



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 08:29 AM EDT

PDB ID : 3C3W / pdb_00003c3w
Title : Crystal Structure of the Mycobacterium tuberculosis Hypoxic Response Regulator DosR
Authors : Wisedchaisri, G.; Wu, M.; Sherman, D.R.; Hol, W.G.J.
Deposited on : 2008-01-28
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

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

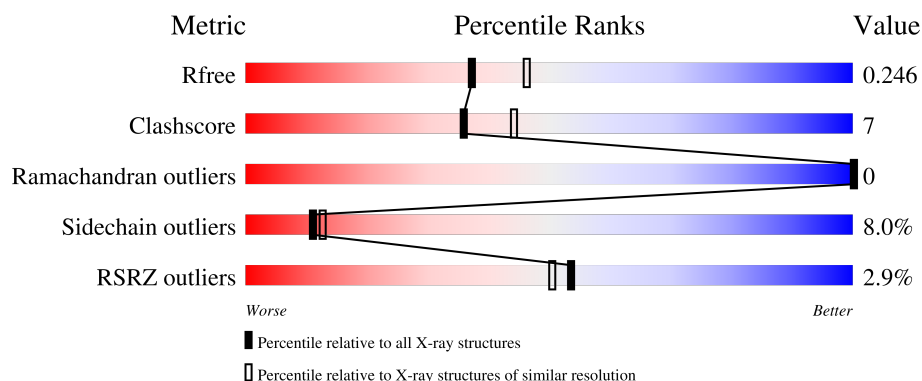
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 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	225	% 79% 12% • 6%
1	B	225	4% 75% 15% • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3243 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 TWO COMPONENT TRANSCRIPTIONAL REGULATORY PROTEIN DEVR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1573	979	283	301	10			
1	B	210	Total	C	N	O	S	0	1	0
			1574	980	284	299	11			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	218	LEU	-	expression tag	UNP P95193
A	219	GLU	-	expression tag	UNP P95193
A	220	HIS	-	expression tag	UNP P95193
A	221	HIS	-	expression tag	UNP P95193
A	222	HIS	-	expression tag	UNP P95193
A	223	HIS	-	expression tag	UNP P95193
A	224	HIS	-	expression tag	UNP P95193
A	225	HIS	-	expression tag	UNP P95193
B	218	LEU	-	expression tag	UNP P95193
B	219	GLU	-	expression tag	UNP P95193
B	220	HIS	-	expression tag	UNP P95193
B	221	HIS	-	expression tag	UNP P95193
B	222	HIS	-	expression tag	UNP P95193
B	223	HIS	-	expression tag	UNP P95193
B	224	HIS	-	expression tag	UNP P95193
B	225	HIS	-	expression tag	UNP P95193

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



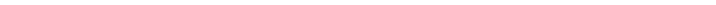
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

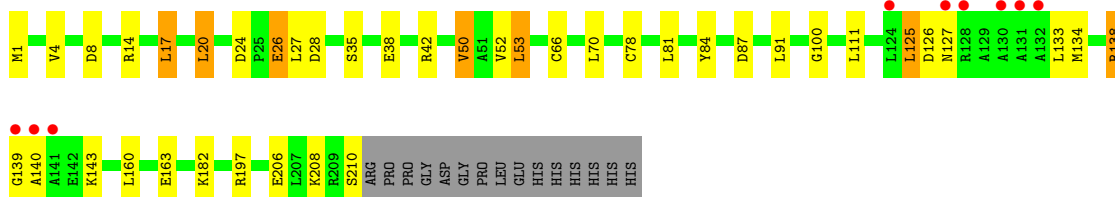
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	33	Total	O	0	0
			33	33		
3	B	58	Total	O	0	0
			58	58		

- Molecule 1: TWO COMPONENT TRANSCRIPTIONAL REGULATORY PROTEIN DEVR

E195	L207	K208	R211	PR0	GLY	GLY	PR0	LEU	GLU	HIS	HIS	HIS	HIS	HIS	HIS	HIS	W1	L20	D24	P26	E26	L27	D28	V43	V52	L53	D54	V55	M73	D87	L91	L96	Y101	K104	L111	V118	R122	S123	L124	L125	R128	A129	A130	M134	L137	A140	L150	E154	D172	R173	S186	R187	L188	L189
------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------

Chain B:  4% 75% 15% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.11Å 98.08Å 53.71Å 90.00° 93.81° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	90.1 (50.00-2.20) 90.3 (50.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.4.0066	Depositor
R, R_{free}	0.186 , 0.249 0.187 , 0.246	Depositor DCC
R_{free} test set	1172 reflections (4.62%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3243	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.18% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.67	1/1585 (0.1%)	0.89	3/2139 (0.1%)
1	B	0.65	0/1589	0.91	1/2143 (0.0%)
All	All	0.66	1/3174 (0.0%)	0.90	4/4282 (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
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	43	VAL	CA-CB	6.71	1.58	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	MET	CA-C-N	5.59	125.43	119.28
1	A	73	MET	C-N-CA	5.59	125.43	119.28
1	B	42	ARG	N-CA-C	5.06	117.92	111.69
1	A	173	ARG	NE-CZ-NH1	-5.05	116.45	121.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	52	VAL	Peptide
1	B	52	VAL	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	1573	0	1627	23	0
1	B	1574	0	1641	29	0
2	A	5	0	0	0	0
3	A	33	0	0	4	0
3	B	58	0	0	1	0
All	All	3243	0	3268	47	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 7.

All (47) 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:125:LEU:H	1:B:125:LEU:HD23	1.31	0.95
1:B:125:LEU:HD23	1:B:125:LEU:N	1.92	0.84
1:B:20:LEU:HD11	1:B:87:ASP:HB3	1.60	0.83
1:A:128:ARG:HH11	1:A:128:ARG:HB2	1.44	0.83
1:A:125:LEU:CD2	1:B:140:ALA:HB1	2.16	0.75
1:B:134:MET:O	1:B:138:ARG:HG3	1.88	0.74
1:A:1:MET:HE2	1:A:28:ASP:HB2	1.78	0.66
1:A:128:ARG:HB2	1:A:128:ARG:NH1	2.11	0.66
1:A:173:ARG:NH1	1:B:163:GLU:O	2.29	0.64
1:B:20:LEU:HD22	1:B:91:LEU:HD21	1.79	0.64
1:B:182:LYS:HE3	3:B:237:HOH:O	1.98	0.63
1:A:195:GLU:HG2	3:A:233:HOH:O	2.02	0.59
1:A:140:ALA:HB1	1:B:125:LEU:HD13	1.84	0.59
1:B:139:GLY:O	1:B:143:LYS:HG2	2.03	0.58
1:B:1:MET:HB2	1:B:26:GLU:O	2.04	0.58
1:B:53:LEU:HD11	1:B:66[A]:CYS:SG	2.44	0.57
1:A:54:ASP:OD2	3:A:229:HOH:O	2.17	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:TYR:O	1:B:197:ARG:HG2	2.05	0.56
1:A:26:GLU:HG3	1:A:26:GLU:O	2.06	0.55
1:A:172:ASP:HB3	3:A:251:HOH:O	2.08	0.54
1:A:134:MET:HE3	1:A:134:MET:HA	1.89	0.54
1:B:26:GLU:O	1:B:26:GLU:HG3	2.07	0.53
1:A:101:TYR:O	1:A:104:LYS:HD3	2.10	0.52
1:B:66[B]:CYS:SG	1:B:70:LEU:HD12	2.50	0.52
1:B:125:LEU:N	1:B:125:LEU:CD2	2.65	0.51
1:A:20:LEU:HD21	1:A:87:ASP:HB3	1.93	0.51
1:A:125:LEU:HD23	1:B:140:ALA:HB1	1.91	0.51
1:A:208:LYS:HD3	3:A:239:HOH:O	2.10	0.51
1:B:78:CYS:O	1:B:100:GLY:HA3	2.14	0.48
1:A:55:VAL:HG11	1:A:189:LEU:CD1	2.44	0.47
1:A:118:VAL:O	1:A:122:ARG:HB3	2.15	0.47
1:A:154:GLU:HB3	1:A:188:LEU:HD13	1.97	0.46
1:B:35:SER:OG	1:B:38:GLU:HG3	2.16	0.46
1:B:66[B]:CYS:SG	1:B:70:LEU:CD1	3.04	0.45
1:A:1:MET:CE	1:A:28:ASP:HB2	2.45	0.45
1:A:101:TYR:HD1	1:A:104:LYS:HE2	1.82	0.44
1:B:1:MET:SD	1:B:28:ASP:HB2	2.58	0.43
1:B:87:ASP:OD2	1:B:206:GLU:OE2	2.36	0.43
1:B:17:LEU:HD21	1:B:81:LEU:HD21	2.00	0.42
1:A:137:LEU:HG	1:B:133:LEU:HD11	2.01	0.42
1:B:4:VAL:HG22	1:B:50:VAL:HG13	2.02	0.42
1:B:24:ASP:OD1	1:B:26:GLU:HB3	2.19	0.41
1:A:24:ASP:OD1	1:A:26:GLU:HB3	2.19	0.41
1:B:208:LYS:C	1:B:210:SER:H	2.28	0.41
1:A:208:LYS:HA	1:A:208:LYS:HD2	1.92	0.40
1:B:8:ASP:O	1:B:14:ARG:NH1	2.50	0.40
1:B:53:LEU:CD1	1:B:66[A]:CYS:SG	3.09	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	209/225 (93%)	208 (100%)	1 (0%)	0	100	100
1	B	209/225 (93%)	204 (98%)	5 (2%)	0	100	100
All	All	418/450 (93%)	412 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	163/179 (91%)	149 (91%)	14 (9%)	10	10
1	B	165/179 (92%)	153 (93%)	12 (7%)	13	15
All	All	328/358 (92%)	302 (92%)	26 (8%)	11	13

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	GLU
1	A	53	LEU
1	A	54	ASP
1	A	55	VAL
1	A	91	LEU
1	A	95	LEU
1	A	104	LYS
1	A	111	LEU
1	A	125	LEU
1	A	128	ARG
1	A	150	LEU
1	A	186	SER
1	A	207	LEU
1	A	208	LYS
1	B	17	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	20	LEU
1	B	26	GLU
1	B	27	LEU
1	B	50	VAL
1	B	53	LEU
1	B	111	LEU
1	B	125	LEU
1	B	126	ASP
1	B	127	ASN
1	B	138	ARG
1	B	160	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	GLN
1	A	167	ASN
1	A	199	GLN
1	B	144	GLN
1	B	167	ASN
1	B	183	ASN

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](#)

1 ligand is modelled in this entry.

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	$\# Z > 2$
2	SO4	A	226	-	4,4,4	0.24	0	6,6,6	0.23	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.

No monomer is involved in short contacts.

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, 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	211/225 (93%)	-0.98	3 (1%) 73 71	27, 37, 48, 59	0
1	B	210/225 (93%)	-0.80	9 (4%) 40 36	24, 37, 48, 54	1 (0%)
All	All	421/450 (93%)	-0.89	12 (2%) 53 50	24, 37, 48, 59	1 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	128	ARG	3.7
1	B	132	ALA	3.0
1	B	140	ALA	2.9
1	A	130	ALA	2.8
1	B	130	ALA	2.8
1	A	123	SER	2.5
1	B	124	LEU	2.4
1	B	131	ALA	2.3
1	A	124	LEU	2.3
1	B	139	GLY	2.3
1	B	141	ALA	2.3
1	B	127	ASN	2.2

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

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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å ²)	Q<0.9
2	SO4	A	226	5/5	0.99	0.03	45,46,49,49	0

6.5 Other polymers

There are no such residues in this entry.