



Preliminary Full wwPDB X-ray Structure Validation Report ⓘ

Aug 15, 2021 – 12:30 PM JST

Deposition ID : D_1300024008

This is a Preliminary Full wwPDB X-ray Structure Validation Report.

This report is produced by the wwPDB Deposition System during initial deposition but before annotation of the structure.

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

MolProbity	:	4.02b-467
Xtrriage (Phenix)	:	1.13
EDS	:	2.23.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.1

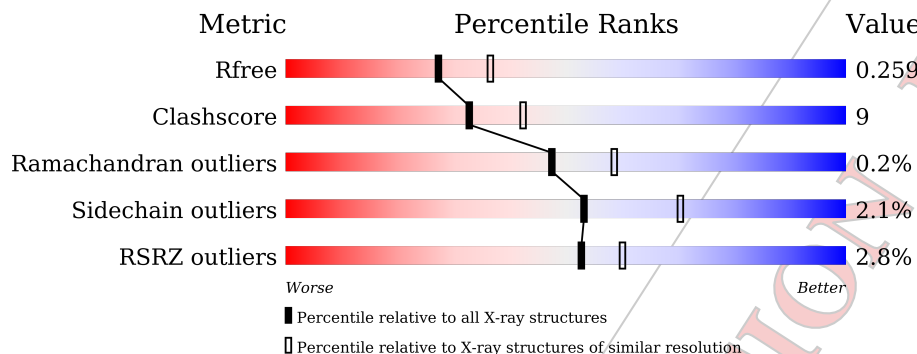
1 Overall quality at a glance i

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))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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	164	 2% 81% 16% ...
1	B	164	 4% 84% 14% ..
1	C	164	 4% 81% 17% ..
1	H	164	 % 85% 15% .

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 10779 atoms, of which 5222 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 *Thermotoga maritima* ferritin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	H	164	2686	874	1319	223	264	6	0	0	0
1	A	163	2666	869	1307	222	263	5	0	0	0
1	B	162	2649	864	1298	221	262	4	0	0	0
1	C	162	2649	864	1298	221	262	4	0	0	0

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	S	121	Total 121	O 121	0	0

PRELIMINARY VALIDATION REPORT

4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.81Å 94.50Å 108.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.69 – 2.30 38.69 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.7 (38.69-2.30) 98.7 (38.69-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.14 3260	Depositor
R, R_{free}	0.201 , 0.259 0.201 , 0.259	Depositor DCC
R_{free} test set	1998 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	43.6	Xtrriage
Anisotropy	0.521	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10779	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

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	0.78	4/1388 (0.3%)	0.73	1/1866 (0.1%)
1	B	0.65	0/1380	0.74	2/1856 (0.1%)
1	C	0.62	0/1380	0.68	0/1856
1	H	0.67	0/1396	0.75	1/1876 (0.1%)
All	All	0.68	4/5544 (0.1%)	0.73	4/7454 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	114	LYS	CE-NZ	9.15	1.72	1.49
1	A	94	LYS	CD-CE	8.93	1.73	1.51
1	A	94	LYS	CE-NZ	-7.90	1.29	1.49
1	A	94	LYS	CG-CD	5.13	1.70	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	MET	CG-SD-CE	7.39	112.02	100.20
1	H	1	MET	CG-SD-CE	-7.15	88.76	100.20
1	A	94	LYS	CD-CE-NZ	6.73	127.18	111.70
1	B	65	GLU	CA-CB-CG	-5.96	100.29	113.40

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	1359	1307	1307	31	0
1	B	1351	1298	1298	19	0
1	C	1351	1298	1298	30	0
1	H	1367	1319	1319	19	0
2	D	8	0	0	0	0
3	S	121	0	0	7	0
All	All	5557	5222	5222	93	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 9.

All (93) 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:114:LYS:CE	1:A:114:LYS:NZ	1.71	1.48
1:C:3:VAL:CG1	1:C:66:ARG:HH21	1.63	1.12
1:A:98:PHE:HB3	3:S:75:HOH:O	1.53	1.07
1:B:61:GLU:O	1:B:65:GLU:OE1	1.73	1.06
1:H:47:LYS:NZ	1:H:164:GLU:OXT	1.88	1.06
1:C:3:VAL:HG13	1:C:66:ARG:NH2	1.70	1.06
1:C:3:VAL:HG13	1:C:66:ARG:HH21	0.87	1.01
1:A:94:LYS:O	3:S:75:HOH:O	1.94	0.84
1:A:149:GLN:O	1:A:153:ILE:HD12	1.80	0.81
1:H:47:LYS:HZ1	1:H:164:GLU:C	1.89	0.74
1:A:26:TYR:CD1	1:A:45:MET:HE1	2.22	0.73
1:A:98:PHE:CB	3:S:75:HOH:O	2.20	0.73
1:B:61:GLU:C	1:B:65:GLU:OE1	2.27	0.73
1:A:4:ILE:HD12	1:A:9:ARG:HB2	1.72	0.72
1:C:3:VAL:CG1	1:C:66:ARG:NH2	2.42	0.71
1:A:29:MET:HB2	1:A:45:MET:HE2	1.72	0.70
1:C:4:ILE:HG13	1:C:66:ARG:HB3	1.76	0.68
1:A:10:LYS:NZ	3:S:107:HOH:O	2.26	0.68
1:B:57:MET:HE1	1:B:60:TYR:CD2	2.31	0.66
1:A:114:LYS:NZ	1:A:114:LYS:CD	2.57	0.66
1:A:29:MET:CB	1:A:45:MET:HE2	2.26	0.65
1:C:3:VAL:CG1	1:C:66:ARG:HE	2.09	0.65
1:B:142:LEU:O	3:S:51:HOH:O	2.14	0.65
1:H:7:LYS:O	3:S:83:HOH:O	2.15	0.64
1:H:117:ALA:O	1:A:155:GLN:HG2	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:TYR:HA	1:A:45:MET:HE3	1.81	0.62
1:C:4:ILE:HG13	1:C:66:ARG:CB	2.29	0.62
1:A:127:ASP:OD2	3:S:60:HOH:O	2.16	0.62
1:A:38:PHE:HB3	1:A:41:PHE:HD2	1.69	0.57
1:B:41:PHE:CD1	1:B:160:LEU:HD12	2.39	0.57
1:B:4:ILE:HG23	1:B:66:ARG:HB3	1.86	0.57
1:B:102:SER:O	1:B:106:ILE:HG13	2.05	0.56
1:A:26:TYR:CA	1:A:45:MET:HE3	2.36	0.56
1:H:41:PHE:CE1	1:H:160:LEU:HD12	2.40	0.56
1:B:41:PHE:CE1	1:B:160:LEU:HD12	2.41	0.55
1:A:26:TYR:HA	1:A:45:MET:CE	2.36	0.55
1:C:100:THR:HG21	1:C:133:GLU:HG3	1.88	0.55
1:B:57:MET:HA	1:B:57:MET:HE2	1.89	0.54
1:A:31:THR:HG22	1:B:60:TYR:HE2	1.71	0.54
1:A:31:THR:HG22	1:B:60:TYR:CE2	2.43	0.54
1:C:40:GLY:HA3	1:C:157:ASP:O	2.08	0.53
1:A:29:MET:HB2	1:A:45:MET:CE	2.37	0.52
1:C:90:GLU:HA	1:C:140:LEU:HD11	1.92	0.51
1:H:41:PHE:CD1	1:H:160:LEU:HD12	2.46	0.51
1:H:4:ILE:HG23	1:H:66:ARG:HB3	1.91	0.50
1:B:109:LEU:O	1:B:113:GLU:HG3	2.11	0.50
1:H:3:VAL:HG13	1:H:66:ARG:HE	1.75	0.50
1:A:4:ILE:HG23	1:A:66:ARG:CZ	2.42	0.50
1:C:7:LYS:HE2	1:C:113:GLU:OE1	2.11	0.50
1:B:3:VAL:HG23	1:B:66:ARG:HE	1.76	0.50
1:A:34:ASP:OD1	1:A:39:LYS:HD2	2.10	0.50
1:H:38:PHE:CE2	1:H:85:ILE:HD11	2.48	0.49
1:C:41:PHE:CE2	1:C:157:ASP:HB2	2.48	0.49
1:C:26:TYR:HD1	1:C:45:MET:SD	2.36	0.49
1:C:103:ILE:HG22	1:C:126:VAL:HG12	1.94	0.49
1:H:79:PRO:HB2	1:H:82:TRP:CZ2	2.48	0.48
1:H:158:ARG:HD3	1:C:39:LYS:HG2	1.95	0.48
1:B:57:MET:HE1	1:B:60:TYR:CG	2.49	0.48
1:C:3:VAL:CG1	1:C:66:ARG:NE	2.77	0.48
1:A:77:LYS:NZ	1:B:71:GLU:OE2	2.42	0.48
1:C:3:VAL:HG11	1:C:66:ARG:HE	1.79	0.47
1:H:38:PHE:CZ	1:H:85:ILE:HD11	2.50	0.47
1:A:107:LEU:O	1:A:107:LEU:HD23	2.15	0.46
1:H:15:GLN:NE2	1:H:19:GLU:OE2	2.47	0.45
1:C:128:GLU:O	1:C:131:GLU:HG2	2.16	0.45
1:A:46:LYS:O	1:A:50:GLN:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ILE:HD13	1:A:160:LEU:HD11	1.99	0.45
1:A:4:ILE:CD1	1:A:9:ARG:HB2	2.45	0.45
1:C:103:ILE:CG2	1:C:126:VAL:HG12	2.47	0.45
1:C:122:LEU:O	1:C:126:VAL:HG13	2.17	0.45
1:H:15:GLN:HB2	1:H:106:ILE:HD11	1.99	0.44
1:A:26:TYR:HD1	1:A:45:MET:HE1	1.75	0.44
1:H:44:TRP:HZ2	1:H:135:GLN:HG2	1.83	0.44
1:B:158:ARG:NH2	1:C:101:GLN:HG2	2.33	0.44
1:H:3:VAL:CG1	1:H:66:ARG:HE	2.32	0.43
1:H:29:MET:HG2	1:H:82:TRP:CH2	2.53	0.43
1:A:2:MET:O	1:A:2:MET:HG3	2.18	0.43
1:C:16:LEU:C	1:C:16:LEU:HD23	2.38	0.43
1:B:103:ILE:HD13	1:B:125:PHE:HB3	2.01	0.43
1:C:3:VAL:HG12	1:C:66:ARG:HE	1.83	0.42
1:C:124:TRP:HZ3	1:C:128:GLU:OE1	2.03	0.42
1:C:61:GLU:OE2	1:C:61:GLU:HA	2.19	0.42
1:C:125:PHE:O	1:C:128:GLU:N	2.53	0.41
1:C:33:PHE:CD2	1:C:41:PHE:HB3	2.55	0.41
1:B:3:VAL:CG2	1:B:66:ARG:HE	2.33	0.41
1:C:3:VAL:CG1	1:C:66:ARG:CZ	2.98	0.41
1:H:131:GLU:O	1:H:135:GLN:HB2	2.21	0.41
1:B:59:PHE:O	1:B:63:ILE:HG13	2.21	0.41
1:H:97:GLU:OE2	1:H:137:ARG:NH2	2.54	0.40
1:A:38:PHE:CD2	1:A:85:ILE:HD11	2.56	0.40
1:C:90:GLU:HG3	1:C:140:LEU:HD21	2.04	0.40
1:A:86:LYS:O	1:A:90:GLU:HG3	2.21	0.40
1:C:149:GLN:O	1:C:153:ILE:HG12	2.21	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	161/164 (98%)	158 (98%)	3 (2%)	0	100	100
1	B	160/164 (98%)	154 (96%)	6 (4%)	0	100	100
1	C	160/164 (98%)	154 (96%)	5 (3%)	1 (1%)	25	31
1	H	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
All	All	643/656 (98%)	625 (97%)	17 (3%)	1 (0%)	47	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	4	ILE

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	143/144 (99%)	139 (97%)	4 (3%)	43	60
1	B	142/144 (99%)	137 (96%)	5 (4%)	36	50
1	C	142/144 (99%)	141 (99%)	1 (1%)	84	92
1	H	144/144 (100%)	142 (99%)	2 (1%)	67	81
All	All	571/576 (99%)	559 (98%)	12 (2%)	53	70

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	MET
1	H	111	SER
1	A	9	ARG
1	A	94	LYS
1	A	98	PHE
1	A	137	ARG
1	B	9	ARG
1	B	24	TYR
1	B	94	LYS
1	B	138	GLU

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Mol	Chain	Res	Type
1	B	147	ASN
1	C	108	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	162	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 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.

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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	163/164 (99%)	0.41	4 (2%) 57 64	36, 49, 67, 85	0
1	B	162/164 (98%)	0.47	6 (3%) 41 48	35, 51, 68, 82	0
1	C	162/164 (98%)	0.29	6 (3%) 41 48	35, 56, 76, 93	0
1	H	164/164 (100%)	0.09	2 (1%) 79 83	37, 46, 62, 95	0
All	All	651/656 (99%)	0.31	18 (2%) 53 60	35, 50, 71, 95	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	MET	8.1
1	H	1	MET	7.7
1	H	2	MET	5.2
1	C	124	TRP	4.1
1	A	130	VAL	3.9
1	A	126	VAL	3.5
1	B	126	VAL	3.2
1	B	104	TYR	3.2
1	C	107	LEU	2.7
1	B	106	ILE	2.6
1	C	3	VAL	2.5
1	B	107	LEU	2.5
1	C	125	PHE	2.4
1	B	111	SER	2.4
1	B	130	VAL	2.4
1	C	117	ALA	2.3
1	C	156	LEU	2.3
1	A	56	ALA	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 monosaccharides 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	D	8	1/?	0.89	0.14	96,96,96,96	0
2	FE	D	7	1/?	0.98	0.15	66,66,66,66	0
2	FE	D	6	1/?	0.98	0.15	64,64,64,64	0
2	FE	D	4	1/?	0.99	0.18	47,47,47,47	0
2	FE	D	5	1/?	0.99	0.16	50,50,50,50	0
2	FE	D	1	1/?	0.99	0.20	50,50,50,50	0
2	FE	D	2	1/?	0.99	0.18	50,50,50,50	0
2	FE	D	3	1/?	0.99	0.13	64,64,64,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.