



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 08:44 AM UTC

PDB ID : 5U1A / pdb_00005u1a
Title : Ferritin with Gc MtrE loop 1 inserted at His34
Authors : Wang, S.
Deposited on : 2016-11-28
Resolution : 2.00 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

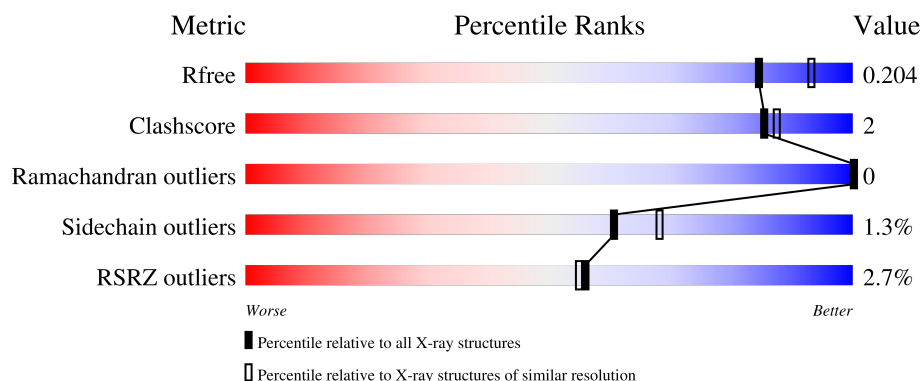
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	182	
1	B	182	
1	C	182	
1	D	182	
1	E	182	

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Mol	Chain	Length	Quality of chain
1	F	182	3% 84% 9% 8%
1	G	182	3% 80% 14% 7%
1	H	182	3% 77% 13% 8%
1	I	182	2% 84% 8% 7%
1	J	182	2% 83% 9% 8%
1	K	182	2% 87% 5% 8%
1	L	182	2% 77% 14% 9%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17961 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 Ferritin,MtrE protein chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	S	0	3	0
			1392	888	227	271	6			
1	B	167	Total	C	N	O	S	0	0	0
			1360	866	224	265	5			
1	C	168	Total	C	N	O	S	0	0	0
			1369	871	225	267	6			
1	D	169	Total	C	N	O	S	0	0	0
			1378	875	229	269	5			
1	E	169	Total	C	N	O	S	0	1	0
			1379	877	228	269	5			
1	F	168	Total	C	N	O	S	0	0	0
			1368	871	225	266	6			
1	G	170	Total	C	N	O	S	0	0	0
			1380	877	227	270	6			
1	H	167	Total	C	N	O	S	0	0	0
			1360	866	224	265	5			
1	I	169	Total	C	N	O	S	0	0	0
			1375	874	226	269	6			
1	J	168	Total	C	N	O	S	0	0	0
			1366	869	225	267	5			
1	K	168	Total	C	N	O	S	0	0	0
			1366	869	225	267	5			
1	L	166	Total	C	N	O	S	0	0	0
			1354	863	223	263	5			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	GLY	HIS	conflict	UNP A0A0B2EHC8
A	54	CYS	SER	conflict	UNP A0A0B2EHC8
A	56	THR	LEU	conflict	UNP A0A0B2EHC8
A	140	SER	ALA	conflict	UNP A0A0B2EHC8
B	34	GLY	HIS	conflict	UNP A0A0B2EHC8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	54	CYS	SER	conflict	UNP A0A0B2EHC8
B	56	THR	LEU	conflict	UNP A0A0B2EHC8
B	140	SER	ALA	conflict	UNP A0A0B2EHC8
C	34	GLY	HIS	conflict	UNP A0A0B2EHC8
C	54	CYS	SER	conflict	UNP A0A0B2EHC8
C	56	THR	LEU	conflict	UNP A0A0B2EHC8
C	140	SER	ALA	conflict	UNP A0A0B2EHC8
D	34	GLY	HIS	conflict	UNP A0A0B2EHC8
D	54	CYS	SER	conflict	UNP A0A0B2EHC8
D	56	THR	LEU	conflict	UNP A0A0B2EHC8
D	140	SER	ALA	conflict	UNP A0A0B2EHC8
E	34	GLY	HIS	conflict	UNP A0A0B2EHC8
E	54	CYS	SER	conflict	UNP A0A0B2EHC8
E	56	THR	LEU	conflict	UNP A0A0B2EHC8
E	140	SER	ALA	conflict	UNP A0A0B2EHC8
F	34	GLY	HIS	conflict	UNP A0A0B2EHC8
F	54	CYS	SER	conflict	UNP A0A0B2EHC8
F	56	THR	LEU	conflict	UNP A0A0B2EHC8
F	140	SER	ALA	conflict	UNP A0A0B2EHC8
G	34	GLY	HIS	conflict	UNP A0A0B2EHC8
G	54	CYS	SER	conflict	UNP A0A0B2EHC8
G	56	THR	LEU	conflict	UNP A0A0B2EHC8
G	140	SER	ALA	conflict	UNP A0A0B2EHC8
H	34	GLY	HIS	conflict	UNP A0A0B2EHC8
H	54	CYS	SER	conflict	UNP A0A0B2EHC8
H	56	THR	LEU	conflict	UNP A0A0B2EHC8
H	140	SER	ALA	conflict	UNP A0A0B2EHC8
I	34	GLY	HIS	conflict	UNP A0A0B2EHC8
I	54	CYS	SER	conflict	UNP A0A0B2EHC8
I	56	THR	LEU	conflict	UNP A0A0B2EHC8
I	140	SER	ALA	conflict	UNP A0A0B2EHC8
J	34	GLY	HIS	conflict	UNP A0A0B2EHC8
J	54	CYS	SER	conflict	UNP A0A0B2EHC8
J	56	THR	LEU	conflict	UNP A0A0B2EHC8
J	140	SER	ALA	conflict	UNP A0A0B2EHC8
K	34	GLY	HIS	conflict	UNP A0A0B2EHC8
K	54	CYS	SER	conflict	UNP A0A0B2EHC8
K	56	THR	LEU	conflict	UNP A0A0B2EHC8
K	140	SER	ALA	conflict	UNP A0A0B2EHC8
L	34	GLY	HIS	conflict	UNP A0A0B2EHC8
L	54	CYS	SER	conflict	UNP A0A0B2EHC8
L	56	THR	LEU	conflict	UNP A0A0B2EHC8

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Chain	Residue	Modelled	Actual	Comment	Reference
L	140	SER	ALA	conflict	UNP A0A0B2EHC8

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Na 2 2	0	0
2	B	2	Total Na 2 2	0	0
2	C	1	Total Na 1 1	0	0
2	D	1	Total Na 1 1	0	0
2	E	3	Total Na 3 3	0	0
2	F	4	Total Na 4 4	0	0
2	G	3	Total Na 3 3	0	0
2	H	4	Total Na 4 4	0	0
2	I	2	Total Na 2 2	0	0
2	J	1	Total Na 1 1	0	0
2	K	2	Total Na 2 2	0	0
2	L	5	Total Na 5 5	0	0

- Molecule 3 is FE (III) ION (CCD ID: FE) (formula: Fe).

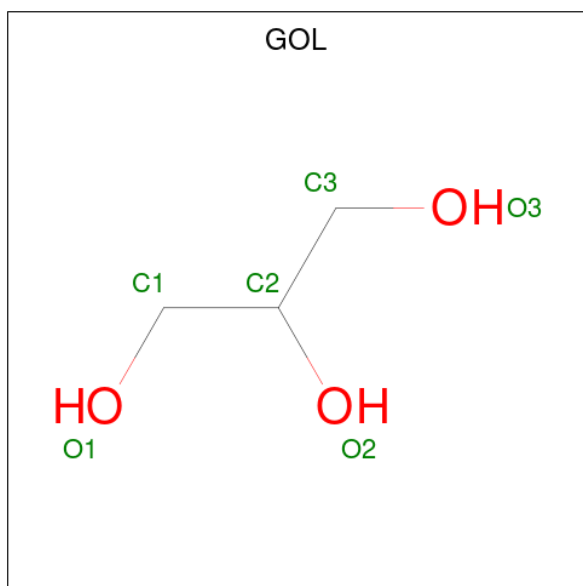
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Fe 1 1	0	0
3	E	1	Total Fe 1 1	0	0
3	F	1	Total Fe 1 1	0	0
3	G	1	Total Fe 1 1	0	0
3	K	1	Total Fe 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	1	Total	Fe	0	0
			1	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		
4	I	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	I	1	Total	C	O	0	0
			6	3	3		
4	L	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Cl	0	0
			1	1		
5	D	1	Total	Cl	0	0
			1	1		
5	F	1	Total	Cl	0	0
			1	1		
5	G	1	Total	Cl	0	0
			1	1		
5	K	1	Total	Cl	0	0
			1	1		
5	L	3	Total	Cl	0	0
			3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	96	Total	O	0	0
			96	96		
6	B	106	Total	O	0	0
			106	106		
6	C	106	Total	O	0	0
			106	106		
6	D	95	Total	O	0	0
			95	95		
6	E	110	Total	O	0	0
			110	110		
6	F	146	Total	O	0	0
			146	146		
6	G	133	Total	O	0	0
			133	133		
6	H	102	Total	O	0	0
			102	102		
6	I	119	Total	O	0	0
			119	119		

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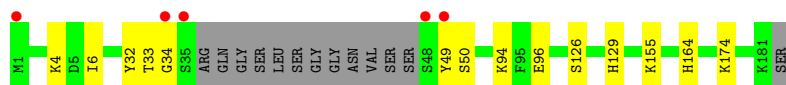
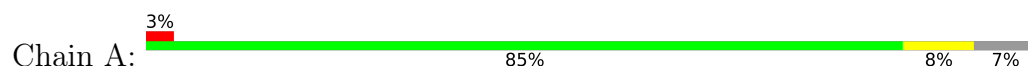
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	J	102	Total 102	O 102	0	0
6	K	137	Total 137	O 137	0	0
6	L	146	Total 146	O 146	0	0

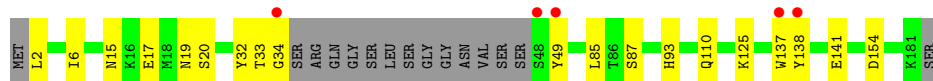
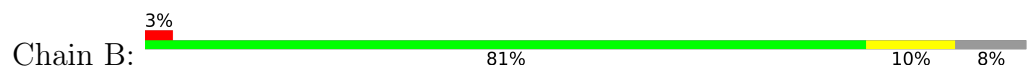
3 Residue-property plots [i](#)

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.

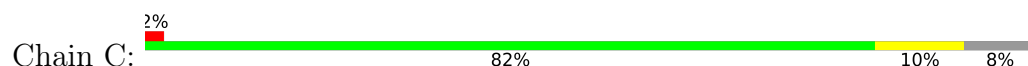
- Molecule 1: Ferritin,MtrE protein chimera



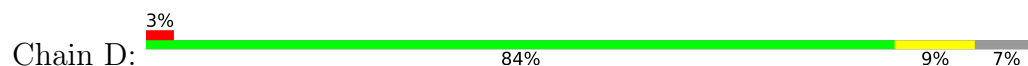
- Molecule 1: Ferritin,MtrE protein chimera



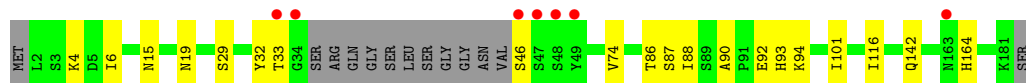
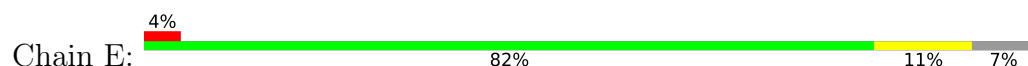
- Molecule 1: Ferritin,MtrE protein chimera



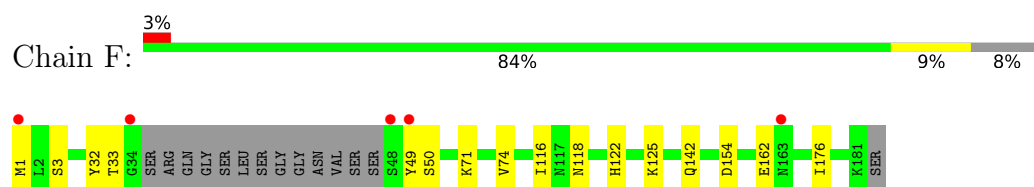
- Molecule 1: Ferritin,MtrE protein chimera



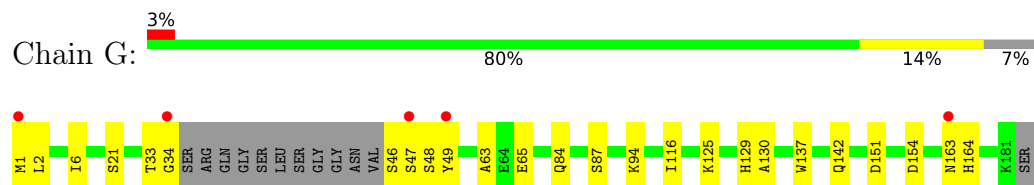
- Molecule 1: Ferritin,MtrE protein chimera



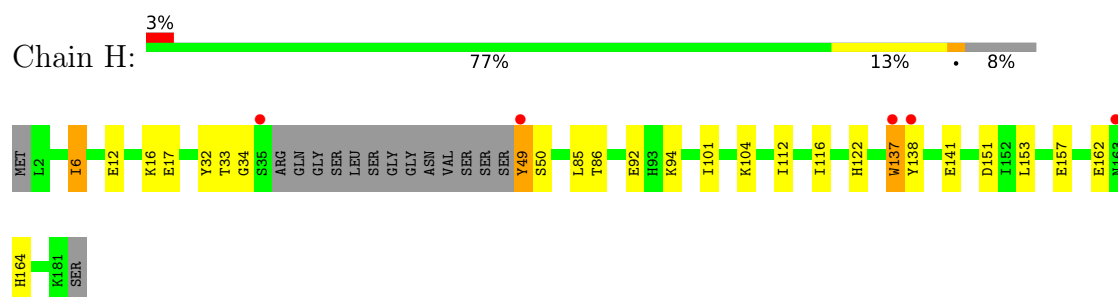
- Molecule 1: Ferritin,MtrE protein chimera



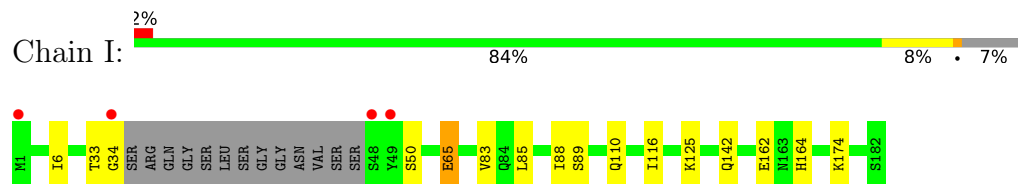
- Molecule 1: Ferritin,MtrE protein chimera



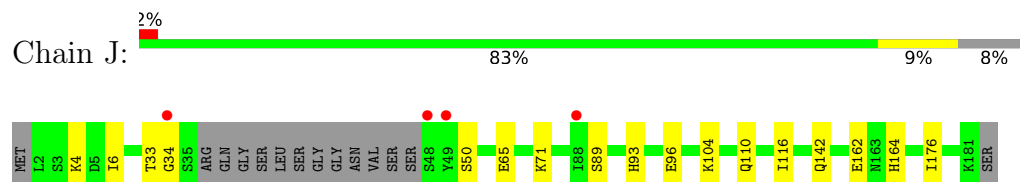
- Molecule 1: Ferritin,MtrE protein chimera



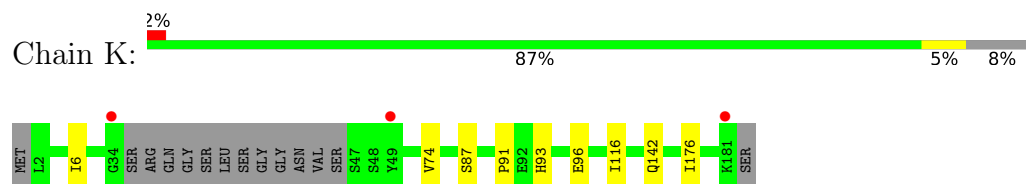
- Molecule 1: Ferritin,MtrE protein chimera



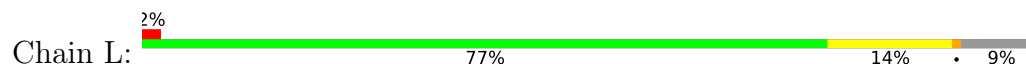
- Molecule 1: Ferritin,MtrE protein chimera

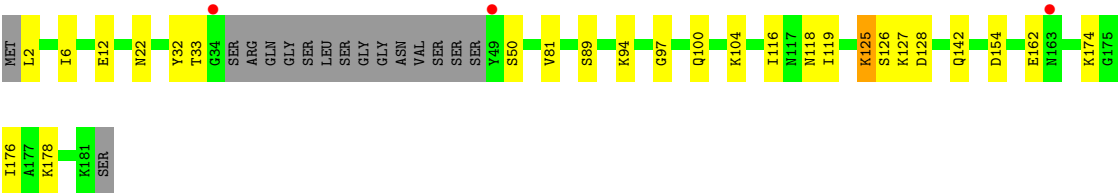


- Molecule 1: Ferritin,MtrE protein chimera



- Molecule 1: Ferritin,MtrE protein chimera





4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	128.41Å 128.41Å 165.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-2.00) 99.7 (20.00-2.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.168 , 0.196 0.178 , 0.204	Depositor DCC
R_{free} test set	9002 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 51.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17961	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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 28.59 % 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 1.7938e-03. 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 ⓘ

5.1 Standard geometry ⓘ

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

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	1.47	6/1434 (0.4%)	1.21	6/1933 (0.3%)
1	B	1.49	4/1389 (0.3%)	1.27	6/1872 (0.3%)
1	C	1.43	6/1398 (0.4%)	1.21	6/1882 (0.3%)
1	D	1.44	8/1407 (0.6%)	1.20	4/1894 (0.2%)
1	E	1.44	5/1412 (0.4%)	1.15	2/1903 (0.1%)
1	F	1.48	7/1397 (0.5%)	1.20	2/1882 (0.1%)
1	G	1.49	11/1409 (0.8%)	1.22	3/1898 (0.2%)
1	H	1.55	9/1389 (0.6%)	1.26	6/1872 (0.3%)
1	I	1.45	9/1404 (0.6%)	1.19	4/1890 (0.2%)
1	J	1.44	7/1395 (0.5%)	1.18	3/1880 (0.2%)
1	K	1.46	2/1395 (0.1%)	1.17	2/1880 (0.1%)
1	L	1.54	12/1383 (0.9%)	1.17	5/1864 (0.3%)
All	All	1.47	86/16812 (0.5%)	1.20	49/22650 (0.2%)

All (86) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	88	ILE	C-O	-7.82	1.16	1.24
1	H	34	GLY	C-O	7.31	1.33	1.23
1	D	87	SER	C-O	-7.29	1.14	1.23
1	I	89	SER	C-O	-6.97	1.15	1.23
1	G	2	LEU	C-O	-6.84	1.15	1.23
1	K	91	PRO	C-O	-6.59	1.15	1.23
1	L	94	LYS	C-O	-6.49	1.16	1.24
1	L	33	THR	CA-C	6.48	1.60	1.52
1	G	63	ALA	C-O	-6.48	1.16	1.24
1	G	137	TRP	C-O	-6.35	1.16	1.24
1	I	33	THR	CA-C	6.29	1.60	1.52
1	E	90	ALA	C-O	-6.12	1.16	1.24
1	E	32	TYR	C-O	-6.09	1.16	1.24
1	H	12	GLU	C-O	-6.03	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	88	ILE	N-CA	6.00	1.53	1.46
1	J	89	SER	C-O	-5.96	1.16	1.23
1	A	34	GLY	C-O	5.90	1.31	1.23
1	A	49	TYR	C-O	5.89	1.31	1.23
1	K	87	SER	N-CA	-5.85	1.38	1.45
1	H	137	TRP	C-O	-5.83	1.17	1.24
1	B	19	ASN	CA-C	5.79	1.60	1.52
1	I	125	LYS	C-O	-5.70	1.17	1.24
1	L	127	LYS	CA-CB	5.68	1.62	1.53
1	C	174	LYS	C-O	-5.68	1.17	1.24
1	A	94	LYS	C-O	-5.63	1.17	1.24
1	J	65	GLU	CD-OE1	-5.55	1.14	1.25
1	F	50	SER	N-CA	-5.54	1.39	1.46
1	F	33	THR	CA-C	5.53	1.59	1.52
1	L	97	GLY	N-CA	5.51	1.51	1.45
1	D	33	THR	CA-C	5.50	1.59	1.52
1	H	32	TYR	C-O	-5.47	1.17	1.24
1	L	118	ASN	CG-OD1	5.45	1.33	1.23
1	C	65	GLU	CD-OE1	-5.45	1.15	1.25
1	I	50	SER	C-O	-5.42	1.17	1.24
1	F	118	ASN	CG-ND2	-5.39	1.22	1.33
1	J	33	THR	CA-C	5.39	1.59	1.52
1	G	151	ASP	CB-CG	5.38	1.65	1.52
1	F	71	LYS	C-O	-5.36	1.17	1.24
1	L	178	LYS	C-O	-5.35	1.17	1.24
1	B	32	TYR	C-O	-5.34	1.17	1.24
1	E	164	HIS	C-O	-5.34	1.16	1.24
1	D	34	GLY	C-O	5.33	1.31	1.23
1	H	151	ASP	CB-CG	5.33	1.65	1.52
1	F	3	SER	CB-OG	5.32	1.52	1.42
1	A	126	SER	C-O	-5.31	1.17	1.24
1	F	32	TYR	C-O	-5.29	1.17	1.24
1	G	130	ALA	C-O	-5.27	1.18	1.24
1	L	22	ASN	C-O	-5.25	1.18	1.24
1	E	88	ILE	C-O	-5.25	1.18	1.24
1	E	87	SER	CB-OG	-5.25	1.31	1.42
1	B	110	GLN	CD-NE2	-5.24	1.22	1.33
1	H	92	GLU	C-O	-5.23	1.17	1.23
1	D	32	TYR	C-O	-5.21	1.17	1.24
1	A	164	HIS	C-O	-5.20	1.17	1.24
1	H	50	SER	N-CA	-5.20	1.39	1.46
1	C	157	GLU	C-O	-5.20	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	87	SER	C-O	-5.20	1.17	1.23
1	H	33	THR	CA-C	5.19	1.59	1.52
1	J	93	HIS	C-O	-5.18	1.17	1.23
1	L	6	ILE	C-O	-5.18	1.18	1.24
1	B	87	SER	C-O	-5.18	1.17	1.23
1	C	32	TYR	C-O	-5.16	1.17	1.24
1	J	71	LYS	C-O	-5.15	1.18	1.24
1	I	110	GLN	CD-NE2	-5.15	1.22	1.33
1	I	83	VAL	C-O	-5.14	1.18	1.24
1	C	50	SER	C-O	-5.13	1.17	1.24
1	D	126	SER	C-O	-5.13	1.17	1.24
1	F	122	HIS	C-O	-5.12	1.18	1.24
1	D	65	GLU	CD-OE1	5.12	1.35	1.25
1	G	65	GLU	C-O	-5.11	1.18	1.24
1	J	50	SER	C-O	-5.11	1.17	1.24
1	J	110	GLN	CD-NE2	-5.11	1.22	1.33
1	H	164	HIS	C-O	-5.11	1.17	1.24
1	G	125	LYS	C-O	-5.09	1.18	1.24
1	G	164	HIS	C-O	-5.09	1.17	1.24
1	D	89	SER	N-CA	5.08	1.52	1.46
1	G	94	LYS	CA-C	-5.08	1.46	1.52
1	L	32	TYR	C-O	-5.07	1.17	1.24
1	D	125	LYS	C-O	-5.06	1.18	1.24
1	L	100	GLN	C-O	-5.06	1.18	1.24
1	A	32	TYR	C-O	-5.05	1.17	1.24
1	I	164	HIS	C-O	-5.04	1.17	1.24
1	I	65	GLU	C-O	-5.04	1.18	1.24
1	L	12	GLU	C-O	-5.04	1.18	1.24
1	G	49	TYR	N-CA	5.03	1.52	1.45
1	L	119	ILE	C-O	-5.03	1.18	1.24

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	49	TYR	CA-C-O	-7.29	113.12	121.72
1	D	33	THR	CB-CA-C	-6.89	98.09	110.35
1	I	33	THR	CB-CA-C	-6.80	98.24	110.35
1	J	33	THR	CB-CA-C	-6.70	98.43	110.35
1	C	33	THR	CB-CA-C	-6.69	98.45	110.35
1	G	129	HIS	N-CA-C	6.66	119.55	111.82
1	F	33	THR	CB-CA-C	-6.60	98.60	110.35
1	G	33	THR	CB-CA-C	-6.59	98.62	110.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	33	THR	CB-CA-C	-6.54	98.70	110.35
1	B	33	THR	CB-CA-C	-6.50	98.77	110.35
1	L	33	THR	CB-CA-C	-6.43	98.90	110.35
1	L	125	LYS	CB-CG-CD	6.25	125.67	111.30
1	J	96	GLU	N-CA-C	6.20	119.02	111.82
1	H	33	THR	CB-CA-C	-6.16	99.38	110.35
1	A	50	SER	N-CA-C	6.13	118.18	108.67
1	I	85	LEU	CD1-CG-CD2	-5.93	97.75	110.80
1	D	96	GLU	N-CA-C	5.88	117.77	111.36
1	E	101	ILE	CB-CA-C	5.85	119.46	111.97
1	C	50	SER	N-CA-C	5.81	117.68	108.67
1	H	101	ILE	CB-CA-C	5.79	119.38	111.97
1	I	89	SER	CA-CB-OG	-5.67	99.77	111.10
1	I	174	LYS	CD-CE-NZ	5.54	129.62	111.90
1	G	49	TYR	CA-C-O	-5.37	114.92	121.36
1	C	181	LYS	N-CA-C	-5.36	105.88	112.90
1	E	32	TYR	N-CA-C	5.34	117.18	111.36
1	B	34	GLY	CA-C-O	-5.32	109.62	120.80
1	D	176	ILE	CB-CA-C	-5.30	105.19	111.97
1	K	96	GLU	N-CA-C	5.28	117.12	111.36
1	L	176	ILE	CB-CA-C	-5.28	105.22	111.97
1	J	176	ILE	CB-CA-C	-5.27	105.22	111.97
1	H	49	TYR	CA-C-N	5.27	130.27	123.00
1	H	49	TYR	C-N-CA	5.27	130.27	123.00
1	L	50	SER	N-CA-C	5.26	116.83	108.67
1	A	129	HIS	N-CA-C	5.26	117.92	111.82
1	B	85	LEU	CD1-CG-CD2	-5.25	99.25	110.80
1	F	176	ILE	CB-CA-C	-5.24	105.27	111.97
1	B	125	LYS	CB-CG-CD	5.19	123.23	111.30
1	C	125	LYS	CB-CG-CD	5.19	123.23	111.30
1	B	93	HIS	CA-CB-CG	5.18	118.98	113.80
1	H	86	THR	OG1-CB-CG2	-5.18	98.94	109.30
1	D	101	ILE	CB-CA-C	5.16	119.11	112.14
1	C	88	ILE	CA-C-O	-5.12	114.93	120.67
1	H	94	LYS	CB-CG-CD	5.12	123.08	111.30
1	K	176	ILE	CB-CA-C	-5.11	105.43	111.97
1	C	34	GLY	CA-C-O	-5.11	110.08	120.80
1	A	96	GLU	CB-CG-CD	5.08	121.23	112.60
1	A	174	LYS	CD-CE-NZ	5.03	127.98	111.90
1	L	89	SER	CA-C-O	-5.01	116.23	121.55
1	B	20	SER	CA-CB-OG	-5.01	101.08	111.10

There are no chirality outliers.

There are no planarity outliers.

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	1392	0	1342	1	0
1	B	1360	0	1308	6	0
1	C	1369	0	1320	4	0
1	D	1378	0	1326	4	0
1	E	1379	0	1325	8	0
1	F	1368	0	1320	8	0
1	G	1380	0	1330	7	0
1	H	1360	0	1307	14	0
1	I	1375	0	1325	4	0
1	J	1366	0	1313	4	0
1	K	1366	0	1313	5	0
1	L	1354	0	1301	9	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	3	0	0	0	0
2	F	4	0	0	0	0
2	G	3	0	0	0	0
2	H	4	0	0	0	0
2	I	2	0	0	0	0
2	J	1	0	0	0	0
2	K	2	0	0	0	0
2	L	5	0	0	0	0
3	A	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
4	B	6	0	8	2	0
4	C	12	0	16	0	0
4	D	12	0	16	1	0
4	F	18	0	23	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	6	0	8	2	0
4	I	12	0	16	1	0
4	L	6	0	8	0	0
5	C	1	0	0	0	0
5	D	1	0	0	1	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
5	K	1	0	0	0	0
5	L	3	0	0	2	0
6	A	96	0	0	0	1
6	B	106	0	0	2	0
6	C	106	0	0	1	0
6	D	95	0	0	0	0
6	E	110	0	0	2	0
6	F	146	0	0	2	1
6	G	133	0	0	3	0
6	H	102	0	0	1	0
6	I	119	0	0	2	0
6	J	102	0	0	1	0
6	K	137	0	0	0	0
6	L	146	0	0	3	0
All	All	17961	0	15925	71	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 2.

All (71) 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:L:126:SER:OG	1:L:128:ASP:OD2	1.58	1.19
1:E:15:ASN:HB2	6:E:386:HOH:O	1.82	0.79
1:J:34:GLY:O	6:J:301:HOH:O	2.02	0.78
1:H:17:GLU:OE2	1:H:138:TYR:OH	2.02	0.77
1:G:34:GLY:HA3	1:G:47:SER:O	1.90	0.71
1:I:34:GLY:O	6:I:301:HOH:O	2.11	0.67
1:H:153:LEU:O	1:H:157:GLU:HG3	1.96	0.64
1:E:29:SER:O	1:E:33:THR:HG23	1.99	0.62
1:G:21:SER:OG	6:G:301:HOH:O	2.17	0.60
1:D:93:HIS:HD2	4:D:203:GOL:O2	1.86	0.59
1:H:49:TYR:O	1:H:49:TYR:CD1	2.55	0.59
1:H:49:TYR:CD1	1:H:49:TYR:C	2.79	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:85:LEU:H	1:K:93:HIS:HE1	1.54	0.56
1:B:49:TYR:C	1:B:49:TYR:CD1	2.83	0.56
1:B:15:ASN:HB3	6:B:369:HOH:O	2.05	0.56
1:F:74:VAL:HG22	4:F:201:GOL:H32	1.88	0.56
4:H:204:GOL:H32	1:K:74:VAL:HG22	1.88	0.55
1:I:162:GLU:H	1:I:162:GLU:CD	2.13	0.55
1:C:87:SER:OG	5:D:204:CL:CL	2.61	0.54
1:C:162:GLU:H	1:C:162:GLU:CD	2.16	0.53
1:H:16:LYS:HE3	1:H:112:ILE:HD13	1.91	0.53
1:E:15:ASN:ND2	1:E:86:THR:OG1	2.42	0.53
1:F:49:TYR:C	1:F:49:TYR:CD1	2.85	0.52
1:E:92:GLU:HG3	6:E:382:HOH:O	2.09	0.52
1:H:49:TYR:C	1:H:49:TYR:HD1	2.17	0.52
1:H:49:TYR:O	1:H:49:TYR:HD1	1.93	0.51
1:H:6:ILE:HG12	1:H:122:HIS:HD2	1.74	0.50
1:H:162:GLU:H	1:H:162:GLU:CD	2.19	0.50
1:H:116:ILE:HG12	1:H:138:TYR:CE2	2.46	0.50
1:L:162:GLU:H	1:L:162:GLU:CD	2.21	0.49
4:B:203:GOL:H11	1:E:74:VAL:HG22	1.95	0.48
1:L:104:LYS:HE3	5:L:209:CL:CL	2.51	0.48
1:G:84:GLN:NE2	6:G:307:HOH:O	2.47	0.47
1:F:162:GLU:HG2	6:F:369:HOH:O	2.13	0.47
1:G:1:MET:HE2	1:G:1:MET:N	2.30	0.47
1:B:49:TYR:C	1:B:49:TYR:HD1	2.22	0.47
1:D:162:GLU:H	1:D:162:GLU:CD	2.21	0.46
1:A:155:LYS:NZ	1:D:52:ASP:OD1	2.45	0.46
1:H:104:LYS:HE3	6:H:330:HOH:O	2.15	0.46
1:B:154:ASP:HB2	6:B:313:HOH:O	2.16	0.46
1:H:85:LEU:H	1:K:93:HIS:CE1	2.34	0.46
1:L:125:LYS:NZ	6:L:305:HOH:O	2.46	0.46
1:F:154:ASP:OD2	6:F:301:HOH:O	2.20	0.46
1:C:111:HIS:CE1	6:C:307:HOH:O	2.69	0.45
1:E:93:HIS:CE1	1:E:94:LYS:HG3	2.51	0.45
1:I:162:GLU:OE1	1:I:162:GLU:N	2.36	0.45
1:L:128:ASP:HB3	5:L:210:CL:CL	2.54	0.45
1:F:49:TYR:CD1	1:F:49:TYR:O	2.70	0.45
1:F:116:ILE:HD12	1:F:142:GLN:HG2	1.99	0.45
1:G:154:ASP:HB2	6:G:314:HOH:O	2.16	0.45
1:B:137:TRP:CZ2	1:B:141:GLU:HG3	2.52	0.44
1:F:49:TYR:C	1:F:49:TYR:HD1	2.25	0.44
4:H:204:GOL:C3	1:K:74:VAL:HG22	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:74:VAL:HG22	4:F:201:GOL:C3	2.48	0.44
1:E:116:ILE:HD12	1:E:142:GLN:HG2	2.00	0.43
4:I:204:GOL:C3	6:I:302:HOH:O	2.66	0.43
1:K:116:ILE:HD12	1:K:142:GLN:HG2	2.01	0.43
1:L:116:ILE:HD12	1:L:142:GLN:HG2	2.01	0.43
1:I:116:ILE:HD12	1:I:142:GLN:HG2	2.01	0.43
1:C:116:ILE:HD12	1:C:142:GLN:HG2	2.00	0.42
1:B:17:GLU:OE2	1:B:138:TYR:OH	2.18	0.42
1:D:116:ILE:HD12	1:D:142:GLN:HG2	2.02	0.42
1:G:163:ASN:OD1	1:J:164:HIS:HE1	2.03	0.42
1:H:137:TRP:CE2	1:H:141:GLU:HG3	2.55	0.42
1:J:116:ILE:HD12	1:J:142:GLN:HG2	2.01	0.41
4:B:203:GOL:C1	1:E:74:VAL:HG22	2.50	0.41
1:G:116:ILE:HD12	1:G:142:GLN:HG2	2.02	0.41
1:L:154:ASP:OD1	6:L:302:HOH:O	2.22	0.41
1:L:2:LEU:HD12	1:L:81:VAL:HG21	2.03	0.41
1:L:174:LYS:HD3	6:L:313:HOH:O	2.20	0.41
1:J:104:LYS:HE2	1:J:104:LYS:HB2	1.80	0.40

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)
6:A:377:HOH:O	6:F:384:HOH:O[3_765]	2.13	0.07

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	168/182 (92%)	165 (98%)	3 (2%)	0	100	100
1	B	163/182 (90%)	163 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	164/182 (90%)	163 (99%)	1 (1%)	0	100	100
1	D	165/182 (91%)	164 (99%)	1 (1%)	0	100	100
1	E	166/182 (91%)	164 (99%)	2 (1%)	0	100	100
1	F	164/182 (90%)	163 (99%)	1 (1%)	0	100	100
1	G	166/182 (91%)	164 (99%)	2 (1%)	0	100	100
1	H	163/182 (90%)	162 (99%)	1 (1%)	0	100	100
1	I	165/182 (91%)	164 (99%)	1 (1%)	0	100	100
1	J	164/182 (90%)	162 (99%)	2 (1%)	0	100	100
1	K	164/182 (90%)	161 (98%)	3 (2%)	0	100	100
1	L	162/182 (89%)	161 (99%)	1 (1%)	0	100	100
All	All	1974/2184 (90%)	1956 (99%)	18 (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	156/163 (96%)	154 (99%)	2 (1%)	61	68
1	B	151/163 (93%)	149 (99%)	2 (1%)	61	68
1	C	152/163 (93%)	149 (98%)	3 (2%)	48	54
1	D	153/163 (94%)	152 (99%)	1 (1%)	76	82
1	E	154/163 (94%)	150 (97%)	4 (3%)	40	44
1	F	152/163 (93%)	150 (99%)	2 (1%)	61	68
1	G	154/163 (94%)	151 (98%)	3 (2%)	50	56
1	H	151/163 (93%)	150 (99%)	1 (1%)	76	82
1	I	153/163 (94%)	151 (99%)	2 (1%)	61	68
1	J	152/163 (93%)	149 (98%)	3 (2%)	48	54
1	K	152/163 (93%)	151 (99%)	1 (1%)	76	82

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	150/163 (92%)	150 (100%)	0	100	100
All	All	1830/1956 (94%)	1806 (99%)	24 (1%)	61	68

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS
1	A	6	ILE
1	B	2	LEU
1	B	6	ILE
1	C	1	MET
1	C	6	ILE
1	C	80	ASN
1	D	6	ILE
1	E	4	LYS
1	E	6	ILE
1	E	19	ASN
1	E	46	SER
1	F	1	MET
1	F	125	LYS
1	G	6	ILE
1	G	46	SER
1	G	48	SER
1	H	6	ILE
1	I	6	ILE
1	I	65	GLU
1	J	4	LYS
1	J	6	ILE
1	J	162	GLU
1	K	6	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	80	ASN
1	B	129	HIS
1	C	13	GLN
1	C	93	HIS
1	D	13	GLN
1	D	19	ASN

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Mol	Chain	Res	Type
1	D	22	ASN
1	D	80	ASN
1	D	93	HIS
1	E	13	GLN
1	E	15	ASN
1	E	80	ASN
1	F	13	GLN
1	F	80	ASN
1	G	13	GLN
1	H	80	ASN
1	H	84	GLN
1	H	122	HIS
1	H	129	HIS
1	I	13	GLN
1	I	111	HIS
1	J	13	GLN
1	K	13	GLN
1	K	93	HIS
1	K	122	HIS
1	L	13	GLN
1	L	93	HIS
1	L	118	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](#)

Of 56 ligands modelled in this entry, 44 are monoatomic - leaving 12 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	$\# Z > 2$
4	GOL	F	208	-	5,5,5	0.41	0	5,5,5	0.75	0
4	GOL	L	207	-	5,5,5	0.91	0	5,5,5	1.24	0
4	GOL	C	203	-	5,5,5	0.91	0	5,5,5	0.86	0
4	GOL	D	203	-	5,5,5	0.52	0	5,5,5	0.92	0
4	GOL	B	203	-	5,5,5	0.72	0	5,5,5	1.08	0
4	GOL	D	202	-	5,5,5	0.51	0	5,5,5	0.62	0
4	GOL	I	203	-	5,5,5	0.47	0	5,5,5	1.04	0
4	GOL	I	204	-	5,5,5	0.49	0	5,5,5	0.83	0
4	GOL	H	204	-	5,5,5	0.49	0	5,5,5	0.91	0
4	GOL	F	201	2	5,5,5	0.43	0	5,5,5	1.06	0
4	GOL	F	207	2	5,5,5	0.44	0	5,5,5	0.60	0
4	GOL	C	202	-	5,5,5	0.35	0	5,5,5	0.87	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	F	208	-	-	0/4/4/4	-
4	GOL	L	207	-	-	2/4/4/4	-
4	GOL	C	203	-	-	0/4/4/4	-
4	GOL	D	203	-	-	0/4/4/4	-
4	GOL	B	203	-	-	2/4/4/4	-
4	GOL	D	202	-	-	0/4/4/4	-
4	GOL	I	203	-	-	3/4/4/4	-
4	GOL	I	204	-	-	1/4/4/4	-
4	GOL	H	204	-	-	2/4/4/4	-
4	GOL	F	201	2	-	2/4/4/4	-
4	GOL	F	207	2	-	2/4/4/4	-
4	GOL	C	202	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	203	GOL	C1-C2-C3-O3
4	C	202	GOL	C1-C2-C3-O3
4	F	201	GOL	O1-C1-C2-O2
4	F	201	GOL	O1-C1-C2-C3
4	H	204	GOL	O1-C1-C2-C3
4	I	203	GOL	C1-C2-C3-O3
4	L	207	GOL	C1-C2-C3-O3
4	L	207	GOL	O2-C2-C3-O3
4	F	207	GOL	O1-C1-C2-C3
4	B	203	GOL	O2-C2-C3-O3
4	H	204	GOL	O1-C1-C2-O2
4	C	202	GOL	O2-C2-C3-O3
4	I	203	GOL	O1-C1-C2-O2
4	F	207	GOL	O1-C1-C2-O2
4	I	204	GOL	O2-C2-C3-O3
4	I	203	GOL	O2-C2-C3-O3

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	203	GOL	1	0
4	B	203	GOL	2	0
4	I	204	GOL	1	0
4	H	204	GOL	2	0
4	F	201	GOL	2	0

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	169/182 (92%)	-0.17	5 (2%)	52	51	15, 24, 47, 66	3 (1%)
1	B	167/182 (91%)	-0.09	5 (2%)	52	51	14, 26, 47, 61	0
1	C	168/182 (92%)	-0.19	4 (2%)	59	59	16, 24, 46, 81	0
1	D	169/182 (92%)	-0.08	5 (2%)	52	51	16, 27, 48, 79	0
1	E	169/182 (92%)	-0.15	7 (4%)	41	40	15, 24, 47, 55	1 (0%)
1	F	168/182 (92%)	-0.29	5 (2%)	52	51	15, 22, 43, 76	0
1	G	170/182 (93%)	-0.24	5 (2%)	53	52	13, 23, 45, 65	0
1	H	167/182 (91%)	-0.12	5 (2%)	52	51	13, 24, 48, 62	0
1	I	169/182 (92%)	-0.23	4 (2%)	59	59	14, 23, 44, 77	0
1	J	168/182 (92%)	-0.11	4 (2%)	59	59	14, 25, 47, 64	0
1	K	168/182 (92%)	-0.32	3 (1%)	67	67	14, 23, 47, 63	0
1	L	166/182 (91%)	-0.34	3 (1%)	67	67	14, 21, 41, 59	0
All	All	2018/2184 (92%)	-0.19	55 (2%)	56	55	13, 24, 47, 81	4 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	49	TYR	4.8
1	I	49	TYR	4.4
1	C	1	MET	4.3
1	C	49	TYR	4.3
1	L	49	TYR	4.3
1	F	48	SER	4.1
1	K	49	TYR	4.0
1	C	34	GLY	4.0
1	D	49	TYR	4.0
1	I	1	MET	3.8
1	J	34	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	H	49	TYR	3.6
1	F	49	TYR	3.5
1	J	88	ILE	3.5
1	L	34	GLY	3.5
1	A	49	TYR	3.5
1	A	35	SER	3.4
1	K	34	GLY	3.4
1	B	49	TYR	3.3
1	E	33	THR	3.2
1	D	34	GLY	3.2
1	F	34	GLY	3.2
1	E	34	GLY	3.1
1	I	34	GLY	3.1
1	E	49	TYR	3.0
1	I	48	SER	3.0
1	B	48	SER	2.9
1	H	35	SER	2.9
1	G	1	MET	2.7
1	G	34	GLY	2.7
1	L	163	ASN	2.7
1	E	163	ASN	2.7
1	A	34	GLY	2.6
1	A	1	MET	2.6
1	B	34	GLY	2.5
1	H	138	TYR	2.5
1	E	48	SER	2.5
1	B	137	TRP	2.5
1	A	48	SER	2.5
1	D	35	SER	2.5
1	B	138	TYR	2.3
1	D	163	ASN	2.3
1	G	47	SER	2.3
1	G	163	ASN	2.2
1	F	1	MET	2.2
1	H	137	TRP	2.2
1	E	46	SER	2.2
1	E	47	SER	2.2
1	H	163	ASN	2.2
1	G	49	TYR	2.2
1	J	48	SER	2.1
1	F	163	ASN	2.1
1	D	182	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	162	GLU	2.0
1	K	181	LYS	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](#)

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
3	FE	E	204	1/1	0.75	0.14	37,37,37,37	1
4	GOL	I	203	6/6	0.78	0.14	36,38,40,43	0
4	GOL	D	203	6/6	0.82	0.16	43,46,47,49	0
3	FE	K	203	1/1	0.84	0.11	38,38,38,38	1
2	NA	H	205	1/1	0.85	0.15	49,49,49,49	0
4	GOL	C	202	6/6	0.85	0.15	50,52,59,59	0
4	GOL	F	201	6/6	0.86	0.13	38,44,49,50	0
4	GOL	B	203	6/6	0.86	0.13	32,45,46,54	0
3	FE	L	206	1/1	0.87	0.09	42,42,42,42	1
2	NA	G	205	1/1	0.87	0.11	44,44,44,44	0
4	GOL	F	208	6/6	0.87	0.13	40,46,53,54	0
4	GOL	H	204	6/6	0.87	0.13	30,39,42,45	0
2	NA	F	205	1/1	0.87	0.12	53,53,53,53	0
2	NA	L	204	1/1	0.88	0.10	52,52,52,52	0
4	GOL	L	207	6/6	0.88	0.12	33,44,46,50	0
2	NA	B	202	1/1	0.89	0.10	46,46,46,46	0
2	NA	I	202	1/1	0.89	0.12	51,51,51,51	0
4	GOL	I	204	6/6	0.89	0.12	36,48,51,53	0
4	GOL	D	202	6/6	0.89	0.13	39,46,55,58	0
4	GOL	F	207	6/6	0.91	0.12	34,43,46,50	0
2	NA	A	202	1/1	0.91	0.08	41,41,41,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	C	203	6/6	0.91	0.13	32,35,36,38	0
2	NA	E	203	1/1	0.91	0.21	46,46,46,46	0
2	NA	H	202	1/1	0.91	0.13	49,49,49,49	0
2	NA	L	205	1/1	0.91	0.20	43,43,43,43	0
2	NA	H	203	1/1	0.93	0.20	38,38,38,38	0
2	NA	F	203	1/1	0.93	0.12	40,40,40,40	0
3	FE	F	206	1/1	0.94	0.06	37,37,37,37	1
2	NA	H	201	1/1	0.95	0.05	25,25,25,25	0
2	NA	F	204	1/1	0.96	0.09	23,23,23,23	0
2	NA	E	202	1/1	0.96	0.11	29,29,29,29	0
2	NA	K	202	1/1	0.96	0.09	46,46,46,46	0
2	NA	A	201	1/1	0.97	0.04	22,22,22,22	0
2	NA	G	201	1/1	0.97	0.07	22,22,22,22	0
2	NA	G	202	1/1	0.97	0.13	36,36,36,36	0
2	NA	E	201	1/1	0.97	0.05	22,22,22,22	0
2	NA	I	201	1/1	0.97	0.08	23,23,23,23	0
5	CL	D	204	1/1	0.97	0.14	34,34,34,34	1
3	FE	A	203	1/1	0.98	0.03	41,41,41,41	0
2	NA	D	201	1/1	0.98	0.03	26,26,26,26	0
2	NA	L	202	1/1	0.98	0.03	26,26,26,26	0
2	NA	L	203	1/1	0.98	0.10	29,29,29,29	0
2	NA	C	201	1/1	0.98	0.06	23,23,23,23	0
2	NA	J	201	1/1	0.98	0.05	24,24,24,24	0
5	CL	G	204	1/1	0.98	0.08	34,34,34,34	0
5	CL	L	209	1/1	0.98	0.05	35,35,35,35	0
5	CL	L	210	1/1	0.98	0.05	35,35,35,35	0
2	NA	F	202	1/1	0.99	0.03	22,22,22,22	0
3	FE	G	203	1/1	0.99	0.03	39,39,39,39	0
5	CL	F	209	1/1	0.99	0.04	22,22,22,22	0
2	NA	L	201	1/1	0.99	0.03	19,19,19,19	0
5	CL	K	204	1/1	0.99	0.15	32,32,32,32	0
5	CL	L	208	1/1	0.99	0.06	21,21,21,21	0
2	NA	B	201	1/1	0.99	0.04	23,23,23,23	0
2	NA	K	201	1/1	0.99	0.03	24,24,24,24	0
5	CL	C	204	1/1	1.00	0.01	24,24,24,24	1

6.5 Other polymers

There are no such residues in this entry.