



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 11:17 AM UTC

PDB ID : 1OVB / pdb_00001ovb
Title : THE MECHANISM OF IRON UPTAKE BY TRANSFERRINS: THE
STRUCTURE OF AN 18KD NII-DOMAIN FRAGMENT AT 2.3
ANGSTROMS RESOLUTION
Authors : Kuser, P.; Lindley, P.; Sarra, R.
Deposited on : 1992-10-05
Resolution : 2.30 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

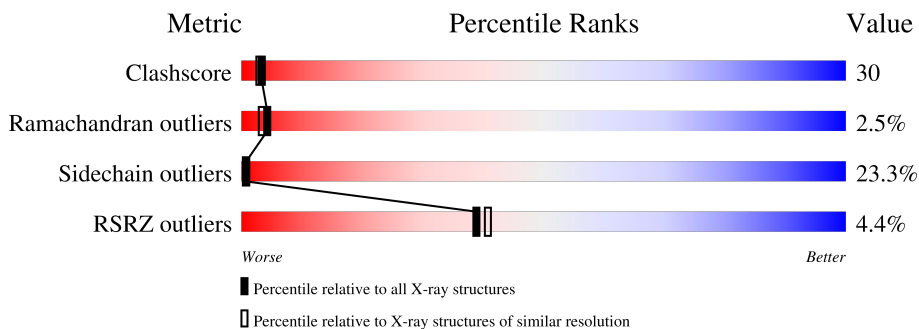
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	159	4% 44% 35% 18% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 1331 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 OVOTRANSFERRIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	0	0
			1220	762	213	236	9			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	145	SER	ILE	conflict	UNP P56410
A	146	VAL	SER	conflict	UNP P56410
A	183	ALA	GLY	conflict	UNP P56410

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		

- Molecule 3 is CARBONATE ION (CCD ID: CO3) (formula: CO₃).



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

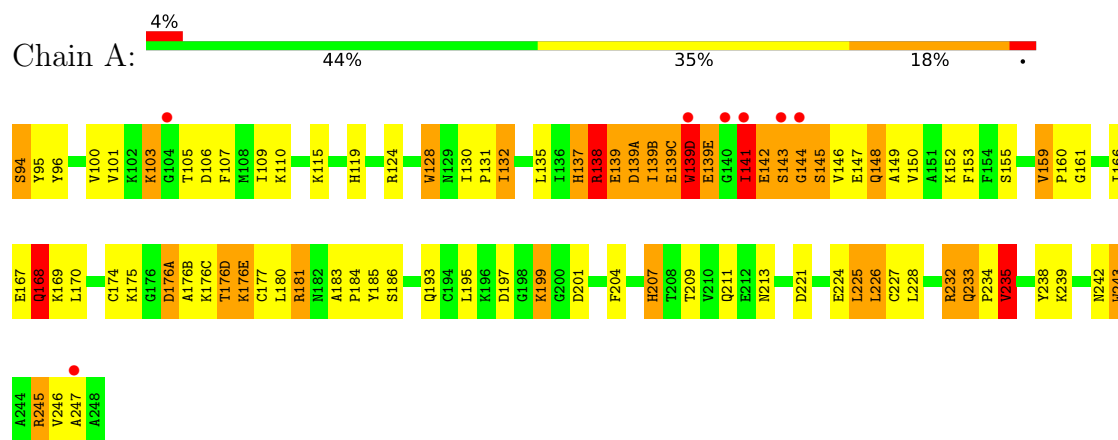
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	106	Total	O	0	0
			106	106		

3 Residue-property plots

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: OVOTRANSFERRIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	41.50Å 41.50Å 81.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	7.00 – 2.30 7.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (7.00-2.30) 91.8 (7.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.70 (at 2.30Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.195 , (Not available) 0.257 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.489	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 78.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.008 for -h,-k,l 0.059 for h,-h-k,-l 0.036 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	1331	wwPDB-VP
Average B, all atoms (Å ²)	0.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.53% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FE, CO3

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	1.20	15/1244 (1.2%)	1.42	11/1676 (0.7%)

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	6

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	137	HIS	CE1-NE2	7.47	1.40	1.32
1	A	137	HIS	ND1-CE1	7.18	1.39	1.32
1	A	207	HIS	CE1-NE2	6.98	1.39	1.32
1	A	119	HIS	ND1-CE1	6.96	1.39	1.32
1	A	207	HIS	ND1-CE1	6.80	1.39	1.32
1	A	119	HIS	CE1-NE2	6.78	1.39	1.32
1	A	128	TRP	NE1-CE2	-6.01	1.30	1.37
1	A	139(D)	TRP	NE1-CE2	-5.58	1.31	1.37
1	A	243	TRP	NE1-CE2	-5.57	1.31	1.37
1	A	168	GLN	CD-OE1	5.51	1.34	1.23
1	A	233	GLN	CD-OE1	5.34	1.33	1.23
1	A	193	GLN	CD-OE1	5.31	1.33	1.23
1	A	211	GLN	CD-OE1	5.19	1.33	1.23
1	A	148	GLN	CD-OE1	5.13	1.33	1.23
1	A	242	ASN	CG-OD1	5.01	1.33	1.23

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	105	THR	N-CA-C	-7.99	98.03	110.42
1	A	159	VAL	CA-CB-CG2	5.82	120.30	110.40
1	A	207	HIS	CA-CB-CG	-5.71	108.09	113.80
1	A	243	TRP	N-CA-CB	5.65	119.03	110.28
1	A	153	PHE	CA-CB-CG	-5.62	108.18	113.80
1	A	247	ALA	CA-C-N	5.58	131.75	121.70
1	A	247	ALA	C-N-CA	5.58	131.75	121.70
1	A	100	VAL	CA-CB-CG2	5.48	119.72	110.40
1	A	235	VAL	CA-CB-CG2	5.14	119.15	110.40
1	A	101	VAL	CA-CB-CG2	5.02	118.93	110.40
1	A	181	ARG	CD-NE-CZ	-5.00	117.40	124.40

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	124	ARG	Sidechain
1	A	138	ARG	Sidechain
1	A	141	ILE	Peptide
1	A	142	GLU	Peptide
1	A	232	ARG	Sidechain
1	A	245	ARG	Sidechain

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	1220	0	1183	73	4
2	A	1	0	0	0	0
3	A	4	0	0	0	0
4	A	106	0	0	6	4
All	All	1331	0	1183	73	4

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 30.

All (73) 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:139:GLU:O	1:A:139:GLU:HG3	1.58	1.01
1:A:131:PRO:HA	1:A:243:TRP:CE3	2.10	0.87
1:A:147:GLU:HG2	1:A:170:LEU:HD12	1.58	0.84
1:A:147:GLU:HG2	1:A:170:LEU:CD1	2.08	0.83
1:A:141:ILE:CG2	1:A:141:ILE:O	2.30	0.79
1:A:131:PRO:HA	1:A:243:TRP:CZ3	2.22	0.74
1:A:197:ASP:HB3	1:A:199:LYS:HE2	1.68	0.74
1:A:166:ILE:HG13	1:A:167:GLU:HG2	1.68	0.73
1:A:174:CYS:HB2	1:A:176(E):LYS:HD3	1.72	0.71
1:A:166:ILE:CG1	1:A:167:GLU:HG2	2.22	0.70
1:A:130:ILE:HD11	1:A:246:VAL:HG22	1.74	0.69
1:A:128:TRP:O	1:A:132:ILE:HB	1.93	0.68
1:A:197:ASP:HB3	1:A:199:LYS:CE	2.22	0.68
1:A:139(C):GLU:O	1:A:139(E):GLU:HG2	1.96	0.66
1:A:139:GLU:O	1:A:139:GLU:CG	2.40	0.64
1:A:139(E):GLU:HG3	1:A:149:ALA:HA	1.79	0.63
1:A:174:CYS:HB2	1:A:176(E):LYS:CD	2.27	0.63
1:A:225:LEU:HD22	1:A:235:VAL:HA	1.81	0.63
1:A:175:LYS:HB2	1:A:185:TYR:CZ	2.35	0.62
1:A:141:ILE:O	1:A:141:ILE:HG22	1.99	0.62
1:A:146:VAL:HG23	4:A:463:HOH:O	2.02	0.59
1:A:137:HIS:O	1:A:137:HIS:CG	2.54	0.59
1:A:146:VAL:O	1:A:150:VAL:HG23	2.03	0.58
1:A:131:PRO:HB3	1:A:204:PHE:CD1	2.38	0.58
1:A:130:ILE:CD1	1:A:246:VAL:HG22	2.34	0.58
1:A:176(A):ASP:HB2	1:A:176(D):THR:OG1	2.04	0.57
1:A:131:PRO:HA	1:A:243:TRP:HE3	1.67	0.57
1:A:168:GLN:HE21	1:A:168:GLN:HA	1.70	0.57
1:A:175:LYS:O	1:A:184:PRO:HG2	2.05	0.57
1:A:130:ILE:N	1:A:131:PRO:HD2	2.19	0.56
1:A:185:TYR:HB2	4:A:444:HOH:O	2.04	0.56
1:A:144:GLY:HA3	4:A:499:HOH:O	2.05	0.56
1:A:95:TYR:CB	1:A:207:HIS:HE1	2.19	0.55
1:A:141:ILE:O	1:A:141:ILE:HG23	2.05	0.55
1:A:238:TYR:CD1	1:A:238:TYR:C	2.85	0.54
1:A:138:ARG:HG3	1:A:139(A):ASP:OD1	2.09	0.53
1:A:160:PRO:HG2	1:A:186:SER:HA	1.90	0.52
1:A:176(D):THR:O	1:A:180:LEU:HG	2.09	0.52
1:A:183:ALA:HB1	1:A:184:PRO:HD2	1.92	0.52
1:A:139(E):GLU:H	1:A:149:ALA:HB2	1.76	0.51
1:A:233:GLN:HB3	1:A:234:PRO:HD2	1.91	0.51
1:A:161:GLY:HA3	1:A:181:ARG:HD3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:TYR:HA	1:A:207:HIS:CE1	2.47	0.50
1:A:147:GLU:HG2	1:A:170:LEU:HD11	1.92	0.49
1:A:174:CYS:O	1:A:176(E):LYS:HD3	2.13	0.48
1:A:96:TYR:H	1:A:207:HIS:CE1	2.30	0.48
1:A:110:LYS:HG2	4:A:477:HOH:O	2.14	0.48
1:A:107:PHE:CZ	1:A:201:ASP:HB3	2.48	0.48
1:A:135:LEU:HB3	1:A:139(B):ILE:HG13	1.96	0.47
1:A:175:LYS:HB2	1:A:185:TYR:CE2	2.50	0.47
1:A:227:CYS:SG	1:A:233:GLN:HG3	2.56	0.46
1:A:94:SER:HA	1:A:246:VAL:O	2.15	0.46
1:A:131:PRO:CB	1:A:204:PHE:CD1	2.99	0.46
1:A:141:ILE:HD13	1:A:141:ILE:HA	1.83	0.45
1:A:95:TYR:CB	1:A:207:HIS:CE1	2.99	0.45
1:A:174:CYS:CB	1:A:176(E):LYS:HD3	2.42	0.45
1:A:238:TYR:CD1	1:A:239:LYS:N	2.84	0.45
1:A:137:HIS:O	1:A:137:HIS:ND1	2.50	0.45
1:A:145:SER:HB2	1:A:148:GLN:HB3	2.00	0.44
1:A:115:LYS:HD3	4:A:466:HOH:O	2.17	0.44
1:A:209:THR:O	1:A:213:ASN:HB2	2.18	0.44
1:A:185:TYR:OH	1:A:199:LYS:HE3	2.18	0.43
1:A:174:CYS:HA	1:A:185:TYR:HD2	1.83	0.43
1:A:139(D):TRP:HB3	4:A:478:HOH:O	2.18	0.43
1:A:197:ASP:HB3	1:A:199:LYS:HE3	1.98	0.43
1:A:207:HIS:ND1	1:A:207:HIS:N	2.64	0.43
1:A:137:HIS:ND1	1:A:137:HIS:C	2.75	0.42
1:A:148:GLN:HG2	1:A:152:LYS:HE3	2.00	0.42
1:A:226:LEU:HD12	1:A:226:LEU:HA	1.81	0.42
1:A:224:GLU:CD	1:A:232:ARG:HD3	2.46	0.41
1:A:128:TRP:CZ3	1:A:159:VAL:HG21	2.55	0.41
1:A:103:LYS:HE2	1:A:221:ASP:O	2.21	0.41
1:A:168:GLN:HE21	1:A:168:GLN:CA	2.31	0.40

All (4) 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:107:PHE:CA	4:A:447:HOH:O[3_464]	1.62	0.58
1:A:106:ASP:O	4:A:447:HOH:O[3_464]	1.91	0.29
1:A:107:PHE:C	4:A:447:HOH:O[3_464]	2.07	0.13
1:A:107:PHE:N	4:A:447:HOH:O[3_464]	2.11	0.09

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	157/159 (99%)	141 (90%)	12 (8%)	4 (2%)	4 3

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	SER
1	A	141	ILE
1	A	144	GLY
1	A	176(B)	ALA

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	129/129 (100%)	99 (77%)	30 (23%)	1 1

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	SER
1	A	103	LYS
1	A	109	ILE
1	A	132	ILE
1	A	138	ARG
1	A	139	GLU
1	A	139(A)	ASP

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Mol	Chain	Res	Type
1	A	139(B)	ILE
1	A	139(C)	GLU
1	A	139(D)	TRP
1	A	139(E)	GLU
1	A	141	ILE
1	A	142	GLU
1	A	143	SER
1	A	145	SER
1	A	155	SER
1	A	168	GLN
1	A	169	LYS
1	A	176(A)	ASP
1	A	176(C)	LYS
1	A	176(D)	THR
1	A	176(E)	LYS
1	A	177	CYS
1	A	195	LEU
1	A	199	LYS
1	A	225	LEU
1	A	226	LEU
1	A	228	LEU
1	A	235	VAL
1	A	245	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	A	168	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is 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	$\# Z > 2$
3	CO3	A	400	2	3,3,3	0.85	0	2,3,3	0.07	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	159/159 (100%)	0.31	7 (4%)	39 41	0, 0, 1, 1	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	143	SER	5.2
1	A	139(D)	TRP	4.2
1	A	144	GLY	4.0
1	A	140	GLY	3.9
1	A	141	ILE	3.9
1	A	104	GLY	2.4
1	A	247	ALA	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(Å ²)	Q<0.9
2	FE	A	300	1/1	0.97	0.03	0,0,0,0	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CO3	A	400	4/4	0.97	0.07	0,0,0,0	0

6.5 Other polymers [i](#)

There are no such residues in this entry.