



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:18 AM GMT

PDB ID : 3HXK
Title : Crystal Structure of a sugar hydrolase (YeeB) from Lactococcus lactis, Northeast Structural Genomics Consortium Target KR108
Authors : Forouhar,F.; Su, M.; Seetharaman, J.; Mao,M.; Xiao,R.; Ciccianti,C.; Foote,E.L.; Maglaqui,M.; Zhao,L.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong,L.; Hunt,J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2009-06-20
Resolution : 3.20 Å(reported)

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

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

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

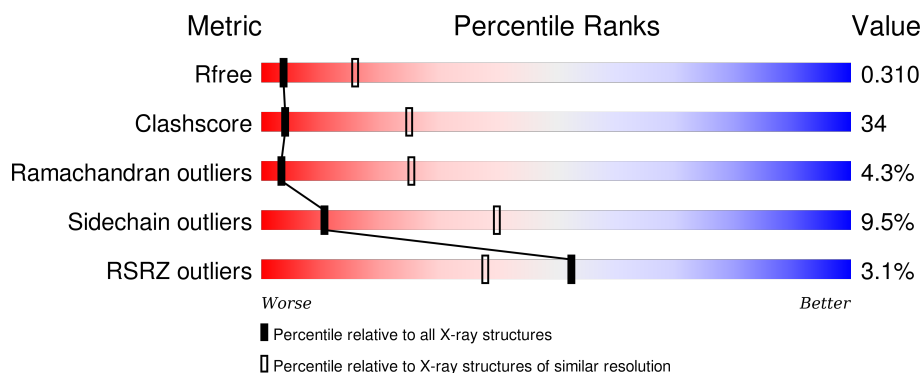
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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

Mol	Chain	Length	Quality of chain
1	A	276	<div> <div>2%</div> <div>37% 46% 7% 9%</div> </div>
1	B	276	<div> <div>3%</div> <div>36% 47% 8% 9%</div> </div>
1	C	276	<div> <div>2%</div> <div>33% 49% 8% 9%</div> </div>
1	D	276	<div> <div>4%</div> <div>38% 46% 7% 9%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8141 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 Sugar hydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	250	Total 2020	C 1300	N 337	O 377	S 5	Se 1	0	0	0
1	B	250	Total 2020	C 1300	N 337	O 377	S 5	Se 1	0	0	0
1	C	250	Total 2020	C 1300	N 337	O 377	S 5	Se 1	0	0	0
1	D	250	Total 2020	C 1300	N 337	O 377	S 5	Se 1	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	269	LEU	-	expression tag	UNP Q9CID4
A	270	GLU	-	expression tag	UNP Q9CID4
A	271	HIS	-	expression tag	UNP Q9CID4
A	272	HIS	-	expression tag	UNP Q9CID4
A	273	HIS	-	expression tag	UNP Q9CID4
A	274	HIS	-	expression tag	UNP Q9CID4
A	275	HIS	-	expression tag	UNP Q9CID4
A	276	HIS	-	expression tag	UNP Q9CID4
B	269	LEU	-	expression tag	UNP Q9CID4
B	270	GLU	-	expression tag	UNP Q9CID4
B	271	HIS	-	expression tag	UNP Q9CID4
B	272	HIS	-	expression tag	UNP Q9CID4
B	273	HIS	-	expression tag	UNP Q9CID4
B	274	HIS	-	expression tag	UNP Q9CID4
B	275	HIS	-	expression tag	UNP Q9CID4
B	276	HIS	-	expression tag	UNP Q9CID4
C	269	LEU	-	expression tag	UNP Q9CID4
C	270	GLU	-	expression tag	UNP Q9CID4
C	271	HIS	-	expression tag	UNP Q9CID4
C	272	HIS	-	expression tag	UNP Q9CID4
C	273	HIS	-	expression tag	UNP Q9CID4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	274	HIS	-	expression tag	UNP Q9CID4
C	275	HIS	-	expression tag	UNP Q9CID4
C	276	HIS	-	expression tag	UNP Q9CID4
D	269	LEU	-	expression tag	UNP Q9CID4
D	270	GLU	-	expression tag	UNP Q9CID4
D	271	HIS	-	expression tag	UNP Q9CID4
D	272	HIS	-	expression tag	UNP Q9CID4
D	273	HIS	-	expression tag	UNP Q9CID4
D	274	HIS	-	expression tag	UNP Q9CID4
D	275	HIS	-	expression tag	UNP Q9CID4
D	276	HIS	-	expression tag	UNP Q9CID4

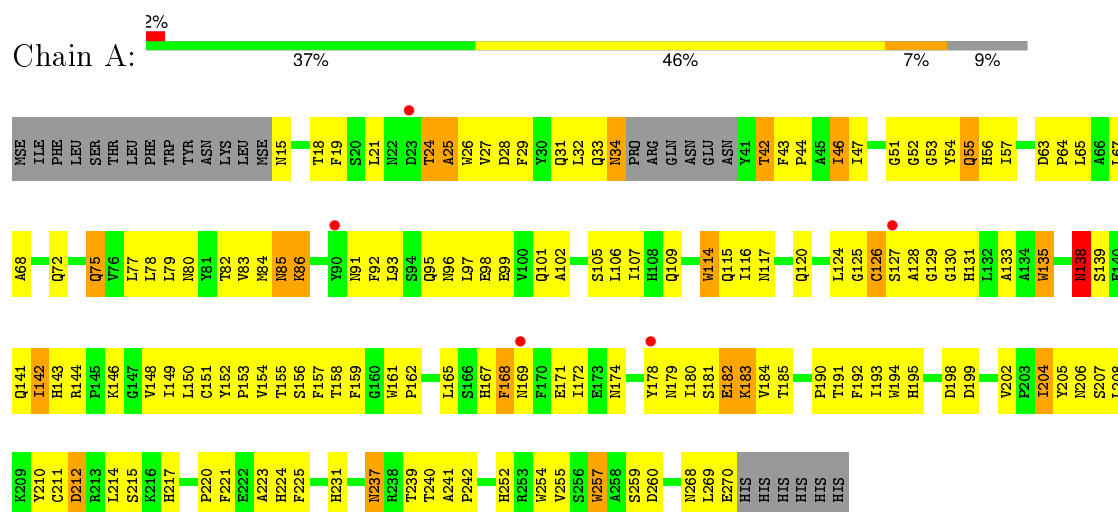
- Molecule 2 is water.

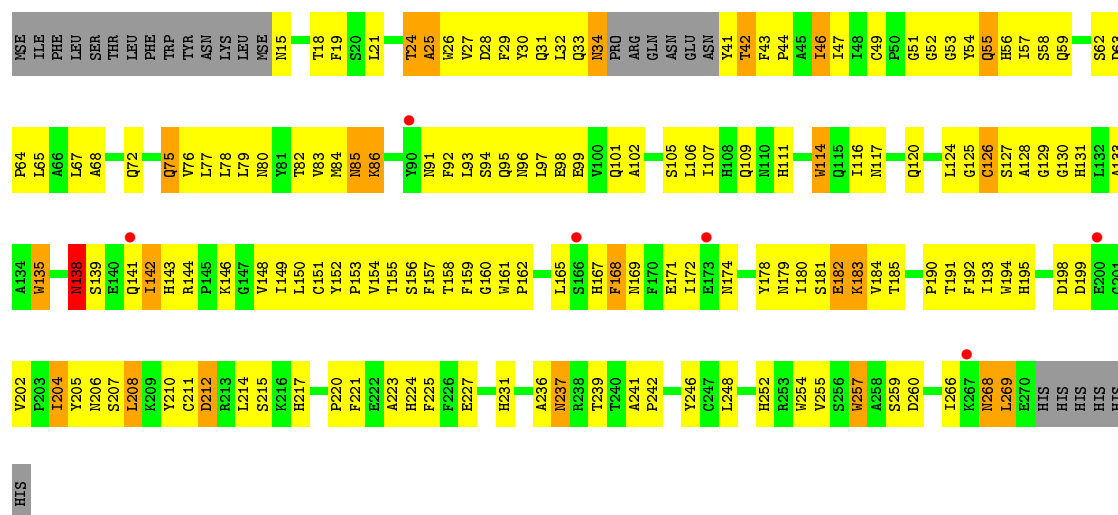
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	9	Total O 9 9	0	0
2	B	13	Total O 13 13	0	0
2	C	14	Total O 14 14	0	0
2	D	25	Total O 25 25	0	0

3 Residue-property plots [i](#)

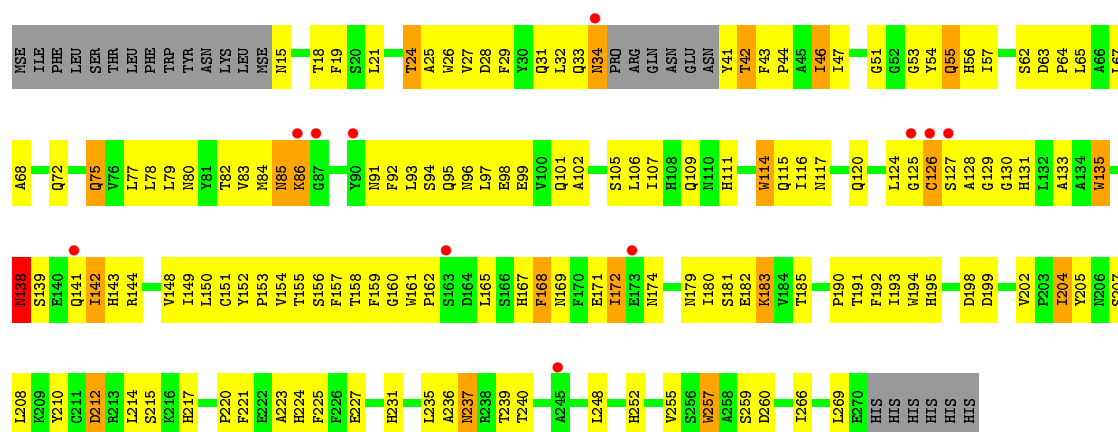
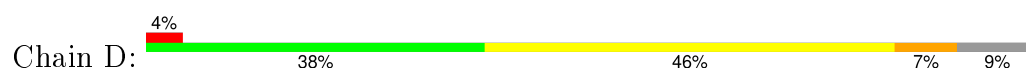
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Sugar hydrolase





• Molecule 1: Sugar hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	68.39Å 150.43Å 218.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 3.20 28.63 – 3.08	Depositor EDS
% Data completeness (in resolution range)	91.7 (19.96-3.20) 91.7 (28.63-3.08)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 3.05Å)	Xtriage
Refinement program	CNS 1.2 & XtalView	Depositor
R, R_{free}	0.256 , 0.296 0.273 , 0.310	Depositor DCC
R_{free} test set	1731 reflections (9.48%)	DCC
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.442	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 47.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	1 of 36062 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	8141	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.50	0/2085	0.58	0/2845
1	B	0.49	0/2085	0.58	0/2845
1	C	0.52	0/2085	0.59	0/2845
1	D	0.51	0/2085	0.58	0/2845
All	All	0.51	0/8340	0.58	0/11380

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2020	0	1897	136	0
1	B	2020	0	1897	134	0
1	C	2020	0	1897	147	0
1	D	2020	0	1897	140	0
2	A	9	0	0	3	0
2	B	13	0	0	3	0
2	C	14	0	0	6	0
2	D	25	0	0	8	0
All	All	8141	0	7588	534	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 34.

All (534) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:ILE:HB	2:A:283:HOH:O	1.50	1.12
1:D:126:CYS:HB3	1:D:151:CYS:HB2	1.35	1.02
1:B:126:CYS:HB3	1:B:151:CYS:HB2	1.39	1.01
1:A:126:CYS:HB3	1:A:151:CYS:HB2	1.45	0.99
1:C:126:CYS:HB3	1:C:151:CYS:HB2	1.43	0.96
1:A:199:ASP:OD2	1:A:231:HIS:HB2	1.71	0.90
1:D:148:VAL:HG22	1:D:191:THR:HG22	1.54	0.89
1:B:199:ASP:OD2	1:B:231:HIS:HB2	1.74	0.88
1:C:199:ASP:OD2	1:C:231:HIS:HB2	1.74	0.88
1:D:199:ASP:OD2	1:D:231:HIS:HB2	1.74	0.88
1:A:148:VAL:HG22	1:A:191:THR:HG22	1.54	0.88
1:C:148:VAL:HG22	1:C:191:THR:HG22	1.56	0.85
1:A:168:PHE:HB2	2:A:282:HOH:O	1.76	0.84
1:C:208:LEU:HD23	1:D:208:LEU:HD23	1.61	0.81
1:D:26:TRP:HB2	1:D:80:ASN:HB3	1.62	0.81
1:C:26:TRP:HB2	1:C:80:ASN:HB3	1.62	0.80
1:A:215:SER:OG	1:B:204:ILE:HD11	1.82	0.80
1:B:26:TRP:HB2	1:B:80:ASN:HB3	1.63	0.80
1:A:26:TRP:HB2	1:A:80:ASN:HB3	1.63	0.80
1:B:124:LEU:HD13	1:B:149:ILE:HB	1.64	0.79
1:B:148:VAL:HG22	1:B:191:THR:HG22	1.65	0.78
1:D:124:LEU:HD13	1:D:149:ILE:HB	1.66	0.78
1:A:131:HIS:HB2	1:A:154:VAL:HB	1.67	0.75
1:D:124:LEU:HD12	1:D:125:GLY:H	1.51	0.75
1:C:124:LEU:HD13	1:C:149:ILE:HB	1.68	0.75
1:A:257:TRP:CH2	1:B:257:TRP:CH2	2.74	0.74
1:A:124:LEU:HD13	1:A:149:ILE:HB	1.67	0.74
1:C:131:HIS:HB2	1:C:154:VAL:HB	1.70	0.74
1:C:41:TYR:HA	2:C:289:HOH:O	1.87	0.74
1:B:32:LEU:HG	1:B:33:GLN:H	1.54	0.73
1:D:131:HIS:HB2	1:D:154:VAL:HB	1.71	0.72
1:A:34:ASN:ND2	2:A:277:HOH:O	2.22	0.72
1:A:185:THR:HA	1:A:217:HIS:HD2	1.54	0.72
1:D:202:VAL:HG12	1:D:231:HIS:HD2	1.53	0.72
1:C:150:LEU:HB2	1:C:193:ILE:HG13	1.72	0.72
1:C:202:VAL:HG12	1:C:231:HIS:HD2	1.55	0.71
1:D:32:LEU:HG	1:D:33:GLN:H	1.55	0.71
1:B:202:VAL:HG12	1:B:231:HIS:HD2	1.54	0.71
1:A:124:LEU:HD12	1:A:125:GLY:H	1.55	0.71
1:A:202:VAL:HG12	1:A:231:HIS:HD2	1.53	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:32:LEU:HG	1:C:33:GLN:H	1.56	0.70
1:B:131:HIS:HB2	1:B:154:VAL:HB	1.71	0.70
1:D:150:LEU:HB2	1:D:193:ILE:HG13	1.73	0.70
1:A:150:LEU:HB2	1:A:193:ILE:HG13	1.73	0.70
1:C:185:THR:HA	1:C:217:HIS:HD2	1.55	0.70
1:B:124:LEU:HD12	1:B:125:GLY:H	1.56	0.69
1:D:97:LEU:HD11	1:D:135:TRP:NE1	2.08	0.69
1:A:32:LEU:HG	1:A:33:GLN:H	1.58	0.68
1:A:185:THR:HA	1:A:217:HIS:CD2	2.28	0.68
1:D:33:GLN:N	2:D:296:HOH:O	2.27	0.68
1:C:102:ALA:O	1:C:106:LEU:HD23	1.93	0.68
1:A:97:LEU:HD11	1:A:135:TRP:NE1	2.09	0.68
1:C:185:THR:HA	1:C:217:HIS:CD2	2.29	0.68
1:D:135:TRP:HZ3	1:D:183:LYS:HG2	1.60	0.67
1:D:102:ALA:O	1:D:106:LEU:HD23	1.95	0.67
1:B:97:LEU:HD11	1:B:135:TRP:NE1	2.10	0.67
1:C:97:LEU:HD11	1:C:135:TRP:NE1	2.09	0.67
1:B:185:THR:HA	1:B:217:HIS:HD2	1.60	0.66
1:C:124:LEU:HD12	1:C:125:GLY:H	1.60	0.66
1:D:162:PRO:HB3	1:D:168:PHE:HE1	1.62	0.65
1:B:54:TYR:O	1:B:83:VAL:HB	1.96	0.65
1:D:172:ILE:HA	2:D:278:HOH:O	1.95	0.65
1:D:185:THR:HA	1:D:217:HIS:HD2	1.61	0.65
1:D:192:PHE:C	1:D:193:ILE:HD12	2.17	0.65
1:A:167:HIS:C	1:A:169:ASN:H	2.00	0.64
1:D:185:THR:HA	1:D:217:HIS:CD2	2.33	0.64
1:D:54:TYR:O	1:D:83:VAL:HB	1.98	0.64
1:B:150:LEU:HB2	1:B:193:ILE:HG13	1.78	0.64
1:C:167:HIS:C	1:C:169:ASN:H	2.00	0.64
1:B:204:ILE:HD12	1:B:205:TYR:N	2.13	0.64
1:C:162:PRO:HB3	1:C:168:PHE:HE1	1.62	0.63
1:B:185:THR:HA	1:B:217:HIS:CD2	2.33	0.63
1:B:167:HIS:C	1:B:169:ASN:H	2.00	0.63
1:B:192:PHE:C	1:B:193:ILE:HD12	2.18	0.63
1:C:257:TRP:CH2	1:D:257:TRP:CH2	2.87	0.63
1:D:167:HIS:C	1:D:169:ASN:H	2.01	0.63
1:A:208:LEU:HD23	1:B:208:LEU:HD23	1.81	0.63
1:D:33:GLN:HB3	2:D:296:HOH:O	1.99	0.63
1:C:46:ILE:HD11	1:C:124:LEU:HB2	1.81	0.62
1:B:202:VAL:HG12	1:B:231:HIS:CD2	2.35	0.62
1:A:162:PRO:HB3	1:A:168:PHE:HE1	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:PRO:HB3	1:B:168:PHE:HE1	1.64	0.62
1:B:204:ILE:HD13	1:B:208:LEU:HD12	1.81	0.62
1:A:102:ALA:O	1:A:106:LEU:HD23	1.99	0.62
1:C:194:TRP:HA	1:C:224:HIS:O	2.00	0.62
1:A:46:ILE:HD11	1:A:124:LEU:HB2	1.82	0.62
1:A:202:VAL:HG12	1:A:231:HIS:CD2	2.35	0.62
1:B:204:ILE:HD12	1:B:205:TYR:H	1.65	0.62
1:B:135:TRP:HZ3	1:B:183:LYS:HG2	1.64	0.62
1:D:42:THR:HG22	1:D:116:ILE:C	2.19	0.62
1:D:46:ILE:HD11	1:D:124:LEU:HB2	1.82	0.61
1:D:156:SER:HB2	1:D:179:ASN:OD1	2.00	0.61
1:B:42:THR:HG22	1:B:116:ILE:C	2.19	0.61
1:D:204:ILE:HD13	1:D:208:LEU:HD12	1.82	0.61
1:C:29:PHE:CZ	1:C:31:GLN:HB2	2.36	0.61
1:A:54:TYR:O	1:A:83:VAL:HB	2.00	0.61
1:C:204:ILE:HD12	1:C:205:TYR:H	1.66	0.61
1:C:204:ILE:HD13	1:C:208:LEU:HD12	1.81	0.61
1:C:204:ILE:HD12	1:C:205:TYR:N	2.16	0.60
1:A:135:TRP:HZ3	1:A:183:LYS:HG2	1.65	0.60
1:A:192:PHE:C	1:A:193:ILE:HD12	2.21	0.60
1:A:204:ILE:HD12	1:A:205:TYR:N	2.16	0.60
1:A:204:ILE:HD11	1:B:215:SER:OG	2.01	0.60
1:C:135:TRP:HZ3	1:C:183:LYS:HG2	1.64	0.60
1:D:204:ILE:HD12	1:D:205:TYR:N	2.17	0.60
1:C:54:TYR:O	1:C:83:VAL:HB	2.01	0.60
1:C:82:THR:HG22	1:C:99:GLU:OE2	2.01	0.60
1:A:67:LEU:HD21	2:D:297:HOH:O	2.02	0.60
1:D:252:HIS:O	1:D:255:VAL:HG12	2.01	0.60
1:C:42:THR:HG22	1:C:117:ASN:N	2.16	0.60
1:B:156:SER:HB2	1:B:179:ASN:OD1	2.02	0.60
1:D:42:THR:HG22	1:D:117:ASN:N	2.17	0.60
1:C:192:PHE:C	1:C:193:ILE:HD12	2.21	0.60
1:A:42:THR:HG22	1:A:116:ILE:C	2.22	0.60
1:C:149:ILE:O	1:C:150:LEU:HD23	2.02	0.59
1:B:64:PRO:HB3	1:B:239:THR:HG22	1.84	0.59
1:C:64:PRO:HB3	1:C:239:THR:HG22	1.84	0.59
1:A:252:HIS:O	1:A:255:VAL:HG12	2.02	0.59
1:B:46:ILE:HD11	1:B:124:LEU:HB2	1.83	0.59
1:D:194:TRP:HA	1:D:224:HIS:O	2.01	0.59
1:A:146:LYS:HG2	1:A:269:LEU:HD21	1.84	0.59
1:B:194:TRP:HA	1:B:224:HIS:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:155:THR:HG22	1:D:181:SER:HB2	1.83	0.59
1:C:202:VAL:HG12	1:C:231:HIS:CD2	2.38	0.59
1:B:149:ILE:O	1:B:150:LEU:HD23	2.02	0.59
1:C:252:HIS:O	1:C:255:VAL:HG12	2.03	0.59
1:C:59:GLN:NE2	2:C:287:HOH:O	2.34	0.59
1:D:204:ILE:HD12	1:D:205:TYR:H	1.66	0.59
1:A:204:ILE:HD12	1:A:205:TYR:H	1.66	0.59
1:C:156:SER:HB2	1:C:179:ASN:OD1	2.03	0.59
1:D:202:VAL:HG12	1:D:231:HIS:CD2	2.36	0.59
1:B:42:THR:HG22	1:B:117:ASN:N	2.17	0.59
1:D:63:ASP:HB3	1:D:64:PRO:HD3	1.84	0.58
1:B:29:PHE:CZ	1:B:31:GLN:HB2	2.37	0.58
1:A:29:PHE:CZ	1:A:31:GLN:HB2	2.39	0.58
1:A:156:SER:HB2	1:A:179:ASN:OD1	2.03	0.58
1:C:162:PRO:HB3	1:C:168:PHE:CE1	2.38	0.58
1:C:63:ASP:HB3	1:C:64:PRO:HD3	1.86	0.58
1:C:42:THR:HG22	1:C:116:ILE:C	2.24	0.58
1:D:162:PRO:HB3	1:D:168:PHE:CE1	2.38	0.58
1:A:63:ASP:HB3	1:A:64:PRO:HD3	1.86	0.58
1:A:82:THR:HG22	1:A:99:GLU:OE2	2.03	0.57
1:A:162:PRO:HB3	1:A:168:PHE:CE1	2.39	0.57
1:A:57:ILE:HG13	1:A:80:ASN:OD1	2.03	0.57
1:B:155:THR:HG22	1:B:181:SER:HB2	1.86	0.57
1:D:29:PHE:CZ	1:D:31:GLN:HB2	2.39	0.57
1:B:102:ALA:O	1:B:106:LEU:HD23	2.05	0.57
1:A:204:ILE:HD13	1:A:208:LEU:HD12	1.86	0.57
1:B:82:THR:HG22	1:B:99:GLU:OE2	2.04	0.57
1:D:82:THR:HG22	1:D:99:GLU:OE2	2.05	0.57
1:D:25:ALA:HB2	1:D:99:GLU:OE1	2.04	0.57
1:B:162:PRO:HB3	1:B:168:PHE:CE1	2.40	0.57
1:A:155:THR:HG22	1:A:181:SER:HB2	1.86	0.57
1:D:64:PRO:HB3	1:D:239:THR:HG22	1.87	0.56
1:D:57:ILE:HG13	1:D:80:ASN:OD1	2.05	0.56
1:A:194:TRP:HA	1:A:224:HIS:O	2.05	0.56
1:C:155:THR:HG22	1:C:181:SER:HB2	1.85	0.56
1:A:64:PRO:HB3	1:A:239:THR:HG22	1.87	0.56
1:C:57:ILE:HG13	1:C:80:ASN:OD1	2.04	0.56
1:D:26:TRP:HB2	2:D:301:HOH:O	2.06	0.56
1:C:93:LEU:HA	1:C:96:ASN:HD22	1.71	0.56
1:B:252:HIS:O	1:B:255:VAL:HG12	2.06	0.56
1:B:144:ARG:HH11	1:B:144:ARG:HG3	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:LEU:HA	1:A:96:ASN:HD22	1.71	0.55
1:D:151:CYS:O	1:D:152:TYR:C	2.44	0.55
1:C:180:ILE:HD12	1:C:180:ILE:C	2.27	0.55
1:C:34:ASN:ND2	2:C:285:HOH:O	2.27	0.55
1:D:156:SER:C	1:D:158:THR:H	2.11	0.55
1:D:141:GLN:C	1:D:143:HIS:H	2.11	0.55
1:A:25:ALA:HB2	1:A:99:GLU:OE1	2.07	0.54
1:A:42:THR:HG22	1:A:117:ASN:N	2.21	0.54
1:A:141:GLN:C	1:A:143:HIS:H	2.11	0.54
1:B:193:ILE:HG21	1:B:207:SER:HB3	1.87	0.54
1:C:144:ARG:HG3	1:C:144:ARG:HH11	1.71	0.54
1:C:165:LEU:HD23	2:C:277:HOH:O	2.08	0.54
1:B:141:GLN:C	1:B:143:HIS:H	2.11	0.54
1:A:151:CYS:O	1:A:152:TYR:C	2.46	0.54
1:C:151:CYS:O	1:C:152:TYR:C	2.46	0.54
1:B:57:ILE:HG13	1:B:80:ASN:OD1	2.06	0.54
1:A:157:PHE:HA	1:A:161:TRP:O	2.07	0.54
1:D:26:TRP:HB2	1:D:80:ASN:CB	2.37	0.54
1:B:63:ASP:HB3	1:B:64:PRO:HD3	1.89	0.54
1:C:135:TRP:CZ3	1:C:183:LYS:HG2	2.43	0.54
1:B:155:THR:HB	1:B:159:PHE:CD1	2.43	0.54
1:D:135:TRP:CZ3	1:D:183:LYS:HG2	2.41	0.53
1:D:155:THR:CG2	1:D:181:SER:HB2	2.39	0.53
1:D:93:LEU:HA	1:D:96:ASN:HD22	1.73	0.53
1:C:156:SER:C	1:C:158:THR:H	2.11	0.53
1:D:85:ASN:C	1:D:86:LYS:HD2	2.29	0.53
1:B:25:ALA:HB2	1:B:99:GLU:OE1	2.08	0.53
1:D:72:GLN:HG2	1:D:259:SER:OG	2.08	0.53
1:B:44:PRO:HB3	1:B:120:GLN:HB3	1.90	0.53
1:D:157:PHE:HA	1:D:161:TRP:O	2.09	0.52
1:B:180:ILE:HD12	1:B:180:ILE:C	2.29	0.52
1:C:25:ALA:HB2	1:C:99:GLU:OE1	2.08	0.52
1:C:237:ASN:C	1:C:239:THR:H	2.11	0.52
1:D:53:GLY:C	1:D:55:GLN:H	2.12	0.52
1:A:149:ILE:O	1:A:150:LEU:HD23	2.09	0.52
1:B:91:ASN:ND2	2:B:288:HOH:O	2.42	0.52
1:A:269:LEU:O	1:A:270:GLU:HG2	2.10	0.52
1:B:15:ASN:HA	1:B:31:GLN:O	2.10	0.52
1:A:242:PRO:O	1:D:41:TYR:HE1	1.92	0.52
1:C:141:GLN:C	1:C:143:HIS:H	2.12	0.52
1:A:144:ARG:HH11	1:A:144:ARG:HG3	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:204:ILE:O	1:D:205:TYR:C	2.48	0.52
1:A:135:TRP:CZ3	1:A:183:LYS:HG2	2.44	0.52
1:A:180:ILE:C	1:A:180:ILE:HD12	2.30	0.52
1:D:141:GLN:HG3	1:D:142:ILE:HG13	1.92	0.52
1:C:105:SER:O	1:C:109:GLN:HG2	2.09	0.52
1:C:155:THR:HB	1:C:159:PHE:CD1	2.45	0.52
1:B:141:GLN:HG3	1:B:142:ILE:HG13	1.92	0.52
1:B:53:GLY:C	1:B:55:GLN:H	2.11	0.52
1:C:56:HIS:O	1:C:57:ILE:HD13	2.10	0.51
1:A:26:TRP:HB2	1:A:80:ASN:CB	2.39	0.51
1:D:33:GLN:NE2	1:D:34:ASN:N	2.58	0.51
1:D:124:LEU:HD12	1:D:125:GLY:N	2.22	0.51
1:D:144:ARG:HH11	1:D:144:ARG:HG3	1.75	0.51
1:B:26:TRP:HB2	1:B:80:ASN:CB	2.38	0.51
1:C:15:ASN:HA	1:C:31:GLN:O	2.10	0.51
1:B:156:SER:C	1:B:158:THR:H	2.13	0.51
1:D:237:ASN:C	1:D:239:THR:H	2.14	0.51
1:B:151:CYS:O	1:B:152:TYR:C	2.49	0.51
1:D:149:ILE:O	1:D:150:LEU:HD23	2.10	0.51
1:B:135:TRP:CZ3	1:B:183:LYS:HG2	2.43	0.51
1:C:208:LEU:HD23	1:D:208:LEU:CD2	2.36	0.51
1:B:93:LEU:HA	1:B:96:ASN:HD22	1.76	0.51
1:A:167:HIS:C	1:A:169:ASN:N	2.64	0.51
1:C:167:HIS:C	1:C:169:ASN:N	2.64	0.51
1:A:141:GLN:HG3	1:A:142:ILE:HG13	1.92	0.51
1:C:72:GLN:HG2	1:C:259:SER:OG	2.10	0.51
1:C:44:PRO:HB3	1:C:120:GLN:HB3	1.94	0.50
1:C:32:LEU:HG	1:C:33:GLN:N	2.25	0.50
1:C:51:GLY:HA2	1:C:129:GLY:H	1.76	0.50
1:C:141:GLN:HG3	1:C:142:ILE:HG13	1.92	0.50
1:D:114:TRP:HB3	1:D:116:ILE:HG13	1.93	0.50
1:C:24:THR:O	1:C:82:THR:HB	2.11	0.50
1:B:167:HIS:C	1:B:169:ASN:N	2.64	0.50
1:A:156:SER:C	1:A:158:THR:H	2.14	0.50
1:A:15:ASN:HA	1:A:31:GLN:O	2.10	0.50
1:B:105:SER:O	1:B:109:GLN:HG2	2.12	0.50
1:A:124:LEU:HD12	1:A:125:GLY:N	2.25	0.50
1:B:32:LEU:HG	1:B:33:GLN:N	2.25	0.50
1:B:33:GLN:NE2	1:B:34:ASN:N	2.59	0.50
1:B:157:PHE:HA	1:B:161:TRP:O	2.12	0.50
1:B:114:TRP:HB3	1:B:116:ILE:HG13	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:LYS:HE3	2:B:289:HOH:O	2.11	0.50
1:A:105:SER:O	1:A:109:GLN:HG2	2.12	0.49
1:D:127:SER:HA	1:D:152:TYR:O	2.11	0.49
1:B:266:ILE:O	1:B:269:LEU:HB2	2.12	0.49
1:B:193:ILE:CG2	1:B:207:SER:HB3	2.42	0.49
1:B:77:LEU:CD1	1:B:107:ILE:HD11	2.43	0.49
1:B:24:THR:O	1:B:82:THR:HB	2.13	0.49
1:B:212:ASP:C	1:B:214:LEU:H	2.16	0.49
1:C:130:GLY:O	1:C:133:ALA:HB3	2.13	0.49
1:C:208:LEU:CD2	1:D:208:LEU:HD23	2.37	0.49
1:B:85:ASN:C	1:B:86:LYS:HD2	2.33	0.49
1:C:83:VAL:HG23	1:C:85:ASN:HB2	1.94	0.49
1:D:27:VAL:HG22	1:D:79:LEU:CD1	2.43	0.49
1:C:27:VAL:HG22	1:C:79:LEU:CD1	2.43	0.49
1:B:204:ILE:O	1:B:205:TYR:C	2.50	0.49
1:D:97:LEU:HD11	1:D:135:TRP:CE2	2.48	0.49
1:A:53:GLY:C	1:A:55:GLN:H	2.15	0.49
1:A:72:GLN:HG2	1:A:259:SER:OG	2.13	0.49
1:C:26:TRP:HB2	1:C:80:ASN:CB	2.38	0.49
1:D:167:HIS:C	1:D:169:ASN:N	2.66	0.49
1:A:155:THR:HA	1:A:180:ILE:HD11	1.95	0.49
1:D:180:ILE:HD12	1:D:180:ILE:C	2.34	0.48
1:A:114:TRP:HB3	1:A:116:ILE:HG13	1.93	0.48
1:A:44:PRO:HB3	1:A:120:GLN:HB3	1.95	0.48
1:D:105:SER:O	1:D:109:GLN:HG2	2.12	0.48
1:A:51:GLY:HA2	1:A:129:GLY:H	1.78	0.48
1:B:124:LEU:HD12	1:B:125:GLY:N	2.26	0.48
1:D:83:VAL:HG23	1:D:85:ASN:HB2	1.95	0.48
1:C:157:PHE:HA	1:C:161:TRP:O	2.13	0.48
1:B:96:ASN:HA	1:B:99:GLU:OE2	2.13	0.48
1:C:114:TRP:HB3	1:C:116:ILE:HG13	1.95	0.48
1:A:237:ASN:C	1:A:239:THR:H	2.15	0.48
1:D:24:THR:O	1:D:82:THR:HB	2.14	0.48
1:D:21:LEU:HD22	1:D:27:VAL:CG2	2.44	0.48
1:A:211:CYS:HB2	1:B:208:LEU:HD21	1.96	0.48
1:D:86:LYS:HD2	1:D:86:LYS:N	2.29	0.48
1:A:85:ASN:C	1:A:86:LYS:HD2	2.34	0.48
1:D:15:ASN:HA	1:D:31:GLN:O	2.14	0.48
1:D:212:ASP:C	1:D:214:LEU:H	2.17	0.48
1:C:190:PRO:HG3	1:C:220:PRO:HG2	1.95	0.48
1:A:165:LEU:O	1:A:169:ASN:ND2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:THR:O	1:A:82:THR:HB	2.13	0.48
1:A:208:LEU:HD21	1:B:211:CYS:HB2	1.94	0.48
1:C:124:LEU:HD12	1:C:125:GLY:N	2.28	0.48
1:C:127:SER:HA	1:C:152:TYR:O	2.13	0.47
1:B:42:THR:HA	1:B:115:GLN:O	2.14	0.47
1:C:83:VAL:C	1:C:85:ASN:H	2.18	0.47
1:C:204:ILE:O	1:C:205:TYR:C	2.52	0.47
1:C:86:LYS:HD2	1:C:86:LYS:N	2.29	0.47
1:D:51:GLY:HA2	1:D:129:GLY:H	1.78	0.47
1:C:53:GLY:C	1:C:55:GLN:H	2.16	0.47
1:B:237:ASN:C	1:B:239:THR:H	2.17	0.47
1:B:150:LEU:HB3	1:B:153:PRO:HB3	1.96	0.47
1:C:150:LEU:HB3	1:C:153:PRO:HB3	1.95	0.47
1:D:114:TRP:CE3	1:D:114:TRP:HA	2.50	0.47
1:C:237:ASN:C	1:C:239:THR:N	2.68	0.47
1:B:47:ILE:N	1:B:47:ILE:HD12	2.29	0.47
1:A:212:ASP:C	1:A:214:LEU:H	2.16	0.47
1:C:85:ASN:C	1:C:86:LYS:HD2	2.35	0.47
1:C:114:TRP:CE3	1:C:114:TRP:HA	2.49	0.47
1:A:237:ASN:HD22	1:A:237:ASN:HA	1.55	0.47
1:A:155:THR:CG2	1:A:181:SER:HB2	2.44	0.47
1:D:236:ALA:HB1	1:D:248:LEU:O	2.14	0.47
1:B:90:TYR:HB3	2:B:278:HOH:O	2.15	0.47
1:A:97:LEU:HD11	1:A:135:TRP:CE2	2.49	0.47
1:B:27:VAL:HG22	1:B:79:LEU:CD1	2.44	0.47
1:D:148:VAL:HG22	1:D:191:THR:CG2	2.37	0.47
1:B:86:LYS:N	1:B:86:LYS:HD2	2.30	0.47
1:D:83:VAL:C	1:D:85:ASN:H	2.19	0.47
1:D:155:THR:HB	1:D:159:PHE:CD1	2.50	0.47
1:D:126:CYS:HA	1:D:151:CYS:O	2.15	0.47
1:D:150:LEU:HB3	1:D:153:PRO:HB3	1.96	0.47
1:D:44:PRO:HB3	1:D:120:GLN:HB3	1.95	0.47
1:D:56:HIS:O	1:D:57:ILE:HD13	2.15	0.46
1:D:96:ASN:HA	1:D:99:GLU:OE2	2.15	0.46
1:D:190:PRO:HG3	1:D:220:PRO:HG2	1.97	0.46
1:C:212:ASP:C	1:C:214:LEU:H	2.18	0.46
1:B:97:LEU:HD11	1:B:135:TRP:CE2	2.50	0.46
1:A:96:ASN:HA	1:A:99:GLU:OE2	2.15	0.46
1:B:155:THR:CG2	1:B:181:SER:HB2	2.45	0.46
1:A:21:LEU:HD22	1:A:27:VAL:CG2	2.44	0.46
1:A:98:GLU:O	1:A:101:GLN:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:LEU:HB3	1:A:153:PRO:HB3	1.97	0.46
1:B:72:GLN:HG2	1:B:259:SER:OG	2.15	0.46
1:A:83:VAL:HG23	1:A:85:ASN:HB2	1.98	0.46
1:C:114:TRP:HA	1:C:114:TRP:HE3	1.80	0.46
1:B:43:PHE:HD2	1:B:75:GLN:HE22	1.63	0.46
1:D:114:TRP:HE3	1:D:114:TRP:HA	1.81	0.46
1:C:156:SER:OG	1:C:179:ASN:HB2	2.16	0.46
1:C:77:LEU:CD1	1:C:107:ILE:HD11	2.45	0.46
1:D:26:TRP:CB	2:D:301:HOH:O	2.63	0.46
1:B:156:SER:OG	1:B:179:ASN:HB2	2.15	0.46
1:A:193:ILE:HG21	1:A:207:SER:HB3	1.97	0.46
1:D:165:LEU:O	1:D:169:ASN:ND2	2.48	0.46
1:A:86:LYS:HD2	1:A:86:LYS:N	2.31	0.46
1:A:254:TRP:CG	1:A:255:VAL:N	2.84	0.46
1:B:210:TYR:O	1:B:214:LEU:HB2	2.15	0.46
1:C:21:LEU:HD22	1:C:27:VAL:CG2	2.46	0.46
1:B:67:LEU:O	1:B:68:ALA:C	2.52	0.46
1:A:221:PHE:O	1:B:227:GLU:HA	2.16	0.46
1:A:97:LEU:HD21	1:A:135:TRP:CD1	2.51	0.46
1:C:97:LEU:HD11	1:C:135:TRP:CE2	2.50	0.46
1:D:47:ILE:N	1:D:47:ILE:HD12	2.31	0.46
1:B:56:HIS:O	1:B:57:ILE:HD13	2.15	0.46
1:A:33:GLN:NE2	1:A:34:ASN:N	2.64	0.46
1:C:33:GLN:NE2	1:C:34:ASN:N	2.64	0.46
1:D:155:THR:HA	1:D:180:ILE:HD11	1.97	0.46
1:B:114:TRP:HA	1:B:114:TRP:CE3	2.51	0.46
1:A:127:SER:HA	1:A:152:TYR:O	2.16	0.45
1:C:155:THR:CG2	1:C:181:SER:HB2	2.45	0.45
1:C:96:ASN:HA	1:C:99:GLU:OE2	2.16	0.45
1:A:27:VAL:HG22	1:A:79:LEU:CD1	2.45	0.45
1:A:56:HIS:O	1:A:57:ILE:HD13	2.16	0.45
1:B:83:VAL:HG23	1:B:85:ASN:HB2	1.97	0.45
1:C:155:THR:HA	1:C:180:ILE:HD11	1.97	0.45
1:D:43:PHE:HB2	1:D:75:GLN:NE2	2.31	0.45
1:A:114:TRP:HA	1:A:114:TRP:CE3	2.51	0.45
1:B:268:ASN:O	1:B:269:LEU:HD12	2.16	0.45
1:D:130:GLY:O	1:D:133:ALA:HB3	2.17	0.45
1:A:130:GLY:O	1:A:133:ALA:HB3	2.16	0.45
1:A:155:THR:HB	1:A:159:PHE:CD1	2.51	0.45
1:C:97:LEU:HD21	1:C:135:TRP:CD1	2.52	0.45
1:C:193:ILE:HG21	1:C:207:SER:HB3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:PHE:CE1	1:A:178:TYR:HB3	2.51	0.45
1:B:43:PHE:HB2	1:B:75:GLN:NE2	2.32	0.45
1:B:114:TRP:HA	1:B:114:TRP:HE3	1.82	0.45
1:A:190:PRO:HG3	1:A:220:PRO:HG2	1.97	0.45
1:C:241:ALA:HB2	1:C:246:TYR:O	2.17	0.45
1:D:266:ILE:O	1:D:269:LEU:HB2	2.16	0.45
1:B:51:GLY:HA2	1:B:129:GLY:H	1.81	0.45
1:A:223:ALA:HB3	1:B:225:PHE:HB2	1.98	0.45
1:A:43:PHE:HD2	1:A:75:GLN:HE22	1.65	0.45
1:C:204:ILE:HD11	1:D:215:SER:OG	2.16	0.45
1:A:83:VAL:C	1:A:85:ASN:H	2.20	0.45
1:B:98:GLU:O	1:B:101:GLN:HB3	2.16	0.45
1:D:237:ASN:C	1:D:239:THR:N	2.70	0.44
1:D:33:GLN:CD	1:D:34:ASN:N	2.71	0.44
1:D:97:LEU:HD21	1:D:135:TRP:CD1	2.52	0.44
1:A:67:LEU:O	1:A:68:ALA:C	2.54	0.44
1:A:47:ILE:N	1:A:47:ILE:HD12	2.32	0.44
1:A:210:TYR:O	1:A:214:LEU:HB2	2.17	0.44
1:B:91:ASN:C	1:B:93:LEU:H	2.21	0.44
1:D:43:PHE:HD2	1:D:75:GLN:HE22	1.65	0.44
1:D:128:ALA:HA	1:D:154:VAL:CG2	2.48	0.44
1:B:33:GLN:CD	1:B:34:ASN:N	2.70	0.44
1:C:165:LEU:O	1:C:169:ASN:ND2	2.50	0.44
1:A:156:SER:OG	1:A:179:ASN:HB2	2.17	0.44
1:A:32:LEU:HG	1:A:33:GLN:N	2.29	0.44
1:D:32:LEU:HG	1:D:33:GLN:N	2.26	0.44
1:C:168:PHE:HZ	2:C:290:HOH:O	2.00	0.44
1:A:93:LEU:HD12	1:A:178:TYR:CE1	2.53	0.44
1:D:192:PHE:O	1:D:193:ILE:HD12	2.18	0.44
1:A:128:ALA:HA	1:A:154:VAL:CG2	2.48	0.44
1:D:27:VAL:HG22	1:D:79:LEU:HD12	2.00	0.44
1:C:98:GLU:O	1:C:101:GLN:HB3	2.18	0.44
1:D:195:HIS:CE1	1:D:204:ILE:HA	2.52	0.44
1:C:236:ALA:HB1	1:C:248:LEU:O	2.17	0.44
1:B:138:ASN:O	1:B:139:SER:HB2	2.17	0.44
1:A:114:TRP:HA	1:A:114:TRP:HE3	1.82	0.43
1:C:43:PHE:HB2	1:C:75:GLN:NE2	2.33	0.43
1:C:128:ALA:HA	1:C:154:VAL:CG2	2.48	0.43
1:B:21:LEU:HD22	1:B:27:VAL:CG2	2.48	0.43
1:A:204:ILE:O	1:A:205:TYR:C	2.56	0.43
1:C:223:ALA:HB3	1:D:225:PHE:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:THR:HG22	1:D:19:PHE:N	2.33	0.43
1:A:77:LEU:CD1	1:A:107:ILE:HD11	2.49	0.43
1:D:77:LEU:CD1	1:D:107:ILE:HD11	2.48	0.43
1:A:18:THR:HG22	1:A:19:PHE:N	2.33	0.43
1:C:202:VAL:O	1:C:202:VAL:HG13	2.18	0.43
1:A:202:VAL:HG23	1:A:206:ASN:ND2	2.33	0.43
1:C:241:ALA:HA	1:C:242:PRO:HD2	1.86	0.43
1:C:67:LEU:O	1:C:68:ALA:C	2.56	0.43
1:A:237:ASN:C	1:A:239:THR:N	2.72	0.43
1:A:241:ALA:HA	1:A:242:PRO:HD2	1.86	0.43
1:B:165:LEU:O	1:B:169:ASN:ND2	2.51	0.43
1:C:237:ASN:HA	1:C:237:ASN:HD22	1.51	0.43
1:C:43:PHE:HD2	1:C:75:GLN:HE22	1.65	0.43
1:D:67:LEU:O	1:D:68:ALA:C	2.55	0.43
1:D:34:ASN:CG	1:D:34:ASN:O	2.57	0.43
1:C:144:ARG:HG3	1:C:144:ARG:NH1	2.34	0.43
1:B:28:ASP:HB2	1:B:78:LEU:HB3	2.00	0.43
1:C:18:THR:HG22	1:C:19:PHE:N	2.33	0.43
1:A:202:VAL:CG2	1:A:206:ASN:ND2	2.82	0.43
1:D:67:LEU:HD21	2:D:286:HOH:O	2.18	0.43
1:B:182:GLU:C	1:B:184:VAL:H	2.21	0.43
1:C:49:CYS:N	1:C:124:LEU:O	2.52	0.42
1:C:92:PHE:C	1:C:94:SER:N	2.72	0.42
1:D:98:GLU:O	1:D:101:GLN:HB3	2.19	0.42
1:A:28:ASP:HB2	1:A:78:LEU:HB3	2.01	0.42
1:D:138:ASN:O	1:D:139:SER:HB2	2.18	0.42
1:B:92:PHE:O	1:B:95:GLN:N	2.52	0.42
1:A:126:CYS:O	1:A:127:SER:C	2.58	0.42
1:B:155:THR:HA	1:B:180:ILE:HD11	2.01	0.42
1:C:27:VAL:HG22	1:C:79:LEU:HD12	2.01	0.42
1:C:138:ASN:O	1:C:139:SER:HB2	2.19	0.42
1:A:42:THR:HA	1:A:115:GLN:O	2.20	0.42
1:B:237:ASN:C	1:B:239:THR:N	2.72	0.42
1:B:18:THR:HG22	1:B:19:PHE:N	2.35	0.42
1:B:117:ASN:HD22	1:B:120:GLN:HB2	1.85	0.42
1:C:91:ASN:C	1:C:93:LEU:H	2.22	0.42
1:C:107:ILE:O	1:C:111:HIS:N	2.52	0.42
1:B:198:ASP:OD1	1:B:229:GLY:N	2.52	0.42
1:C:127:SER:HG	1:C:231:HIS:CE1	2.37	0.42
1:C:268:ASN:ND2	1:C:268:ASN:O	2.52	0.42
1:A:138:ASN:O	1:A:139:SER:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:ASP:HB2	1:C:78:LEU:HB3	2.02	0.42
1:D:156:SER:HB3	1:D:159:PHE:CD1	2.54	0.42
1:B:190:PRO:HG3	1:B:220:PRO:HG2	2.01	0.42
1:C:195:HIS:CE1	1:C:204:ILE:HA	2.55	0.42
1:C:33:GLN:CD	1:C:34:ASN:N	2.73	0.42
1:D:42:THR:HA	1:D:115:GLN:O	2.20	0.42
1:B:130:GLY:O	1:B:133:ALA:HB3	2.19	0.42
1:B:235:LEU:HB2	1:B:240:THR:OG1	2.19	0.42
1:D:128:ALA:C	2:D:290:HOH:O	2.58	0.42
1:B:202:VAL:O	1:B:202:VAL:HG13	2.20	0.42
1:C:92:PHE:O	1:C:95:GLN:N	2.53	0.42
1:B:127:SER:HA	1:B:152:TYR:O	2.19	0.41
1:B:193:ILE:O	1:B:223:ALA:HA	2.20	0.41
1:C:161:TRP:HA	1:C:162:PRO:C	2.40	0.41
1:D:210:TYR:O	1:D:214:LEU:HB2	2.20	0.41
1:B:83:VAL:C	1:B:85:ASN:H	2.22	0.41
1:C:210:TYR:O	1:C:214:LEU:HB2	2.20	0.41
1:B:97:LEU:HD21	1:B:135:TRP:CD1	2.54	0.41
1:C:165:LEU:HA	2:C:277:HOH:O	2.20	0.41
1:A:225:PHE:HD1	1:B:223:ALA:HB3	1.85	0.41
1:A:83:VAL:HG22	1:A:86:LYS:HE3	2.01	0.41
1:C:254:TRP:CG	1:C:255:VAL:N	2.88	0.41
1:C:221:PHE:O	1:D:227:GLU:HA	2.19	0.41
1:A:32:LEU:HD23	1:D:239:THR:HG23	2.02	0.41
1:C:157:PHE:CE1	1:C:178:TYR:HB3	2.56	0.41
1:C:225:PHE:HB2	1:D:223:ALA:HB3	2.03	0.41
1:C:30:TYR:HB2	1:C:76:VAL:HB	2.01	0.41
1:C:126:CYS:HA	1:C:151:CYS:O	2.20	0.41
1:C:205:TYR:O	1:C:206:ASN:C	2.58	0.41
1:A:212:ASP:HA	1:B:208:LEU:HD13	2.02	0.41
1:D:21:LEU:HD22	1:D:27:VAL:HG21	2.01	0.41
1:C:47:ILE:HD12	1:C:47:ILE:N	2.36	0.41
1:C:127:SER:OG	1:C:231:HIS:CE1	2.73	0.41
1:D:193:ILE:HG21	1:D:207:SER:HB3	2.03	0.41
1:D:83:VAL:HG22	1:D:86:LYS:HE3	2.02	0.41
1:C:227:GLU:OE2	1:D:220:PRO:HB2	2.21	0.41
1:A:193:ILE:O	1:A:223:ALA:HA	2.20	0.41
1:B:132:LEU:O	1:B:135:TRP:HB3	2.19	0.41
1:C:159:PHE:C	1:C:161:TRP:H	2.24	0.41
1:B:157:PHE:CE1	1:B:178:TYR:HB3	2.56	0.41
1:D:159:PHE:C	1:D:161:TRP:H	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:GLY:C	1:B:55:GLN:N	2.74	0.41
1:D:191:THR:O	1:D:221:PHE:HB2	2.21	0.41
1:A:195:HIS:CE1	1:A:204:ILE:HA	2.55	0.41
1:B:188:THR:HG23	1:B:189:PRO:HD2	2.02	0.41
1:C:193:ILE:O	1:C:223:ALA:HA	2.21	0.41
1:B:93:LEU:HD12	1:B:178:TYR:CE1	2.56	0.41
1:C:62:SER:O	1:C:63:ASP:C	2.58	0.41
1:B:144:ARG:NH1	1:B:144:ARG:HG3	2.35	0.41
1:A:27:VAL:HG22	1:A:79:LEU:HD12	2.03	0.41
1:D:92:PHE:O	1:D:95:GLN:N	2.54	0.41
1:C:182:GLU:C	1:C:184:VAL:H	2.24	0.41
1:D:28:ASP:HB2	1:D:78:LEU:HB3	2.03	0.41
1:B:59:GLN:C	1:B:61:GLU:H	2.25	0.41
1:D:62:SER:O	1:D:63:ASP:C	2.59	0.41
1:C:83:VAL:HG22	1:C:86:LYS:HE3	2.03	0.41
1:D:91:ASN:C	1:D:93:LEU:H	2.23	0.41
1:D:92:PHE:C	1:D:94:SER:N	2.74	0.41
1:C:127:SER:C	1:C:154:VAL:HG23	2.42	0.40
1:C:193:ILE:CG2	1:C:207:SER:HB3	2.51	0.40
1:B:135:TRP:CE3	1:B:135:TRP:HA	2.56	0.40
1:D:53:GLY:C	1:D:55:GLN:N	2.75	0.40
1:C:211:CYS:O	1:C:214:LEU:HB3	2.21	0.40
1:A:182:GLU:C	1:A:184:VAL:H	2.25	0.40
1:A:92:PHE:O	1:A:95:GLN:N	2.54	0.40
1:D:127:SER:C	1:D:154:VAL:HG23	2.41	0.40
1:A:151:CYS:SG	1:A:194:TRP:NE1	2.91	0.40
1:C:215:SER:OG	1:D:204:ILE:HD11	2.20	0.40
1:A:33:GLN:CD	1:A:34:ASN:N	2.75	0.40
1:A:135:TRP:HA	1:A:135:TRP:HE3	1.86	0.40
1:B:161:TRP:HA	1:B:162:PRO:C	2.42	0.40
1:A:117:ASN:HD22	1:A:120:GLN:HB2	1.86	0.40
1:A:240:THR:O	1:A:241:ALA:HB2	2.21	0.40
1:D:235:LEU:HB2	1:D:240:THR:OG1	2.20	0.40
1:C:266:ILE:O	1:C:269:LEU:HB2	2.21	0.40
1:B:189:PRO:O	1:B:191:THR:HG23	2.21	0.40
1:B:21:LEU:HD21	1:B:99:GLU:O	2.21	0.40
1:A:91:ASN:C	1:A:93:LEU:H	2.23	0.40
1:B:254:TRP:CG	1:B:255:VAL:N	2.90	0.40
1:D:107:ILE:O	1:D:111:HIS:N	2.55	0.40
1:C:221:PHE:N	1:D:227:GLU:HG3	2.37	0.40
1:D:156:SER:OG	1:D:179:ASN:HB2	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:PHE:C	1:A:161:TRP:H	2.24	0.40
1:C:146:LYS:N	1:C:146:LYS:HD2	2.37	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	246/276 (89%)	180 (73%)	56 (23%)	10 (4%)	3	27
1	B	246/276 (89%)	179 (73%)	56 (23%)	11 (4%)	3	24
1	C	246/276 (89%)	182 (74%)	52 (21%)	12 (5%)	3	22
1	D	246/276 (89%)	179 (73%)	58 (24%)	9 (4%)	4	29
All	All	984/1104 (89%)	720 (73%)	222 (23%)	42 (4%)	3	25

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	MSE
1	A	171	GLU
1	A	174	ASN
1	A	198	ASP
1	B	171	GLU
1	B	174	ASN
1	B	198	ASP
1	C	84	MSE
1	C	171	GLU
1	C	174	ASN
1	C	198	ASP
1	D	171	GLU
1	D	174	ASN

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Mol	Chain	Res	Type
1	D	198	ASP
1	A	183	LYS
1	B	84	MSE
1	B	183	LYS
1	C	183	LYS
1	D	84	MSE
1	D	183	LYS
1	B	25	ALA
1	C	25	ALA
1	C	138	ASN
1	A	25	ALA
1	A	138	ASN
1	A	142	ILE
1	A	172	ILE
1	B	138	ASN
1	B	142	ILE
1	B	172	ILE
1	B	213	ARG
1	C	58	SER
1	C	142	ILE
1	C	172	ILE
1	D	138	ASN
1	D	142	ILE
1	D	172	ILE
1	D	160	GLY
1	A	52	GLY
1	B	52	GLY
1	C	52	GLY
1	C	160	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	221/244 (91%)	200 (90%)	21 (10%)	11	40
1	B	221/244 (91%)	201 (91%)	20 (9%)	12	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	221/244 (91%)	198 (90%)	23 (10%)	9	35
1	D	221/244 (91%)	201 (91%)	20 (9%)	12	43
All	All	884/976 (91%)	800 (90%)	84 (10%)	11	40

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	THR
1	A	34	ASN
1	A	42	THR
1	A	46	ILE
1	A	55	GLN
1	A	65	LEU
1	A	75	GLN
1	A	85	ASN
1	A	86	LYS
1	A	114	TRP
1	A	126	CYS
1	A	135	TRP
1	A	138	ASN
1	A	168	PHE
1	A	182	GLU
1	A	204	ILE
1	A	212	ASP
1	A	237	ASN
1	A	257	TRP
1	A	260	ASP
1	A	268	ASN
1	B	24	THR
1	B	34	ASN
1	B	42	THR
1	B	46	ILE
1	B	55	GLN
1	B	65	LEU
1	B	75	GLN
1	B	85	ASN
1	B	86	LYS
1	B	114	TRP
1	B	126	CYS
1	B	135	TRP
1	B	138	ASN

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Mol	Chain	Res	Type
1	B	168	PHE
1	B	182	GLU
1	B	204	ILE
1	B	212	ASP
1	B	237	ASN
1	B	257	TRP
1	B	260	ASP
1	C	24	THR
1	C	34	ASN
1	C	42	THR
1	C	46	ILE
1	C	55	GLN
1	C	65	LEU
1	C	75	GLN
1	C	85	ASN
1	C	86	LYS
1	C	114	TRP
1	C	126	CYS
1	C	135	TRP
1	C	138	ASN
1	C	168	PHE
1	C	182	GLU
1	C	204	ILE
1	C	208	LEU
1	C	212	ASP
1	C	237	ASN
1	C	257	TRP
1	C	260	ASP
1	C	268	ASN
1	C	269	LEU
1	D	24	THR
1	D	34	ASN
1	D	42	THR
1	D	46	ILE
1	D	55	GLN
1	D	65	LEU
1	D	75	GLN
1	D	85	ASN
1	D	86	LYS
1	D	114	TRP
1	D	126	CYS
1	D	135	TRP

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Mol	Chain	Res	Type
1	D	138	ASN
1	D	168	PHE
1	D	182	GLU
1	D	204	ILE
1	D	212	ASP
1	D	237	ASN
1	D	257	TRP
1	D	260	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	33	GLN
1	A	34	ASN
1	A	95	GLN
1	A	96	ASN
1	A	117	ASN
1	A	195	HIS
1	A	217	HIS
1	A	237	ASN
1	A	265	GLN
1	A	268	ASN
1	B	31	GLN
1	B	33	GLN
1	B	91	ASN
1	B	95	GLN
1	B	96	ASN
1	B	108	HIS
1	B	117	ASN
1	B	217	HIS
1	B	237	ASN
1	B	265	GLN
1	C	31	GLN
1	C	33	GLN
1	C	95	GLN
1	C	96	ASN
1	C	117	ASN
1	C	195	HIS
1	C	217	HIS
1	C	237	ASN
1	C	265	GLN

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Mol	Chain	Res	Type
1	C	268	ASN
1	D	31	GLN
1	D	33	GLN
1	D	85	ASN
1	D	91	ASN
1	D	95	GLN
1	D	96	ASN
1	D	117	ASN
1	D	217	HIS
1	D	237	ASN
1	D	265	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	249/276 (90%)	0.06	5 (2%)	68	54	28, 68, 106, 123	1 (0%)
1	B	249/276 (90%)	0.03	9 (3%)	46	31	33, 69, 105, 121	1 (0%)
1	C	249/276 (90%)	0.22	6 (2%)	62	47	34, 72, 105, 122	1 (0%)
1	D	249/276 (90%)	0.15	11 (4%)	38	24	36, 69, 106, 120	1 (0%)
All	All	996/1104 (90%)	0.12	31 (3%)	52	38	28, 70, 106, 123	4 (0%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	141	GLN	6.0
1	C	90	TYR	4.8
1	D	127	SER	4.7
1	C	166	SER	4.1
1	B	90	TYR	3.7
1	B	173	GLU	3.4
1	D	87	GLY	3.0
1	A	127	SER	3.0
1	B	20	SER	2.9
1	B	178	TYR	2.8
1	D	173	GLU	2.7
1	A	90	TYR	2.6
1	B	125	GLY	2.5
1	C	173	GLU	2.5
1	D	126	CYS	2.5
1	D	90	TYR	2.4
1	D	141	GLN	2.4
1	B	164	ASP	2.4
1	A	169	ASN	2.4
1	D	86	LYS	2.4
1	B	165	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	26	TRP	2.3
1	C	200	GLU	2.3
1	B	127	SER	2.3
1	D	125	GLY	2.3
1	D	34	ASN	2.2
1	C	267	LYS	2.2
1	A	178	TYR	2.2
1	D	163	SER	2.2
1	A	23	ASP	2.1
1	D	245	ALA	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.