



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:19 PM UTC

PDB ID : 6OFU / pdb_00006ofu
Title : X-ray crystal structure of the YdjI aldolase from Escherichia coli K12
Authors : Dopkins, B.J.; Thoden, J.B.; Huddleston, J.P.; Narindoshvili, T.; Fose, B.; Rachel, F.M.; Holden, H.M.
Deposited on : 2019-04-01
Resolution : 1.75 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

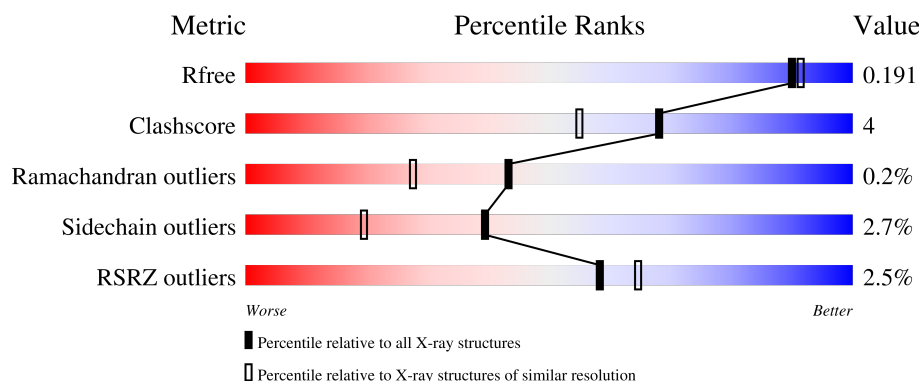
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	286	2% 82% 7% • 9%
1	B	286	2% 84% 8% • 8%
1	C	286	3% 85% 5% •• 8%
1	D	286	2% 79% 10% • 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9148 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 YdjI aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	2	0
			2046	1292	358	384	12			
1	B	264	Total	C	N	O	S	0	0	0
			2061	1301	354	394	12			
1	C	262	Total	C	N	O	S	0	1	0
			2044	1290	352	390	12			
1	D	256	Total	C	N	O	S	0	1	0
			1991	1259	342	378	12			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	279	LEU	-	expression tag	UNP P77704
A	280	GLU	-	expression tag	UNP P77704
A	281	HIS	-	expression tag	UNP P77704
A	282	HIS	-	expression tag	UNP P77704
A	283	HIS	-	expression tag	UNP P77704
A	284	HIS	-	expression tag	UNP P77704
A	285	HIS	-	expression tag	UNP P77704
A	286	HIS	-	expression tag	UNP P77704
B	279	LEU	-	expression tag	UNP P77704
B	280	GLU	-	expression tag	UNP P77704
B	281	HIS	-	expression tag	UNP P77704
B	282	HIS	-	expression tag	UNP P77704
B	283	HIS	-	expression tag	UNP P77704
B	284	HIS	-	expression tag	UNP P77704
B	285	HIS	-	expression tag	UNP P77704
B	286	HIS	-	expression tag	UNP P77704
C	279	LEU	-	expression tag	UNP P77704
C	280	GLU	-	expression tag	UNP P77704
C	281	HIS	-	expression tag	UNP P77704
C	282	HIS	-	expression tag	UNP P77704
C	283	HIS	-	expression tag	UNP P77704

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Chain	Residue	Modelled	Actual	Comment	Reference
C	284	HIS	-	expression tag	UNP P77704
C	285	HIS	-	expression tag	UNP P77704
C	286	HIS	-	expression tag	UNP P77704
D	279	LEU	-	expression tag	UNP P77704
D	280	GLU	-	expression tag	UNP P77704
D	281	HIS	-	expression tag	UNP P77704
D	282	HIS	-	expression tag	UNP P77704
D	283	HIS	-	expression tag	UNP P77704
D	284	HIS	-	expression tag	UNP P77704
D	285	HIS	-	expression tag	UNP P77704
D	286	HIS	-	expression tag	UNP P77704

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	2	Total Zn 2 2	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	B	1	Total Cl 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	232	Total O 232 232	0	0
4	B	274	Total O 274 274	0	0
4	C	261	Total O 261 261	0	0

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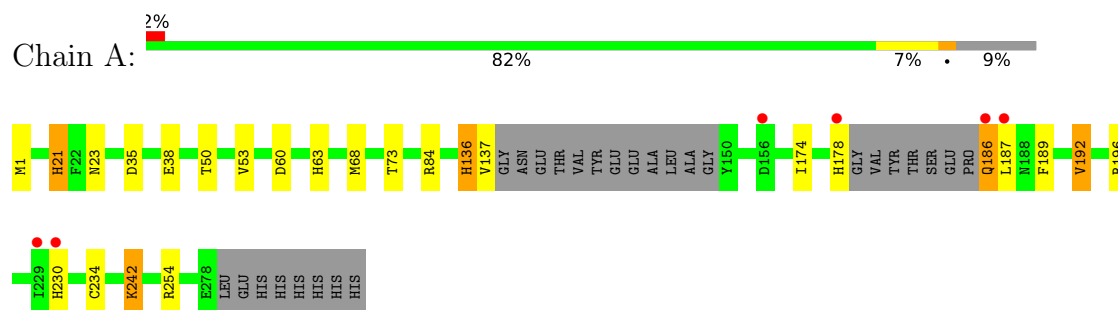
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	232	Total	O	0	0
			232	232		

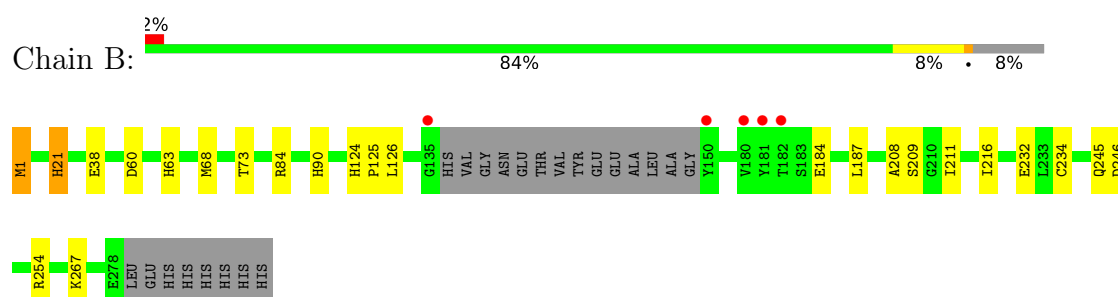
3 Residue-property plots [i](#)

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.

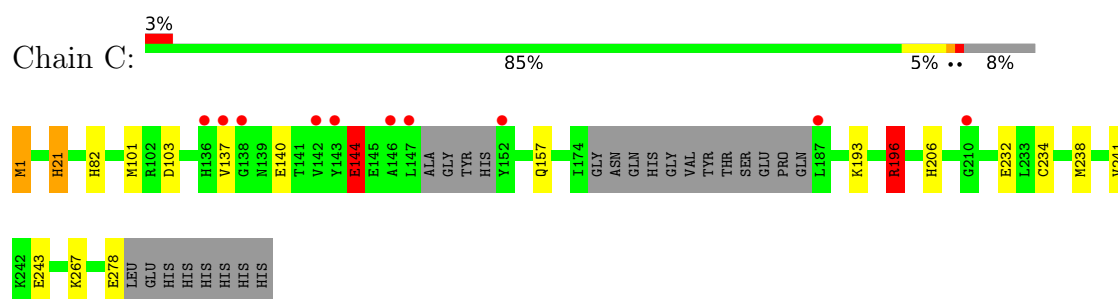
• Molecule 1: YdjI aldolase



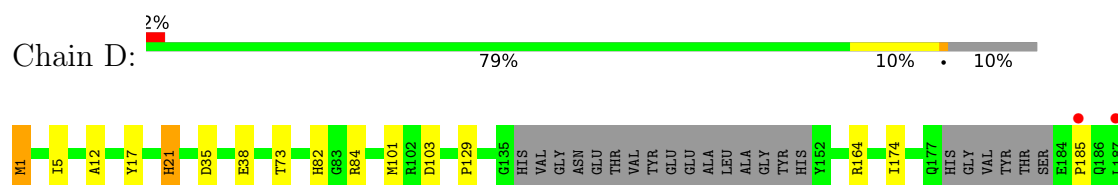
• Molecule 1: YdjI aldolase

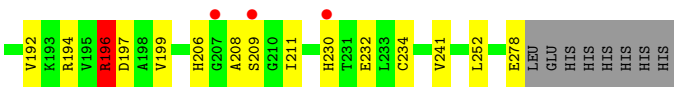


• Molecule 1: YdjI aldolase



• Molecule 1: YdjI aldolase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.76Å 85.66Å 84.27Å 90.00° 109.88° 90.00°	Depositor
Resolution (Å)	34.74 – 1.75 34.74 – 1.75	Depositor EDS
% Data completeness (in resolution range)	96.5 (34.74-1.75) 96.5 (34.74-1.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.39 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.141 , 0.182 0.153 , 0.191	Depositor DCC
R_{free} test set	5200 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.019 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9148	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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: OCS, ZN, 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	0.84	0/2086	1.00	1/2819 (0.0%)
1	B	0.83	0/2095	0.97	0/2835
1	C	0.81	0/2079	1.00	1/2812 (0.0%)
1	D	0.85	0/2024	0.99	2/2737 (0.1%)
All	All	0.83	0/8284	0.99	4/11203 (0.0%)

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	4
1	B	0	1
1	C	0	1
1	D	0	3
All	All	0	9

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	35	ASP	CA-CB-CG	6.59	119.19	112.60
1	D	35	ASP	CA-CB-CG	6.26	118.86	112.60
1	D	197	ASP	CB-CA-C	5.19	120.96	110.38
1	C	144	GLU	CB-CG-CD	5.03	121.15	112.60

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	196	ARG	Sidechain
1	A	254[A]	ARG	Sidechain
1	A	254[B]	ARG	Sidechain
1	A	84	ARG	Sidechain
1	B	84	ARG	Sidechain
1	C	196	ARG	Sidechain
1	D	164	ARG	Sidechain
1	D	196	ARG	Sidechain
1	D	84	ARG	Sidechain

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	2046	0	1990	16	0
1	B	2061	0	1997	17	0
1	C	2044	0	1975	14	0
1	D	1991	0	1939	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	232	0	0	2	0
4	B	274	0	0	2	0
4	C	261	0	0	1	0
4	D	232	0	0	4	0
All	All	9148	0	7901	58	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 4.

All (58) 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:D:82:HIS:HE1	4:D:505:HOH:O	1.26	1.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:194:ARG:NH2	4:D:401:HOH:O	1.80	1.11
1:D:196:ARG:HD3	1:D:196:ARG:O	1.82	0.80
1:B:254:ARG:NH2	4:B:401:HOH:O	2.17	0.77
1:D:82:HIS:CE1	4:D:505:HOH:O	2.09	0.75
1:A:136[B]:HIS:HE1	4:A:1278:HOH:O	1.69	0.73
1:C:140:GLU:HB2	1:C:144:GLU:HG2	1.73	0.70
1:D:38:GLU:OE1	1:D:73:THR:OG1	2.13	0.65
1:C:82[B]:HIS:HD2	1:C:103:ASP:OD2	1.82	0.63
1:B:90:HIS:HD2	4:B:491:HOH:O	1.80	0.62
1:A:242:LYS:NZ	1:B:246:ASP:OD1	2.32	0.61
1:A:136[A]:HIS:HE1	4:A:1218:HOH:O	1.84	0.59
1:C:196:ARG:O	1:C:196:ARG:HD3	2.03	0.58
1:C:232:GLU:HG3	4:C:429:HOH:O	2.04	0.57
1:B:232:GLU:OE2	1:B:267:LYS:HE3	2.05	0.57
1:C:232:GLU:OE2	1:C:267:LYS:NZ	2.39	0.56
1:D:196:ARG:HD3	1:D:196:ARG:C	2.27	0.55
1:D:241:VAL:HA	1:D:252[A]:LEU:HD11	1.88	0.55
1:A:186:GLN:HG3	1:A:187:LEU:H	1.73	0.53
1:A:137:VAL:HG23	1:A:178:HIS:CE1	2.43	0.53
1:A:242:LYS:HE3	1:B:245:GLN:OE1	2.11	0.51
1:A:174:ILE:HG21	1:A:192:VAL:HG22	1.91	0.51
1:A:21:HIS:C	1:A:21:HIS:CD2	2.89	0.50
1:B:208:ALA:HB1	1:B:216:ILE:HD11	1.92	0.50
1:D:232:GLU:HG3	4:D:545:HOH:O	2.12	0.50
1:C:137:VAL:HB	1:C:157:GLN:OE1	2.12	0.49
1:D:208:ALA:HA	1:D:211:ILE:HD12	1.95	0.49
1:A:187:LEU:HD23	1:A:189:PHE:CZ	2.48	0.47
1:C:238:MET:SD	1:D:252[B]:LEU:HD13	2.54	0.47
1:C:21:HIS:C	1:C:21:HIS:CD2	2.92	0.47
1:C:196:ARG:HD3	1:C:196:ARG:C	2.38	0.47
1:B:232:GLU:OE2	1:B:267:LYS:CE	2.64	0.46
1:D:174:ILE:HG21	1:D:192:VAL:HG22	1.98	0.45
1:D:12:ALA:HA	1:D:17:TYR:CE2	2.51	0.45
1:A:60:ASP:CG	1:B:68:MET:HG2	2.42	0.45
1:A:23:ASN:ND2	1:A:230:HIS:HB2	2.32	0.45
1:B:38:GLU:OE1	1:B:73:THR:OG1	2.25	0.45
1:B:124:HIS:N	1:B:125:PRO:CD	2.80	0.45
1:D:101:MET:HE1	1:D:206:HIS:HE1	1.82	0.45
1:D:82:HIS:HD2	1:D:103:ASP:OD2	1.99	0.45
1:C:241:VAL:HG11	1:D:252[B]:LEU:HD11	1.98	0.44
1:A:186:GLN:CG	1:A:187:LEU:H	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:241:VAL:HG11	1:D:252[B]:LEU:CD1	2.47	0.43
1:C:1:MET:HE2	1:C:1:MET:HB2	1.93	0.43
1:B:126:LEU:HD21	1:D:1:MET:HG3	2.01	0.43
1:A:50:THR:O	1:A:53:VAL:HG22	2.19	0.42
1:A:63:HIS:HB3	1:B:63:HIS:HB3	2.02	0.42
1:D:194:ARG:HD3	1:D:194:ARG:HA	1.82	0.42
1:B:1:MET:HE2	1:B:1:MET:HB2	1.91	0.41
1:C:232:GLU:OE2	1:C:267:LYS:CE	2.68	0.41
1:A:38:GLU:OE1	1:A:73:THR:OG1	2.30	0.41
1:B:267:LYS:HA	1:B:267:LYS:HD3	1.89	0.41
1:C:101:MET:HE1	1:C:206:HIS:HE1	1.85	0.41
1:B:187:LEU:HD21	1:B:211:ILE:HD13	2.02	0.41
1:B:21:HIS:CD2	1:B:21:HIS:C	2.99	0.41
1:D:21:HIS:CD2	1:D:21:HIS:C	2.98	0.40
1:A:68:MET:HG2	1:B:60:ASP:CG	2.46	0.40
1:D:5:ILE:HG22	1:D:129:PRO:HG3	2.03	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	254/286 (89%)	252 (99%)	2 (1%)	0	100	100
1	B	259/286 (91%)	256 (99%)	3 (1%)	0	100	100
1	C	256/286 (90%)	254 (99%)	2 (1%)	0	100	100
1	D	250/286 (87%)	248 (99%)	0	2 (1%)	16	5
All	All	1019/1144 (89%)	1010 (99%)	7 (1%)	2 (0%)	43	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	230	HIS
1	D	185	PRO

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	215/235 (92%)	208 (97%)	7 (3%)	33	13
1	B	216/235 (92%)	212 (98%)	4 (2%)	50	31
1	C	213/235 (91%)	206 (97%)	7 (3%)	33	13
1	D	208/235 (88%)	202 (97%)	6 (3%)	37	17
All	All	852/940 (91%)	828 (97%)	24 (3%)	39	18

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	21	HIS
1	A	136[A]	HIS
1	A	136[B]	HIS
1	A	186	GLN
1	A	192	VAL
1	A	242	LYS
1	B	1	MET
1	B	21	HIS
1	B	184	GLU
1	B	209	SER
1	C	1	MET
1	C	21	HIS
1	C	144	GLU
1	C	193	LYS
1	C	196	ARG
1	C	243	GLU
1	C	278	GLU
1	D	1	MET
1	D	21	HIS

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Mol	Chain	Res	Type
1	D	196	ARG
1	D	199	VAL
1	D	209	SER
1	D	278	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	177	GLN
1	A	186	GLN
1	A	230	HIS
1	B	10	ASN
1	B	90	HIS
1	B	91	ASN
1	C	206	HIS
1	C	235	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are 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	$\# Z > 2$
1	OCS	A	234	1	6,8,9	2.13	2 (33%)	7,11,13	1.54	2 (28%)
1	OCS	C	234	1	6,8,9	1.79	2 (33%)	7,11,13	1.54	2 (28%)
1	OCS	D	234	1	6,8,9	1.90	2 (33%)	7,11,13	0.90	0
1	OCS	B	234	1	6,8,9	1.56	1 (16%)	7,11,13	1.84	2 (28%)

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
1	OCS	A	234	1	-	0/4/7/9	-
1	OCS	C	234	1	-	0/4/7/9	-
1	OCS	D	234	1	-	0/4/7/9	-
1	OCS	B	234	1	-	0/4/7/9	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	234	OCS	OD1-SG	4.49	1.57	1.45
1	D	234	OCS	OD3-SG	3.87	1.56	1.45
1	C	234	OCS	OD3-SG	3.44	1.54	1.45
1	B	234	OCS	OD2-SG	2.52	1.56	1.47
1	D	234	OCS	OD2-SG	-2.17	1.38	1.47
1	C	234	OCS	CB-CA	2.15	1.55	1.53
1	A	234	OCS	CB-CA	-2.13	1.52	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	234	OCS	OD1-SG-CB	3.59	112.11	106.76
1	C	234	OCS	OD2-SG-OD3	2.69	118.13	111.40
1	A	234	OCS	OD1-SG-CB	2.40	110.34	106.76
1	C	234	OCS	OD2-SG-OD1	-2.29	105.67	111.40
1	A	234	OCS	OD2-SG-CB	2.18	110.19	105.97
1	B	234	OCS	OD2-SG-OD1	2.15	116.78	111.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 7 ligands modelled in this entry, 7 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	258/286 (90%)	-0.22	6 (2%) 61 67	13, 20, 36, 59	2 (0%)
1	B	263/286 (91%)	-0.22	5 (1%) 66 73	15, 20, 39, 80	0
1	C	261/286 (91%)	-0.01	10 (3%) 44 50	12, 22, 46, 65	1 (0%)
1	D	255/286 (89%)	-0.22	5 (1%) 65 71	11, 20, 40, 94	1 (0%)
All	All	1037/1144 (90%)	-0.17	26 (2%) 58 64	11, 21, 41, 94	4 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	137	VAL	5.2
1	C	142	VAL	4.7
1	C	187	LEU	4.7
1	C	152	TYR	4.0
1	B	182	THR	3.9
1	C	147	LEU	3.8
1	C	143	TYR	3.8
1	D	185	PRO	3.5
1	C	146	ALA	3.5
1	C	136	HIS	3.2
1	B	181	TYR	2.8
1	A	230	HIS	2.8
1	B	180	VAL	2.8
1	A	187	LEU	2.6
1	B	135	GLY	2.5
1	D	187	LEU	2.3
1	C	138	GLY	2.3
1	A	178	HIS	2.2
1	A	156	ASP	2.2
1	C	210	GLY	2.2
1	D	207	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	186	GLN	2.1
1	D	230	HIS	2.1
1	D	209	SER	2.1
1	B	150	TYR	2.1
1	A	229	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	OCS	A	234	9/10	0.97	0.07	20,21,28,32	0
1	OCS	B	234	9/10	0.97	0.07	17,19,24,25	0
1	OCS	C	234	9/10	0.98	0.06	17,21,26,28	0
1	OCS	D	234	9/10	0.98	0.06	18,19,24,26	0

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	ZN	D	301	1/1	0.96	0.19	40,40,40,40	0
3	CL	B	302	1/1	0.96	0.10	32,32,32,32	0
2	ZN	C	302	1/1	0.97	0.27	58,58,58,58	0
2	ZN	C	301	1/1	0.98	0.14	60,60,60,60	0
3	CL	A	302	1/1	0.99	0.04	23,23,23,23	0
2	ZN	B	301	1/1	0.99	0.09	26,26,26,26	0
2	ZN	A	301	1/1	1.00	0.09	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.