



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:30 PM UTC

PDB ID : 7E5C / pdb_00007e5c
Title : Bacterial prolidase mutant D45W/L225Y/H226L/H343I
Authors : Yang, J.
Deposited on : 2021-02-18
Resolution : 2.22 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
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

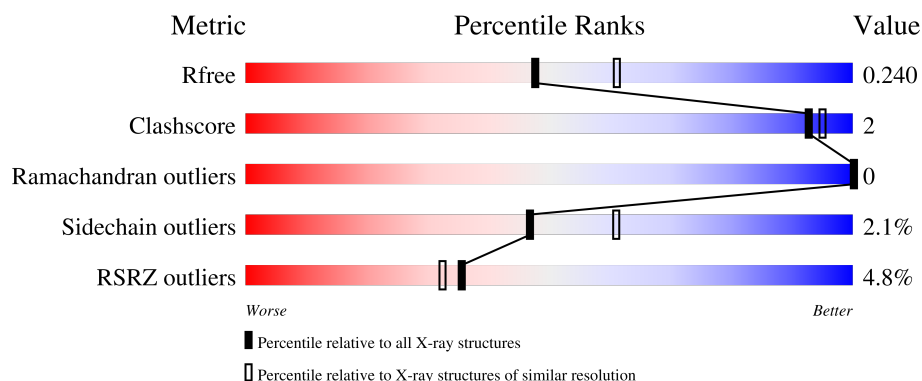
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	448	3% 95% 5% •
1	B	448	4% 93% 5% •
1	C	448	8% 91% 6% •
1	D	448	4% 93% 5% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14635 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 Xaa-Pro dipeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	439	Total	C	N	O	S	0	0	0
			3569	2298	603	654	14			
1	B	439	Total	C	N	O	S	0	0	0
			3569	2298	603	654	14			
1	C	439	Total	C	N	O	S	0	0	0
			3569	2298	603	654	14			
1	D	438	Total	C	N	O	S	0	0	0
			3560	2293	602	651	14			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	TRP	ASP	engineered mutation	UNP A0A1I7CHQ2
A	225	TYR	LEU	engineered mutation	UNP A0A1I7CHQ2
A	226	LEU	HIS	engineered mutation	UNP A0A1I7CHQ2
A	343	ILE	HIS	engineered mutation	UNP A0A1I7CHQ2
A	441	LEU	-	expression tag	UNP A0A1I7CHQ2
A	442	GLU	-	expression tag	UNP A0A1I7CHQ2
A	443	HIS	-	expression tag	UNP A0A1I7CHQ2
A	444	HIS	-	expression tag	UNP A0A1I7CHQ2
A	445	HIS	-	expression tag	UNP A0A1I7CHQ2
A	446	HIS	-	expression tag	UNP A0A1I7CHQ2
A	447	HIS	-	expression tag	UNP A0A1I7CHQ2
A	448	HIS	-	expression tag	UNP A0A1I7CHQ2
B	45	TRP	ASP	engineered mutation	UNP A0A1I7CHQ2
B	225	TYR	LEU	engineered mutation	UNP A0A1I7CHQ2
B	226	LEU	HIS	engineered mutation	UNP A0A1I7CHQ2
B	343	ILE	HIS	engineered mutation	UNP A0A1I7CHQ2
B	441	LEU	-	expression tag	UNP A0A1I7CHQ2
B	442	GLU	-	expression tag	UNP A0A1I7CHQ2
B	443	HIS	-	expression tag	UNP A0A1I7CHQ2
B	444	HIS	-	expression tag	UNP A0A1I7CHQ2
B	445	HIS	-	expression tag	UNP A0A1I7CHQ2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	446	HIS	-	expression tag	UNP A0A1I7CHQ2
B	447	HIS	-	expression tag	UNP A0A1I7CHQ2
B	448	HIS	-	expression tag	UNP A0A1I7CHQ2
C	45	TRP	ASP	engineered mutation	UNP A0A1I7CHQ2
C	225	TYR	LEU	engineered mutation	UNP A0A1I7CHQ2
C	226	LEU	HIS	engineered mutation	UNP A0A1I7CHQ2
C	343	ILE	HIS	engineered mutation	UNP A0A1I7CHQ2
C	441	LEU	-	expression tag	UNP A0A1I7CHQ2
C	442	GLU	-	expression tag	UNP A0A1I7CHQ2
C	443	HIS	-	expression tag	UNP A0A1I7CHQ2
C	444	HIS	-	expression tag	UNP A0A1I7CHQ2
C	445	HIS	-	expression tag	UNP A0A1I7CHQ2
C	446	HIS	-	expression tag	UNP A0A1I7CHQ2
C	447	HIS	-	expression tag	UNP A0A1I7CHQ2
C	448	HIS	-	expression tag	UNP A0A1I7CHQ2
D	45	TRP	ASP	engineered mutation	UNP A0A1I7CHQ2
D	225	TYR	LEU	engineered mutation	UNP A0A1I7CHQ2
D	226	LEU	HIS	engineered mutation	UNP A0A1I7CHQ2
D	343	ILE	HIS	engineered mutation	UNP A0A1I7CHQ2
D	441	LEU	-	expression tag	UNP A0A1I7CHQ2
D	442	GLU	-	expression tag	UNP A0A1I7CHQ2
D	443	HIS	-	expression tag	UNP A0A1I7CHQ2
D	444	HIS	-	expression tag	UNP A0A1I7CHQ2
D	445	HIS	-	expression tag	UNP A0A1I7CHQ2
D	446	HIS	-	expression tag	UNP A0A1I7CHQ2
D	447	HIS	-	expression tag	UNP A0A1I7CHQ2
D	448	HIS	-	expression tag	UNP A0A1I7CHQ2

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	122	Total 122	O 122	0	0
3	B	83	Total 83	O 83	0	0
3	C	61	Total 61	O 61	0	0
3	D	94	Total 94	O 94	0	0

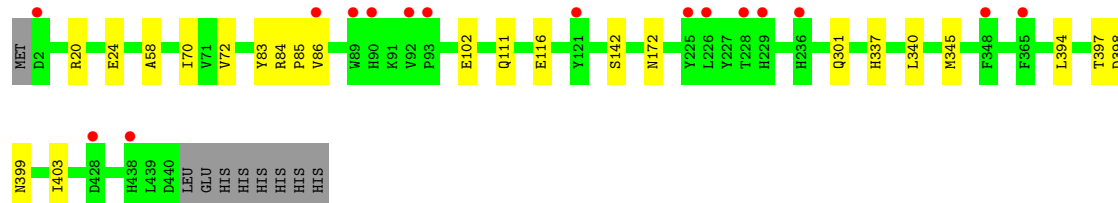
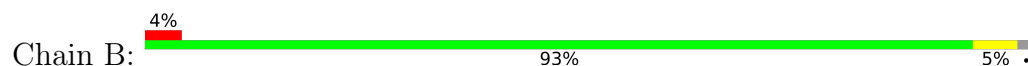
3 Residue-property plots [i](#)

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.

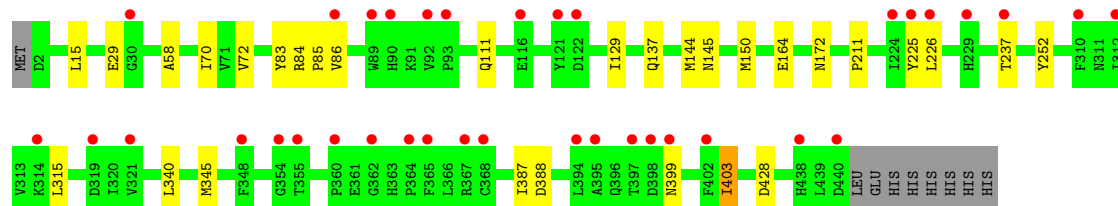
- Molecule 1: Xaa-Pro dipeptidase



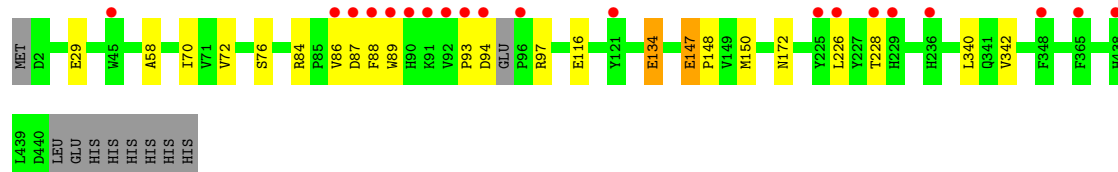
- Molecule 1: Xaa-Pro dipeptidase



- Molecule 1: Xaa-Pro dipeptidase



- Molecule 1: Xaa-Pro dipeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	182.97Å 182.97Å 372.08Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	67.45 – 2.22 67.45 – 2.22	Depositor EDS
% Data completeness (in resolution range)	93.9 (67.45-2.22) 93.9 (67.45-2.22)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.22Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.203 , 0.235 0.211 , 0.240	Depositor DCC
R_{free} test set	5564 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	45.2	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 31.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14635	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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: MN

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	0.97	0/3672	1.12	0/4987
1	B	0.98	0/3672	1.12	2/4987 (0.0%)
1	C	1.06	1/3672 (0.0%)	1.14	2/4987 (0.0%)
1	D	0.98	0/3662	1.14	3/4971 (0.1%)
All	All	1.00	1/14678 (0.0%)	1.13	7/19932 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	211	PRO	C-O	-5.42	1.17	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	84	ARG	CA-C-N	8.24	128.27	119.78
1	B	84	ARG	C-N-CA	8.24	128.27	119.78
1	D	84	ARG	CA-C-N	8.03	128.03	119.76
1	D	84	ARG	C-N-CA	8.03	128.03	119.76
1	C	84	ARG	CA-C-N	7.94	127.96	119.78
1	C	84	ARG	C-N-CA	7.94	127.96	119.78
1	D	147	GLU	CB-CG-CD	5.09	121.26	112.60

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	3569	0	3433	7	0
1	B	3569	0	3433	13	0
1	C	3569	0	3433	14	0
1	D	3560	0	3427	10	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	122	0	0	0	0
3	B	83	0	0	2	2
3	C	61	0	0	0	2
3	D	94	0	0	0	0
All	All	14635	0	13726	43	2

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 (43) 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:C:129:ILE:HD12	1:C:150:MET:CE	1.94	0.98
1:C:129:ILE:HD12	1:C:150:MET:HE1	1.53	0.90
1:C:145:ASN:HA	1:C:150:MET:HE3	1.63	0.81
1:D:150:MET:HA	1:D:150:MET:HE2	1.64	0.79
1:B:394:LEU:O	1:B:397:THR:HG22	1.84	0.78
1:D:134:GLU:H	1:D:134:GLU:CD	2.01	0.69
1:C:129:ILE:HG23	1:C:150:MET:HE1	1.75	0.67
1:C:137:GLN:CG	1:C:144:MET:HE1	2.24	0.66
1:C:387:ILE:HG21	1:D:88:PHE:HB2	1.78	0.64
1:B:20:ARG:O	1:B:24:GLU:HG2	1.98	0.64
1:C:137:GLN:HG3	1:C:144:MET:HE1	1.82	0.62
1:C:58:ALA:HA	1:C:340:LEU:HD11	1.84	0.60
1:B:102:GLU:HG2	3:B:627:HOH:O	2.03	0.57
1:C:137:GLN:HG2	1:C:144:MET:HE1	1.87	0.56
1:A:116:GLU:HG3	1:A:117:LYS:N	2.21	0.56
1:B:86:VAL:HG22	1:B:86:VAL:O	2.06	0.55
1:B:58:ALA:HA	1:B:340:LEU:HD11	1.88	0.55
1:D:87:ASP:OD2	1:D:89:TRP:CZ2	2.59	0.55
1:D:58:ALA:HA	1:D:340:LEU:HD11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:GLU:CG	1:A:117:LYS:N	2.71	0.53
1:A:58:ALA:HA	1:A:340:LEU:HD11	1.90	0.53
1:A:79:LYS:HE2	1:A:107:GLU:HG3	1.93	0.51
1:D:93:PRO:O	1:D:94:ASP:HB3	2.13	0.49
1:A:397:THR:O	1:A:400:LYS:HE2	2.14	0.48
1:C:345:MET:HA	1:C:345:MET:HE2	1.96	0.48
1:D:70:ILE:HG22	1:D:72:VAL:HG13	1.96	0.48
1:C:399:ASN:O	1:C:403:ILE:HD13	2.15	0.47
1:B:20:ARG:O	1:B:24:GLU:CG	2.64	0.46
1:D:147:GLU:N	1:D:148:PRO:CD	2.80	0.45
1:C:83:TYR:CZ	1:C:85:PRO:HD3	2.52	0.45
1:B:345:MET:HA	1:B:345:MET:HE2	1.99	0.44
1:D:93:PRO:O	1:D:94:ASP:CB	2.64	0.44
1:B:83:TYR:CE1	1:B:85:PRO:HD3	2.52	0.44
1:B:394:LEU:HD21	1:B:403:ILE:HD11	1.98	0.44
1:C:70:ILE:HG22	1:C:72:VAL:HG13	2.00	0.43
1:A:86:VAL:O	1:A:86:VAL:HG23	2.19	0.43
1:B:70:ILE:HG22	1:B:72:VAL:HG13	2.00	0.42
1:B:83:TYR:CZ	1:B:85:PRO:HD3	2.55	0.42
1:B:337:HIS:HD2	3:B:644:HOH:O	2.02	0.42
1:D:226:LEU:O	1:D:228:THR:HG23	2.20	0.42
1:C:164:GLU:HG2	1:C:252:TYR:OH	2.21	0.41
1:A:46:MET:HB2	1:A:46:MET:HE2	1.86	0.41
1:B:397:THR:HG23	1:B:399:ASN:H	1.86	0.41

All (2) 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)
3:B:678:HOH:O	3:C:656:HOH:O[8_555]	2.13	0.07
3:B:623:HOH:O	3:C:659:HOH:O[8_555]	2.15	0.05

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	437/448 (98%)	431 (99%)	6 (1%)	0	100	100
1	B	437/448 (98%)	432 (99%)	5 (1%)	0	100	100
1	C	437/448 (98%)	434 (99%)	3 (1%)	0	100	100
1	D	434/448 (97%)	428 (99%)	6 (1%)	0	100	100
All	All	1745/1792 (97%)	1725 (99%)	20 (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	378/387 (98%)	372 (98%)	6 (2%)	55	70
1	B	378/387 (98%)	372 (98%)	6 (2%)	55	70
1	C	378/387 (98%)	366 (97%)	12 (3%)	34	45
1	D	377/387 (97%)	369 (98%)	8 (2%)	47	61
All	All	1511/1548 (98%)	1479 (98%)	32 (2%)	47	61

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	MET
1	A	117	LYS
1	A	142	SER
1	A	172	ASN
1	A	342	VAL
1	A	367	ARG
1	B	111	GLN
1	B	116	GLU
1	B	142	SER
1	B	172	ASN

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Mol	Chain	Res	Type
1	B	301	GLN
1	B	398	ASP
1	C	15	LEU
1	C	29	GLU
1	C	86	VAL
1	C	111	GLN
1	C	172	ASN
1	C	225	TYR
1	C	226	LEU
1	C	237	THR
1	C	315	LEU
1	C	388	ASP
1	C	403	ILE
1	C	428	ASP
1	D	29	GLU
1	D	76	SER
1	D	86	VAL
1	D	97	ARG
1	D	116	GLU
1	D	134	GLU
1	D	172	ASN
1	D	342	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	296	HIS
1	A	305	GLN
1	A	363	HIS
1	A	438	HIS
1	B	55	GLN
1	B	125	ASN
1	B	188	ASN
1	B	305	GLN
1	B	337	HIS
1	C	55	GLN
1	C	239	HIS
1	C	273	GLN
1	C	337	HIS
1	D	111	GLN
1	D	114	GLN
1	D	197	HIS

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Mol	Chain	Res	Type
1	D	239	HIS
1	D	296	HIS
1	D	363	HIS

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 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.

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	439/448 (97%)	0.15	13 (2%) 52 50	37, 44, 70, 118	0
1	B	439/448 (97%)	0.34	16 (3%) 46 43	33, 48, 72, 98	0
1	C	439/448 (97%)	0.66	36 (8%) 17 15	34, 51, 76, 112	0
1	D	438/448 (97%)	0.28	20 (4%) 37 34	38, 46, 72, 131	0
All	All	1755/1792 (97%)	0.36	85 (4%) 35 32	33, 47, 73, 131	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	93	PRO	6.9
1	A	438	HIS	4.9
1	C	92	VAL	4.8
1	C	225	TYR	4.7
1	C	93	PRO	4.5
1	C	226	LEU	4.4
1	D	88	PHE	4.4
1	C	365	PHE	4.4
1	B	93	PRO	4.1
1	C	402	PHE	4.0
1	D	92	VAL	4.0
1	D	225	TYR	3.8
1	D	90	HIS	3.8
1	D	438	HIS	3.7
1	D	89	TRP	3.7
1	C	121	TYR	3.6
1	D	96	PRO	3.6
1	A	89	TRP	3.6
1	C	89	TRP	3.6
1	B	121	TYR	3.5
1	C	90	HIS	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	365	PHE	3.5
1	D	94	ASP	3.4
1	D	365	PHE	3.3
1	B	92	VAL	3.2
1	D	86	VAL	3.2
1	B	229	HIS	3.2
1	B	89	TRP	3.2
1	B	226	LEU	3.1
1	D	226	LEU	3.1
1	C	440	ASP	3.1
1	B	225	TYR	3.1
1	C	397	THR	3.1
1	D	236	HIS	3.1
1	C	398	ASP	3.0
1	B	228	THR	2.9
1	C	438	HIS	2.9
1	A	93	PRO	2.8
1	B	90	HIS	2.8
1	C	367	ARG	2.8
1	B	86	VAL	2.8
1	D	121	TYR	2.8
1	C	86	VAL	2.8
1	A	90	HIS	2.8
1	D	229	HIS	2.7
1	A	229	HIS	2.7
1	D	91	LYS	2.7
1	C	399	ASN	2.7
1	C	116	GLU	2.7
1	C	310	PHE	2.6
1	C	229	HIS	2.6
1	C	394	LEU	2.6
1	A	94	ASP	2.6
1	A	348	PHE	2.6
1	A	121	TYR	2.6
1	B	236	HIS	2.5
1	D	228	THR	2.5
1	D	87	ASP	2.5
1	A	86	VAL	2.4
1	A	92	VAL	2.4
1	D	348	PHE	2.4
1	A	225	TYR	2.4
1	C	362	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	122	ASP	2.4
1	D	45	TRP	2.3
1	C	354	GLY	2.3
1	B	348	PHE	2.3
1	C	312	ILE	2.3
1	C	395	ALA	2.2
1	C	355	THR	2.2
1	A	365	PHE	2.2
1	C	348	PHE	2.2
1	C	314	LYS	2.2
1	A	226	LEU	2.2
1	C	360	PRO	2.2
1	C	321	VAL	2.1
1	C	237	THR	2.1
1	C	30	GLY	2.1
1	B	428	ASP	2.1
1	C	368	CYS	2.1
1	B	2	ASP	2.1
1	C	224	ILE	2.1
1	C	364	PRO	2.1
1	B	438	HIS	2.0
1	C	319	ASP	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(Å ²)	Q<0.9
2	MN	A	502	1/1	0.95	0.08	71,71,71,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MN	C	502	1/1	0.95	0.08	85,85,85,85	0
2	MN	C	501	1/1	0.96	0.08	85,85,85,85	0
2	MN	D	502	1/1	0.96	0.07	67,67,67,67	0
2	MN	B	502	1/1	0.97	0.04	71,71,71,71	0
2	MN	D	501	1/1	0.97	0.06	63,63,63,63	0
2	MN	B	501	1/1	0.97	0.07	68,68,68,68	0
2	MN	A	501	1/1	0.98	0.05	71,71,71,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.