



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:33 AM GMT

PDB ID : 3F5S
Title : CRYSTAL STRUCTURE OF putative short chain dehydrogenase from
Shigella flexneri 2a str. 301
Authors : Malashkevich, V.N.; Toro, R.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New
York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2008-11-04
Resolution : 1.36 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

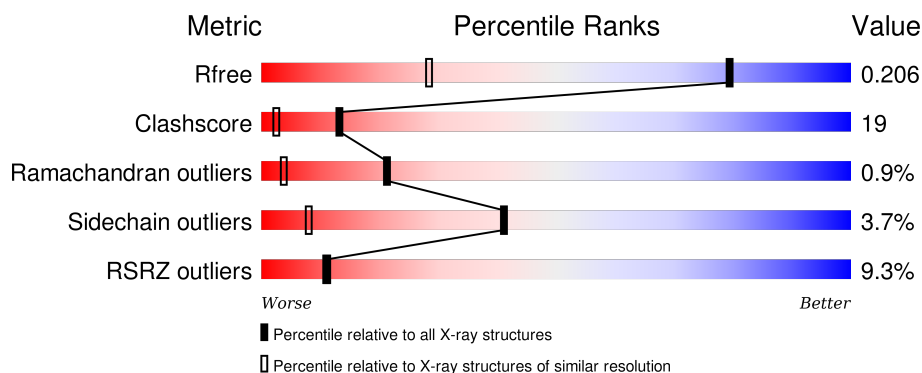
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.36 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	2199 (1.40-1.32)
Clashscore	102246	2337 (1.40-1.32)
Ramachandran outliers	100387	2280 (1.40-1.32)
Sidechain outliers	100360	2279 (1.40-1.32)
RSRZ outliers	91569	2199 (1.40-1.32)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	255	<div> <div>9%</div> <div>70%</div> <div>15%</div> <div>•</div> <div>11%</div> </div>
1	B	255	<div> <div>8%</div> <div>75%</div> <div>11%</div> <div>•</div> <div>11%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3837 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 dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	4	0
			1786	1114	324	335	13			
1	B	227	Total	C	N	O	S	0	5	0
			1800	1125	326	336	13			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q83RM3
A	2	SER	-	expression tag	UNP Q83RM3
A	3	LEU	-	expression tag	UNP Q83RM3
A	248	GLU	-	expression tag	UNP Q83RM3
A	249	GLY	-	expression tag	UNP Q83RM3
A	250	HIS	-	expression tag	UNP Q83RM3
A	251	HIS	-	expression tag	UNP Q83RM3
A	252	HIS	-	expression tag	UNP Q83RM3
A	253	HIS	-	expression tag	UNP Q83RM3
A	254	HIS	-	expression tag	UNP Q83RM3
A	255	HIS	-	expression tag	UNP Q83RM3
B	1	MET	-	expression tag	UNP Q83RM3
B	2	SER	-	expression tag	UNP Q83RM3
B	3	LEU	-	expression tag	UNP Q83RM3
B	248	GLU	-	expression tag	UNP Q83RM3
B	249	GLY	-	expression tag	UNP Q83RM3
B	250	HIS	-	expression tag	UNP Q83RM3
B	251	HIS	-	expression tag	UNP Q83RM3
B	252	HIS	-	expression tag	UNP Q83RM3
B	253	HIS	-	expression tag	UNP Q83RM3
B	254	HIS	-	expression tag	UNP Q83RM3
B	255	HIS	-	expression tag	UNP Q83RM3

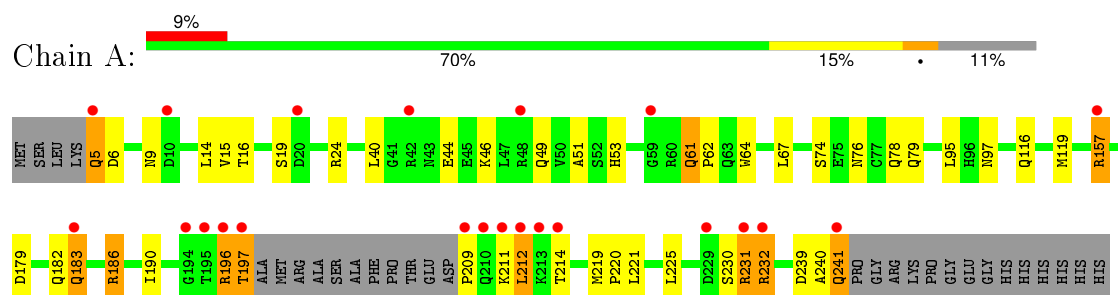
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	251	Total 251	O 251	0	0

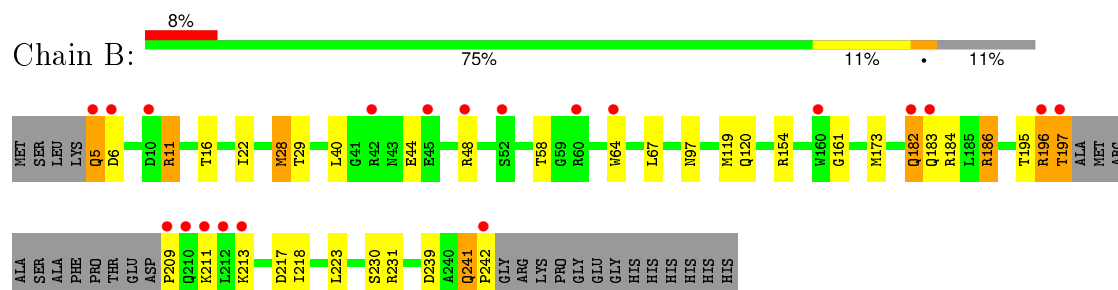
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: dehydrogenase



• Molecule 1: dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.99Å 67.41Å 65.03Å 90.00° 108.69° 90.00°	Depositor
Resolution (Å)	7.99 – 1.36 7.97 – 1.36	Depositor EDS
% Data completeness (in resolution range)	100.0 (7.99-1.36) 94.1 (7.97-1.36)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 1.36Å)	Xtriage
Refinement program	REFMAC 5.4.0069	Depositor
R, R_{free}	0.184 , 0.208 0.183 , 0.206	Depositor DCC
R_{free} test set	4452 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	11.4	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.62 , 68.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 88705 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3837	wwPDB-VP
Average B, all atoms (Å ²)	15.0	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 41.36 % 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 2.3866e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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 [i](#)

5.1 Standard geometry [i](#)

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	0.55	0/1823	0.70	1/2467 (0.0%)
1	B	0.56	0/1841	0.77	2/2492 (0.1%)
All	All	0.56	0/3664	0.73	3/4959 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	186	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	B	28	MET	CG-SD-CE	5.74	109.38	100.20
1	B	173	MET	CA-CB-CG	5.54	122.72	113.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	232	ARG	Peptide

5.2 Too-close contacts [i](#)

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	1786	0	1792	77	0
1	B	1800	0	1813	64	0
2	B	251	0	0	16	2
All	All	3837	0	3605	136	2

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 19.

All (136) 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:231:ARG:NH1	1:A:232:ARG:NH1	1.70	1.38
1:B:44:GLU:HG2	1:B:64:TRP:CH2	1.70	1.27
1:B:28:MET:CE	1:B:58:THR:CG2	2.19	1.19
1:B:11[A]:ARG:HG3	1:B:11[A]:ARG:HH11	1.03	1.17
1:A:231:ARG:HH11	1:A:232:ARG:NH1	1.32	1.13
1:B:241:GLN:CB	1:B:242:PRO:HD3	1.78	1.13
1:B:44:GLU:CG	1:B:64:TRP:CH2	2.32	1.11
1:A:239:ASP:HB3	1:A:241:GLN:HB2	1.12	1.10
1:A:231:ARG:NH1	1:A:232:ARG:CZ	2.14	1.10
1:A:231:ARG:HH21	1:A:231:ARG:CG	1.67	1.07
1:B:28:MET:HE1	1:B:58:THR:CG2	1.80	1.05
1:A:231:ARG:HG2	1:A:231:ARG:HH21	0.90	1.05
1:B:186[B]:ARG:HD2	2:B:483:HOH:O	1.58	1.03
1:B:28:MET:HE2	1:B:58:THR:CG2	1.88	1.03
1:B:28:MET:HE1	1:B:58:THR:HG23	1.38	1.01
1:B:28:MET:CE	1:B:58:THR:HG21	1.90	0.97
1:A:239:ASP:CB	1:A:241:GLN:HB2	1.94	0.97
1:B:241:GLN:HB2	1:B:242:PRO:HD3	1.47	0.97
1:A:231:ARG:HH11	1:A:232:ARG:HH11	1.01	0.97
1:A:157[A]:ARG:CG	1:A:157[A]:ARG:HH11	1.78	0.96
1:A:231:ARG:HD2	1:A:232:ARG:H	1.30	0.96
1:A:182:GLN:O	1:A:183:GLN:HG2	1.65	0.95
1:A:231:ARG:HG2	1:A:231:ARG:NH2	1.74	0.95
1:A:241:GLN:O	1:A:241:GLN:HG3	1.63	0.95
1:B:241:GLN:HB3	1:B:242:PRO:HD3	1.45	0.94
1:B:11[A]:ARG:HG3	1:B:11[A]:ARG:NH1	1.84	0.89
1:A:241:GLN:O	1:A:241:GLN:CG	2.21	0.88
1:B:48:ARG:HH21	1:B:48:ARG:HG2	1.38	0.87
1:A:157[A]:ARG:HG3	1:A:157[A]:ARG:HH11	1.40	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:GLN:HA	2:B:286:HOH:O	1.75	0.86
1:B:241:GLN:CB	1:B:242:PRO:CD	2.52	0.86
1:A:239:ASP:HB3	1:A:241:GLN:CB	2.04	0.86
1:B:241:GLN:HB2	1:B:242:PRO:CD	2.06	0.85
1:A:182:GLN:O	1:A:183:GLN:CG	2.24	0.85
1:A:231:ARG:HD2	1:A:232:ARG:N	1.95	0.81
1:B:154[A]:ARG:NH2	1:B:239:ASP:OD1	2.15	0.81
1:A:119:MET:CE	1:B:119:MET:CE	2.61	0.79
1:A:196:ARG:HD2	1:A:212:LEU:CD1	2.13	0.79
1:B:44:GLU:HG2	1:B:64:TRP:CZ3	2.18	0.79
1:B:11[A]:ARG:NH1	2:B:265:HOH:O	2.16	0.78
1:B:28:MET:CE	1:B:58:THR:HG23	1.98	0.77
1:B:44:GLU:HG3	1:B:64:TRP:CH2	2.18	0.77
1:B:28:MET:HE2	1:B:58:THR:HG22	1.66	0.77
1:A:231:ARG:HH12	1:A:232:ARG:NH1	1.84	0.76
1:A:231:ARG:CG	1:A:231:ARG:NH2	2.35	0.74
1:B:11[A]:ARG:CG	1:B:11[A]:ARG:HH11	1.91	0.74
1:B:213:LYS:HE3	1:B:217:ASP:HB3	1.71	0.73
1:B:44:GLU:CG	1:B:64:TRP:HH2	2.01	0.73
1:B:186[A]:ARG:CD	2:B:505:HOH:O	2.36	0.72
1:B:44:GLU:HG3	1:B:64:TRP:HH2	1.54	0.72
1:A:196:ARG:HD2	1:A:212:LEU:HD12	1.70	0.72
1:A:119:MET:HE3	1:B:119:MET:CE	2.21	0.70
1:B:196:ARG:HD2	1:B:209:PRO:N	2.06	0.69
1:A:119:MET:CE	1:B:119:MET:HE3	2.24	0.68
1:B:28:MET:HE1	1:B:58:THR:HG21	1.60	0.68
1:B:5:GLN:HG2	1:B:6:ASP:OD1	1.95	0.67
1:B:195:THR:O	1:B:197:THR:N	2.24	0.66
1:B:197:THR:OG1	1:B:197:THR:O	2.13	0.66
1:B:186[A]:ARG:HD3	2:B:505:HOH:O	1.93	0.66
1:A:40:LEU:HD21	1:A:67:LEU:HB3	1.77	0.65
1:B:186[B]:ARG:NH1	2:B:502:HOH:O	2.10	0.65
1:B:48:ARG:HG2	1:B:48:ARG:NH2	2.05	0.65
1:A:241:GLN:HA	2:B:462:HOH:O	1.97	0.64
1:B:195:THR:C	1:B:197:THR:H	2.00	0.63
1:A:231:ARG:CZ	1:A:232:ARG:CZ	2.75	0.63
1:A:6:ASP:HB2	1:A:9:ASN:ND2	2.13	0.63
1:A:44:GLU:HG3	1:A:64:TRP:CH2	2.35	0.61
1:A:197:THR:HG23	2:B:276:HOH:O	2.01	0.61
1:A:240:ALA:N	1:A:241:GLN:HA	2.15	0.60
1:B:5:GLN:HG2	1:B:6:ASP:H	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157[A]:ARG:HG2	1:A:157[A]:ARG:HH11	1.63	0.59
1:A:196:ARG:HA	1:A:196:ARG:NE	2.16	0.58
1:A:196:ARG:HA	1:A:196:ARG:HE	1.69	0.58
1:B:44:GLU:HG2	1:B:64:TRP:CZ2	2.35	0.57
1:B:184:ARG:O	2:B:416:HOH:O	2.18	0.57
1:B:183:GLN:N	2:B:286:HOH:O	2.38	0.56
1:A:239:ASP:C	1:A:241:GLN:CB	2.74	0.56
1:A:211:LYS:HA	1:A:211:LYS:HE3	1.87	0.56
1:A:157[A]:ARG:CG	1:A:157[A]:ARG:NH1	2.50	0.55
1:B:6:ASP:OD1	1:B:6:ASP:N	2.40	0.55
1:A:44:GLU:CG	1:A:64:TRP:CH2	2.89	0.55
1:A:221:LEU:HD22	1:A:240:ALA:HA	1.89	0.55
1:A:239:ASP:O	1:A:241:GLN:HB3	2.08	0.54
1:A:196:ARG:N	1:A:196:ARG:HD2	2.22	0.54
1:A:197:THR:O	1:A:197:THR:CG2	2.54	0.54
1:B:242:PRO:HD2	2:B:274:HOH:O	2.07	0.54
1:B:40:LEU:HD21	1:B:67:LEU:HB3	1.91	0.52
1:A:241:GLN:CA	2:B:462:HOH:O	2.58	0.51
1:B:182:GLN:CA	2:B:286:HOH:O	2.46	0.51
1:A:196:ARG:HG2	1:A:214:THR:HA	1.93	0.51
1:A:239:ASP:CB	1:A:241:GLN:CB	2.76	0.51
1:A:157[A]:ARG:NH1	1:A:157[A]:ARG:HG3	2.18	0.50
1:B:28:MET:O	1:B:28:MET:HE3	2.11	0.50
1:A:157[B]:ARG:CZ	2:B:264:HOH:O	2.60	0.50
1:A:49:GLN:NE2	2:B:291:HOH:O	2.45	0.50
1:A:116[B]:GLN:NE2	1:B:120:GLN:HE22	2.10	0.49
1:A:219:MET:N	1:A:220:PRO:CD	2.76	0.49
1:B:213:LYS:HE3	1:B:217:ASP:CB	2.41	0.49
1:A:196:ARG:N	1:A:212:LEU:HD13	2.28	0.49
1:A:221:LEU:CD2	1:A:240:ALA:HA	2.43	0.49
1:A:240:ALA:N	1:A:241:GLN:CA	2.76	0.49
1:A:51:ALA:HB1	1:A:61:GLN:OE1	2.12	0.48
1:A:190:ILE:HD11	1:A:225:LEU:HD12	1.94	0.48
1:A:196:ARG:N	1:A:212:LEU:CD1	2.77	0.48
1:A:76:ASN:HA	1:A:79:GLN:HE21	1.80	0.47
1:A:231:ARG:CZ	1:A:232:ARG:NE	2.77	0.46
1:B:195:THR:HG23	1:B:197:THR:N	2.31	0.46
1:B:29:THR:HG22	1:B:223:LEU:HD21	1.98	0.46
1:A:44:GLU:HG3	1:A:64:TRP:CZ3	2.50	0.45
1:A:5:GLN:HB2	1:A:6:ASP:H	1.53	0.45
1:B:186[B]:ARG:HD3	1:B:230:SER:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186[A]:ARG:HD2	2:B:505:HOH:O	2.10	0.45
1:A:95:LEU:HD23	1:A:97[B]:ASN:HD21	1.81	0.45
1:A:74:SER:O	1:A:78:GLN:HG2	2.17	0.44
1:B:186[B]:ARG:HD3	1:B:231:ARG:HA	1.99	0.44
1:A:196:ARG:HD2	1:A:212:LEU:HD11	1.95	0.44
1:A:51:ALA:HA	1:A:62:PRO:HG2	1.99	0.44
1:A:24[A]:ARG:NH2	1:A:53:HIS:ND1	2.65	0.44
1:A:119:MET:HE3	1:B:119:MET:HE2	1.96	0.44
1:A:196:ARG:H	1:A:196:ARG:HD2	1.82	0.44
1:B:196:ARG:CD	1:B:209:PRO:N	2.79	0.44
1:A:19:SER:HB2	1:A:46:LYS:HB3	1.99	0.43
1:A:196:ARG:NH2	1:A:209:PRO:HB2	2.33	0.43
1:B:22[A]:ILE:HD12	1:B:218:ILE:HD11	1.99	0.43
1:A:95:LEU:CD2	1:A:97[B]:ASN:HD21	2.32	0.42
1:B:16:THR:O	1:B:97:ASN:HB3	2.19	0.42
1:A:186:ARG:HD3	1:A:230:SER:O	2.19	0.42
1:A:179:ASP:O	1:A:182:GLN:NE2	2.53	0.42
1:A:157[A]:ARG:NH1	1:A:157[A]:ARG:HG2	2.27	0.41
1:B:48:ARG:HH21	1:B:48:ARG:CG	2.20	0.41
1:B:44:GLU:CG	1:B:64:TRP:CZ3	2.91	0.41
1:A:14:LEU:HD23	1:A:14:LEU:C	2.41	0.41
1:B:48:ARG:CG	1:B:48:ARG:NH2	2.80	0.41
1:A:16:THR:O	1:A:97[B]:ASN:HB2	2.21	0.41
1:A:15:VAL:HG13	1:A:97[B]:ASN:ND2	2.36	0.41
1:B:5:GLN:HB3	1:B:5:GLN:HE21	1.57	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:313:HOH:O	2:B:340:HOH:O[2_646]	1.40	0.80
2:B:270:HOH:O	2:B:474:HOH:O[1_554]	1.99	0.21

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	226/255 (89%)	217 (96%)	8 (4%)	1 (0%)	39	14
1	B	228/255 (89%)	221 (97%)	4 (2%)	3 (1%)	15	2
All	All	454/510 (89%)	438 (96%)	12 (3%)	4 (1%)	21	4

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	241	GLN
1	B	196	ARG
1	A	196	ARG
1	B	161	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	193/212 (91%)	184 (95%)	9 (5%)	32	4
1	B	195/212 (92%)	187 (96%)	8 (4%)	37	6
All	All	388/424 (92%)	371 (96%)	17 (4%)	41	5

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	61	GLN
1	A	157[A]	ARG
1	A	157[B]	ARG
1	A	183	GLN
1	A	197	THR
1	A	212	LEU
1	A	231	ARG
1	A	241	GLN

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Mol	Chain	Res	Type
1	B	5	GLN
1	B	11[A]	ARG
1	B	11[B]	ARG
1	B	182	GLN
1	B	186[A]	ARG
1	B	186[B]	ARG
1	B	197	THR
1	B	211	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	79	GLN
1	A	155	GLN
1	A	159	ASN
1	B	5	GLN
1	B	63	GLN
1	B	87	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	226/255 (88%)	0.50	22 (9%)	10 10	8, 13, 30, 42	0
1	B	227/255 (89%)	0.39	20 (8%)	12 12	8, 12, 29, 42	0
All	All	453/510 (88%)	0.44	42 (9%)	11 11	8, 13, 30, 42	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	242	PRO	11.1
1	B	197	THR	8.4
1	A	197	THR	8.4
1	B	209	PRO	7.0
1	B	211	LYS	6.7
1	A	232	ARG	6.6
1	A	241	GLN	5.8
1	A	212	LEU	4.8
1	A	213	LYS	4.4
1	A	211	LYS	4.2
1	A	196	ARG	4.2
1	B	196	ARG	3.9
1	A	183	GLN	3.8
1	A	5	GLN	3.7
1	B	183	GLN	3.2
1	A	42	ARG	3.1
1	B	5	GLN	3.0
1	B	210	GLN	2.9
1	B	213	LYS	2.8
1	A	20	ASP	2.8
1	A	157[A]	ARG	2.7
1	B	60	ARG	2.6
1	B	42	ARG	2.6
1	A	59	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	212	LEU	2.6
1	A	214	THR	2.6
1	A	195	THR	2.5
1	B	160	TRP	2.4
1	A	48	ARG	2.4
1	B	64	TRP	2.3
1	B	182	GLN	2.3
1	A	210	GLN	2.3
1	A	231	ARG	2.3
1	B	10[A]	ASP	2.3
1	A	10	ASP	2.3
1	A	194	GLY	2.2
1	A	209	PRO	2.2
1	B	45	GLU	2.1
1	B	52	SER	2.1
1	A	229	ASP	2.0
1	B	6	ASP	2.0
1	B	48	ARG	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.