



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 01:06 am GMT

PDB ID : 1RSV  
Title : azide complex of the diferrous E238A mutant R2 subunit of ribonucleotide reductase  
Authors : Assarsson, M.; Andersson, M.E.; Hogbom, M.; Persson, B.O.; Sahlin, M.; Barra, A.L.; Sjoberg, B.M.; Nordlund, P.; Graslund, A.  
Deposited on : 2003-12-10  
Resolution : 2.20 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](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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

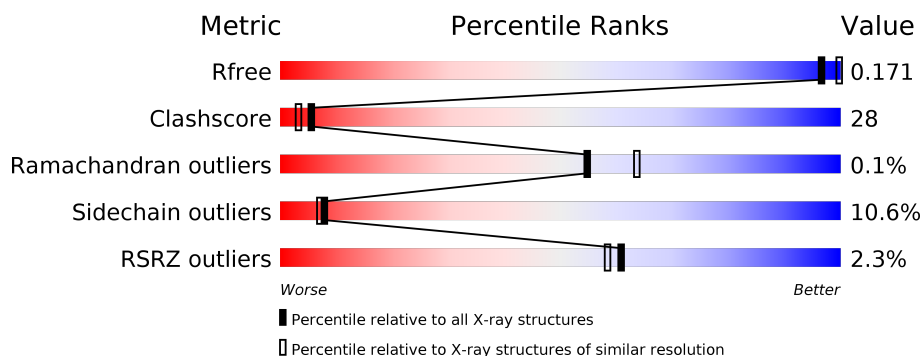
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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	375	<div> <div></div> <div>56%</div> <div>29%</div> <div>6%</div> <div>9%</div> </div>
1	B	375	<div> <div>3%</div> <div>48%</div> <div>33%</div> <div>10%</div> <div>9%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	AZI	A	901	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5919 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 Ribonucleoside-diphosphate reductase 1 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2783	1782	464	524	13			
1	B	341	Total	C	N	O	S	0	0	0
			2789	1785	465	526	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	122	PHE	TYR	ENGINEERED	UNP P69924
A	238	ALA	GLU	ENGINEERED	UNP P69924
B	122	PHE	TYR	ENGINEERED	UNP P69924
B	238	ALA	GLU	ENGINEERED	UNP P69924

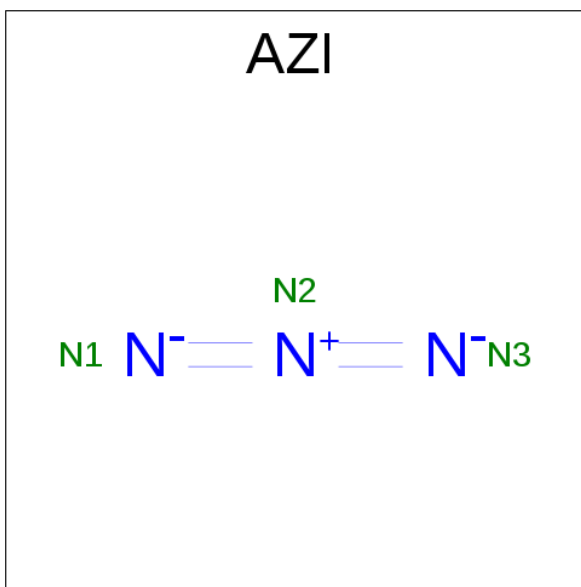
- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Fe	0	0
			2	2		
2	A	2	Total	Fe	0	0
			2	2		

- Molecule 3 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	6	Total	Hg	0	0
			6	6		
3	A	6	Total	Hg	0	0
			6	6		

- Molecule 4 is AZIDE ION (three-letter code: AZI) (formula: N<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total N 3 3	0	0

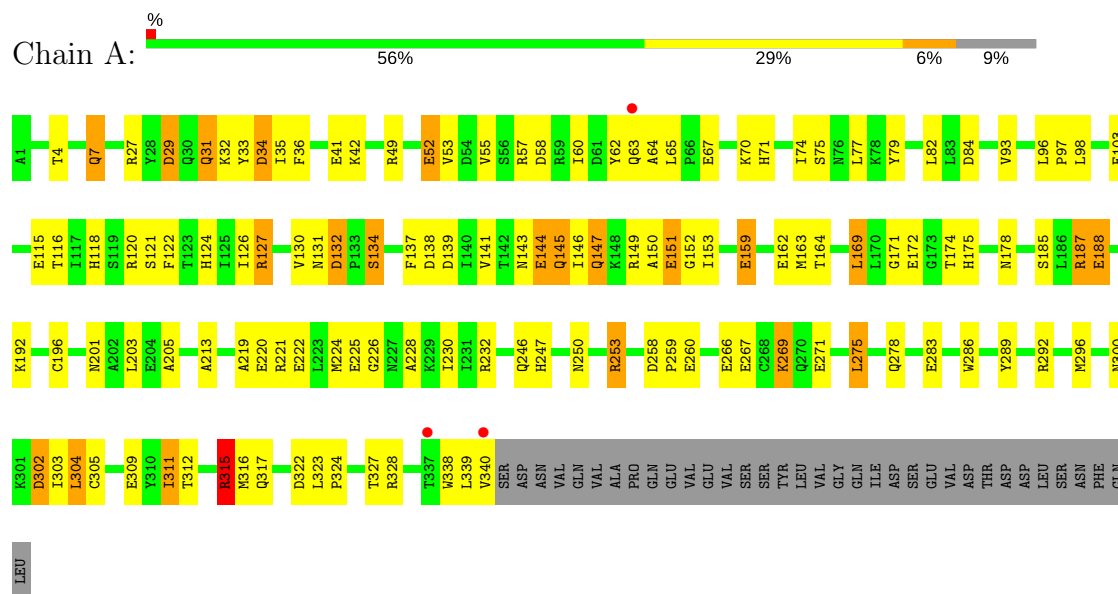
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	161	Total O 161 161	0	0
5	B	167	Total O 167 167	0	0

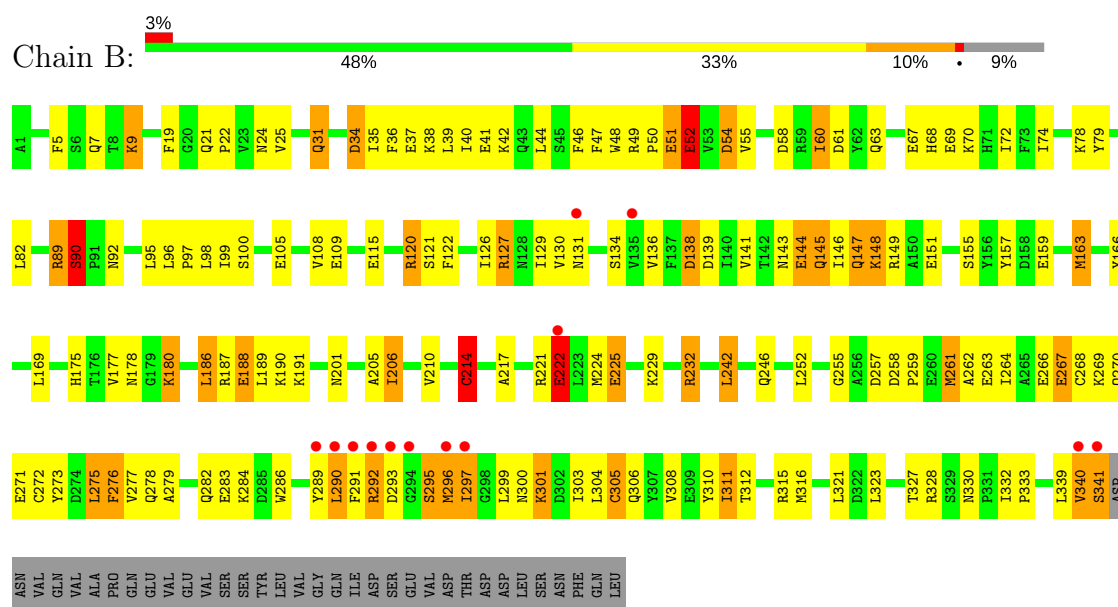
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Ribonucleoside-diphosphate reductase 1 beta chain



- Molecule 1: Ribonucleoside-diphosphate reductase 1 beta chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.89Å 84.50Å 113.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.00 – 2.20 16.74 – 2.19	Depositor EDS
% Data completeness (in resolution range)	(Not available) (17.00-2.20) 97.2 (16.74-2.19)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.28 (at 2.18Å)	Xtriage
Refinement program	TNT	Depositor
R, $R_{free}$	0.182 , 0.256 0.174 , 0.171	Depositor DCC
$R_{free}$ test set	1786 reflections (5.24%)	DCC
Wilson B-factor (Å <sup>2</sup> )	17.4	Xtriage
Anisotropy	0.489	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 79.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5919	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AZI, FE, HG

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	1.04	16/2847 (0.6%)	1.11	9/3862 (0.2%)
1	B	1.04	15/2853 (0.5%)	1.11	9/3870 (0.2%)
All	All	1.04	31/5700 (0.5%)	1.11	18/7732 (0.2%)

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	214	CYS	CB-SG	8.64	1.97	1.82
1	B	222	GLU	CD-OE2	7.24	1.33	1.25
1	A	41	GLU	CD-OE2	6.99	1.33	1.25
1	B	271	GLU	CD-OE2	6.96	1.33	1.25
1	B	188	GLU	CD-OE2	6.94	1.33	1.25
1	A	172	GLU	CD-OE2	6.76	1.33	1.25
1	B	51	GLU	CD-OE2	6.70	1.33	1.25
1	B	159	GLU	CD-OE2	6.41	1.32	1.25
1	A	267	GLU	CD-OE2	6.32	1.32	1.25
1	A	222	GLU	CD-OE2	6.28	1.32	1.25
1	A	144	GLU	CD-OE2	6.20	1.32	1.25
1	B	214	CYS	CA-CB	6.20	1.67	1.53
1	A	115	GLU	CD-OE2	6.18	1.32	1.25
1	A	271	GLU	CD-OE2	6.06	1.32	1.25
1	A	159	GLU	CD-OE2	6.05	1.32	1.25
1	A	225	GLU	CD-OE2	6.05	1.32	1.25
1	A	52	GLU	CD-OE2	6.04	1.32	1.25
1	A	162	GLU	CD-OE2	6.03	1.32	1.25
1	B	52	GLU	CD-OE2	6.00	1.32	1.25
1	B	266	GLU	CD-OE2	5.95	1.32	1.25
1	A	260	GLU	CD-OE2	5.76	1.31	1.25
1	B	225	GLU	CD-OE2	5.65	1.31	1.25
1	B	263	GLU	CD-OE2	5.65	1.31	1.25
1	B	67	GLU	CD-OE2	5.56	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	67	GLU	CD-OE2	5.44	1.31	1.25
1	A	151	GLU	CD-OE2	5.39	1.31	1.25
1	A	220	GLU	CD-OE2	5.39	1.31	1.25
1	A	103	GLU	CD-OE2	5.29	1.31	1.25
1	B	267	GLU	CD-OE2	5.27	1.31	1.25
1	B	115	GLU	CD-OE1	-5.24	1.19	1.25
1	B	144	GLU	CD-OE2	5.15	1.31	1.25

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	214	CYS	CA-CB-SG	10.29	132.53	114.00
1	B	214	CYS	CB-CA-C	7.28	124.96	110.40
1	B	89	ARG	NE-CZ-NH2	-7.07	116.77	120.30
1	A	132	ASP	CB-CG-OD2	-6.60	112.36	118.30
1	B	54	ASP	CB-CG-OD2	-6.57	112.39	118.30
1	B	54	ASP	CB-CG-OD1	6.26	123.94	118.30
1	A	315	ARG	NE-CZ-NH2	6.22	123.41	120.30
1	B	138	ASP	CB-CG-OD1	5.95	123.66	118.30
1	B	257	ASP	CB-CG-OD2	-5.58	113.28	118.30
1	A	139	ASP	CB-CG-OD1	5.55	123.29	118.30
1	A	169	LEU	CB-CG-CD1	-5.44	101.76	111.00
1	A	34	ASP	CB-CG-OD1	5.39	123.16	118.30
1	A	187	ARG	NE-CZ-NH2	-5.33	117.63	120.30
1	A	302	ASP	CB-CG-OD2	-5.11	113.70	118.30
1	B	90	SER	N-CA-C	5.10	124.78	111.00
1	B	232	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	A	29	ASP	CB-CG-OD2	-5.03	113.78	118.30
1	A	29	ASP	CB-CG-OD1	5.01	122.81	118.30

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	2783	0	2732	121	0
1	B	2789	0	2735	208	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	6	0	0	4	0
3	B	6	0	0	1	0
4	A	3	0	0	1	0
5	A	161	0	0	17	0
5	B	167	0	0	28	0
All	All	5919	0	5467	311	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 28.

All (311) 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)
3:B:2004:HG:HG	5:B:2175:HOH:O	1.15	1.18
3:A:2009:HG:HG	5:A:2165:HOH:O	1.34	1.05
3:A:2012:HG:HG	5:A:2151:HOH:O	1.32	1.03
1:A:205:ALA:HB1	1:A:315:ARG:HD3	1.36	1.02
3:A:2007:HG:HG	5:A:2127:HOH:O	1.37	1.00
1:B:301:LYS:HE3	1:B:301:LYS:H	1.23	0.99
1:A:138:ASP:HB3	1:B:9:LYS:HZ3	1.30	0.97
1:B:252:LEU:HD22	1:B:261:MET:HG3	1.49	0.95
1:A:7:GLN:HE22	1:B:147:GLN:HE21	1.12	0.92
1:B:149:ARG:HD3	1:B:286:TRP:HB2	1.51	0.91
1:B:296:MET:HG2	1:B:297:ILE:H	1.36	0.91
1:B:299:LEU:HD11	1:B:304:LEU:HD13	1.53	0.90
1:A:7:GLN:NE2	1:B:147:GLN:HE21	1.71	0.88
1:B:190:LYS:HB3	1:B:261:MET:HE2	1.56	0.87
1:A:205:ALA:HB1	1:A:315:ARG:CD	2.04	0.87
1:B:72:ILE:HG12	1:B:290:LEU:CD1	2.06	0.84
1:B:284:LYS:HE2	1:B:305:CYS:SG	2.17	0.84
1:B:300:ASN:HB2	1:B:301:LYS:HE3	1.60	0.83
1:A:62:TYR:HB2	1:A:224:MET:HE1	1.61	0.83
1:A:7:GLN:HE22	1:B:147:GLN:NE2	1.75	0.83
1:B:72:ILE:HA	1:B:290:LEU:CD2	2.09	0.83
1:A:145:GLN:HG3	1:A:289:TYR:HB2	1.58	0.82
1:B:301:LYS:H	1:B:301:LYS:CE	1.91	0.82
1:A:178:ASN:HD21	1:B:175:HIS:CD2	1.96	0.82
1:B:301:LYS:HE3	1:B:301:LYS:N	1.96	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:PHE:HD2	1:B:295:SER:HA	1.46	0.80
1:B:286:TRP:CH2	1:B:290:LEU:HG	2.17	0.80
1:B:120:ARG:HB2	5:B:2166:HOH:O	1.81	0.79
1:B:255:GLY:HA2	1:B:262:ALA:HB2	1.64	0.78
1:B:72:ILE:HA	1:B:290:LEU:HD22	1.66	0.78
1:A:127:ARG:HD3	5:A:2173:HOH:O	1.84	0.78
1:A:145:GLN:HG3	1:A:289:TYR:CB	2.13	0.77
1:A:60:ILE:O	1:A:63:GLN:HG2	1.84	0.77
1:B:275:LEU:HD12	1:B:276:PHE:N	2.00	0.76
1:A:7:GLN:NE2	1:B:147:GLN:NE2	2.32	0.76
1:B:292:ARG:HE	1:B:293:ASP:H	1.34	0.76
1:B:149:ARG:HD3	1:B:286:TRP:CB	2.16	0.76
1:B:296:MET:HG2	1:B:297:ILE:N	2.02	0.74
1:A:150:ALA:O	1:A:153:ILE:HG13	1.88	0.74
1:A:62:TYR:HB2	1:A:224:MET:CE	2.16	0.73
1:B:290:LEU:HD12	1:B:291:PHE:CD1	2.23	0.73
1:B:48:TRP:HA	5:B:2068:HOH:O	1.86	0.73
1:B:145:GLN:HA	1:B:145:GLN:NE2	2.02	0.73
1:B:177:VAL:O	1:B:180:LYS:HE2	1.89	0.73
1:B:50:PRO:HG2	1:B:120:ARG:HG2	1.71	0.72
1:B:31:GLN:HG3	1:B:34:ASP:HA	1.71	0.72
1:B:292:ARG:HE	1:B:293:ASP:N	1.89	0.71
1:B:190:LYS:HB3	1:B:261:MET:CE	2.20	0.71
1:B:72:ILE:HG12	1:B:290:LEU:HD13	1.72	0.71
1:B:96:LEU:N	5:B:2169:HOH:O	2.23	0.71
1:B:60:ILE:HG12	1:B:61:ASP:N	2.04	0.70
1:A:175:HIS:HD2	1:B:178:ASN:HD21	1.37	0.70
1:B:79:TYR:CZ	1:B:149:ARG:HG2	2.26	0.70
1:B:72:ILE:HG12	1:B:290:LEU:HD11	1.73	0.70
1:A:143:ASN:O	1:A:147:GLN:HG3	1.93	0.69
1:A:253:ARG:HG3	1:A:253:ARG:HH11	1.56	0.69
1:B:300:ASN:HB2	1:B:301:LYS:CE	2.22	0.69
1:B:51:GLU:CD	1:B:51:GLU:H	1.96	0.69
1:B:95:LEU:N	5:B:2169:HOH:O	2.25	0.69
1:B:272:CYS:O	1:B:275:LEU:HG	1.92	0.69
1:A:145:GLN:HG3	1:A:289:TYR:CG	2.28	0.68
1:B:291:PHE:HB3	1:B:301:LYS:HD3	1.74	0.68
1:A:145:GLN:HA	1:A:145:GLN:NE2	2.08	0.68
1:B:295:SER:HB3	1:B:301:LYS:HE2	1.77	0.67
1:A:138:ASP:HB3	1:B:9:LYS:NZ	2.07	0.67
1:A:178:ASN:HD21	1:B:175:HIS:HD2	1.41	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:ILE:HD13	1:B:297:ILE:N	2.11	0.66
1:B:323:LEU:HD23	1:B:323:LEU:N	2.11	0.66
1:B:290:LEU:HD12	1:B:291:PHE:CE1	2.31	0.66
1:B:92:ASN:O	5:B:2169:HOH:O	2.14	0.66
1:A:169:LEU:HD22	1:B:166:TYR:CE2	2.31	0.66
1:A:253:ARG:NH2	1:A:266:GLU:OE1	2.29	0.66
1:A:57:ARG:HB2	5:A:2164:HOH:O	1.96	0.65
5:A:2112:HOH:O	1:B:147:GLN:HG2	1.96	0.65
1:B:187:ARG:NH2	1:B:267:GLU:OE1	2.29	0.65
1:B:300:ASN:HB2	1:B:301:LYS:HZ1	1.61	0.65
1:B:190:LYS:O	1:B:261:MET:HE1	1.97	0.65
1:A:27:ARG:NH2	1:A:29:ASP:OD2	2.29	0.65
1:A:32:LYS:HE2	1:A:33:TYR:CZ	2.32	0.65
1:A:205:ALA:O	1:A:315:ARG:HD2	1.97	0.64
1:B:141:VAL:N	5:B:2073:HOH:O	2.30	0.64
1:B:52:GLU:HG2	1:B:52:GLU:O	1.96	0.64
1:B:47:PHE:HE2	1:B:49:ARG:CZ	2.10	0.64
1:A:130:VAL:HG12	1:A:132:ASP:H	1.63	0.63
1:A:185:SER:HB3	1:A:188:GLU:HB2	1.79	0.63
1:B:286:TRP:CZ2	1:B:290:LEU:HD23	2.34	0.63
1:B:305:CYS:HB2	5:B:2168:HOH:O	1.98	0.63
1:B:78:LYS:HE2	1:B:139:ASP:HB3	1.78	0.63
1:B:49:ARG:HB3	1:B:51:GLU:OE1	1.97	0.63
1:B:255:GLY:CA	1:B:262:ALA:HB2	2.27	0.63
1:B:286:TRP:CZ3	1:B:290:LEU:HG	2.34	0.62
1:B:143:ASN:OD1	1:B:145:GLN:HB2	1.99	0.62
1:B:268:CYS:HB3	5:B:2141:HOH:O	1.98	0.62
1:A:63:GLN:HG3	1:A:64:ALA:N	2.14	0.62
1:A:31:GLN:HG3	1:A:34:ASP:HA	1.82	0.61
1:B:139:ASP:O	1:B:143:ASN:HB2	1.98	0.61
1:B:222:GLU:O	1:B:222:GLU:HG2	2.00	0.61
1:A:98:LEU:HD21	1:A:164:THR:HG23	1.81	0.61
1:B:138:ASP:O	5:B:2073:HOH:O	2.16	0.61
1:B:291:PHE:CB	1:B:301:LYS:HD3	2.30	0.61
1:B:312:THR:N	5:B:2167:HOH:O	2.32	0.61
1:B:272:CYS:HA	1:B:275:LEU:HG	1.82	0.60
1:B:72:ILE:HA	1:B:290:LEU:HD21	1.81	0.60
1:B:305:CYS:SG	5:B:2135:HOH:O	2.53	0.60
1:B:79:TYR:CE1	1:B:149:ARG:HG2	2.37	0.60
1:A:175:HIS:CD2	1:B:178:ASN:HD21	2.18	0.60
1:B:54:ASP:HB3	5:B:2153:HOH:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:GLU:O	1:B:148:LYS:HB2	2.02	0.59
1:B:295:SER:CB	1:B:301:LYS:HE2	2.32	0.59
1:B:163:MET:HE2	1:B:189:LEU:HD13	1.85	0.59
1:B:300:ASN:CB	1:B:301:LYS:HE3	2.30	0.59
1:A:196:CYS:HG	3:A:2008:HG:HG	1.42	0.58
1:A:31:GLN:NE2	5:A:2116:HOH:O	2.35	0.58
1:B:82:LEU:HD22	1:B:146:ILE:HG22	1.83	0.58
1:B:143:ASN:HB3	1:B:146:ILE:HG12	1.86	0.58
1:B:273:TYR:O	1:B:277:VAL:HG23	2.03	0.58
1:A:63:GLN:HA	1:A:70:LYS:HZ2	1.68	0.58
1:A:221:ARG:CZ	1:A:296:MET:HE3	2.34	0.58
1:A:269:LYS:NZ	1:A:322:ASP:H	2.01	0.58
1:A:305:CYS:HB3	5:A:2151:HOH:O	2.04	0.58
1:A:122:PHE:O	1:A:126:ILE:HG13	2.03	0.58
1:A:55:VAL:HG12	1:A:226:GLY:HA3	1.84	0.57
1:B:143:ASN:CG	1:B:146:ILE:HG12	2.24	0.57
1:A:144:GLU:HA	1:A:147:GLN:NE2	2.20	0.57
1:A:32:LYS:HE2	1:A:33:TYR:CE2	2.38	0.57
1:B:229:LYS:HB3	5:B:2034:HOH:O	2.04	0.57
1:A:327:THR:HG22	1:A:328:ARG:N	2.19	0.57
1:B:300:ASN:HB2	1:B:301:LYS:NZ	2.21	0.56
1:B:312:THR:O	1:B:316:MET:HG3	2.04	0.56
1:A:96:LEU:HB2	1:A:97:PRO:HD3	1.87	0.56
1:A:27:ARG:HH21	1:A:29:ASP:CG	2.07	0.56
1:A:253:ARG:NH1	1:A:253:ARG:HG3	2.15	0.56
1:B:50:PRO:HG3	1:B:121:SER:HB3	1.87	0.56
1:B:306:GLN:OE1	1:B:328:ARG:NH1	2.34	0.56
1:A:315:ARG:NH2	5:A:2148:HOH:O	2.25	0.56
1:B:295:SER:OG	1:B:301:LYS:NZ	2.37	0.56
1:B:31:GLN:NE2	5:B:2084:HOH:O	2.38	0.55
1:A:116:THR:O	1:A:120:ARG:HG3	2.07	0.55
1:B:301:LYS:H	1:B:301:LYS:CD	2.13	0.55
1:B:99:ILE:HD11	1:B:108:VAL:HG21	1.87	0.55
1:B:5:PHE:HD1	1:B:7:GLN:HE21	1.54	0.55
1:A:57:ARG:HG2	1:A:57:ARG:O	2.06	0.55
1:A:304:LEU:HD21	5:A:2149:HOH:O	2.06	0.55
1:A:159:GLU:OE1	1:A:192:LYS:NZ	2.40	0.54
1:A:219:ALA:HB1	1:A:338:TRP:CH2	2.42	0.54
1:B:163:MET:HB3	1:B:189:LEU:HD13	1.89	0.54
1:A:232:ARG:NH1	1:A:338:TRP:HA	2.23	0.54
1:A:149:ARG:NH2	1:A:283:GLU:OE1	2.36	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:ASP:OD1	1:A:259:PRO:HD2	2.07	0.54
1:B:37:GLU:OE2	1:B:41:GLU:OE2	2.26	0.54
1:B:127:ARG:NH1	1:B:127:ARG:HG3	2.23	0.54
1:B:178:ASN:HB2	1:B:180:LYS:CE	2.38	0.54
1:B:295:SER:HB3	1:B:301:LYS:CE	2.38	0.54
1:B:311:ILE:HG23	5:B:2167:HOH:O	2.08	0.54
1:A:62:TYR:HD1	1:A:224:MET:HE2	1.73	0.54
1:B:300:ASN:ND2	1:B:303:ILE:HG13	2.23	0.54
1:A:228:ALA:O	1:A:232:ARG:HG3	2.08	0.53
1:A:118:HIS:CE1	4:A:901:AZI:N1	2.75	0.53
1:B:149:ARG:NE	1:B:286:TRP:CG	2.77	0.53
1:B:35:ILE:HG23	1:B:36:PHE:N	2.24	0.53
1:A:149:ARG:HD2	1:A:286:TRP:HB2	1.91	0.52
1:B:129:ILE:HG13	1:B:130:VAL:HG13	1.91	0.52
1:A:201:ASN:HD21	1:A:246:GLN:HG3	1.74	0.52
1:A:126:ILE:HD12	1:A:137:PHE:CZ	2.45	0.52
1:A:53:VAL:HG11	1:A:230:ILE:HG13	1.92	0.52
1:B:300:ASN:CA	1:B:301:LYS:HE3	2.40	0.52
1:B:69:GLU:HG2	1:B:296:MET:HG3	1.92	0.51
1:B:40:ILE:O	1:B:44:LEU:HG	2.10	0.51
1:A:178:ASN:ND2	1:B:175:HIS:HD2	2.08	0.51
1:B:306:GLN:HG3	1:B:328:ARG:NH1	2.26	0.51
1:B:97:PRO:HD3	5:B:2169:HOH:O	2.10	0.51
1:A:134:SER:O	1:A:138:ASP:HB2	2.10	0.51
1:B:177:VAL:O	1:B:180:LYS:HG2	2.10	0.51
1:B:340:VAL:HG22	5:B:2025:HOH:O	2.10	0.51
1:A:213:ALA:HB3	5:A:2165:HOH:O	2.09	0.51
1:B:68:HIS:HE1	1:B:295:SER:O	1.94	0.51
1:B:311:ILE:HD13	5:B:2167:HOH:O	2.11	0.51
1:A:159:GLU:O	1:A:163:MET:HG3	2.11	0.50
1:B:127:ARG:CG	1:B:127:ARG:HH11	2.23	0.50
1:B:78:LYS:HE3	1:B:136:VAL:HG13	1.93	0.50
1:A:93:VAL:HG13	5:A:2099:HOH:O	2.10	0.50
1:A:138:ASP:C	1:B:9:LYS:HZ1	2.14	0.50
1:B:339:LEU:O	1:B:341:SER:O	2.29	0.50
1:B:55:VAL:HA	5:B:2059:HOH:O	2.10	0.50
1:A:205:ALA:CB	1:A:315:ARG:HD3	2.26	0.50
1:B:278:GLN:O	1:B:282:GLN:HG3	2.11	0.50
1:A:32:LYS:HE2	1:A:33:TYR:OH	2.12	0.49
1:A:203:LEU:HD23	1:A:203:LEU:C	2.32	0.49
1:B:58:ASP:HB3	1:B:224:MET:CE	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:ARG:NH1	5:A:2148:HOH:O	2.29	0.49
1:B:143:ASN:O	1:B:147:GLN:HG3	2.12	0.49
1:B:143:ASN:CB	1:B:146:ILE:HG12	2.41	0.49
1:A:145:GLN:CG	1:A:289:TYR:CD1	2.96	0.49
1:B:31:GLN:HG3	1:B:34:ASP:CA	2.42	0.49
1:A:145:GLN:HG2	1:A:289:TYR:CD1	2.47	0.48
1:A:275:LEU:HD12	1:A:275:LEU:C	2.34	0.48
1:B:332:ILE:HG13	1:B:332:ILE:O	2.13	0.48
1:B:105:GLU:O	1:B:109:GLU:HG3	2.13	0.48
1:A:309:GLU:OE1	1:A:328:ARG:NH1	2.47	0.48
1:B:225:GLU:O	1:B:229:LYS:HG3	2.13	0.48
1:B:290:LEU:HD13	1:B:290:LEU:C	2.34	0.48
1:B:284:LYS:CE	1:B:305:CYS:SG	2.95	0.48
1:B:339:LEU:O	1:B:341:SER:N	2.47	0.48
1:A:62:TYR:O	1:A:70:LYS:NZ	2.46	0.48
1:A:96:LEU:N	1:A:97:PRO:CD	2.77	0.48
1:B:286:TRP:CH2	1:B:290:LEU:CG	2.94	0.48
1:A:63:GLN:CG	1:A:64:ALA:N	2.77	0.47
1:A:79:TYR:CE2	1:A:149:ARG:CZ	2.98	0.47
1:B:290:LEU:HD13	1:B:290:LEU:O	2.13	0.47
1:B:310:TYR:CZ	1:B:330:ASN:HB2	2.49	0.47
1:B:79:TYR:CD2	1:B:149:ARG:CZ	2.98	0.47
1:A:145:GLN:CG	1:A:289:TYR:HB2	2.38	0.46
1:B:232:ARG:CZ	1:B:340:VAL:HG13	2.45	0.46
1:B:306:GLN:CD	1:B:328:ARG:HH11	2.17	0.46
1:B:58:ASP:HB3	1:B:224:MET:HE1	1.96	0.46
1:A:312:THR:O	1:A:316:MET:HG3	2.14	0.46
1:B:122:PHE:O	1:B:126:ILE:HD12	2.15	0.46
1:B:178:ASN:HB2	1:B:180:LYS:HE2	1.97	0.46
1:B:19:PHE:CE1	1:B:98:LEU:HD22	2.51	0.46
1:B:286:TRP:CH2	1:B:290:LEU:CD2	2.98	0.46
1:B:311:ILE:O	1:B:315:ARG:HG2	2.15	0.46
1:B:49:ARG:O	1:B:52:GLU:OE1	2.33	0.46
1:B:131:ASN:N	5:B:2049:HOH:O	2.48	0.46
1:B:145:GLN:HG3	1:B:289:TYR:CD2	2.51	0.46
1:B:232:ARG:NH2	1:B:340:VAL:CG1	2.79	0.46
1:B:232:ARG:NH2	1:B:340:VAL:HG13	2.31	0.46
1:B:332:ILE:N	1:B:333:PRO:HD3	2.30	0.46
1:A:138:ASP:CB	1:B:9:LYS:NZ	2.77	0.46
1:B:291:PHE:HB3	1:B:301:LYS:CD	2.45	0.46
1:A:79:TYR:CD2	1:A:149:ARG:CZ	2.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:GLN:CG	1:A:289:TYR:CG	2.98	0.46
1:B:178:ASN:HB2	1:B:180:LYS:NZ	2.29	0.46
1:B:24:ASN:OD1	1:B:25:VAL:N	2.42	0.45
1:A:35:ILE:HG23	1:A:36:PHE:N	2.30	0.45
1:B:300:ASN:CG	1:B:303:ILE:HG13	2.36	0.45
1:A:62:TYR:O	1:A:65:LEU:HG	2.16	0.45
1:B:191:LYS:HE3	5:B:2141:HOH:O	2.16	0.45
1:A:152:GLY:N	5:A:2038:HOH:O	2.46	0.45
1:A:311:ILE:CG2	1:A:312:THR:N	2.80	0.45
1:A:327:THR:CG2	1:A:328:ARG:N	2.79	0.45
1:B:272:CYS:HB2	5:B:2171:HOH:O	2.16	0.45
1:B:304:LEU:O	1:B:308:VAL:HG23	2.17	0.45
1:B:327:THR:HA	5:B:2093:HOH:O	2.17	0.45
1:A:144:GLU:HA	1:A:147:GLN:CD	2.38	0.45
1:A:250:ASN:ND2	5:A:2111:HOH:O	2.50	0.45
1:B:39:LEU:HA	1:B:39:LEU:HD23	1.74	0.45
1:B:120:ARG:HG3	5:B:2060:HOH:O	2.17	0.44
1:B:286:TRP:CH2	1:B:290:LEU:HD23	2.51	0.44
1:A:62:TYR:CD1	1:A:224:MET:HE2	2.51	0.44
1:A:49:ARG:O	1:A:52:GLU:HB3	2.17	0.44
1:A:74:ILE:CG2	1:A:75:SER:N	2.80	0.44
1:B:149:ARG:NH1	1:B:283:GLU:OE1	2.38	0.44
1:B:217:ALA:HB2	1:B:299:LEU:HD22	2.00	0.44
1:A:4:THR:O	1:B:89:ARG:HD2	2.17	0.44
1:B:187:ARG:HH21	1:B:267:GLU:CD	2.19	0.44
1:B:46:PHE:O	1:B:48:TRP:HD1	2.01	0.44
1:A:82:LEU:CD1	1:B:5:PHE:CE2	3.01	0.44
1:A:7:GLN:CD	1:B:147:GLN:HE21	2.19	0.43
1:B:186:LEU:O	1:B:190:LYS:HG3	2.17	0.43
1:B:21:GLN:HG2	1:B:22:PRO:O	2.18	0.43
1:B:214:CYS:SG	1:B:286:TRP:HZ3	2.42	0.43
1:B:279:ALA:O	1:B:283:GLU:HG2	2.18	0.43
1:B:47:PHE:CE2	1:B:49:ARG:CZ	2.96	0.43
1:B:55:VAL:HG22	5:B:2038:HOH:O	2.17	0.43
1:B:306:GLN:CG	1:B:328:ARG:NH1	2.82	0.43
1:A:138:ASP:CB	1:B:9:LYS:HZ3	2.15	0.43
1:A:171:GLY:O	1:A:175:HIS:HE1	2.01	0.43
1:B:178:ASN:CB	1:B:180:LYS:NZ	2.82	0.43
1:B:242:LEU:HD22	1:B:242:LEU:O	2.19	0.43
1:A:304:LEU:HA	1:A:304:LEU:HD22	1.56	0.43
1:A:323:LEU:HA	1:A:324:PRO:HD3	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:LEU:HD12	1:B:5:PHE:CE2	2.54	0.43
1:B:258:ASP:O	1:B:261:MET:HB2	2.19	0.43
1:B:63:GLN:HA	1:B:70:LYS:HE3	2.01	0.43
1:A:300:ASN:ND2	1:A:303:ILE:HD12	2.35	0.42
1:B:31:GLN:CG	1:B:34:ASP:HB3	2.49	0.42
1:A:130:VAL:HG12	1:A:131:ASN:N	2.34	0.42
1:A:317:GLN:HB2	1:A:323:LEU:HD21	2.01	0.42
1:A:141:VAL:HG22	1:B:25:VAL:HG23	2.01	0.42
1:B:304:LEU:HA	1:B:304:LEU:HD12	1.51	0.42
1:B:292:ARG:HH21	1:B:293:ASP:CG	2.23	0.42
1:B:69:GLU:CD	1:B:221:ARG:HH22	2.23	0.42
1:B:201:ASN:HD21	1:B:246:GLN:HG3	1.85	0.42
1:A:62:TYR:HA	1:A:65:LEU:HG	2.01	0.41
1:B:205:ALA:HB2	1:B:242:LEU:HD21	2.01	0.41
1:A:247:HIS:ND1	5:A:2019:HOH:O	2.29	0.41
1:B:308:VAL:O	5:B:2167:HOH:O	2.22	0.41
1:A:84:ASP:OD2	1:A:122:PHE:HE1	2.04	0.41
1:B:232:ARG:CZ	1:B:340:VAL:CG1	2.99	0.41
1:B:272:CYS:CA	1:B:275:LEU:HG	2.49	0.41
1:A:82:LEU:CD1	1:B:5:PHE:CZ	3.03	0.41
1:B:258:ASP:HA	1:B:259:PRO:HD3	1.85	0.41
1:B:47:PHE:HE2	1:B:49:ARG:NH2	2.16	0.41
1:B:296:MET:HB3	1:B:296:MET:HE3	1.74	0.41
1:B:284:LYS:CD	1:B:305:CYS:SG	3.09	0.41
1:A:143:ASN:CB	1:A:146:ILE:HD12	2.50	0.41
1:A:311:ILE:O	1:A:315:ARG:HG2	2.21	0.41
1:A:4:THR:HG21	1:B:157:TYR:HB3	2.03	0.41
1:B:316:MET:HB3	1:B:321:LEU:HB2	2.02	0.41
1:B:190:LYS:HB2	1:B:264:ILE:HD13	2.03	0.41
1:B:272:CYS:HA	1:B:275:LEU:CD2	2.51	0.41
1:A:65:LEU:HD12	1:A:70:LYS:HG2	2.01	0.41
1:A:96:LEU:HD23	1:A:96:LEU:HA	1.67	0.41
1:B:206:ILE:O	1:B:210:VAL:HG23	2.21	0.40
1:A:121:SER:O	1:A:124:HIS:HB3	2.21	0.40
1:B:163:MET:HE2	1:B:189:LEU:HB2	2.03	0.40
1:B:89:ARG:CG	1:B:90:SER:N	2.84	0.40
1:A:130:VAL:CG1	1:A:131:ASN:N	2.84	0.40
1:A:71:HIS:HD2	5:A:2161:HOH:O	2.03	0.40
1:B:163:MET:HB3	1:B:163:MET:HE2	1.88	0.40
1:B:78:LYS:NZ	5:B:2065:HOH:O	2.54	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	338/375 (90%)	324 (96%)	14 (4%)	0	100	100
1	B	339/375 (90%)	323 (95%)	15 (4%)	1 (0%)	44	49
All	All	677/750 (90%)	647 (96%)	29 (4%)	1 (0%)	55	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	340	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	306/339 (90%)	282 (92%)	24 (8%)	15	15
1	B	307/339 (91%)	266 (87%)	41 (13%)	4	3
All	All	613/678 (90%)	548 (89%)	65 (11%)	8	7

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	31	GLN
1	A	42	LYS
1	A	58	ASP
1	A	77	LEU

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Mol	Chain	Res	Type
1	A	127	ARG
1	A	134	SER
1	A	145	GLN
1	A	147	GLN
1	A	151	GLU
1	A	174	THR
1	A	187	ARG
1	A	188	GLU
1	A	253	ARG
1	A	269	LYS
1	A	275	LEU
1	A	278	GLN
1	A	292	ARG
1	A	302	ASP
1	A	304	LEU
1	A	311	ILE
1	A	315	ARG
1	A	339	LEU
1	A	340	VAL
1	B	9	LYS
1	B	31	GLN
1	B	34	ASP
1	B	38	LYS
1	B	42	LYS
1	B	52	GLU
1	B	60	ILE
1	B	74	ILE
1	B	90	SER
1	B	100	SER
1	B	120	ARG
1	B	127	ARG
1	B	134	SER
1	B	145	GLN
1	B	147	GLN
1	B	148	LYS
1	B	151	GLU
1	B	155	SER
1	B	163	MET
1	B	169	LEU
1	B	180	LYS
1	B	186	LEU
1	B	188	GLU

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Mol	Chain	Res	Type
1	B	206	ILE
1	B	214	CYS
1	B	222	GLU
1	B	242	LEU
1	B	261	MET
1	B	269	LYS
1	B	270	GLN
1	B	275	LEU
1	B	276	PHE
1	B	290	LEU
1	B	292	ARG
1	B	295	SER
1	B	296	MET
1	B	297	ILE
1	B	301	LYS
1	B	305	CYS
1	B	311	ILE
1	B	341	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	68	HIS
1	A	71	HIS
1	A	147	GLN
1	A	175	HIS
1	A	201	ASN
1	A	246	GLN
1	A	250	ASN
1	A	278	GLN
1	B	7	GLN
1	B	12	GLN
1	B	30	GLN
1	B	31	GLN
1	B	68	HIS
1	B	76	ASN
1	B	80	GLN
1	B	87	GLN
1	B	128	ASN
1	B	131	ASN
1	B	145	GLN

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Mol	Chain	Res	Type
1	B	147	GLN
1	B	168	HIS
1	B	175	HIS
1	B	201	ASN
1	B	247	HIS

### 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](#)

Of 17 ligands modelled in this entry, 16 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	AZI	A	901	2	0,2,2	0.00	-	0,1,1	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	AZI	A	901	2	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	901	AZI	1	0

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	340/375 (90%)	-0.51	3 (0%) 84 83	14, 25, 51, 73	0
1	B	341/375 (90%)	-0.26	13 (3%) 41 39	12, 28, 58, 84	0
All	All	681/750 (90%)	-0.38	16 (2%) 61 58	12, 27, 56, 84	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	340	VAL	6.4
1	B	297	ILE	6.3
1	B	289	TYR	4.9
1	B	291	PHE	4.1
1	B	341	SER	3.6
1	B	292	ARG	3.5
1	B	293	ASP	3.1
1	B	290	LEU	2.8
1	A	340	VAL	2.5
1	A	63	GLN	2.5
1	B	294	GLY	2.4
1	A	337	THR	2.4
1	B	296	MET	2.2
1	B	135	VAL	2.2
1	B	222	GLU	2.2
1	B	131	ASN	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	AZI	A	901	3/3	0.91	0.23	5.76	22,22,30,39	3
3	HG	B	2006	1/1	0.97	0.10	-0.78	35,35,35,35	1
3	HG	B	2013	1/1	0.98	0.05	-2.20	26,26,26,26	1
3	HG	A	2001	1/1	1.00	0.02	-2.27	24,24,24,24	1
3	HG	B	2002	1/1	0.97	0.06	-2.42	39,39,39,39	1
2	FE	A	1003	1/1	0.99	0.04	-2.55	22,22,22,22	1
3	HG	A	2007	1/1	1.00	0.02	-2.62	24,24,24,24	1
2	FE	B	1002	1/1	0.99	0.06	-3.18	24,24,24,24	0
3	HG	B	2010	1/1	0.99	0.03	-3.25	42,42,42,42	1
3	HG	A	2012	1/1	1.00	0.04	-3.38	31,31,31,31	1
3	HG	A	2008	1/1	0.97	0.06	-4.01	37,37,37,37	1
3	HG	B	2004	1/1	1.00	0.03	-4.07	23,23,23,23	1
2	FE	A	1004	1/1	1.00	0.03	-4.48	17,17,17,17	0
3	HG	B	2003	1/1	0.97	0.04	-4.58	49,49,49,49	1
2	FE	B	1001	1/1	0.98	0.04	-4.93	41,41,41,41	1
3	HG	A	2011	1/1	0.93	0.18	-	40,40,40,40	1
3	HG	A	2009	1/1	1.00	0.03	-	24,24,24,24	1

### 6.5 Other polymers [i](#)

There are no such residues in this entry.