



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 08:57 AM UTC

PDB ID : 3CDL / pdb_00003cdl
Title : Crystal structure of a TetR family transcriptional regulator from *Pseudomonas syringae* pv. tomato str. DC3000
Authors : Tan, K.; Bigelow, L.; Gu, M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2008-02-27
Resolution : 2.36 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

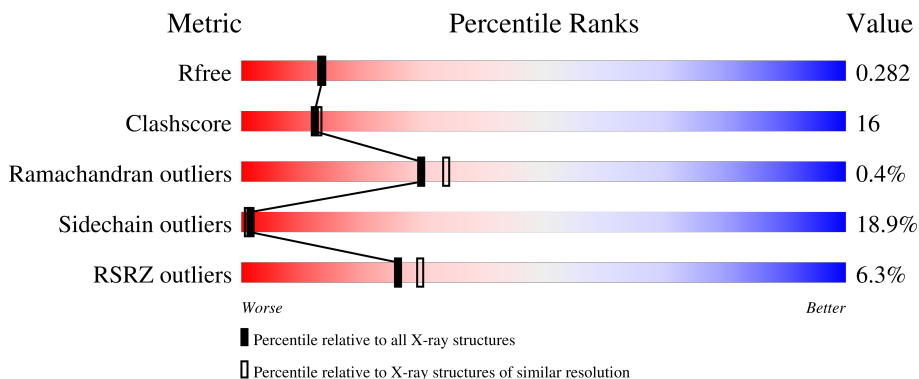
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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}	180053	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	203	
1	B	203	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3068 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 Transcriptional regulator AefR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	187	1507	967	261	273	1	5	0	0	0
1	B	188	1520	974	266	273	1	6	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q87Z81
A	-1	ASN	-	expression tag	UNP Q87Z81
A	0	ALA	-	expression tag	UNP Q87Z81
B	-2	SER	-	expression tag	UNP Q87Z81
B	-1	ASN	-	expression tag	UNP Q87Z81
B	0	ALA	-	expression tag	UNP Q87Z81

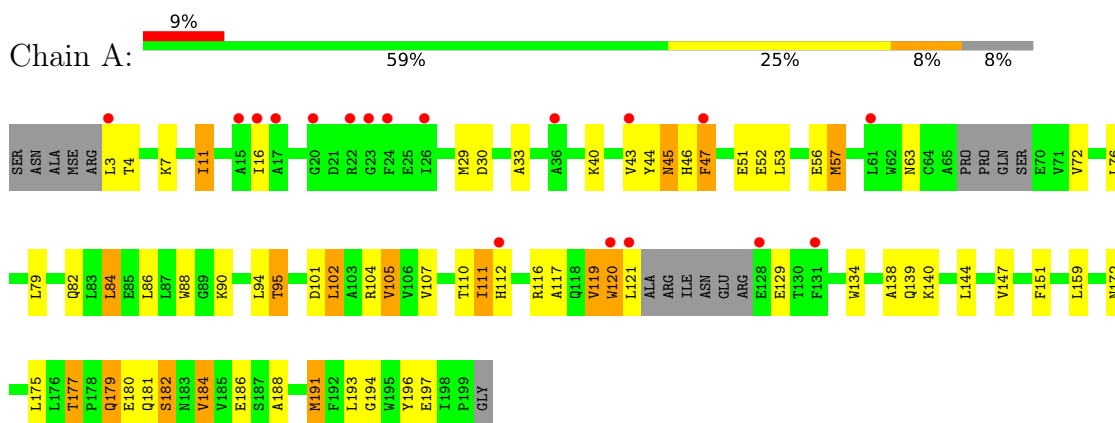
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	24	Total 24	O 24	0	0
2	B	17	Total 17	O 17	0	0

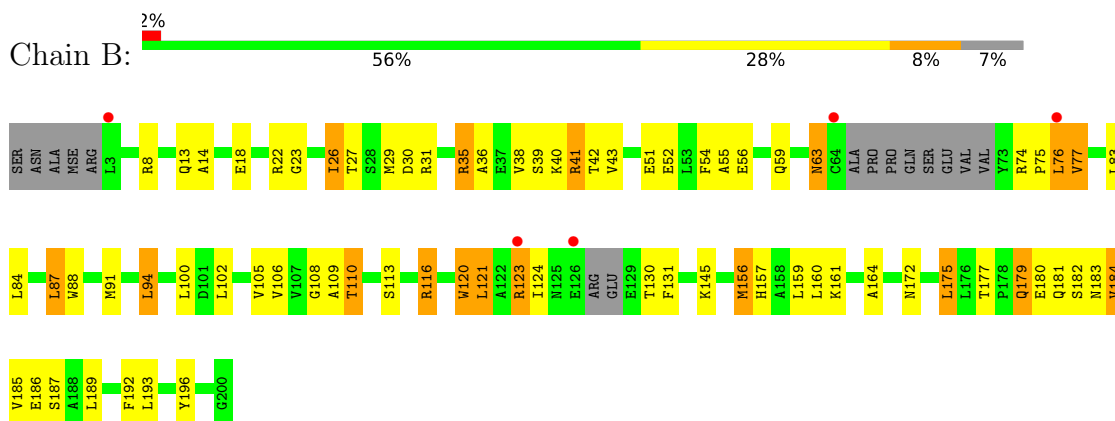
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Transcriptional regulator AefR



- Molecule 1: Transcriptional regulator AefR



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	132.31Å 53.41Å 55.83Å 90.00° 113.25° 90.00°	Depositor
Resolution (Å)	33.26 – 2.36 33.26 – 2.36	Depositor EDS
% Data completeness (in resolution range)	96.3 (33.26-2.36) 96.5 (33.26-2.36)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.36Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.207 , 0.275 (Not available) , 0.282	Depositor DCC
R_{free} test set	728 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	43.3	Xtrriage
Anisotropy	0.072	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.025 for -h+k-1,-l,-k 0.000 for -h-k-1,l,k 0.029 for -h-2*1,-k,l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3068	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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	0.90	2/1536 (0.1%)	1.13	6/2073 (0.3%)
1	B	0.78	2/1552 (0.1%)	1.07	4/2092 (0.2%)
All	All	0.84	4/3088 (0.1%)	1.10	10/4165 (0.2%)

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	105	VAL	CA-CB	6.97	1.62	1.54
1	B	41	ARG	NE-CZ	6.51	1.40	1.33
1	B	41	ARG	CZ-NH1	5.74	1.40	1.32
1	A	63	ASN	CG-OD1	5.51	1.34	1.23

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	74	ARG	CA-C-N	6.80	128.34	119.84
1	B	74	ARG	C-N-CA	6.80	128.34	119.84
1	A	120	TRP	N-CA-C	6.62	119.98	112.97
1	A	177	THR	CA-C-N	6.11	127.48	119.84
1	A	177	THR	C-N-CA	6.11	127.48	119.84
1	A	105	VAL	N-CA-CB	5.88	117.43	110.55
1	A	47	PHE	CA-C-N	5.71	126.98	119.84
1	A	47	PHE	C-N-CA	5.71	126.98	119.84
1	B	77	VAL	N-CA-C	-5.13	101.01	109.12
1	B	36	ALA	N-CA-C	-5.08	106.92	113.01

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	76	LEU	Peptide

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	1507	0	1493	50	0
1	B	1520	0	1511	50	0
2	A	24	0	0	0	0
2	B	17	0	0	2	0
All	All	3068	0	3004	94	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 16.

All (94) 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:179:GLN:H	1:B:179:GLN:HE21	1.17	0.93
1:A:16:ILE:HA	1:A:57:MSE:HE1	1.50	0.92
1:A:7:LYS:O	1:A:11:ILE:HG23	1.71	0.90
1:A:111:ILE:HD11	1:B:108:GLY:HA2	1.52	0.89
1:B:63:ASN:HB3	2:B:207:HOH:O	1.80	0.82
1:A:182:SER:O	1:A:186:GLU:HG2	1.82	0.78
1:B:38:VAL:HG22	1:B:42:THR:OG1	1.83	0.77
1:B:110:THR:HG21	1:B:120:TRP:HD1	1.49	0.75
1:B:83:LEU:HD11	1:B:156[B]:MSE:HE3	1.67	0.74
1:A:45:ASN:C	1:A:45:ASN:HD22	1.96	0.74
1:A:111:ILE:CD1	1:B:108:GLY:HA2	2.19	0.71
1:A:191:MSE:HE3	1:B:196:TYR:CE2	2.26	0.70
1:B:106:VAL:O	1:B:110:THR:HG23	1.93	0.69
1:A:16:ILE:HG12	1:A:57:MSE:HE1	1.76	0.67
1:B:113:SER:OG	1:B:116:ARG:HB2	1.95	0.67
1:A:191:MSE:HE3	1:B:196:TYR:HE2	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:GLN:H	1:B:179:GLN:NE2	1.91	0.65
1:A:79:LEU:HD12	1:A:197:GLU:HG3	1.80	0.64
1:B:182:SER:O	1:B:186:GLU:HB2	1.96	0.64
1:B:83:LEU:HD21	1:B:156[B]:MSE:CE	2.28	0.63
1:A:180:GLU:O	1:A:184:VAL:HG13	2.00	0.62
1:B:55:ALA:O	1:B:59:GLN:HG2	1.99	0.61
1:B:51:GLU:HG3	2:B:216:HOH:O	2.01	0.61
1:B:116:ARG:HH11	1:B:116:ARG:HG2	1.64	0.61
1:A:51:GLU:HA	1:A:51:GLU:OE1	1.99	0.60
1:A:57:MSE:HE3	1:A:102:LEU:HD13	1.84	0.59
1:B:14:ALA:HB1	1:B:35:ARG:HG2	1.85	0.58
1:B:59:GLN:O	1:B:63:ASN:HB2	2.03	0.58
1:A:119:VAL:HG22	1:A:120:TRP:CD1	2.39	0.58
1:A:116:ARG:HA	1:A:119:VAL:HG13	1.87	0.56
1:A:45:ASN:HD22	1:A:46:HIS:N	2.05	0.55
1:B:38:VAL:CG1	1:B:43:VAL:HG23	2.37	0.55
1:A:16:ILE:HG12	1:A:57:MSE:CE	2.38	0.53
1:B:87:LEU:HB3	1:B:185:VAL:HG13	1.91	0.53
1:B:83:LEU:HD21	1:B:156[B]:MSE:HE2	1.91	0.53
1:B:59:GLN:HE22	1:B:123:ARG:HD3	1.74	0.53
1:A:139:GLN:NE2	1:A:147:VAL:HG12	2.24	0.52
1:A:151:PHE:HZ	1:B:187:SER:HG	1.58	0.52
1:B:88:TRP:HZ2	1:B:182:SER:HB3	1.75	0.52
1:A:79:LEU:CD1	1:A:197:GLU:HG3	2.38	0.52
1:B:26:ILE:O	1:B:31:ARG:NH1	2.42	0.52
1:A:16:ILE:CA	1:A:57:MSE:HE1	2.31	0.51
1:A:144:LEU:HD22	1:A:196:TYR:HB3	1.92	0.51
1:B:177:THR:OG1	1:B:180:GLU:HG3	2.10	0.51
1:B:23:GLY:O	1:B:27:THR:OG1	2.23	0.51
1:A:16:ILE:CG1	1:A:57:MSE:HE1	2.39	0.51
1:A:179:GLN:HE21	1:A:179:GLN:H	1.58	0.50
1:A:16:ILE:HA	1:A:57:MSE:CE	2.32	0.50
1:B:110:THR:HG21	1:B:120:TRP:CD1	2.39	0.50
1:A:144:LEU:HD23	1:A:197:GLU:HA	1.93	0.49
1:A:11:ILE:HD13	1:A:43:VAL:HG13	1.95	0.49
1:B:109:ALA:O	1:B:116:ARG:NH1	2.46	0.49
1:B:179:GLN:HE21	1:B:179:GLN:N	1.99	0.49
1:A:95:THR:CG2	1:A:95:THR:O	2.61	0.48
1:B:91:MSE:HE2	1:B:160:LEU:HG	1.94	0.48
1:A:30:ASP:HA	1:A:33:ALA:HB3	1.95	0.47
1:B:180:GLU:O	1:B:184:VAL:HG13	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:GLN:HG3	1:A:134:TRP:HZ2	1.78	0.47
1:A:30:ASP:HB3	1:A:40:LYS:HG3	1.97	0.47
1:A:107:VAL:O	1:A:111:ILE:HG23	2.16	0.46
1:B:100:LEU:HD12	1:B:175:LEU:HD11	1.98	0.45
1:B:38:VAL:CG2	1:B:42:THR:OG1	2.59	0.45
1:A:101:ASP:OD1	1:A:104:ARG:NH2	2.50	0.45
1:B:54:PHE:CE2	1:B:116:ARG:HG3	2.51	0.45
1:A:191:MSE:HE1	1:B:192:PHE:CD1	2.52	0.45
1:A:193:LEU:O	1:A:194:GLY:C	2.60	0.44
1:A:40:LYS:HE2	1:A:44:TYR:HE1	1.82	0.44
1:B:94:LEU:HD23	1:B:164:ALA:HB3	1.99	0.44
1:A:177:THR:O	1:A:181:GLN:HG3	2.18	0.44
1:B:29:MSE:HE3	1:B:43:VAL:HG12	1.99	0.43
1:B:39:SER:O	1:B:40:LYS:C	2.60	0.43
1:A:82:GLN:HG3	1:A:134:TRP:CZ2	2.53	0.43
1:B:18:GLU:HG2	1:B:35:ARG:HD3	2.00	0.43
1:B:193:LEU:HD23	1:B:193:LEU:HA	1.79	0.43
1:A:179:GLN:H	1:A:179:GLN:NE2	2.16	0.43
1:A:84:LEU:HD22	1:A:88:TRP:HD1	1.84	0.43
1:B:131:PHE:CD1	1:B:157:HIS:CE1	3.07	0.43
1:A:11:ILE:HD11	1:A:47:PHE:CD2	2.54	0.42
1:A:138:ALA:HB3	1:A:144:LEU:HD12	2.00	0.42
1:A:159:LEU:HD13	1:A:188:ALA:HA	2.02	0.42
1:A:111:ILE:HG12	1:A:112:HIS:N	2.34	0.42
1:B:181:GLN:O	1:B:185:VAL:HG23	2.20	0.41
1:B:38:VAL:HG13	1:B:43:VAL:HG23	2.01	0.41
1:B:54:PHE:CZ	1:B:116:ARG:HG3	2.56	0.41
1:A:95:THR:O	1:A:95:THR:HG22	2.20	0.41
1:A:110:THR:HB	1:A:117:ALA:HB2	2.03	0.41
1:A:111:ILE:CG1	1:A:112:HIS:N	2.84	0.41
1:B:22:ARG:HB3	1:B:26:ILE:HD11	2.03	0.41
1:B:121:LEU:HA	1:B:124:ILE:HG22	2.02	0.40
1:A:116:ARG:O	1:A:120:TRP:HD1	2.05	0.40
1:B:87:LEU:HD21	1:B:156[B]:MSE:CE	2.51	0.40
1:A:151:PHE:CD2	1:A:151:PHE:C	3.00	0.40
1:A:30:ASP:HB3	1:A:40:LYS:CG	2.52	0.40
1:B:39:SER:C	1:B:41:ARG:N	2.78	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	45/203 (22%)	42 (93%)	3 (7%)	0	100	100
1	B	183/203 (90%)	174 (95%)	8 (4%)	1 (0%)	24	27
All	All	228/406 (56%)	216 (95%)	11 (5%)	1 (0%)	30	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	75	PRO

5.3.2 Protein sidechains [i](#)

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	159/166 (96%)	130 (82%)	29 (18%)	2	1
1	B	160/166 (96%)	128 (80%)	32 (20%)	1	1
All	All	319/332 (96%)	258 (81%)	61 (19%)	1	1

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	4	THR
1	A	11	ILE
1	A	29	MSE
1	A	45	ASN

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Mol	Chain	Res	Type
1	A	52	GLU
1	A	53	LEU
1	A	56	GLU
1	A	57	MSE
1	A	72	VAL
1	A	76	LEU
1	A	84	LEU
1	A	86	LEU
1	A	90	LYS
1	A	94	LEU
1	A	95	THR
1	A	102	LEU
1	A	105	VAL
1	A	111	ILE
1	A	119	VAL
1	A	121	LEU
1	A	129	GLU
1	A	140	LYS
1	A	172	ASN
1	A	175	LEU
1	A	179	GLN
1	A	182	SER
1	A	184	VAL
1	A	191	MSE
1	B	8	ARG
1	B	13	GLN
1	B	26	ILE
1	B	30	ASP
1	B	35	ARG
1	B	52	GLU
1	B	56	GLU
1	B	63	ASN
1	B	76	LEU
1	B	77	VAL
1	B	84	LEU
1	B	87	LEU
1	B	94	LEU
1	B	102	LEU
1	B	105	VAL
1	B	110	THR
1	B	116	ARG
1	B	120	TRP

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Mol	Chain	Res	Type
1	B	121	LEU
1	B	123	ARG
1	B	130	THR
1	B	145	LYS
1	B	156[A]	MSE
1	B	156[B]	MSE
1	B	159	LEU
1	B	161	LYS
1	B	172	ASN
1	B	175	LEU
1	B	179	GLN
1	B	183	ASN
1	B	184	VAL
1	B	189	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	155	GLN
1	A	179	GLN
1	A	183	ASN
1	B	45	ASN
1	B	118	GLN
1	B	139	GLN
1	B	157	HIS
1	B	172	ASN
1	B	179	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	182/203 (89%)	0.72	18 (9%) 13 14	40, 50, 63, 74	0
1	B	183/203 (90%)	0.49	5 (2%) 56 63	41, 49, 66, 83	0
All	All	365/406 (89%)	0.60	23 (6%) 26 29	40, 50, 64, 83	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3	LEU	5.9
1	A	3	LEU	5.0
1	A	120	TRP	3.5
1	A	26	ILE	3.5
1	A	15	ALA	2.7
1	A	23	GLY	2.6
1	A	17	ALA	2.6
1	A	16	ILE	2.5
1	A	24	PHE	2.5
1	B	76	LEU	2.4
1	B	64	CYS	2.4
1	A	43	VAL	2.4
1	A	121	LEU	2.3
1	A	36	ALA	2.3
1	A	22	ARG	2.3
1	A	131	PHE	2.2
1	A	20	GLY	2.2
1	A	61	LEU	2.1
1	A	128	GLU	2.1
1	B	126	GLU	2.1
1	B	123	ARG	2.1
1	A	112	HIS	2.0
1	A	47	PHE	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 oligosaccharides 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.