



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:17 PM UTC

PDB ID : 3QDN / pdb_00003qdn
Title : Putative thioredoxin protein from Salmonella typhimurium
Authors : Osipiuk, J.; Zhou, M.; Kwon, K.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2011-01-18
Resolution : 2.09 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

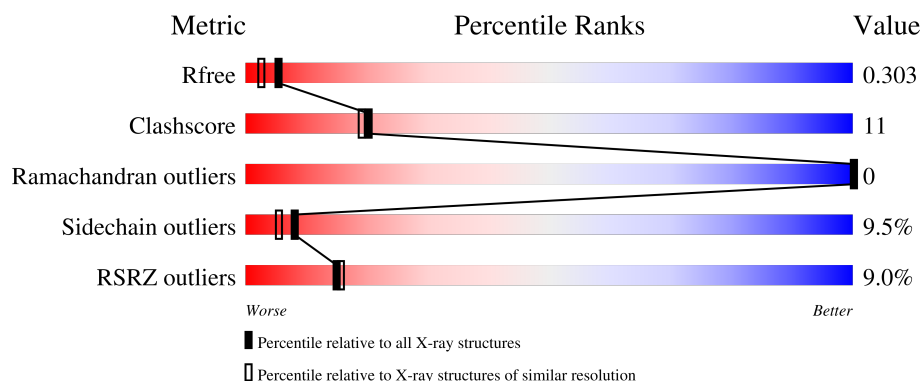
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (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 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	287	
1	B	287	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4710 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 Putative thioredoxin protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	Se	0	17	0
			2338	1479	397	453	2	7			
1	B	280	Total	C	N	O	S	Se	0	12	0
			2276	1437	391	442	2	4			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q8ZR92
A	-1	ASN	-	expression tag	UNP Q8ZR92
A	0	ALA	-	expression tag	UNP Q8ZR92
B	-2	SER	-	expression tag	UNP Q8ZR92
B	-1	ASN	-	expression tag	UNP Q8ZR92
B	0	ALA	-	expression tag	UNP Q8ZR92

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		

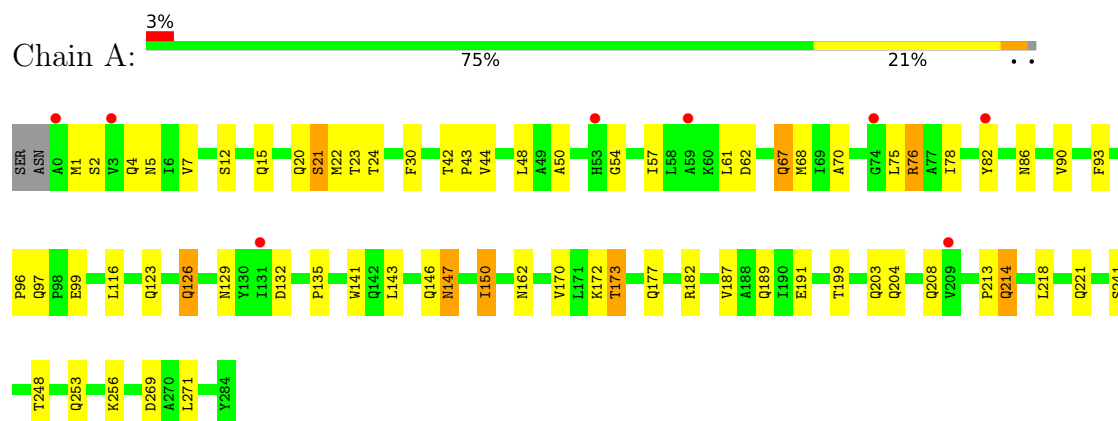
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total	O	0	0
			54	54		
3	B	40	Total	O	0	0
			40	40		

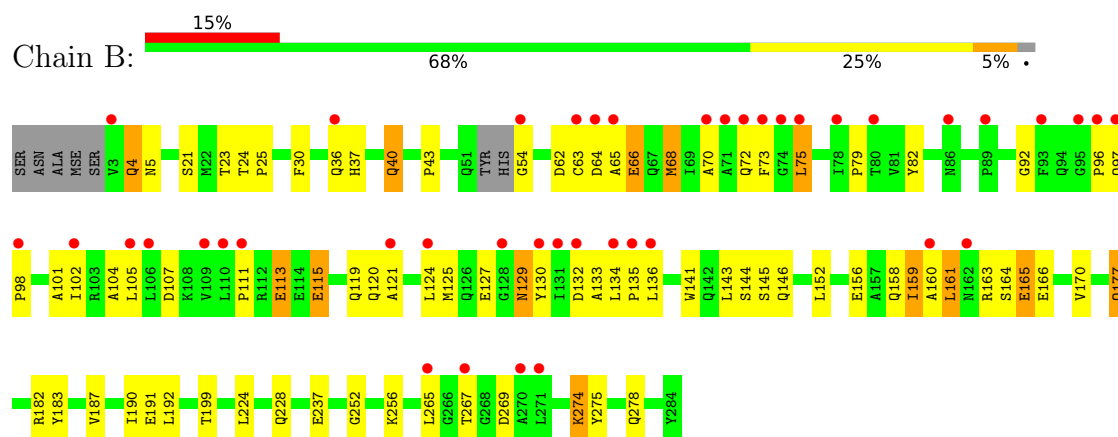
3 Residue-property plots [i](#)

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: Putative thioredoxin protein



- Molecule 1: Putative thioredoxin protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.23Å 62.55Å 79.52Å 90.00° 98.66° 90.00°	Depositor
Resolution (Å)	35.70 – 2.09 35.70 – 2.09	Depositor EDS
% Data completeness (in resolution range)	96.3 (35.70-2.09) 96.3 (35.70-2.09)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.242 , 0.302 0.246 , 0.303	Depositor DCC
R_{free} test set	1629 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	39.8	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4710	wwPDB-VP
Average B, all atoms (Å ²)	45.0	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 56.96 % 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 2.5359e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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](#)

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

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	0.91	0/2418	1.03	1/3267 (0.0%)
1	B	0.86	0/2338	1.03	3/3156 (0.1%)
All	All	0.89	0/4756	1.03	4/6423 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	66	GLU	N-CA-C	-7.69	99.00	109.54
1	B	104	ALA	N-CA-C	-6.37	104.45	112.93
1	B	129	ASN	N-CA-C	5.55	116.36	108.54
1	A	90	VAL	N-CA-C	-5.14	107.14	111.56

There are no chirality outliers.

There are no planarity outliers.

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	2338	0	2373	46	1
1	B	2276	0	2309	64	1
2	A	2	0	0	0	0
3	A	54	0	0	3	0
3	B	40	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4710	0	4682	105	1

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 11.

All (105) 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:158:GLN:HE21	1:B:166:GLU:HB3	1.29	0.97
1:B:158:GLN:NE2	1:B:166:GLU:HB3	1.83	0.93
1:B:130:TYR:HB2	1:B:161:LEU:HD11	1.49	0.92
1:B:252:GLY:O	1:B:256[B]:LYS:HG3	1.76	0.85
1:A:170:VAL:O	1:A:173:THR:HB	1.79	0.83
1:A:126:GLN:HE21	1:A:126:GLN:HA	1.43	0.82
1:A:20:GLN:HG2	3:A:304:HOH:O	1.80	0.81
1:B:113:GLU:HB2	3:B:309:HOH:O	1.80	0.80
1:B:274[A]:LYS:HE2	1:B:278:GLN:OE1	1.83	0.77
1:A:126:GLN:HA	1:A:126:GLN:NE2	1.99	0.77
1:B:265:LEU:O	1:B:269:ASP:HB2	1.85	0.76
1:B:63:CYS:O	1:B:65:ALA:HA	1.87	0.74
1:A:67[A]:GLN:HE21	1:A:67[A]:GLN:H	1.33	0.74
1:B:119:GLN:HA	1:B:119:GLN:OE1	1.85	0.74
1:B:97:GLN:HG2	1:B:98:PRO:HD2	1.71	0.72
1:B:165:GLU:HB2	3:B:311:HOH:O	1.90	0.71
1:A:187:VAL:O	1:A:191[B]:GLU:HG2	1.90	0.71
1:A:141:TRP:HZ2	1:A:173:THR:HG21	1.58	0.68
1:B:120:GLN:HB3	1:B:136:LEU:HD21	1.74	0.67
1:B:5:ASN:OD1	1:B:54:GLY:HA2	1.94	0.66
1:B:130:TYR:CB	1:B:161:LEU:HD11	2.27	0.63
1:A:1:MSE:SE	1:B:73:PHE:HD1	2.32	0.62
1:A:218:LEU:HA	1:A:221:GLN:HE21	1.64	0.62
1:B:4:GLN:HE21	1:B:4:GLN:HA	1.64	0.62
1:B:265:LEU:O	1:B:269:ASP:CB	2.49	0.61
1:A:93:PHE:CD1	1:A:97:GLN:HG3	2.36	0.60
1:A:147:ASN:OD1	1:A:150:ILE:HG12	2.00	0.60
1:B:121:ALA:O	1:B:125:MSE:HG3	2.00	0.60
1:B:21:SER:HA	1:B:24:THR:O	2.02	0.60
1:A:269:ASP:OD1	1:A:271:LEU:HB3	2.01	0.59
1:B:97:GLN:CD	1:B:101:ALA:CB	2.75	0.59
1:A:132:ASP:O	1:A:135:PRO:HD2	2.04	0.58
1:B:63:CYS:C	1:B:65:ALA:HA	2.28	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:LEU:HD13	1:B:182[A]:ARG:HD3	1.85	0.57
1:A:44:VAL:O	1:A:48:LEU:HG	2.05	0.56
3:A:311:HOH:O	1:B:72:GLN:HA	2.06	0.56
1:B:115:GLU:O	1:B:119:GLN:HG2	2.06	0.56
1:B:161:LEU:N	1:B:161:LEU:CD1	2.68	0.56
1:A:22[B]:MSE:SE	1:A:86:ASN:HA	2.56	0.56
1:A:93:PHE:CE1	1:A:97:GLN:HG3	2.41	0.56
1:B:134:LEU:HB3	1:B:135:PRO:HD3	1.88	0.55
1:A:15[B]:GLN:H	1:A:15[B]:GLN:CD	2.15	0.54
1:B:111:PRO:CB	1:B:115:GLU:HG2	2.37	0.54
1:A:67[A]:GLN:H	1:A:67[A]:GLN:NE2	2.04	0.54
1:B:111:PRO:HB3	1:B:115:GLU:HG2	1.89	0.53
1:B:37:HIS:CE1	1:B:96:PRO:HG3	2.44	0.53
1:A:30:PHE:HD1	1:A:61:LEU:HD23	1.75	0.52
1:B:166:GLU:O	1:B:170:VAL:HG23	2.10	0.52
1:B:92:GLY:HA2	1:B:105:LEU:HD21	1.92	0.51
1:B:224:LEU:O	1:B:228:GLN:HG2	2.11	0.51
1:B:164:SER:HB2	1:B:190:ILE:HG23	1.93	0.50
1:A:123:GLN:O	1:A:126:GLN:HB2	2.12	0.50
1:A:76:ARG:HG2	3:A:308:HOH:O	2.12	0.49
1:A:4:GLN:HG3	1:A:50:ALA:HA	1.94	0.48
1:B:183:TYR:C	1:B:183:TYR:CD1	2.92	0.48
1:B:62:ASP:OD1	1:B:64:ASP:C	2.57	0.48
1:A:213:PRO:O	1:A:214:GLN:HG2	2.13	0.47
1:B:129:ASN:ND2	1:B:132:ASP:OD2	2.43	0.47
1:B:30:PHE:O	1:B:79:PRO:HA	2.14	0.47
1:B:37:HIS:ND1	1:B:96:PRO:HG3	2.29	0.47
1:A:62:ASP:OD1	1:A:62:ASP:C	2.57	0.47
1:A:1:MSE:HB2	1:B:73:PHE:O	2.15	0.46
1:A:70:ALA:HB1	1:A:75:LEU:HD22	1.98	0.46
1:B:187:VAL:O	1:B:191[A]:GLU:HG2	2.15	0.46
1:B:70:ALA:O	1:B:75:LEU:HD11	2.15	0.46
1:A:147:ASN:OD1	1:A:150:ILE:CG1	2.64	0.46
1:A:208:GLN:O	1:A:218:LEU:HD23	2.15	0.46
1:B:119:GLN:OE1	1:B:119:GLN:CA	2.56	0.46
1:B:68:MSE:HE3	1:B:68:MSE:HA	1.97	0.45
1:B:4:GLN:HE21	1:B:4:GLN:CA	2.30	0.45
1:B:75:LEU:HD12	1:B:75:LEU:H	1.80	0.45
1:B:237:GLU:OE2	1:B:275:TYR:OH	2.31	0.45
1:A:141:TRP:CD1	1:A:146[A]:GLN:HA	2.52	0.45
1:A:253:GLN:HE22	1:A:256:LYS:HD3	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:GLN:H	1:B:177:GLN:HG3	1.19	0.44
1:A:1:MSE:SE	1:B:73:PHE:HB3	2.68	0.44
1:B:121:ALA:HB1	1:B:133:ALA:HB1	2.00	0.44
1:A:182:ARG:HA	1:A:182:ARG:HD2	1.65	0.44
1:B:159:ILE:HG22	1:B:160:ALA:N	2.32	0.43
1:B:121:ALA:HB2	1:B:136:LEU:HD23	2.00	0.43
1:A:2:SER:HB2	1:A:5:ASN:HA	2.01	0.43
1:A:30:PHE:CD1	1:A:61:LEU:HD23	2.52	0.43
1:A:129:ASN:ND2	1:A:132:ASP:OD2	2.46	0.43
1:A:21:SER:HA	1:A:24:THR:O	2.19	0.43
1:A:253:GLN:NE2	1:A:256:LYS:HD3	2.34	0.42
1:A:42:THR:HB	1:A:43:PRO:HD3	2.02	0.42
1:A:218:LEU:HA	1:A:221:GLN:NE2	2.34	0.42
1:A:1:MSE:SE	1:B:73:PHE:O	2.87	0.42
1:B:125:MSE:HE1	1:B:156:GLU:HB3	2.00	0.42
1:B:163:ARG:HB3	1:B:166:GLU:HB2	2.00	0.42
1:A:5:ASN:ND2	1:A:54:GLY:N	2.67	0.42
1:A:48:LEU:HD21	1:A:99:GLU:HG3	2.01	0.42
1:B:121:ALA:HA	1:B:124:LEU:HD12	2.01	0.42
1:A:1:MSE:SE	1:B:73:PHE:CD1	3.19	0.42
1:B:40:GLN:O	1:B:43:PRO:HD2	2.20	0.42
1:B:62:ASP:O	1:B:66:GLU:N	2.52	0.42
1:A:96:PRO:C	1:A:97:GLN:HG2	2.45	0.42
1:A:75:LEU:HD23	1:A:78:ILE:HD11	2.02	0.41
1:B:144:SER:O	1:B:145[B]:SER:HB3	2.20	0.41
1:B:141:TRP:CD1	1:B:146:GLN:HA	2.55	0.41
1:B:182[A]:ARG:NH1	3:B:286:HOH:O	2.54	0.41
1:B:161:LEU:N	1:B:161:LEU:HD13	2.36	0.41
1:A:141:TRP:CD1	1:A:146[B]:GLN:HA	2.56	0.41
1:B:97:GLN:CD	1:B:101:ALA:HB1	2.45	0.41

All (1) 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)
1:A:177:GLN:NE2	1:B:160:ALA:O[1_455]	2.08	0.12

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	300/287 (104%)	289 (96%)	11 (4%)	0	100	100
1	B	288/287 (100%)	271 (94%)	17 (6%)	0	100	100
All	All	588/574 (102%)	560 (95%)	28 (5%)	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	255/235 (108%)	227 (89%)	28 (11%)	6	3
1	B	246/235 (105%)	224 (91%)	22 (9%)	9	6
All	All	501/470 (107%)	451 (90%)	50 (10%)	8	5

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	12	SER
1	A	21	SER
1	A	23	THR
1	A	57	ILE
1	A	67[A]	GLN
1	A	67[B]	GLN
1	A	68[A]	MSE

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Mol	Chain	Res	Type
1	A	68[B]	MSE
1	A	76	ARG
1	A	82	TYR
1	A	116	LEU
1	A	126	GLN
1	A	143	LEU
1	A	147	ASN
1	A	150	ILE
1	A	162	ASN
1	A	172	LYS
1	A	173	THR
1	A	189	GLN
1	A	199[A]	THR
1	A	199[B]	THR
1	A	203	GLN
1	A	204[A]	GLN
1	A	204[B]	GLN
1	A	214	GLN
1	A	241	SER
1	A	248	THR
1	B	4	GLN
1	B	23	THR
1	B	36	GLN
1	B	40	GLN
1	B	68	MSE
1	B	75	LEU
1	B	82	TYR
1	B	102	ILE
1	B	107	ASP
1	B	113	GLU
1	B	115	GLU
1	B	127	GLU
1	B	143	LEU
1	B	159	ILE
1	B	161	LEU
1	B	165	GLU
1	B	177	GLN
1	B	192	LEU
1	B	199	THR
1	B	267	THR
1	B	274[A]	LYS
1	B	274[B]	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	126	GLN
1	A	142	GLN
1	A	184	GLN
1	A	189	GLN
1	A	214	GLN
1	A	221	GLN
1	A	253	GLN
1	B	4	GLN
1	B	40	GLN
1	B	85	GLN
1	B	97	GLN
1	B	146	GLN
1	B	189	GLN
1	B	228	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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

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	280/287 (97%)	0.35	8 (2%) 53 56	19, 40, 56, 65	15 (5%)
1	B	276/287 (96%)	0.88	42 (15%) 5 5	17, 47, 85, 98	12 (4%)
All	All	556/574 (96%)	0.61	50 (8%) 15 16	17, 42, 78, 98	27 (4%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	75	LEU	4.6
1	B	93	PHE	4.5
1	B	71	ALA	4.1
1	B	130	TYR	3.9
1	A	0	ALA	3.7
1	B	267	THR	3.6
1	B	65	ALA	3.3
1	B	162	ASN	3.3
1	B	124	LEU	3.2
1	B	74	GLY	3.2
1	B	73	PHE	3.0
1	B	105	LEU	3.0
1	B	131	ILE	2.9
1	A	3	VAL	2.8
1	B	86	ASN	2.8
1	B	54	GLY	2.8
1	B	3	VAL	2.8
1	B	132	ASP	2.7
1	B	36	GLN	2.7
1	B	102	ILE	2.7
1	B	95	GLY	2.7
1	B	160	ALA	2.6
1	A	53	HIS	2.6
1	B	80	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	96	PRO	2.6
1	B	109	VAL	2.6
1	B	63	CYS	2.6
1	B	70	ALA	2.6
1	B	98	PRO	2.6
1	B	128	GLY	2.6
1	A	131[A]	ILE	2.5
1	B	106	LEU	2.5
1	A	82	TYR	2.4
1	B	78	ILE	2.4
1	B	136	LEU	2.4
1	B	89	PRO	2.4
1	B	121	ALA	2.3
1	B	111	PRO	2.3
1	B	110	LEU	2.3
1	B	134	LEU	2.3
1	B	265	LEU	2.2
1	B	270	ALA	2.1
1	B	64	ASP	2.1
1	B	72	GLN	2.1
1	A	209	VAL	2.1
1	B	135	PRO	2.1
1	A	74	GLY	2.1
1	A	59	ALA	2.1
1	B	271	LEU	2.1
1	B	97	GLN	2.1

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](#)

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($\text{\AA}^2$)	Q<0.9
2	CA	A	501	1/1	0.97	0.10	41,41,41,41	0
2	CA	A	502	1/1	0.98	0.10	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.