



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:17 PM GMT

PDB ID : 1JLH  
Title : Human Glucose-6-phosphate Isomerase  
Authors : Cordeiro, A.T.  
Deposited on : 2001-07-16  
Resolution : 2.10 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](mailto: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.

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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

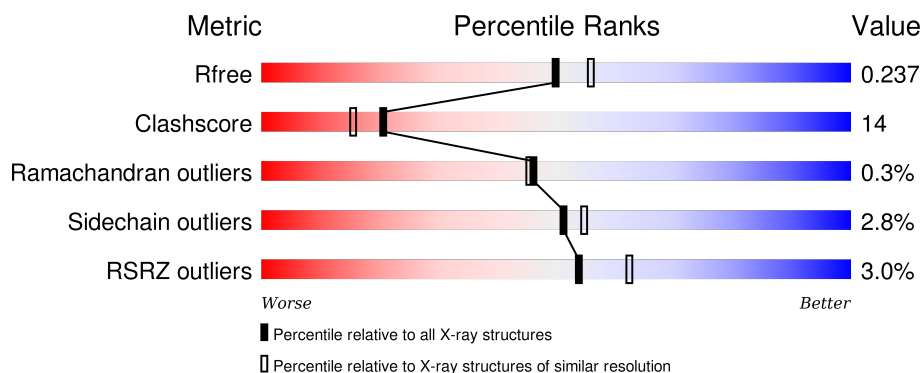
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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  $\geq 3$ , 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	558	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 73%, yellow 73%, yellow 98%, red 98%, red 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>73%</span> <span>25%</span> </div> </div>
1	B	558	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 2%, green 2%, green 71%, yellow 71%, yellow 98%, red 98%, red 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>71%</span> <span>26%</span> </div> </div>
1	C	558	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 5%, green 5%, green 69%, yellow 69%, yellow 98%, red 98%, red 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>69%</span> <span>29%</span> </div> </div>
1	D	558	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, green 3%, green 75%, yellow 75%, yellow 98%, red 98%, red 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>75%</span> <span>23%</span> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 19037 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 phosphoglucose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	556	Total	C	N	O	S	0	0	0
			4437	2826	781	812	18			
1	B	556	Total	C	N	O	S	0	0	0
			4437	2826	781	812	18			
1	C	556	Total	C	N	O	S	0	0	0
			4437	2826	781	812	18			
1	D	556	Total	C	N	O	S	0	0	0
			4437	2826	781	812	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	399	GLU	LYS	see remark 999	UNP P06744
B	399	GLU	LYS	see remark 999	UNP P06744
C	399	GLU	LYS	see remark 999	UNP P06744
D	399	GLU	LYS	see remark 999	UNP P06744

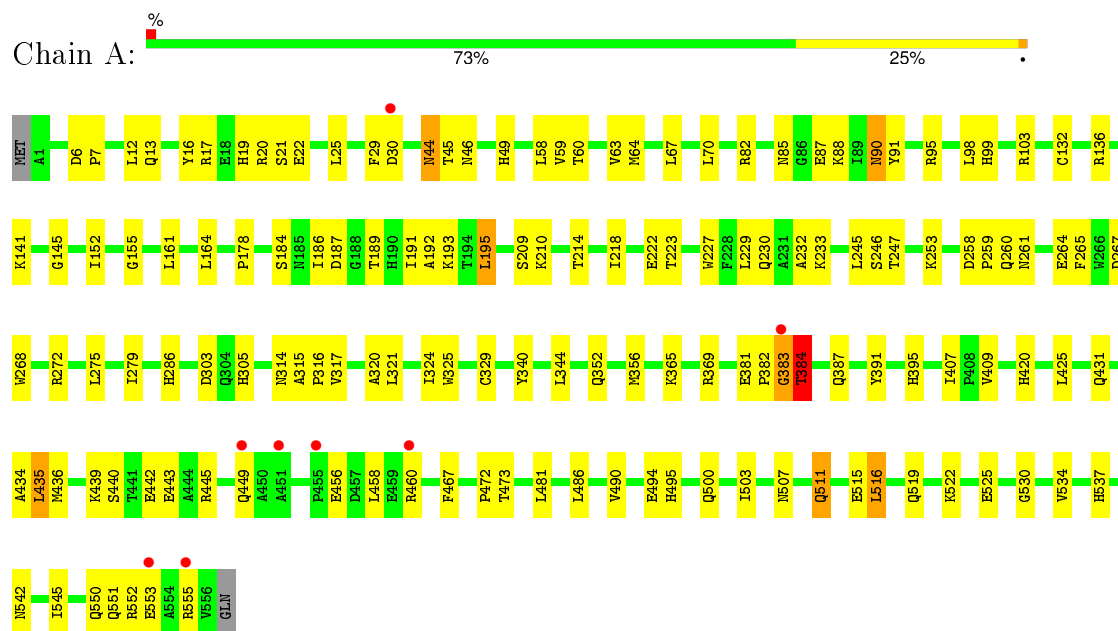
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	350	Total	O	0	0
			350	350		
2	B	352	Total	O	0	0
			352	352		
2	C	275	Total	O	0	0
			275	275		
2	D	312	Total	O	0	0
			312	312		

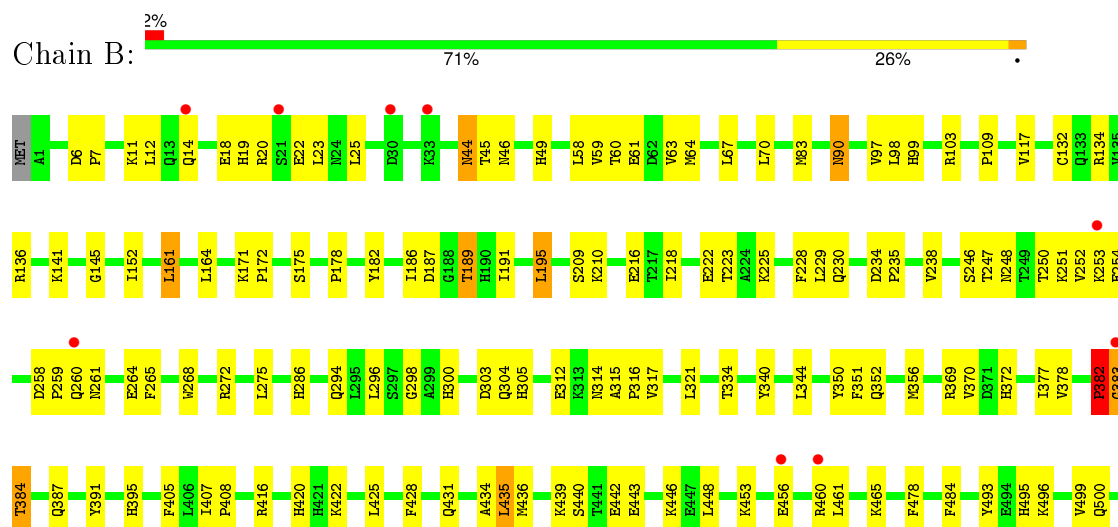
### 3 Residue-property plots

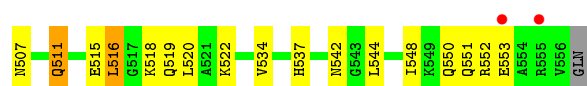
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

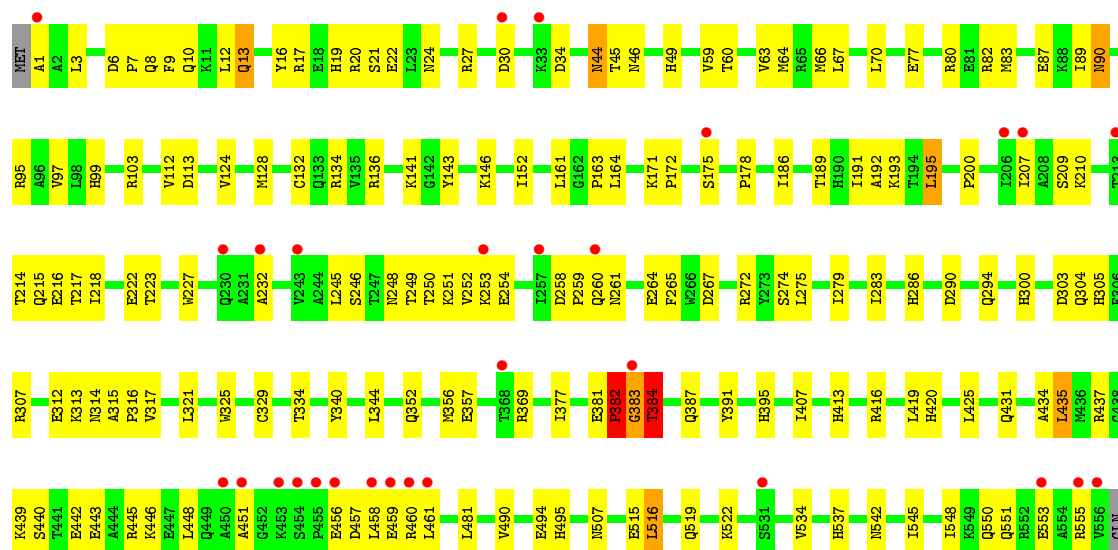


- Molecule 1: phosphoglucose isomerase

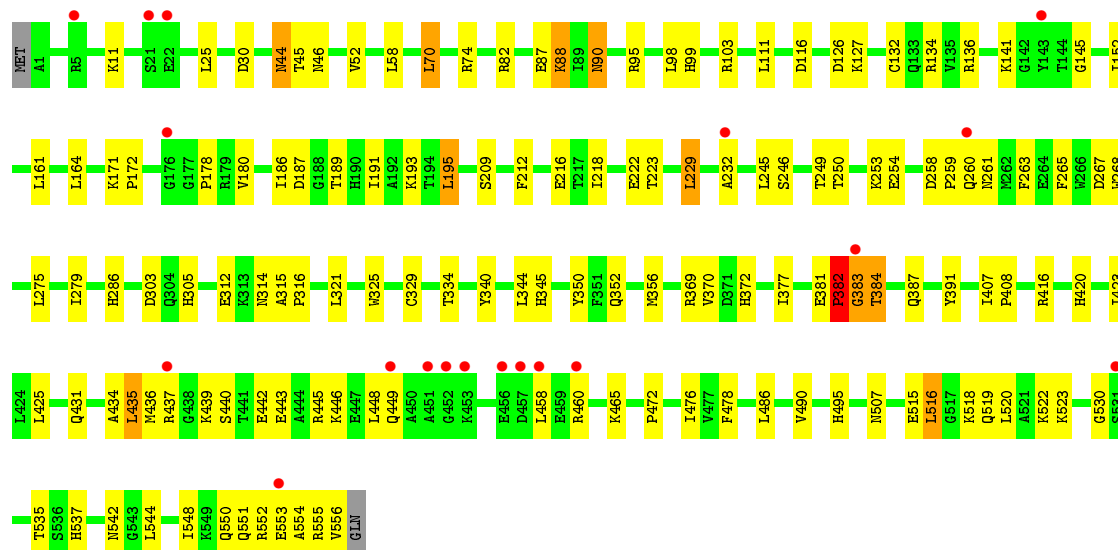
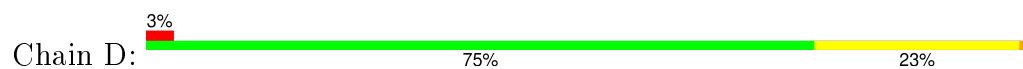




- Molecule 1: phosphoglucose isomerase



- Molecule 1: phosphoglucose isomerase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.75Å 108.02Å 271.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.21 – 2.10 22.21 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.7 (22.21-2.10) 99.7 (22.21-2.10)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.39 (at 2.09Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.197 , 0.237 0.197 , 0.237	Depositor DCC
$R_{free}$ test set	6914 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.8	Xtriage
Anisotropy	0.103	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 54.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	7 of 138364 reflections (0.005%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	19037	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.15 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.3155e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.35	0/4545	0.60	3/6155 (0.0%)
1	B	0.44	2/4545 (0.0%)	0.63	4/6155 (0.1%)
1	C	0.36	2/4545 (0.0%)	0.58	3/6155 (0.0%)
1	D	0.54	1/4545 (0.0%)	0.60	3/6155 (0.0%)
All	All	0.43	5/18180 (0.0%)	0.60	13/24620 (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	B	0	1
1	C	0	1
1	D	0	1
All	All	0	3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	382	PRO	C-O	28.31	1.79	1.23
1	B	382	PRO	CA-C	15.78	1.84	1.52
1	C	382	PRO	CA-C	8.88	1.70	1.52
1	B	382	PRO	C-O	6.66	1.36	1.23
1	C	382	PRO	C-O	6.29	1.35	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	382	PRO	CA-C-N	-10.48	95.23	116.20
1	B	382	PRO	O-C-N	10.07	140.33	123.20
1	C	383	GLY	C-N-CA	7.91	141.47	121.70
1	B	383	GLY	C-N-CA	7.64	140.79	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	382	PRO	CA-C-N	-7.52	101.16	116.20
1	A	383	GLY	C-N-CA	7.51	140.48	121.70
1	D	383	GLY	C-N-CA	7.27	139.88	121.70
1	C	384	THR	N-CA-CB	7.20	123.98	110.30
1	D	384	THR	N-CA-CB	6.79	123.21	110.30
1	A	384	THR	N-CA-CB	6.55	122.75	110.30
1	B	384	THR	N-CA-CB	5.91	121.53	110.30
1	C	382	PRO	CA-C-N	-5.68	104.84	116.20
1	A	383	GLY	CA-C-N	-5.30	105.53	117.20

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	382	PRO	Mainchain
1	C	382	PRO	Mainchain
1	D	382	PRO	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	4437	0	4388	132	0
1	B	4437	0	4388	141	0
1	C	4437	0	4388	156	0
1	D	4437	0	4388	113	0
2	A	350	0	0	13	0
2	B	352	0	0	20	0
2	C	275	0	0	20	0
2	D	312	0	0	13	0
All	All	19037	0	17552	512	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 14.

All (512) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:382:PRO:C	1:B:382:PRO:CA	1.84	1.44
1:D:382:PRO:O	1:D:382:PRO:C	1.79	1.20
1:C:304:GLN:HG3	2:C:558:HOH:O	1.45	1.14
1:B:304:GLN:HG3	2:B:558:HOH:O	1.53	1.05
1:A:550:GLN:HE21	1:A:551:GLN:HE21	1.04	0.97
1:C:134:ARG:HD2	2:C:803:HOH:O	1.67	0.94
1:C:439:LYS:HZ1	1:D:519:GLN:HE22	0.89	0.89
1:B:382:PRO:CA	1:B:383:GLY:N	2.37	0.87
1:A:537:HIS:H	1:A:542:ASN:HD21	1.20	0.87
1:B:537:HIS:H	1:B:542:ASN:HD21	1.23	0.86
1:D:340:TYR:HA	1:D:383:GLY:HA3	1.58	0.85
1:C:305:HIS:HE1	1:C:315:ALA:H	1.25	0.85
1:C:442:GLU:HG2	1:C:446:LYS:HE2	1.60	0.84
1:B:340:TYR:HA	1:B:383:GLY:HA3	1.58	0.84
1:B:11:LYS:HE2	2:B:612:HOH:O	1.75	0.84
1:A:59:VAL:HG21	1:A:64:MET:SD	2.17	0.84
1:C:340:TYR:HA	1:C:383:GLY:HA3	1.59	0.83
1:D:253:LYS:HG2	1:D:259:PRO:HG3	1.61	0.83
1:D:209:SER:O	1:D:246:SER:HB3	1.79	0.81
1:D:550:GLN:HE21	1:D:551:GLN:HE21	1.26	0.80
1:C:439:LYS:HZ1	1:D:519:GLN:NE2	1.75	0.80
1:A:519:GLN:HE22	1:B:439:LYS:NZ	1.80	0.78
1:D:537:HIS:H	1:D:542:ASN:HD21	1.31	0.77
1:D:370:VAL:HG12	1:D:372:HIS:H	1.48	0.76
1:C:90:ASN:ND2	1:C:507:ASN:HD21	1.83	0.76
1:D:250:THR:O	1:D:254:GLU:HG3	1.86	0.76
1:C:59:VAL:HG21	1:C:64:MET:SD	2.25	0.76
1:A:439:LYS:NZ	1:B:519:GLN:HE22	1.84	0.75
1:C:132:CYS:O	1:C:136:ARG:HG3	1.86	0.75
1:C:321:LEU:HG	2:C:639:HOH:O	1.85	0.75
1:C:550:GLN:HE21	1:C:551:GLN:HE21	1.34	0.75
1:B:317:VAL:O	1:B:321:LEU:HD13	1.86	0.75
1:A:272:ARG:HH12	1:A:511:GLN:HE21	1.35	0.75
1:B:132:CYS:O	1:B:136:ARG:HG3	1.87	0.74
1:C:264:GLU:HG2	2:C:564:HOH:O	1.86	0.74
1:B:59:VAL:HG21	1:B:64:MET:SD	2.27	0.74
1:A:59:VAL:HG23	1:A:63:VAL:HB	1.71	0.72
1:A:305:HIS:HE1	1:A:315:ALA:H	1.38	0.71
1:A:90:ASN:ND2	1:A:507:ASN:HD21	1.88	0.71
1:C:209:SER:O	1:C:246:SER:HB3	1.91	0.70
1:C:59:VAL:HG23	1:C:63:VAL:HB	1.71	0.70
1:A:272:ARG:HH12	1:A:511:GLN:NE2	1.89	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:553:GLU:O	1:D:555:ARG:HD2	1.91	0.70
1:B:305:HIS:HE1	1:B:315:ALA:H	1.38	0.70
1:B:315:ALA:HB3	1:B:316:PRO:HD3	1.74	0.70
1:C:537:HIS:H	1:C:542:ASN:HD21	1.38	0.70
1:B:59:VAL:HG23	1:B:63:VAL:HB	1.74	0.69
1:B:550:GLN:HE21	1:B:551:GLN:HE21	1.40	0.69
1:C:317:VAL:O	1:C:321:LEU:HD13	1.92	0.69
1:A:387:GLN:HA	1:A:391:TYR:CD1	2.28	0.69
1:B:300:HIS:O	1:B:304:GLN:HG2	1.93	0.69
1:A:85:ASN:HB3	2:A:703:HOH:O	1.92	0.69
1:A:515:GLU:O	1:A:519:GLN:HG3	1.93	0.68
1:B:90:ASN:ND2	1:B:507:ASN:HD21	1.90	0.68
1:D:90:ASN:ND2	1:D:507:ASN:HD21	1.92	0.68
1:A:445:ARG:O	1:A:449:GLN:HG3	1.93	0.68
1:D:305:HIS:HE1	1:D:315:ALA:H	1.39	0.68
1:A:314:ASN:OD1	1:A:316:PRO:HD2	1.93	0.67
1:C:272:ARG:NH1	1:C:357:GLU:HG2	2.09	0.67
1:A:12:LEU:HD11	1:A:67:LEU:HD23	1.77	0.67
1:C:515:GLU:O	1:C:519:GLN:HG3	1.93	0.67
1:C:387:GLN:HA	1:C:391:TYR:CD1	2.30	0.66
1:A:25:LEU:HB3	1:A:436:MET:HG3	1.77	0.66
1:A:409:VAL:HG22	1:B:548:ILE:HD13	1.76	0.66
1:D:25:LEU:HB3	1:D:436:MET:HG3	1.77	0.66
1:C:459:GLU:HB2	1:C:460:ARG:NH1	2.11	0.66
1:D:387:GLN:HA	1:D:391:TYR:CD1	2.30	0.66
1:A:209:SER:O	1:A:246:SER:HB3	1.97	0.65
1:B:218:ILE:O	1:B:222:GLU:HG3	1.95	0.65
1:A:434:ALA:HB1	1:B:516:LEU:HD13	1.77	0.65
1:B:522:LYS:HE2	1:B:522:LYS:HA	1.79	0.64
1:D:186:ILE:HD13	1:D:216:GLU:HA	1.79	0.64
1:A:340:TYR:HA	1:A:383:GLY:HA3	1.80	0.64
1:D:522:LYS:HE2	1:D:522:LYS:HA	1.77	0.64
1:C:186:ILE:HD13	1:C:216:GLU:HA	1.78	0.64
1:A:522:LYS:HG2	2:A:640:HOH:O	1.96	0.64
1:A:191:ILE:HG13	1:A:195:LEU:HD22	1.79	0.63
1:C:548:ILE:HD13	1:D:425:LEU:HD22	1.80	0.63
1:C:439:LYS:NZ	1:D:519:GLN:HE22	1.78	0.63
1:C:352:GLN:O	1:C:356:MET:HB2	1.98	0.63
1:A:132:CYS:O	1:A:136:ARG:HG3	1.99	0.63
1:D:249:THR:HG22	1:D:253:LYS:HE3	1.79	0.63
1:C:407:ILE:HD13	1:C:425:LEU:HD23	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:522:LYS:HG2	2:B:903:HOH:O	1.96	0.63
1:B:248:ASN:O	1:B:252:VAL:HG23	1.99	0.62
1:D:258:ASP:OD1	1:D:260:GLN:HG2	1.99	0.62
1:C:314:ASN:OD1	1:C:316:PRO:HD2	1.99	0.62
1:B:314:ASN:OD1	1:B:316:PRO:HD2	1.97	0.62
1:A:552:ARG:HB2	2:A:664:HOH:O	1.99	0.62
1:A:439:LYS:HZ3	1:B:519:GLN:HE22	1.46	0.62
1:A:218:ILE:O	1:A:222:GLU:HG3	1.99	0.62
1:C:90:ASN:HD22	1:C:507:ASN:HD21	1.46	0.62
1:C:519:GLN:HE22	1:D:439:LYS:NZ	1.96	0.62
1:B:250:THR:O	1:B:254:GLU:HG3	2.00	0.62
1:B:387:GLN:HA	1:B:391:TYR:CD1	2.34	0.62
1:C:305:HIS:HD2	2:C:560:HOH:O	1.83	0.62
1:A:456:GLU:O	1:A:460:ARG:HG2	2.00	0.62
1:B:161:LEU:HD12	1:B:350:TYR:HA	1.82	0.61
1:C:19:HIS:O	1:C:22:GLU:HB3	2.00	0.61
1:B:61:GLU:HG2	2:B:757:HOH:O	2.00	0.61
1:A:550:GLN:HE21	1:A:551:GLN:NE2	1.87	0.61
1:C:83:MET:HE2	1:C:97:VAL:CG1	2.31	0.61
1:D:431:GLN:O	1:D:435:LEU:HD22	2.00	0.61
1:A:519:GLN:HE22	1:B:439:LYS:HZ2	1.48	0.60
1:D:132:CYS:O	1:D:136:ARG:HG3	2.01	0.60
1:C:45:THR:O	1:C:46:ASN:HB2	2.00	0.60
1:C:315:ALA:HB3	1:C:316:PRO:HD3	1.83	0.60
1:D:134:ARG:HD2	2:D:575:HOH:O	2.01	0.60
1:D:315:ALA:HB3	1:D:316:PRO:HD3	1.84	0.60
1:A:19:HIS:O	1:A:22:GLU:HB3	2.02	0.60
1:A:258:ASP:OD1	1:A:260:GLN:HG2	2.01	0.60
1:C:178:PRO:O	1:C:286:HIS:HD2	1.84	0.60
1:A:95:ARG:HD2	1:A:267:ASP:OD2	2.01	0.60
1:C:522:LYS:HE2	1:C:522:LYS:HA	1.82	0.60
1:C:249:THR:HG22	1:C:253:LYS:HE2	1.83	0.60
1:A:44:ASN:ND2	1:A:46:ASN:H	2.00	0.60
1:D:556:VAL:HG12	1:D:556:VAL:O	2.02	0.60
1:B:191:ILE:HG13	1:B:195:LEU:HD22	1.84	0.59
1:C:434:ALA:HB1	1:D:516:LEU:HD13	1.82	0.59
1:B:407:ILE:HD13	1:B:425:LEU:HD23	1.84	0.59
1:D:45:THR:O	1:D:46:ASN:HB2	2.00	0.59
1:B:352:GLN:O	1:B:356:MET:HB2	2.02	0.59
1:B:370:VAL:HG12	1:B:372:HIS:H	1.68	0.59
1:A:59:VAL:HG23	1:A:63:VAL:CG1	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:THR:OG1	1:B:420:HIS:HE1	1.86	0.59
1:C:59:VAL:HG23	1:C:63:VAL:CG1	2.33	0.59
1:B:544:LEU:O	1:B:548:ILE:HG12	2.02	0.59
1:B:109:PRO:HG2	2:B:614:HOH:O	2.03	0.59
1:D:191:ILE:HG13	1:D:195:LEU:HD22	1.85	0.59
1:D:522:LYS:HG2	2:D:637:HOH:O	2.03	0.58
1:C:124:VAL:HG12	1:C:128:MET:HE2	1.85	0.58
1:A:442:GLU:HG3	2:A:907:HOH:O	2.02	0.58
1:C:12:LEU:HD11	1:C:67:LEU:HD23	1.86	0.58
1:B:59:VAL:HG23	1:B:63:VAL:CG1	2.34	0.58
1:C:253:LYS:HG2	1:C:259:PRO:HG3	1.85	0.58
1:A:230:GLN:O	1:A:233:LYS:HE3	2.03	0.58
1:A:519:GLN:HE22	1:B:439:LYS:HZ1	1.50	0.58
1:C:519:GLN:HE22	1:D:439:LYS:HZ1	1.51	0.57
2:A:561:HOH:O	1:B:49:HIS:HD2	1.87	0.57
1:C:16:TYR:O	1:C:20:ARG:HB2	2.04	0.57
1:B:461:LEU:HG	1:B:465:LYS:HD2	1.85	0.57
1:A:420:HIS:HE1	1:B:223:THR:OG1	1.87	0.57
1:B:442:GLU:HG2	1:B:446:LYS:HE2	1.87	0.57
1:D:265:PHE:CE2	1:D:279:ILE:HD11	2.40	0.57
1:B:45:THR:O	1:B:46:ASN:HB2	2.05	0.57
1:C:305:HIS:CE1	1:C:315:ALA:H	2.15	0.57
1:D:440:SER:OG	1:D:443:GLU:HG3	2.05	0.57
1:A:45:THR:O	1:A:46:ASN:HB2	2.03	0.57
1:A:315:ALA:HB3	1:A:316:PRO:HD3	1.87	0.57
2:A:566:HOH:O	1:B:49:HIS:HE1	1.87	0.57
1:A:440:SER:OG	1:A:443:GLU:HG3	2.05	0.56
1:B:258:ASP:HB3	1:B:261:ASN:ND2	2.19	0.56
1:C:395:HIS:HD2	2:C:820:HOH:O	1.87	0.56
1:A:387:GLN:HA	1:A:391:TYR:CG	2.40	0.56
1:A:553:GLU:O	1:A:555:ARG:HD2	2.05	0.56
1:C:420:HIS:HE1	1:D:223:THR:OG1	1.87	0.56
1:D:442:GLU:HG2	1:D:446:LYS:HE2	1.87	0.56
1:C:59:VAL:HG23	1:C:63:VAL:CB	2.35	0.56
1:A:456:GLU:HG3	1:A:460:ARG:NH2	2.20	0.56
1:B:83:MET:HE2	1:B:97:VAL:HG11	1.88	0.56
1:C:152:ILE:HD12	1:C:152:ILE:N	2.21	0.56
1:B:59:VAL:CG2	1:B:63:VAL:HB	2.35	0.56
1:B:264:GLU:HG2	2:B:833:HOH:O	2.05	0.56
1:B:251:LYS:HD3	1:B:254:GLU:OE1	2.06	0.56
1:D:442:GLU:O	1:D:446:LYS:HG3	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:193:LYS:HE2	2:D:739:HOH:O	2.06	0.55
1:B:210:LYS:O	1:B:246:SER:HB2	2.06	0.55
1:D:515:GLU:O	1:D:519:GLN:HG3	2.05	0.55
1:B:395:HIS:HD2	2:B:791:HOH:O	1.88	0.55
1:A:152:ILE:N	1:A:152:ILE:HD12	2.20	0.55
1:C:395:HIS:NE2	1:C:435:LEU:HD13	2.22	0.55
1:B:59:VAL:HG23	1:B:63:VAL:CB	2.36	0.55
1:A:90:ASN:HD22	1:A:507:ASN:HD21	1.52	0.55
1:A:13:GLN:HG2	1:A:17:ARG:NH1	2.22	0.55
1:B:518:LYS:HE2	2:B:764:HOH:O	2.07	0.55
1:C:13:GLN:HG2	1:C:17:ARG:HH12	1.71	0.55
1:B:258:ASP:HB3	1:B:261:ASN:HD22	1.72	0.55
1:C:248:ASN:O	1:C:252:VAL:HG23	2.07	0.55
1:A:369:ARG:HH11	1:A:369:ARG:HG2	1.72	0.55
1:C:440:SER:OG	1:C:443:GLU:HG3	2.07	0.55
1:D:314:ASN:OD1	1:D:316:PRO:HD2	2.07	0.54
1:B:209:SER:O	1:B:246:SER:CB	2.55	0.54
1:A:59:VAL:CG2	1:A:63:VAL:HB	2.36	0.54
1:B:99:HIS:HD2	2:B:752:HOH:O	1.90	0.54
1:B:228:PHE:HD2	1:B:229:LEU:HD12	1.72	0.54
1:C:413:HIS:HD2	1:D:189:THR:HG22	1.73	0.54
1:C:442:GLU:O	1:C:446:LYS:HG3	2.08	0.54
1:C:210:LYS:HD3	1:C:265:PHE:CE1	2.43	0.54
1:C:124:VAL:HG12	1:C:128:MET:CE	2.37	0.54
1:D:275:LEU:HD12	1:D:275:LEU:C	2.27	0.54
1:A:59:VAL:HG23	1:A:63:VAL:CB	2.36	0.54
1:A:409:VAL:HG22	1:B:548:ILE:CD1	2.38	0.54
1:C:146:LYS:HB2	2:C:618:HOH:O	2.08	0.54
1:A:534:VAL:HG21	1:A:545:ILE:HG21	1.89	0.54
1:D:352:GLN:O	1:D:356:MET:HB2	2.08	0.54
1:B:90:ASN:HD22	1:B:507:ASN:HD21	1.56	0.54
1:D:258:ASP:CG	1:D:260:GLN:HG2	2.28	0.54
1:C:44:ASN:ND2	1:C:46:ASN:H	2.05	0.54
1:A:525:GLU:O	1:B:422:LYS:HE3	2.06	0.54
1:D:535:THR:HG22	2:D:686:HOH:O	2.08	0.54
1:D:189:THR:HG21	2:D:690:HOH:O	2.08	0.54
1:A:516:LEU:HD13	1:B:434:ALA:HB1	1.90	0.54
1:A:522:LYS:HE2	1:A:522:LYS:HA	1.89	0.53
1:B:225:LYS:O	1:B:229:LEU:HD13	2.08	0.53
1:C:191:ILE:HG13	1:C:195:LEU:HD22	1.89	0.53
1:C:59:VAL:CG2	1:C:63:VAL:HB	2.37	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:210:LYS:O	1:C:246:SER:HB2	2.07	0.53
1:B:14:GLN:O	1:B:18:GLU:HG3	2.07	0.53
1:C:209:SER:O	1:C:246:SER:CB	2.56	0.53
1:C:83:MET:HE2	1:C:97:VAL:HG12	1.90	0.53
1:A:91:TYR:CE2	1:A:365:LYS:HA	2.43	0.53
1:A:303:ASP:OD1	1:A:495:HIS:HE1	1.92	0.53
1:B:175:SER:HB3	2:B:638:HOH:O	2.08	0.53
1:D:265:PHE:CZ	1:D:279:ILE:HD11	2.44	0.53
1:A:164:LEU:HD23	1:A:164:LEU:C	2.29	0.53
1:C:258:ASP:HB3	1:C:261:ASN:HD22	1.74	0.53
1:C:381:GLU:HG3	1:C:382:PRO:HD2	1.90	0.53
1:A:253:LYS:NZ	1:A:259:PRO:HG2	2.24	0.53
1:B:44:ASN:ND2	1:B:46:ASN:H	2.06	0.53
1:A:13:GLN:HG2	1:A:17:ARG:HH12	1.72	0.53
1:D:44:ASN:ND2	1:D:46:ASN:H	2.07	0.53
1:C:49:HIS:HE1	2:D:574:HOH:O	1.91	0.53
1:A:209:SER:O	1:A:246:SER:CB	2.57	0.52
1:D:58:LEU:CD2	1:D:472:PRO:HB3	2.39	0.52
1:C:245:LEU:HD13	1:C:279:ILE:HA	1.91	0.52
1:C:200:PRO:HG3	1:C:227:TRP:CZ2	2.45	0.52
1:B:98:LEU:HB2	1:B:268:TRP:CE3	2.45	0.52
1:B:189:THR:HG21	2:B:571:HOH:O	2.09	0.52
1:A:537:HIS:N	1:A:542:ASN:HD21	2.00	0.52
1:A:210:LYS:HD3	1:A:265:PHE:CE1	2.45	0.52
1:D:90:ASN:C	1:D:90:ASN:HD22	2.13	0.52
1:B:369:ARG:HH11	1:B:369:ARG:HG2	1.75	0.52
1:A:550:GLN:NE2	1:A:551:GLN:HE21	1.88	0.52
1:A:352:GLN:O	1:A:356:MET:HB2	2.11	0.51
1:A:264:GLU:HG2	2:A:620:HOH:O	2.10	0.51
1:B:258:ASP:OD1	1:B:260:GLN:HG2	2.09	0.51
1:B:294:GLN:HG3	2:B:650:HOH:O	2.10	0.51
1:B:229:LEU:HD11	1:B:238:VAL:CG2	2.40	0.51
1:C:143:TYR:CZ	1:C:232:ALA:HB2	2.46	0.51
1:D:325:TRP:O	1:D:329:CYS:HB2	2.10	0.51
1:C:294:GLN:HA	1:C:294:GLN:HE21	1.76	0.51
1:B:186:ILE:HD13	1:B:216:GLU:HA	1.93	0.51
1:C:77:GLU:HG3	2:C:731:HOH:O	2.10	0.51
1:A:395:HIS:HE1	1:A:431:GLN:OE1	1.93	0.51
1:C:294:GLN:NE2	1:C:294:GLN:HA	2.25	0.51
1:D:381:GLU:HG3	1:D:382:PRO:HD2	1.91	0.51
1:A:449:GLN:HG2	1:A:458:LEU:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:LEU:CB	1:A:436:MET:HG3	2.40	0.51
1:B:83:MET:HE2	1:B:97:VAL:CG1	2.41	0.51
1:D:218:ILE:O	1:D:222:GLU:HG3	2.11	0.51
1:B:141:LYS:HD3	1:B:145:GLY:O	2.11	0.51
1:C:253:LYS:NZ	1:C:259:PRO:HG2	2.24	0.50
1:B:395:HIS:HE1	1:B:431:GLN:OE1	1.94	0.50
1:C:250:THR:O	1:C:254:GLU:HG3	2.11	0.50
1:B:134:ARG:NH2	2:B:809:HOH:O	2.44	0.50
1:A:434:ALA:CB	1:B:516:LEU:HD13	2.41	0.50
1:C:325:TRP:O	1:C:329:CYS:HB2	2.10	0.50
1:D:229:LEU:HA	1:D:232:ALA:O	2.11	0.50
1:A:16:TYR:O	1:A:20:ARG:HB2	2.11	0.50
1:B:258:ASP:CG	1:B:260:GLN:HG2	2.31	0.50
1:A:395:HIS:NE2	1:A:435:LEU:HD13	2.26	0.50
1:D:258:ASP:HB3	1:D:261:ASN:HD22	1.76	0.50
1:C:6:ASP:O	1:C:10:GLN:HG3	2.12	0.50
1:D:58:LEU:HD22	1:D:472:PRO:HB3	1.93	0.50
1:B:164:LEU:HD13	1:B:164:LEU:C	2.32	0.50
1:D:518:LYS:HG3	2:D:637:HOH:O	2.09	0.50
1:C:334:THR:HG22	1:C:377:ILE:HB	1.94	0.50
1:C:163:PRO:HA	1:C:283:ILE:HD11	1.93	0.50
1:A:439:LYS:HZ1	1:B:519:GLN:HE22	1.57	0.50
1:C:218:ILE:O	1:C:222:GLU:HG3	2.11	0.50
1:D:436:MET:CE	1:D:437:ARG:HG3	2.42	0.50
1:D:209:SER:O	1:D:246:SER:CB	2.57	0.49
1:B:515:GLU:O	1:B:519:GLN:HG3	2.11	0.49
1:D:141:LYS:HD3	1:D:145:GLY:O	2.11	0.49
1:C:258:ASP:HB3	1:C:261:ASN:ND2	2.27	0.49
1:A:58:LEU:CD2	1:A:472:PRO:HB3	2.42	0.49
1:B:209:SER:O	1:B:246:SER:HB2	2.12	0.49
1:C:191:ILE:HG23	1:C:192:ALA:N	2.27	0.49
1:B:230:GLN:HG3	2:B:674:HOH:O	2.12	0.49
1:A:58:LEU:HD22	1:A:472:PRO:HB3	1.95	0.49
1:B:19:HIS:O	1:B:22:GLU:HB3	2.11	0.49
1:D:387:GLN:HA	1:D:391:TYR:CG	2.48	0.49
1:B:25:LEU:CB	1:B:436:MET:HG3	2.42	0.49
1:A:458:LEU:HD23	1:A:458:LEU:C	2.33	0.49
1:B:395:HIS:CD2	1:B:435:LEU:HD13	2.47	0.49
1:B:384:THR:HA	1:B:387:GLN:HB3	1.94	0.49
1:A:325:TRP:O	1:A:329:CYS:HB2	2.12	0.49
1:C:416:ARG:NH2	1:D:222:GLU:HB2	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:303:ASP:OD1	1:D:495:HIS:HE1	1.96	0.49
1:B:275:LEU:C	1:B:275:LEU:HD12	2.32	0.49
1:B:440:SER:OG	1:B:443:GLU:HG3	2.12	0.49
1:B:312:GLU:H	1:B:312:GLU:CD	2.15	0.49
1:C:312:GLU:CD	1:C:312:GLU:H	2.16	0.49
1:D:88:LYS:HD2	1:D:88:LYS:N	2.27	0.49
1:B:314:ASN:HB3	1:B:317:VAL:HB	1.93	0.48
1:A:324:ILE:CD1	1:A:503:ILE:HD12	2.43	0.48
1:C:490:VAL:O	1:C:494:GLU:HG3	2.12	0.48
1:A:90:ASN:HD22	1:A:90:ASN:C	2.16	0.48
1:A:324:ILE:HD13	1:A:503:ILE:HD12	1.95	0.48
1:B:23:LEU:HB3	1:B:58:LEU:HD12	1.94	0.48
1:B:304:GLN:CG	2:B:558:HOH:O	2.34	0.48
1:C:82:ARG:NH2	2:C:788:HOH:O	2.47	0.48
1:B:382:PRO:CA	1:B:383:GLY:H	2.21	0.48
1:C:13:GLN:HG2	1:C:17:ARG:NH1	2.28	0.48
1:B:25:LEU:HB3	1:B:436:MET:HG3	1.95	0.48
1:A:141:LYS:HD3	1:A:145:GLY:O	2.12	0.48
1:B:370:VAL:CG1	1:B:372:HIS:H	2.26	0.48
1:D:445:ARG:O	1:D:449:GLN:HG3	2.14	0.48
1:B:209:SER:O	1:B:246:SER:HB3	2.13	0.48
1:A:132:CYS:O	1:A:136:ARG:CG	2.62	0.48
1:A:384:THR:HA	1:A:387:GLN:HB3	1.96	0.48
1:C:214:THR:O	1:C:218:ILE:HG12	2.14	0.48
1:A:16:TYR:CZ	1:A:20:ARG:HG3	2.49	0.47
1:D:189:THR:O	1:D:193:LYS:HG2	2.15	0.47
1:A:320:ALA:HB1	1:A:500:GLN:HG3	1.97	0.47
1:A:305:HIS:HD2	2:A:867:HOH:O	1.98	0.47
1:D:407:ILE:HD13	1:D:425:LEU:HD23	1.96	0.47
1:C:215:GLN:OE1	1:D:423:ILE:HD12	2.14	0.47
1:C:445:ARG:HD3	2:C:787:HOH:O	2.14	0.47
1:D:161:LEU:HD12	1:D:350:TYR:HA	1.95	0.47
1:A:530:GLY:O	1:A:552:ARG:NH2	2.48	0.47
1:C:313:LYS:HD3	2:C:778:HOH:O	2.15	0.47
1:D:334:THR:HG22	1:D:377:ILE:HB	1.96	0.47
1:B:132:CYS:O	1:B:136:ARG:CG	2.60	0.47
1:C:124:VAL:O	1:C:128:MET:HG3	2.14	0.47
1:B:253:LYS:NZ	1:B:259:PRO:HG2	2.30	0.47
1:B:537:HIS:N	1:B:542:ASN:HD21	2.02	0.47
1:A:59:VAL:HG22	1:A:60:THR:N	2.29	0.47
1:A:99:HIS:HE1	2:A:869:HOH:O	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:ARG:HH12	1:B:511:GLN:HE21	1.63	0.47
1:C:481:LEU:HD12	2:C:593:HOH:O	2.15	0.47
1:D:90:ASN:HD22	1:D:507:ASN:HD21	1.62	0.47
1:B:303:ASP:OD1	1:B:495:HIS:HE1	1.98	0.46
1:A:29:PHE:CD1	1:A:436:MET:CE	2.98	0.46
1:D:132:CYS:O	1:D:136:ARG:CG	2.62	0.46
1:C:200:PRO:HG3	1:C:227:TRP:CH2	2.50	0.46
1:D:164:LEU:HD23	1:D:164:LEU:C	2.36	0.46
1:C:300:HIS:O	1:C:304:GLN:HG2	2.15	0.46
1:B:305:HIS:CE1	1:B:315:ALA:H	2.26	0.46
1:A:481:LEU:HD12	2:A:812:HOH:O	2.15	0.46
1:B:405:PHE:HB3	1:B:428:PHE:CE1	2.51	0.46
1:C:249:THR:O	1:C:253:LYS:HG3	2.16	0.46
1:C:99:HIS:HD2	2:C:683:HOH:O	1.99	0.46
1:B:178:PRO:O	1:B:286:HIS:HD2	1.99	0.46
1:B:228:PHE:CD2	1:B:229:LEU:HD12	2.50	0.46
1:D:95:ARG:HD2	1:D:267:ASP:HB2	1.98	0.46
1:A:258:ASP:HB3	1:A:261:ASN:HD22	1.80	0.46
1:C:303:ASP:OD1	1:C:495:HIS:HE1	1.99	0.46
1:A:222:GLU:HB2	1:B:416:ARG:NH2	2.31	0.46
1:A:189:THR:HG21	2:B:607:HOH:O	2.16	0.46
1:A:486:LEU:O	1:A:490:VAL:HG23	2.16	0.46
1:B:234:ASP:HA	1:B:235:PRO:HD2	1.87	0.46
1:D:111:LEU:HD23	1:D:116:ASP:HA	1.97	0.46
1:D:370:VAL:HG13	1:D:372:HIS:CE1	2.51	0.45
1:D:436:MET:HE3	1:D:437:ARG:HG3	1.98	0.45
1:B:164:LEU:HD23	1:B:182:TYR:CD2	2.51	0.45
1:C:82:ARG:HG2	1:C:87:GLU:CD	2.37	0.45
1:B:456:GLU:O	1:B:460:ARG:HG2	2.16	0.45
1:A:439:LYS:N	1:A:467:PHE:HB2	2.31	0.45
1:B:305:HIS:HD2	2:B:667:HOH:O	1.99	0.45
1:C:395:HIS:HE1	1:C:431:GLN:OE1	1.99	0.45
1:D:70:LEU:HD22	1:D:74:ARG:HG2	1.98	0.45
1:D:126:ASP:HB3	2:D:759:HOH:O	2.16	0.45
1:A:245:LEU:HD13	1:A:279:ILE:HA	1.98	0.45
1:C:80:ARG:HD2	1:C:307:ARG:HA	1.97	0.45
1:C:387:GLN:HA	1:C:391:TYR:CG	2.51	0.45
1:A:247:THR:HG22	1:A:264:GLU:HB2	1.97	0.45
1:D:522:LYS:CE	1:D:522:LYS:HA	2.45	0.45
1:B:187:ASP:OD2	1:B:189:THR:HG23	2.17	0.45
1:D:152:ILE:HD12	1:D:152:ILE:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:460:ARG:HG2	1:C:460:ARG:HH11	1.82	0.45
1:A:395:HIS:HD2	2:A:897:HOH:O	2.00	0.45
1:B:495:HIS:O	1:B:499:VAL:HG23	2.17	0.45
1:D:99:HIS:HE1	2:D:683:HOH:O	1.99	0.45
1:C:6:ASP:OD2	1:C:8:GLN:HB2	2.16	0.45
1:C:522:LYS:CE	1:C:522:LYS:HA	2.46	0.44
1:D:11:LYS:HD3	2:D:860:HOH:O	2.16	0.44
1:D:369:ARG:HG2	1:D:369:ARG:HH11	1.82	0.44
1:C:164:LEU:HD23	1:C:164:LEU:C	2.37	0.44
1:A:20:ARG:C	1:A:22:GLU:H	2.20	0.44
1:A:258:ASP:HB3	1:A:261:ASN:ND2	2.32	0.44
1:B:171:LYS:N	1:B:172:PRO:CD	2.80	0.44
1:C:34:ASP:HA	2:C:694:HOH:O	2.16	0.44
1:D:460:ARG:HG2	1:D:460:ARG:HH11	1.81	0.44
1:C:305:HIS:HE1	1:C:315:ALA:N	2.05	0.44
1:B:229:LEU:HD11	1:B:238:VAL:HG22	1.98	0.44
1:C:95:ARG:HD2	1:C:267:ASP:HB2	1.99	0.44
1:A:214:THR:O	1:A:218:ILE:HG12	2.18	0.44
1:A:186:ILE:O	1:B:420:HIS:HD2	2.00	0.44
1:A:324:ILE:HD13	1:A:503:ILE:CD1	2.48	0.44
1:D:530:GLY:O	1:D:552:ARG:NH2	2.50	0.44
1:B:496:LYS:O	1:B:500:GLN:HG3	2.17	0.44
1:A:98:LEU:HB2	1:A:268:TRP:CE3	2.53	0.44
1:C:516:LEU:HD13	1:D:434:ALA:HB1	2.00	0.44
1:D:312:GLU:CD	1:D:312:GLU:H	2.21	0.44
1:C:303:ASP:HA	1:C:495:HIS:HE1	1.82	0.44
1:D:523:LYS:HE3	2:D:558:HOH:O	2.17	0.44
1:C:251:LYS:HD3	1:C:254:GLU:OE1	2.18	0.44
1:A:320:ALA:O	1:A:324:ILE:HD12	2.17	0.44
1:C:6:ASP:HA	1:C:7:PRO:HD3	1.89	0.43
1:A:253:LYS:HG2	1:A:259:PRO:HG3	2.01	0.43
1:B:334:THR:HG22	1:B:377:ILE:HB	2.01	0.43
1:D:544:LEU:O	1:D:548:ILE:HG13	2.18	0.43
1:C:20:ARG:C	1:C:22:GLU:H	2.20	0.43
1:A:232:ALA:O	1:A:233:LYS:HB2	2.19	0.43
1:A:490:VAL:O	1:A:494:GLU:HG3	2.19	0.43
1:D:458:LEU:HD23	1:D:458:LEU:C	2.38	0.43
1:C:3:LEU:HD11	1:C:9:PHE:CE1	2.54	0.43
1:B:370:VAL:CG1	1:B:372:HIS:O	2.67	0.43
1:C:456:GLU:O	1:C:460:ARG:HG2	2.18	0.43
1:C:171:LYS:N	1:C:172:PRO:CD	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:178:PRO:O	1:D:286:HIS:HD2	2.01	0.43
1:A:82:ARG:HG2	1:A:87:GLU:CD	2.39	0.43
1:C:207:ILE:N	1:C:207:ILE:HD12	2.33	0.43
1:A:275:LEU:HD12	1:A:275:LEU:C	2.37	0.43
1:D:370:VAL:HG13	1:D:372:HIS:CD2	2.53	0.43
1:B:22:GLU:HG2	1:B:22:GLU:O	2.18	0.43
1:C:437:ARG:HB3	1:C:437:ARG:CZ	2.49	0.43
1:C:1:ALA:HB1	2:C:722:HOH:O	2.16	0.43
1:A:305:HIS:CE1	1:A:315:ALA:H	2.27	0.43
1:B:387:GLN:HA	1:B:391:TYR:CG	2.53	0.43
1:A:195:LEU:HG	1:A:227:TRP:CD2	2.54	0.43
1:A:356:MET:HA	2:A:633:HOH:O	2.19	0.43
1:C:207:ILE:N	1:C:207:ILE:CD1	2.81	0.43
1:C:458:LEU:C	1:C:458:LEU:HD23	2.38	0.43
1:C:63:VAL:HA	1:C:66:MET:HE2	2.01	0.43
1:B:448:LEU:O	1:B:453:LYS:HB2	2.19	0.43
1:A:155:GLY:O	1:A:184:SER:HA	2.19	0.43
1:B:382:PRO:HA	1:B:383:GLY:N	2.32	0.42
1:D:381:GLU:CG	1:D:382:PRO:HD2	2.49	0.42
1:D:381:GLU:OE2	1:D:382:PRO:HD2	2.19	0.42
1:A:530:GLY:HA3	2:A:701:HOH:O	2.18	0.42
1:C:141:LYS:HE3	2:C:580:HOH:O	2.19	0.42
1:B:351:PHE:CZ	1:B:493:TYR:HB2	2.54	0.42
1:D:171:LYS:N	1:D:172:PRO:CD	2.82	0.42
1:B:12:LEU:HD11	1:B:67:LEU:HD23	2.01	0.42
1:C:456:GLU:OE1	1:C:456:GLU:HA	2.19	0.42
1:D:486:LEU:O	1:D:490:VAL:HG23	2.19	0.42
1:C:459:GLU:HB2	1:C:460:ARG:HH11	1.82	0.42
1:D:245:LEU:HD13	1:D:279:ILE:HA	2.02	0.42
1:A:189:THR:O	1:A:193:LYS:HG2	2.19	0.42
1:C:553:GLU:O	1:C:555:ARG:HD2	2.19	0.42
1:C:112:VAL:O	1:C:113:ASP:HB2	2.20	0.42
1:C:132:CYS:O	1:C:136:ARG:CG	2.63	0.42
1:A:191:ILE:HG23	1:A:192:ALA:N	2.34	0.42
1:C:381:GLU:HG2	2:D:699:HOH:O	2.18	0.42
1:D:408:PRO:HA	1:D:478:PHE:O	2.19	0.42
1:A:49:HIS:HD2	2:B:587:HOH:O	2.03	0.42
1:C:448:LEU:O	1:C:451:ALA:HB3	2.19	0.42
1:D:356:MET:HA	2:D:641:HOH:O	2.19	0.42
1:B:6:ASP:HA	1:B:7:PRO:HD3	1.85	0.42
1:A:187:ASP:OD2	1:A:189:THR:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:127:LYS:HG2	1:D:263:PHE:CE2	2.55	0.42
1:A:439:LYS:H	1:A:467:PHE:HB2	1.85	0.42
1:C:321:LEU:CD1	2:C:639:HOH:O	2.68	0.42
1:C:253:LYS:HZ2	1:C:259:PRO:HG2	1.85	0.42
1:A:230:GLN:NE2	1:A:230:GLN:HA	2.35	0.42
1:C:195:LEU:HG	1:C:227:TRP:CD2	2.55	0.42
1:D:111:LEU:HD23	1:D:116:ASP:CA	2.50	0.42
1:A:381:GLU:HG3	1:A:382:PRO:HD2	2.02	0.42
1:B:298:GLY:HA3	1:B:484:PHE:O	2.20	0.41
1:D:52:VAL:HG13	1:D:476:ILE:CD1	2.50	0.41
1:C:134:ARG:NH2	2:C:681:HOH:O	2.51	0.41
1:B:59:VAL:HG22	1:B:60:THR:N	2.34	0.41
1:C:223:THR:OG1	1:D:420:HIS:HE1	2.03	0.41
1:C:24:ASN:HB3	1:C:27:ARG:HE	1.85	0.41
1:C:290:ASP:HB2	2:C:714:HOH:O	2.19	0.41
1:C:416:ARG:NH1	1:C:419:LEU:HD23	2.35	0.41
1:C:59:VAL:HG22	1:C:60:THR:N	2.36	0.41
1:C:550:GLN:NE2	1:C:551:GLN:HE21	2.10	0.41
1:C:384:THR:HA	1:C:387:GLN:HB3	2.02	0.41
1:C:274:SER:O	1:C:279:ILE:HB	2.20	0.41
1:B:408:PRO:HA	1:B:478:PHE:O	2.21	0.41
1:D:98:LEU:HB2	1:D:268:TRP:CE3	2.54	0.41
1:D:387:GLN:HA	1:D:391:TYR:CE1	2.55	0.41
1:C:381:GLU:CG	1:C:382:PRO:HD2	2.51	0.41
1:C:457:ASP:O	1:C:461:LEU:HB2	2.21	0.41
1:C:369:ARG:HG2	1:C:369:ARG:HH11	1.85	0.41
1:B:382:PRO:HA	1:B:383:GLY:H	1.84	0.41
1:B:550:GLN:NE2	1:B:551:GLN:HE21	2.14	0.41
1:B:352:GLN:HG3	1:B:378:VAL:O	2.19	0.41
1:A:320:ALA:HB1	1:A:500:GLN:CG	2.50	0.41
1:D:180:VAL:HG11	1:D:286:HIS:HB2	2.01	0.41
1:B:20:ARG:C	1:B:22:GLU:H	2.23	0.41
1:B:408:PRO:HB3	1:B:478:PHE:CE1	2.56	0.41
1:B:370:VAL:HG13	1:B:372:HIS:CD2	2.55	0.41
1:C:258:ASP:OD1	1:C:260:GLN:HG2	2.20	0.41
1:A:6:ASP:HA	1:A:7:PRO:HD3	1.87	0.41
1:A:407:ILE:HD13	1:A:425:LEU:HD23	2.02	0.41
1:D:82:ARG:HG2	1:D:87:GLU:OE1	2.20	0.41
1:C:175:SER:HB3	2:C:582:HOH:O	2.20	0.41
1:A:210:LYS:O	1:A:246:SER:HB2	2.21	0.41
1:C:214:THR:HB	1:C:217:THR:HB	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:THR:HG22	1:B:264:GLU:HB2	2.03	0.40
1:D:187:ASP:OD2	1:D:189:THR:HG23	2.21	0.40
1:D:82:ARG:HG2	1:D:87:GLU:CD	2.41	0.40
1:B:534:VAL:HA	2:B:719:HOH:O	2.21	0.40
1:B:152:ILE:HD12	1:B:152:ILE:N	2.36	0.40
1:C:434:ALA:HB2	1:D:520:LEU:HD12	2.03	0.40
1:C:303:ASP:OD1	1:C:495:HIS:CE1	2.74	0.40
1:C:534:VAL:HG21	1:C:545:ILE:HG21	2.03	0.40
1:C:83:MET:HB2	1:C:89:ILE:HD13	2.02	0.40
1:A:178:PRO:O	1:A:286:HIS:HD2	2.04	0.40
1:D:370:VAL:HG13	1:D:372:HIS:NE2	2.36	0.40
1:B:516:LEU:HD22	1:B:520:LEU:HG	2.02	0.40
1:A:435:LEU:HD23	1:A:473:THR:HG21	2.03	0.40
1:B:552:ARG:HD3	2:B:692:HOH:O	2.21	0.40
1:D:448:LEU:HD21	1:D:465:LYS:HD3	2.02	0.40
1:A:317:VAL:O	1:A:321:LEU:HD13	2.21	0.40
1:D:345:HIS:HA	1:D:382:PRO:HG3	2.03	0.40
1:C:189:THR:O	1:C:193:LYS:HG2	2.21	0.40
1:C:275:LEU:HD12	1:C:275:LEU:C	2.41	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	554/558 (99%)	533 (96%)	19 (3%)	2 (0%)	39	37
1	B	554/558 (99%)	532 (96%)	22 (4%)	0	100	100
1	C	554/558 (99%)	526 (95%)	26 (5%)	2 (0%)	39	37
1	D	554/558 (99%)	524 (95%)	28 (5%)	2 (0%)	39	37
All	All	2216/2232 (99%)	2115 (95%)	95 (4%)	6 (0%)	46	45

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	384	THR
1	A	21	SER
1	C	21	SER
1	D	384	THR
1	A	384	THR
1	D	554	ALA

### 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	475/477 (100%)	462 (97%)	13 (3%)	52	56
1	B	475/477 (100%)	460 (97%)	15 (3%)	46	48
1	C	475/477 (100%)	464 (98%)	11 (2%)	58	62
1	D	475/477 (100%)	461 (97%)	14 (3%)	50	53
All	All	1900/1908 (100%)	1847 (97%)	53 (3%)	51	55

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ASP
1	A	44	ASN
1	A	70	LEU
1	A	88	LYS
1	A	90	ASN
1	A	103	ARG
1	A	161	LEU
1	A	195	LEU
1	A	229	LEU
1	A	344	LEU
1	A	435	LEU
1	A	511	GLN
1	A	516	LEU
1	B	44	ASN

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Mol	Chain	Res	Type
1	B	70	LEU
1	B	90	ASN
1	B	103	ARG
1	B	117	VAL
1	B	161	LEU
1	B	189	THR
1	B	195	LEU
1	B	265	PHE
1	B	296	LEU
1	B	344	LEU
1	B	435	LEU
1	B	511	GLN
1	B	516	LEU
1	B	553	GLU
1	C	13	GLN
1	C	30	ASP
1	C	44	ASN
1	C	70	LEU
1	C	90	ASN
1	C	103	ARG
1	C	161	LEU
1	C	195	LEU
1	C	344	LEU
1	C	435	LEU
1	C	516	LEU
1	D	30	ASP
1	D	44	ASN
1	D	70	LEU
1	D	88	LYS
1	D	90	ASN
1	D	103	ARG
1	D	195	LEU
1	D	212	PHE
1	D	229	LEU
1	D	321	LEU
1	D	344	LEU
1	D	416	ARG
1	D	435	LEU
1	D	516	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (87) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	44	ASN
1	A	49	HIS
1	A	90	ASN
1	A	99	HIS
1	A	153	ASN
1	A	215	GLN
1	A	230	GLN
1	A	261	ASN
1	A	286	HIS
1	A	294	GLN
1	A	304	GLN
1	A	305	HIS
1	A	353	GLN
1	A	387	GLN
1	A	395	HIS
1	A	410	GLN
1	A	420	HIS
1	A	449	GLN
1	A	495	HIS
1	A	511	GLN
1	A	519	GLN
1	A	533	GLN
1	A	542	ASN
1	A	550	GLN
1	B	10	GLN
1	B	14	GLN
1	B	44	ASN
1	B	49	HIS
1	B	90	ASN
1	B	99	HIS
1	B	153	ASN
1	B	230	GLN
1	B	261	ASN
1	B	286	HIS
1	B	294	GLN
1	B	305	HIS
1	B	353	GLN
1	B	387	GLN
1	B	395	HIS
1	B	410	GLN
1	B	420	HIS
1	B	495	HIS

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Mol	Chain	Res	Type
1	B	511	GLN
1	B	519	GLN
1	B	542	ASN
1	B	550	GLN
1	C	10	GLN
1	C	44	ASN
1	C	49	HIS
1	C	90	ASN
1	C	99	HIS
1	C	153	ASN
1	C	261	ASN
1	C	286	HIS
1	C	294	GLN
1	C	305	HIS
1	C	353	GLN
1	C	387	GLN
1	C	395	HIS
1	C	410	GLN
1	C	420	HIS
1	C	495	HIS
1	C	519	GLN
1	C	542	ASN
1	C	550	GLN
1	D	8	GLN
1	D	10	GLN
1	D	44	ASN
1	D	49	HIS
1	D	90	ASN
1	D	99	HIS
1	D	153	ASN
1	D	215	GLN
1	D	261	ASN
1	D	286	HIS
1	D	294	GLN
1	D	305	HIS
1	D	353	GLN
1	D	387	GLN
1	D	395	HIS
1	D	410	GLN
1	D	420	HIS
1	D	495	HIS
1	D	519	GLN

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Mol	Chain	Res	Type
1	D	542	ASN
1	D	550	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	556/558 (99%)	-0.20	8 (1%) 78 82	14, 22, 39, 53	0
1	B	556/558 (99%)	-0.18	11 (1%) 68 73	14, 22, 41, 53	0
1	C	556/558 (99%)	0.18	29 (5%) 31 39	15, 28, 52, 74	0
1	D	556/558 (99%)	-0.02	19 (3%) 49 58	17, 25, 46, 72	0
All	All	2224/2232 (99%)	-0.05	67 (3%) 54 62	14, 24, 44, 74	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	460	ARG	6.1
1	C	232	ALA	5.6
1	C	451	ALA	5.1
1	C	456	GLU	5.0
1	D	232	ALA	3.8
1	C	454	SER	3.8
1	D	460	ARG	3.8
1	D	456	GLU	3.8
1	C	383	GLY	3.5
1	A	553	GLU	3.4
1	C	260	GLN	3.4
1	B	555	ARG	3.4
1	C	553	GLU	3.4
1	A	460	ARG	3.3
1	B	460	ARG	3.3
1	D	22	GLU	3.2
1	C	458	LEU	3.2
1	D	553	GLU	3.2
1	C	1	ALA	3.0
1	D	451	ALA	3.0
1	D	176	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	383	GLY	2.9
1	C	459	GLU	2.8
1	C	453	LYS	2.7
1	A	555	ARG	2.7
1	C	206	ILE	2.7
1	B	553	GLU	2.7
1	C	555	ARG	2.6
1	C	175	SER	2.6
1	B	30	ASP	2.5
1	C	455	PRO	2.5
1	C	213	THR	2.5
1	C	461	LEU	2.5
1	B	260	GLN	2.5
1	B	383	GLY	2.5
1	B	253	LYS	2.5
1	C	243	VAL	2.4
1	D	449	GLN	2.4
1	B	456	GLU	2.4
1	C	30	ASP	2.3
1	C	207	ILE	2.3
1	D	260	GLN	2.3
1	C	368	THR	2.3
1	B	21	SER	2.3
1	D	453	LYS	2.2
1	D	21	SER	2.2
1	D	437	ARG	2.1
1	A	383	GLY	2.1
1	D	452	GLY	2.1
1	C	230	GLN	2.1
1	A	455	PRO	2.1
1	C	556	VAL	2.1
1	D	457	ASP	2.1
1	C	257	ILE	2.1
1	C	33	LYS	2.1
1	A	30	ASP	2.1
1	D	531	SER	2.1
1	C	450	ALA	2.1
1	C	531	SER	2.1
1	A	449	GLN	2.1
1	B	33	LYS	2.1
1	D	5	ARG	2.0
1	B	14	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	458	LEU	2.0
1	A	451	ALA	2.0
1	D	143	TYR	2.0
1	C	253	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.