG:CG	2.85	0.40
1:B:288:ASN:HA	1:B:288:ASN:HD22	1.55	0.40
1:B:56:LYS:HE2	1:B:57:HIS:CE1	2.56	0.40
1:A:116:ILE:HG22	1:A:156:MSE:HE1	2.03	0.40
1:A:259:PHE:CE1	1:A:266:LEU:HD13	2.55	0.40
1:B:51:GLN:HG3	1:B:53:LEU:CD1	2.51	0.40
1:B:299:LYS:NZ	2:B:406:HOH:O	2.54	0.40
1:A:197:ARG:NH1	1:A:255:ASN:OD1	2.55	0.40
1:B:4:VAL:CG2	1:B:4:VAL:O	2.69	0.40
1:B:51:GLN:CG	1:B:53:LEU:HD11	2.51	0.40
1:B:8:GLN:HG3	1:B:49:LEU:HD12	2.03	0.40
1:B:243:LEU:CD2	1:B:243:LEU:H	2.13	0.40
1:A:140:LYS:HZ3	1:A:184:ASN:HB3	1.86	0.40
1:B:248:THR:HG21	2:B:394:HOH:O	2.20	0.40
1:A:68:THR:HB	2:A:424:HOH:O	2.21	0.40
1:A:168:ASN:C	1:A:170:SER:H	2.21	0.40
1:A:79:GLY:HA2	1:A:104:ILE:HG12	2.02	0.40
1:A:300:ILE:HA	1:A:308:CYS:O	2.21	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	312/349 (89%)	268 (86%)	33 (11%)	11 (4%)	6	3
1	B	325/349 (93%)	266 (82%)	41 (13%)	18 (6%)	3	1
All	All	637/698 (91%)	534 (84%)	74 (12%)	29 (5%)	4	1

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	A	113	ASP
1	A	304	ASP
1	B	181	GLU
1	B	204	LYS
1	B	245	LEU
1	B	249	ASN
1	B	262	ARG
1	A	169	ASN
1	B	184	ASN
1	B	234	LYS
1	B	248	THR
1	A	68	THR
1	B	52	SER
1	B	148	LYS
1	B	224	ASP
1	A	150	LYS
1	A	262	ARG
1	A	295	ASP
1	B	88	GLY
1	B	141	ASN
1	A	293	ASN
1	B	183	ASP
1	A	111	GLY
1	A	149	VAL
1	B	295	ASP
1	B	136	VAL

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Mol	Chain	Res	Type
1	B	274	ILE
1	B	297	VAL

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	287/310 (93%)	263 (92%)	24 (8%)	16	17
1	B	296/310 (96%)	262 (88%)	34 (12%)	8	7
All	All	583/620 (94%)	525 (90%)	58 (10%)	11	11

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	23	LYS
1	A	42	ILE
1	A	46	ASN
1	A	53	LEU
1	A	69	ASP
1	A	71	GLN
1	A	86	LEU
1	A	87	ILE
1	A	93	GLN
1	A	109	LYS
1	A	123	LEU
1	A	184	ASN
1	A	186	THR
1	A	196	ILE
1	A	205	GLU
1	A	221	GLU
1	A	258	GLU
1	A	281	GLN
1	A	283	ARG
1	A	304	ASP
1	A	314	ASP
1	A	317	LYS

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Mol	Chain	Res	Type
1	A	327	GLU
1	B	7	GLU
1	B	8	GLN
1	B	22	SER
1	B	23	LYS
1	B	24	SER
1	B	36	THR
1	B	45	LYS
1	B	46	ASN
1	B	49	LEU
1	B	87	ILE
1	B	106	ARG
1	B	109	LYS
1	B	150	LYS
1	B	184	ASN
1	B	186	THR
1	B	188	GLU
1	B	196	ILE
1	B	204	LYS
1	B	221	GLU
1	B	225	ASP
1	B	235	ARG
1	B	243	LEU
1	B	245	LEU
1	B	250	LEU
1	B	274	ILE
1	B	281	GLN
1	B	288	ASN
1	B	295	ASP
1	B	304	ASP
1	B	309	LEU
1	B	311	THR
1	B	314	ASP
1	B	317	LYS
1	B	326	ILE

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	46	ASN
1	A	57	HIS
1	A	71	GLN

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Mol	Chain	Res	Type
1	A	171	GLN
1	A	249	ASN
1	A	263	HIS
1	A	281	GLN
1	A	288	ASN
1	A	329	ASN
1	B	51	GLN
1	B	151	ASN
1	B	171	GLN
1	B	184	ASN
1	B	281	GLN
1	B	288	ASN
1	B	305	ASN
1	B	329	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

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	322/349 (92%)	0.08	3 (0%) 81 84	38, 57, 84, 113	0
1	B	331/349 (94%)	0.06	4 (1%) 75 79	38, 57, 92, 117	0
All	All	653/698 (93%)	0.07	7 (1%) 77 81	38, 57, 90, 117	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	66	ASP	3.6
1	A	180	CYS	3.5
1	A	67	ASN	2.8
1	B	232	SER	2.5
1	B	247	ASP	2.4
1	B	248	THR	2.3
1	A	50	LEU	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.