



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:22 PM UTC

PDB ID : 1RYY / pdb_00001ryy
Title : Acetobacter turbidans alpha-amino acid ester hydrolase Y206A mutant
Authors : Barends, T.R.M.
Deposited on : 2003-12-23
Resolution : 2.80 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

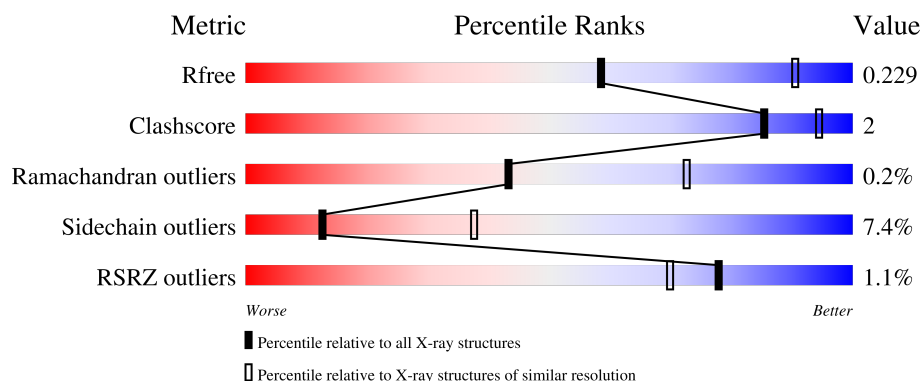
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	652	
1	B	652	
1	C	652	
1	D	652	
1	E	652	

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Mol	Chain	Length	Quality of chain
1	F	652	2% 81% 12% 6%
1	G	652	% 81% 11% 6%
1	H	652	% 80% 13% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 39009 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 alpha-amino acid ester hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			
1	B	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			
1	C	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			
1	D	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			
1	E	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			
1	F	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			
1	G	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			
1	H	612	Total	C	N	O	S	0	0	0
			4836	3081	838	898	19			

There are 208 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	206	ALA	TYR	engineered mutation	UNP Q8VRK8
A	668	LYS	-	expression tag	UNP Q8VRK8
A	669	LEU	-	expression tag	UNP Q8VRK8
A	670	GLY	-	expression tag	UNP Q8VRK8
A	671	PRO	-	expression tag	UNP Q8VRK8
A	672	GLU	-	expression tag	UNP Q8VRK8
A	673	GLN	-	expression tag	UNP Q8VRK8
A	674	LYS	-	expression tag	UNP Q8VRK8
A	675	LEU	-	expression tag	UNP Q8VRK8
A	676	ILE	-	expression tag	UNP Q8VRK8
A	677	SER	-	expression tag	UNP Q8VRK8
A	678	GLU	-	expression tag	UNP Q8VRK8
A	679	GLU	-	expression tag	UNP Q8VRK8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	680	ASP	-	expression tag	UNP Q8VRK8
A	681	LEU	-	expression tag	UNP Q8VRK8
A	682	ASN	-	expression tag	UNP Q8VRK8
A	683	SER	-	expression tag	UNP Q8VRK8
A	684	ALA	-	expression tag	UNP Q8VRK8
A	685	VAL	-	expression tag	UNP Q8VRK8
A	686	ASP	-	expression tag	UNP Q8VRK8
A	687	HIS	-	expression tag	UNP Q8VRK8
A	688	HIS	-	expression tag	UNP Q8VRK8
A	689	HIS	-	expression tag	UNP Q8VRK8
A	690	HIS	-	expression tag	UNP Q8VRK8
A	691	HIS	-	expression tag	UNP Q8VRK8
A	692	HIS	-	expression tag	UNP Q8VRK8
B	206	ALA	TYR	engineered mutation	UNP Q8VRK8
B	668	LYS	-	expression tag	UNP Q8VRK8
B	669	LEU	-	expression tag	UNP Q8VRK8
B	670	GLY	-	expression tag	UNP Q8VRK8
B	671	PRO	-	expression tag	UNP Q8VRK8
B	672	GLU	-	expression tag	UNP Q8VRK8
B	673	GLN	-	expression tag	UNP Q8VRK8
B	674	LYS	-	expression tag	UNP Q8VRK8
B	675	LEU	-	expression tag	UNP Q8VRK8
B	676	ILE	-	expression tag	UNP Q8VRK8
B	677	SER	-	expression tag	UNP Q8VRK8
B	678	GLU	-	expression tag	UNP Q8VRK8
B	679	GLU	-	expression tag	UNP Q8VRK8
B	680	ASP	-	expression tag	UNP Q8VRK8
B	681	LEU	-	expression tag	UNP Q8VRK8
B	682	ASN	-	expression tag	UNP Q8VRK8
B	683	SER	-	expression tag	UNP Q8VRK8
B	684	ALA	-	expression tag	UNP Q8VRK8
B	685	VAL	-	expression tag	UNP Q8VRK8
B	686	ASP	-	expression tag	UNP Q8VRK8
B	687	HIS	-	expression tag	UNP Q8VRK8
B	688	HIS	-	expression tag	UNP Q8VRK8
B	689	HIS	-	expression tag	UNP Q8VRK8
B	690	HIS	-	expression tag	UNP Q8VRK8
B	691	HIS	-	expression tag	UNP Q8VRK8
B	692	HIS	-	expression tag	UNP Q8VRK8
C	206	ALA	TYR	engineered mutation	UNP Q8VRK8
C	668	LYS	-	expression tag	UNP Q8VRK8
C	669	LEU	-	expression tag	UNP Q8VRK8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	670	GLY	-	expression tag	UNP Q8VRK8
C	671	PRO	-	expression tag	UNP Q8VRK8
C	672	GLU	-	expression tag	UNP Q8VRK8
C	673	GLN	-	expression tag	UNP Q8VRK8
C	674	LYS	-	expression tag	UNP Q8VRK8
C	675	LEU	-	expression tag	UNP Q8VRK8
C	676	ILE	-	expression tag	UNP Q8VRK8
C	677	SER	-	expression tag	UNP Q8VRK8
C	678	GLU	-	expression tag	UNP Q8VRK8
C	679	GLU	-	expression tag	UNP Q8VRK8
C	680	ASP	-	expression tag	UNP Q8VRK8
C	681	LEU	-	expression tag	UNP Q8VRK8
C	682	ASN	-	expression tag	UNP Q8VRK8
C	683	SER	-	expression tag	UNP Q8VRK8
C	684	ALA	-	expression tag	UNP Q8VRK8
C	685	VAL	-	expression tag	UNP Q8VRK8
C	686	ASP	-	expression tag	UNP Q8VRK8
C	687	HIS	-	expression tag	UNP Q8VRK8
C	688	HIS	-	expression tag	UNP Q8VRK8
C	689	HIS	-	expression tag	UNP Q8VRK8
C	690	HIS	-	expression tag	UNP Q8VRK8
C	691	HIS	-	expression tag	UNP Q8VRK8
C	692	HIS	-	expression tag	UNP Q8VRK8
D	206	ALA	TYR	engineered mutation	UNP Q8VRK8
D	668	LYS	-	expression tag	UNP Q8VRK8
D	669	LEU	-	expression tag	UNP Q8VRK8
D	670	GLY	-	expression tag	UNP Q8VRK8
D	671	PRO	-	expression tag	UNP Q8VRK8
D	672	GLU	-	expression tag	UNP Q8VRK8
D	673	GLN	-	expression tag	UNP Q8VRK8
D	674	LYS	-	expression tag	UNP Q8VRK8
D	675	LEU	-	expression tag	UNP Q8VRK8
D	676	ILE	-	expression tag	UNP Q8VRK8
D	677	SER	-	expression tag	UNP Q8VRK8
D	678	GLU	-	expression tag	UNP Q8VRK8
D	679	GLU	-	expression tag	UNP Q8VRK8
D	680	ASP	-	expression tag	UNP Q8VRK8
D	681	LEU	-	expression tag	UNP Q8VRK8
D	682	ASN	-	expression tag	UNP Q8VRK8
D	683	SER	-	expression tag	UNP Q8VRK8
D	684	ALA	-	expression tag	UNP Q8VRK8
D	685	VAL	-	expression tag	UNP Q8VRK8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	686	ASP	-	expression tag	UNP Q8VRK8
D	687	HIS	-	expression tag	UNP Q8VRK8
D	688	HIS	-	expression tag	UNP Q8VRK8
D	689	HIS	-	expression tag	UNP Q8VRK8
D	690	HIS	-	expression tag	UNP Q8VRK8
D	691	HIS	-	expression tag	UNP Q8VRK8
D	692	HIS	-	expression tag	UNP Q8VRK8
E	206	ALA	TYR	engineered mutation	UNP Q8VRK8
E	668	LYS	-	expression tag	UNP Q8VRK8
E	669	LEU	-	expression tag	UNP Q8VRK8
E	670	GLY	-	expression tag	UNP Q8VRK8
E	671	PRO	-	expression tag	UNP Q8VRK8
E	672	GLU	-	expression tag	UNP Q8VRK8
E	673	GLN	-	expression tag	UNP Q8VRK8
E	674	LYS	-	expression tag	UNP Q8VRK8
E	675	LEU	-	expression tag	UNP Q8VRK8
E	676	ILE	-	expression tag	UNP Q8VRK8
E	677	SER	-	expression tag	UNP Q8VRK8
E	678	GLU	-	expression tag	UNP Q8VRK8
E	679	GLU	-	expression tag	UNP Q8VRK8
E	680	ASP	-	expression tag	UNP Q8VRK8
E	681	LEU	-	expression tag	UNP Q8VRK8
E	682	ASN	-	expression tag	UNP Q8VRK8
E	683	SER	-	expression tag	UNP Q8VRK8
E	684	ALA	-	expression tag	UNP Q8VRK8
E	685	VAL	-	expression tag	UNP Q8VRK8
E	686	ASP	-	expression tag	UNP Q8VRK8
E	687	HIS	-	expression tag	UNP Q8VRK8
E	688	HIS	-	expression tag	UNP Q8VRK8
E	689	HIS	-	expression tag	UNP Q8VRK8
E	690	HIS	-	expression tag	UNP Q8VRK8
E	691	HIS	-	expression tag	UNP Q8VRK8
E	692	HIS	-	expression tag	UNP Q8VRK8
F	206	ALA	TYR	engineered mutation	UNP Q8VRK8
F	668	LYS	-	expression tag	UNP Q8VRK8
F	669	LEU	-	expression tag	UNP Q8VRK8
F	670	GLY	-	expression tag	UNP Q8VRK8
F	671	PRO	-	expression tag	UNP Q8VRK8
F	672	GLU	-	expression tag	UNP Q8VRK8
F	673	GLN	-	expression tag	UNP Q8VRK8
F	674	LYS	-	expression tag	UNP Q8VRK8
F	675	LEU	-	expression tag	UNP Q8VRK8

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Chain	Residue	Modelled	Actual	Comment	Reference
F	676	ILE	-	expression tag	UNP Q8VRK8
F	677	SER	-	expression tag	UNP Q8VRK8
F	678	GLU	-	expression tag	UNP Q8VRK8
F	679	GLU	-	expression tag	UNP Q8VRK8
F	680	ASP	-	expression tag	UNP Q8VRK8
F	681	LEU	-	expression tag	UNP Q8VRK8
F	682	ASN	-	expression tag	UNP Q8VRK8
F	683	SER	-	expression tag	UNP Q8VRK8
F	684	ALA	-	expression tag	UNP Q8VRK8
F	685	VAL	-	expression tag	UNP Q8VRK8
F	686	ASP	-	expression tag	UNP Q8VRK8
F	687	HIS	-	expression tag	UNP Q8VRK8
F	688	HIS	-	expression tag	UNP Q8VRK8
F	689	HIS	-	expression tag	UNP Q8VRK8
F	690	HIS	-	expression tag	UNP Q8VRK8
F	691	HIS	-	expression tag	UNP Q8VRK8
F	692	HIS	-	expression tag	UNP Q8VRK8
G	206	ALA	TYR	engineered mutation	UNP Q8VRK8
G	668	LYS	-	expression tag	UNP Q8VRK8
G	669	LEU	-	expression tag	UNP Q8VRK8
G	670	GLY	-	expression tag	UNP Q8VRK8
G	671	PRO	-	expression tag	UNP Q8VRK8
G	672	GLU	-	expression tag	UNP Q8VRK8
G	673	GLN	-	expression tag	UNP Q8VRK8
G	674	LYS	-	expression tag	UNP Q8VRK8
G	675	LEU	-	expression tag	UNP Q8VRK8
G	676	ILE	-	expression tag	UNP Q8VRK8
G	677	SER	-	expression tag	UNP Q8VRK8
G	678	GLU	-	expression tag	UNP Q8VRK8
G	679	GLU	-	expression tag	UNP Q8VRK8
G	680	ASP	-	expression tag	UNP Q8VRK8
G	681	LEU	-	expression tag	UNP Q8VRK8
G	682	ASN	-	expression tag	UNP Q8VRK8
G	683	SER	-	expression tag	UNP Q8VRK8
G	684	ALA	-	expression tag	UNP Q8VRK8
G	685	VAL	-	expression tag	UNP Q8VRK8
G	686	ASP	-	expression tag	UNP Q8VRK8
G	687	HIS	-	expression tag	UNP Q8VRK8
G	688	HIS	-	expression tag	UNP Q8VRK8
G	689	HIS	-	expression tag	UNP Q8VRK8
G	690	HIS	-	expression tag	UNP Q8VRK8
G	691	HIS	-	expression tag	UNP Q8VRK8

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Chain	Residue	Modelled	Actual	Comment	Reference
G	692	HIS	-	expression tag	UNP Q8VRK8
H	206	ALA	TYR	engineered mutation	UNP Q8VRK8
H	668	LYS	-	expression tag	UNP Q8VRK8
H	669	LEU	-	expression tag	UNP Q8VRK8
H	670	GLY	-	expression tag	UNP Q8VRK8
H	671	PRO	-	expression tag	UNP Q8VRK8
H	672	GLU	-	expression tag	UNP Q8VRK8
H	673	GLN	-	expression tag	UNP Q8VRK8
H	674	LYS	-	expression tag	UNP Q8VRK8
H	675	LEU	-	expression tag	UNP Q8VRK8
H	676	ILE	-	expression tag	UNP Q8VRK8
H	677	SER	-	expression tag	UNP Q8VRK8
H	678	GLU	-	expression tag	UNP Q8VRK8
H	679	GLU	-	expression tag	UNP Q8VRK8
H	680	ASP	-	expression tag	UNP Q8VRK8
H	681	LEU	-	expression tag	UNP Q8VRK8
H	682	ASN	-	expression tag	UNP Q8VRK8
H	683	SER	-	expression tag	UNP Q8VRK8
H	684	ALA	-	expression tag	UNP Q8VRK8
H	685	VAL	-	expression tag	UNP Q8VRK8
H	686	ASP	-	expression tag	UNP Q8VRK8
H	687	HIS	-	expression tag	UNP Q8VRK8
H	688	HIS	-	expression tag	UNP Q8VRK8
H	689	HIS	-	expression tag	UNP Q8VRK8
H	690	HIS	-	expression tag	UNP Q8VRK8
H	691	HIS	-	expression tag	UNP Q8VRK8
H	692	HIS	-	expression tag	UNP Q8VRK8

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	66	Total O 66 66	0	0
2	B	56	Total O 56 56	0	0
2	C	43	Total O 43 43	0	0
2	D	55	Total O 55 55	0	0
2	E	15	Total O 15 15	0	0
2	F	22	Total O 22 22	0	0

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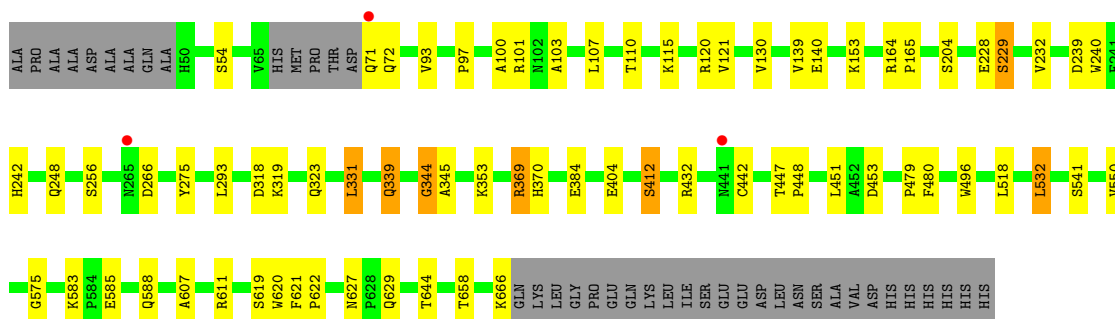
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	32	Total 32	O 32	0	0
2	H	32	Total 32	O 32	0	0

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.

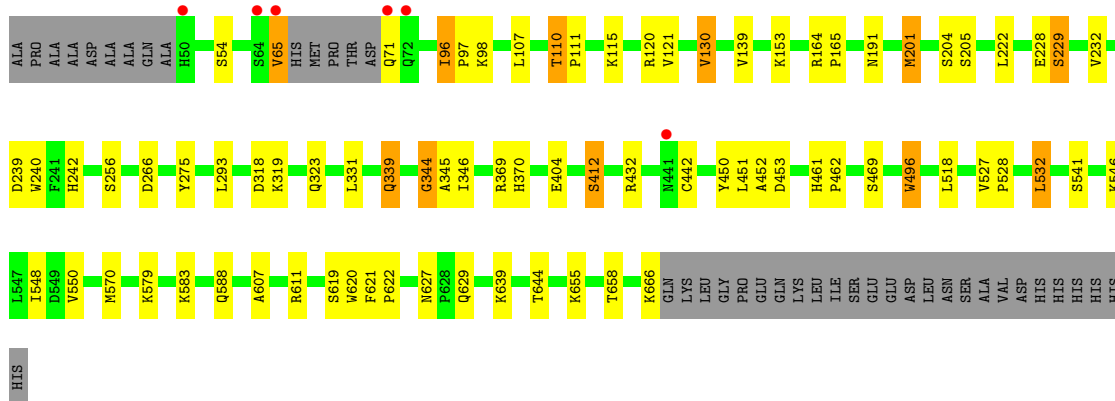
- Molecule 1: alpha-amino acid ester hydrolase

Chain A: 



- Molecule 1: alpha-amino acid ester hydrolase

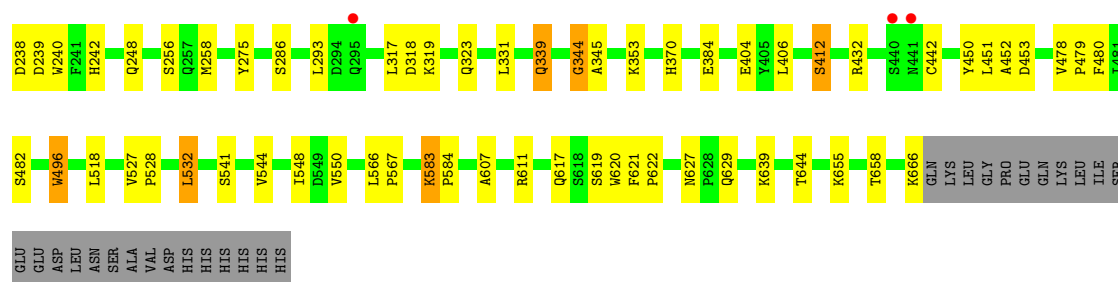
Chain B: 



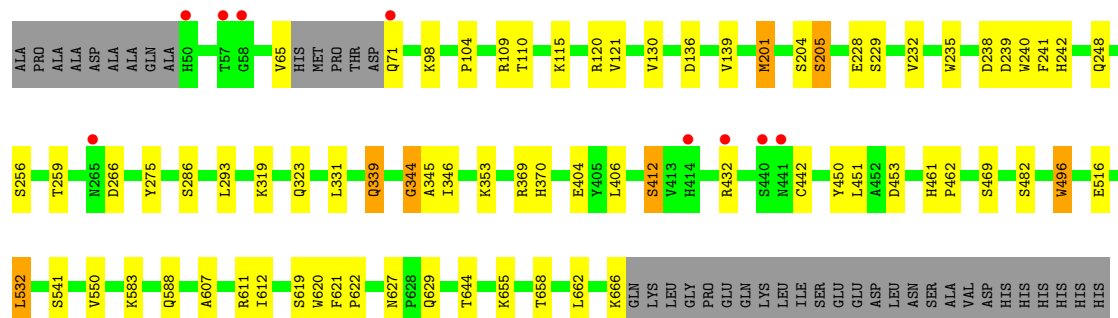
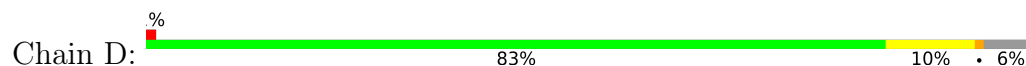
- Molecule 1: alpha-amino acid ester hydrolase

Chain C: 

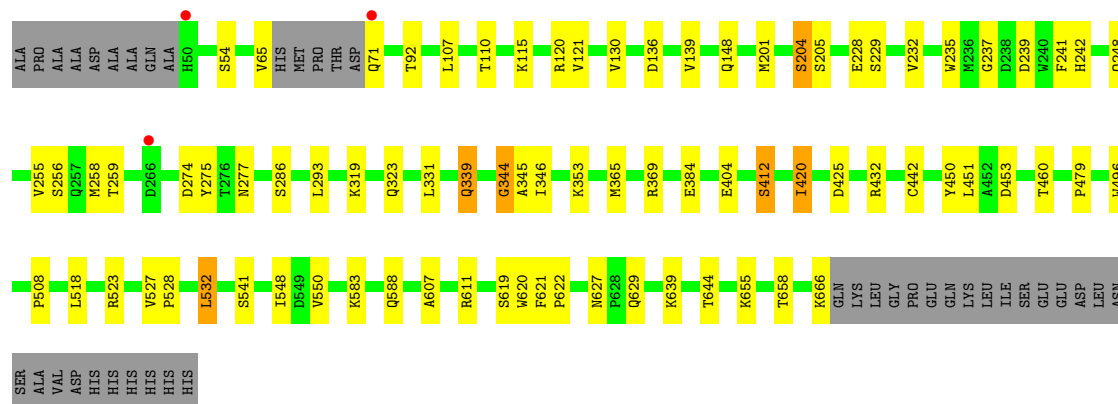
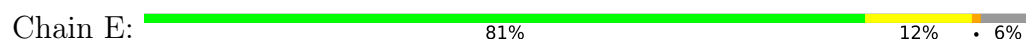




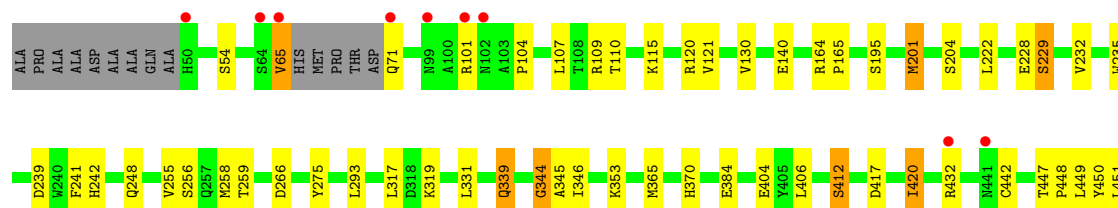
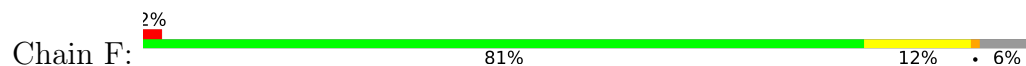
- Molecule 1: alpha-amino acid ester hydrolase



- Molecule 1: alpha-amino acid ester hydrolase

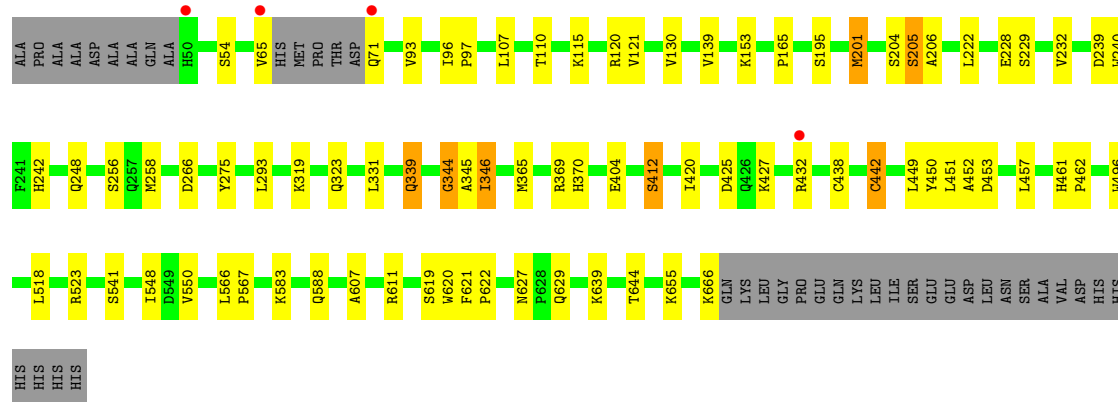
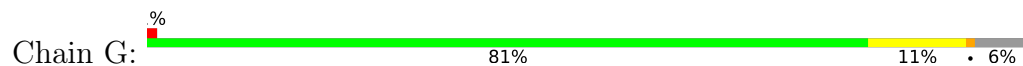


- Molecule 1: alpha-amino acid ester hydrolase

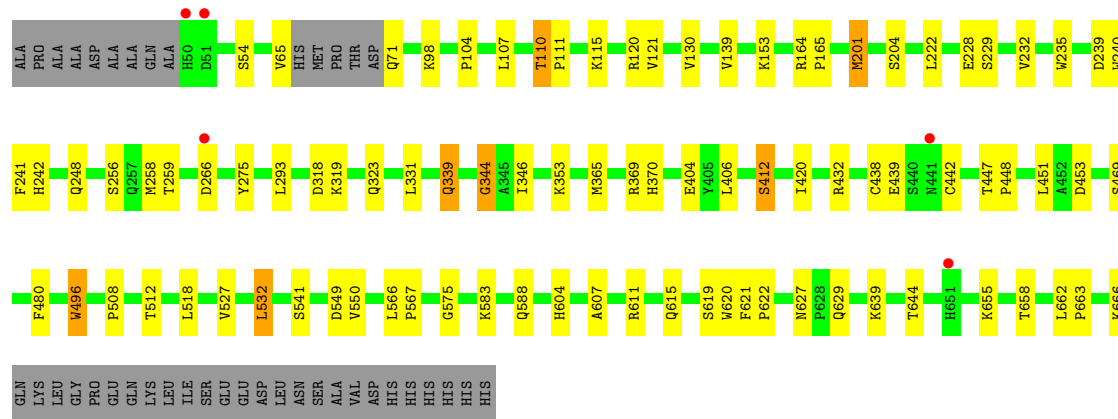
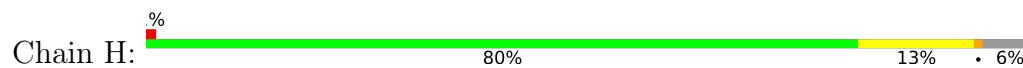




- Molecule 1: alpha-amino acid ester hydrolase



- Molecule 1: alpha-amino acid ester hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.65Å 177.51Å 169.97Å 90.00° 91.03° 90.00°	Depositor
Resolution (Å)	15.00 – 2.80 15.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (15.00-2.80) 98.8 (15.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.60 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.210 , 0.236 0.204 , 0.229	Depositor DCC
R_{free} test set	6623 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.9	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 15.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for -h,-l,-k 0.000 for -h,l,k 0.015 for h,-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	39009	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.80% 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](#)

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	0.31	0/4991	1.01	15/6815 (0.2%)
1	B	0.30	0/4991	1.01	20/6815 (0.3%)
1	C	0.30	0/4991	1.01	21/6815 (0.3%)
1	D	0.30	0/4991	1.01	22/6815 (0.3%)
1	E	0.29	0/4991	0.97	16/6815 (0.2%)
1	F	0.29	0/4991	0.98	14/6815 (0.2%)
1	G	0.29	0/4991	0.99	18/6815 (0.3%)
1	H	0.29	0/4991	0.98	17/6815 (0.2%)
All	All	0.30	0/39928	0.99	143/54520 (0.3%)

There are no bond length outliers.

All (143) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	121	VAL	N-CA-C	-11.44	96.82	108.15
1	A	121	VAL	N-CA-C	-11.06	97.20	108.15
1	D	121	VAL	N-CA-C	-10.90	97.36	108.15
1	C	121	VAL	N-CA-C	-10.81	97.44	108.15
1	D	239	ASP	N-CA-C	10.64	122.63	111.14
1	A	239	ASP	N-CA-C	10.02	121.97	111.14
1	G	121	VAL	N-CA-C	-9.91	98.34	108.15
1	E	121	VAL	N-CA-C	-9.42	98.82	108.15
1	H	121	VAL	N-CA-C	-9.40	98.85	108.15
1	B	239	ASP	N-CA-C	9.04	121.22	111.36
1	C	239	ASP	N-CA-C	8.82	120.67	111.14
1	F	121	VAL	N-CA-C	-8.74	99.49	108.15
1	F	239	ASP	N-CA-C	8.62	122.30	111.69
1	H	110	THR	N-CA-C	7.89	121.02	108.23
1	C	110	THR	N-CA-C	7.79	120.85	108.23
1	A	627	ASN	N-CA-C	-7.70	99.94	109.65
1	D	238	ASP	N-CA-C	7.68	115.08	108.78
1	D	110	THR	N-CA-C	7.64	121.72	108.82
1	E	239	ASP	N-CA-C	7.60	121.04	111.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	110	THR	N-CA-C	7.52	121.53	108.82
1	B	110	THR	N-CA-C	7.45	120.30	108.23
1	E	110	THR	N-CA-C	7.42	120.25	108.23
1	G	239	ASP	N-CA-C	7.35	120.73	111.69
1	D	275	TYR	N-CA-C	-7.30	103.25	111.14
1	G	275	TYR	N-CA-C	-7.30	103.40	111.36
1	H	239	ASP	N-CA-C	7.24	120.23	111.40
1	B	275	TYR	N-CA-C	-6.98	103.75	111.36
1	B	191	ASN	N-CA-C	6.85	121.80	113.17
1	F	275	TYR	N-CA-C	-6.82	103.92	111.36
1	B	627	ASN	N-CA-C	-6.71	101.34	109.72
1	F	345	ALA	N-CA-C	6.66	119.15	111.02
1	D	451	LEU	N-CA-C	-6.64	99.70	110.20
1	A	575	GLY	N-CA-C	6.48	121.89	113.27
1	C	275	TYR	N-CA-C	-6.46	104.32	111.36
1	A	240	TRP	N-CA-C	6.44	119.26	111.40
1	A	248	GLN	N-CA-C	6.27	120.39	112.87
1	A	103	ALA	N-CA-C	6.26	117.69	109.93
1	C	240	TRP	N-CA-C	6.23	119.00	111.40
1	A	275	TYR	N-CA-C	-6.21	104.60	111.36
1	H	275	TYR	N-CA-C	-6.13	104.68	111.36
1	D	345	ALA	N-CA-C	6.11	117.63	110.97
1	E	275	TYR	N-CA-C	-6.11	104.70	111.36
1	C	607	ALA	N-CA-C	6.09	117.74	110.19
1	A	607	ALA	N-CA-C	6.06	117.71	110.19
1	C	345	ALA	N-CA-C	6.04	118.36	111.11
1	A	532	LEU	N-CA-C	6.04	119.08	109.24
1	H	240	TRP	N-CA-C	6.01	118.74	111.40
1	C	238	ASP	N-CA-C	6.00	113.70	108.78
1	D	607	ALA	N-CA-C	6.00	117.63	110.19
1	H	139	VAL	N-CA-C	-6.00	104.46	111.00
1	C	139	VAL	N-CA-C	-5.98	104.68	110.72
1	G	240	TRP	N-CA-C	5.98	118.69	111.40
1	E	248	GLN	N-CA-C	5.96	120.02	112.87
1	E	607	ALA	N-CA-C	5.92	117.96	110.33
1	B	451	LEU	N-CA-C	-5.89	100.89	110.20
1	B	345	ALA	N-CA-C	5.89	117.39	110.97
1	D	248	GLN	N-CA-C	5.87	119.92	112.87
1	F	627	ASN	N-CA-C	-5.86	100.75	109.42
1	C	532	LEU	N-CA-C	5.84	118.76	109.24
1	B	607	ALA	N-CA-C	5.83	117.42	110.19
1	G	139	VAL	N-CA-C	-5.82	104.84	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	248	GLN	N-CA-C	5.82	119.85	112.87
1	D	532	LEU	N-CA-C	5.81	118.70	109.24
1	G	627	ASN	N-CA-C	-5.79	102.35	109.65
1	H	532	LEU	N-CA-C	5.79	118.67	109.24
1	H	627	ASN	N-CA-C	-5.77	101.22	109.24
1	H	607	ALA	N-CA-C	5.77	117.34	110.19
1	A	139	VAL	N-CA-C	-5.73	104.75	111.00
1	C	627	ASN	N-CA-C	-5.72	101.29	109.24
1	E	139	VAL	N-CA-C	-5.70	104.79	111.00
1	F	248	GLN	N-CA-C	5.66	119.66	112.87
1	E	345	ALA	N-CA-C	5.65	117.92	111.02
1	D	136	ASP	N-CA-C	5.61	117.85	111.11
1	E	258	MET	N-CA-C	5.61	119.43	112.59
1	F	451	LEU	N-CA-C	-5.60	101.35	110.20
1	C	518	LEU	N-CA-C	5.59	118.44	110.10
1	B	450	TYR	N-CA-C	5.56	118.51	110.28
1	H	518	LEU	N-CA-C	5.55	118.37	110.10
1	C	482	SER	N-CA-C	5.54	117.44	110.24
1	G	346	ILE	N-CA-C	5.52	115.72	110.42
1	A	518	LEU	N-CA-C	5.52	118.32	110.10
1	D	627	ASN	N-CA-C	-5.51	101.58	109.24
1	H	575	GLY	N-CA-C	5.51	120.60	113.27
1	G	518	LEU	N-CA-C	5.51	118.31	110.10
1	G	548	ILE	N-CA-C	5.48	116.38	108.48
1	G	248	GLN	N-CA-C	5.48	119.45	112.87
1	B	130	VAL	N-CA-C	5.46	117.87	111.05
1	A	345	ALA	N-CA-C	5.45	117.67	111.02
1	D	240	TRP	N-CA-C	5.43	119.86	112.04
1	D	516	GLU	N-CA-C	-5.41	103.55	110.53
1	A	451	LEU	N-CA-C	-5.40	100.89	109.96
1	G	345	ALA	N-CA-C	5.40	117.61	111.02
1	H	248	GLN	N-CA-C	5.38	119.33	112.87
1	B	452	ALA	N-CA-C	5.38	116.82	109.18
1	H	258	MET	N-CA-C	5.38	119.81	112.88
1	D	139	VAL	N-CA-C	-5.37	105.30	110.72
1	C	451	LEU	N-CA-C	-5.35	101.75	110.20
1	B	139	VAL	N-CA-C	-5.35	105.17	111.00
1	G	258	MET	N-CA-C	5.33	119.34	112.41
1	C	258	MET	N-CA-C	5.32	120.44	112.94
1	D	482	SER	N-CA-C	5.30	117.77	110.35
1	H	527	VAL	N-CA-C	5.29	112.79	107.55
1	D	662	LEU	N-CA-C	5.28	116.48	109.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	451	LEU	N-CA-C	-5.28	101.86	110.20
1	F	452	ALA	N-CA-C	5.27	115.24	108.34
1	B	548	ILE	N-CA-C	5.26	116.95	108.85
1	F	482	SER	N-CA-C	5.26	117.65	110.24
1	D	205	SER	CB-CA-C	-5.25	110.10	117.23
1	E	451	LEU	N-CA-C	-5.21	101.96	110.20
1	B	240	TRP	N-CA-C	5.21	119.54	112.04
1	E	532	LEU	N-CA-C	5.20	117.72	109.24
1	B	496	TRP	N-CA-C	5.20	116.95	111.28
1	E	136	ASP	N-CA-C	5.20	117.35	111.11
1	D	450	TYR	N-CA-C	5.19	117.76	110.23
1	G	110	THR	N-CA-C	5.19	121.47	108.40
1	E	548	ILE	N-CA-C	5.18	116.76	108.89
1	D	612	ILE	N-CA-C	-5.17	101.00	108.71
1	H	439	GLU	N-CA-C	5.17	116.60	111.07
1	G	607	ALA	N-CA-C	5.16	116.76	110.41
1	F	258	MET	N-CA-C	5.15	119.11	112.41
1	B	111	PRO	N-CA-C	-5.15	108.09	114.68
1	E	627	ASN	N-CA-C	-5.14	102.09	109.24
1	F	532	LEU	N-CA-C	5.14	117.62	109.24
1	C	452	ALA	N-CA-C	5.14	116.47	109.18
1	B	518	LEU	N-CA-C	5.13	117.75	110.10
1	F	110	THR	N-CA-C	5.12	121.29	108.40
1	F	317	LEU	N-CA-C	5.11	118.61	112.38
1	G	450	TYR	N-CA-C	5.11	117.84	110.28
1	B	532	LEU	N-CA-C	5.10	117.55	109.24
1	D	469	SER	N-CA-C	5.10	117.88	109.72
1	B	96	ILE	N-CA-C	5.09	113.44	107.89
1	C	496	TRP	N-CA-C	5.08	117.56	111.71
1	C	317	LEU	N-CA-C	5.08	117.55	111.71
1	D	496	TRP	N-CA-C	5.06	116.80	111.28
1	H	451	LEU	N-CA-C	-5.06	101.45	109.96
1	F	450	TYR	N-CA-C	5.05	117.75	110.28
1	H	496	TRP	N-CA-C	5.04	116.78	111.28
1	G	452	ALA	N-CA-C	5.04	116.33	109.18
1	C	548	ILE	N-CA-C	5.03	116.60	108.85
1	E	518	LEU	N-CA-C	5.03	117.59	110.10
1	G	165	PRO	O-C-N	5.02	123.62	121.31
1	E	450	TYR	N-CA-C	5.00	117.69	110.28
1	C	450	TYR	N-CA-C	5.00	117.68	110.28

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	4836	0	4609	21	0
1	B	4836	0	4609	22	0
1	C	4836	0	4609	22	0
1	D	4836	0	4609	16	0
1	E	4836	0	4609	18	0
1	F	4836	0	4609	23	0
1	G	4836	0	4609	19	0
1	H	4836	0	4609	28	0
2	A	66	0	0	0	0
2	B	56	0	0	0	0
2	C	43	0	0	0	0
2	D	55	0	0	0	0
2	E	15	0	0	0	0
2	F	22	0	0	0	0
2	G	32	0	0	0	0
2	H	32	0	0	0	0
All	All	39009	0	36872	166	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 2.

All (166) 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:G:550:VAL:HB	1:G:611:ARG:HB2	1.68	0.75
1:C:550:VAL:HB	1:C:611:ARG:HB2	1.69	0.74
1:C:339:GLN:HE21	1:C:339:GLN:H	1.38	0.71
1:A:550:VAL:HB	1:A:611:ARG:HB2	1.73	0.70
1:H:550:VAL:HB	1:H:611:ARG:HB2	1.75	0.69
1:D:550:VAL:HB	1:D:611:ARG:HB2	1.75	0.67
1:B:550:VAL:HB	1:B:611:ARG:HB2	1.76	0.66
1:E:550:VAL:HB	1:E:611:ARG:HB2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:339:GLN:HE22	1:C:370:HIS:H	1.47	0.61
1:F:550:VAL:HB	1:F:611:ARG:HB2	1.82	0.61
1:A:339:GLN:HE22	1:A:370:HIS:H	1.46	0.61
1:A:97:PRO:HG2	1:A:100:ALA:HB2	1.84	0.59
1:A:339:GLN:HE21	1:A:339:GLN:H	1.48	0.59
1:F:339:GLN:HE22	1:F:370:HIS:H	1.51	0.58
1:C:404:GLU:HG3	1:C:412:SER:HA	1.86	0.57
1:D:339:GLN:HE22	1:D:370:HIS:H	1.52	0.56
1:G:339:GLN:HE22	1:G:370:HIS:H	1.53	0.56
1:G:438:CYS:HG	1:G:442:CYS:HG	1.52	0.56
1:F:339:GLN:HE21	1:F:339:GLN:H	1.54	0.55
1:C:339:GLN:H	1:C:339:GLN:NE2	2.04	0.54
1:H:404:GLU:HG3	1:H:412:SER:HA	1.90	0.54
1:F:255:VAL:O	1:F:259:THR:OG1	2.21	0.54
1:D:339:GLN:NE2	1:D:339:GLN:H	2.06	0.53
1:D:204:SER:HA	1:D:228:GLU:O	2.09	0.53
1:D:404:GLU:HG3	1:D:412:SER:HA	1.91	0.53
1:B:65:VAL:HG11	1:D:259:THR:HG23	1.92	0.52
1:C:232:VAL:HB	1:C:344:GLY:HA2	1.90	0.52
1:A:339:GLN:NE2	1:A:370:HIS:H	2.08	0.52
1:E:404:GLU:HG3	1:E:412:SER:HA	1.90	0.51
1:A:232:VAL:HB	1:A:344:GLY:HA2	1.93	0.51
1:H:339:GLN:HE22	1:H:370:HIS:H	1.58	0.51
1:H:232:VAL:HB	1:H:344:GLY:HA2	1.93	0.50
1:C:339:GLN:NE2	1:C:370:HIS:H	2.09	0.50
1:C:532:LEU:HA	1:C:658:THR:HG23	1.94	0.50
1:G:404:GLU:HG3	1:G:412:SER:HA	1.93	0.49
1:H:532:LEU:HA	1:H:658:THR:HG23	1.93	0.49
1:A:339:GLN:H	1:A:339:GLN:NE2	2.11	0.49
1:E:255:VAL:O	1:E:259:THR:OG1	2.27	0.49
1:B:404:GLU:HG3	1:B:412:SER:HA	1.94	0.49
1:B:232:VAL:HB	1:B:344:GLY:HA2	1.94	0.49
1:B:339:GLN:HE22	1:B:370:HIS:H	1.60	0.48
1:E:204:SER:HA	1:E:228:GLU:O	2.12	0.48
1:D:232:VAL:HB	1:D:344:GLY:HA2	1.95	0.48
1:F:65:VAL:HG11	1:H:259:THR:HG23	1.96	0.48
1:G:201:MET:HB2	1:G:222:LEU:HD11	1.95	0.47
1:F:339:GLN:NE2	1:F:370:HIS:H	2.12	0.47
1:A:621:PHE:CG	1:A:622:PRO:HA	2.49	0.47
1:G:425:ASP:OD2	1:G:523:ARG:NH1	2.43	0.47
1:D:339:GLN:H	1:D:339:GLN:HE21	1.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:339:GLN:HE22	1:E:369:ARG:HB2	1.79	0.47
1:E:532:LEU:HA	1:E:658:THR:HG23	1.96	0.47
1:A:318:ASP:OD1	1:A:318:ASP:N	2.47	0.47
1:F:339:GLN:H	1:F:339:GLN:NE2	2.13	0.47
1:H:339:GLN:H	1:H:339:GLN:HE21	1.62	0.46
1:H:621:PHE:CG	1:H:622:PRO:HA	2.50	0.46
1:F:532:LEU:HA	1:F:658:THR:HG23	1.96	0.46
1:D:461:HIS:HA	1:D:462:PRO:HD3	1.78	0.46
1:E:508:PRO:HG2	1:H:508:PRO:HG2	1.98	0.46
1:B:318:ASP:N	1:B:318:ASP:OD1	2.48	0.46
1:C:104:PRO:HB3	1:C:406:LEU:HB3	1.98	0.45
1:H:153:LYS:HA	1:H:153:LYS:HD2	1.77	0.45
1:D:369:ARG:H	1:D:369:ARG:HG3	1.60	0.45
1:F:235:TRP:HA	1:F:241:PHE:HB2	1.98	0.45
1:B:369:ARG:H	1:B:369:ARG:HG3	1.60	0.45
1:B:546:LYS:HA	1:B:570:MET:HB3	1.99	0.45
1:F:164:ARG:HA	1:F:165:PRO:HD3	1.84	0.45
1:G:339:GLN:HE21	1:G:339:GLN:H	1.64	0.45
1:A:532:LEU:HA	1:A:658:THR:HG23	1.98	0.44
1:B:339:GLN:HE21	1:B:339:GLN:H	1.65	0.44
1:E:235:TRP:HA	1:E:241:PHE:HB2	1.98	0.44
1:H:164:ARG:HA	1:H:165:PRO:HD3	1.84	0.44
1:B:339:GLN:HE22	1:B:369:ARG:HB2	1.82	0.44
1:C:204:SER:HA	1:C:228:GLU:O	2.18	0.44
1:F:621:PHE:CG	1:F:622:PRO:HA	2.53	0.44
1:A:404:GLU:HG3	1:A:412:SER:HA	1.98	0.44
1:H:104:PRO:HB3	1:H:406:LEU:HB3	1.99	0.44
1:F:404:GLU:HG3	1:F:412:SER:HA	1.99	0.44
1:E:339:GLN:HE21	1:E:339:GLN:H	1.65	0.44
1:F:201:MET:HB2	1:F:222:LEU:HD11	1.99	0.44
1:D:235:TRP:HA	1:D:241:PHE:HB2	1.98	0.44
1:H:339:GLN:HE22	1:H:369:ARG:HB2	1.82	0.44
1:C:318:ASP:OD1	1:C:318:ASP:N	2.48	0.44
1:F:109:ARG:HG3	1:F:201:MET:HE3	1.98	0.43
1:G:153:LYS:HD2	1:G:153:LYS:HA	1.76	0.43
1:G:204:SER:HA	1:G:228:GLU:O	2.18	0.43
1:B:339:GLN:H	1:B:339:GLN:NE2	2.17	0.43
1:C:621:PHE:CG	1:C:622:PRO:HA	2.53	0.43
1:F:480:PHE:CE1	1:F:622:PRO:HD2	2.53	0.43
1:A:204:SER:HA	1:A:228:GLU:O	2.18	0.43
1:B:153:LYS:HA	1:B:153:LYS:HD2	1.79	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:109:ARG:HG3	1:D:201:MET:HE3	1.99	0.43
1:A:447:THR:HA	1:A:448:PRO:HD3	1.89	0.43
1:C:153:LYS:HA	1:C:153:LYS:HD2	1.67	0.43
1:A:369:ARG:H	1:A:369:ARG:HG3	1.68	0.43
1:D:532:LEU:HA	1:D:658:THR:HG23	2.01	0.43
1:G:232:VAL:HB	1:G:344:GLY:HA2	2.00	0.43
1:H:566:LEU:HA	1:H:567:PRO:HD3	1.91	0.43
1:E:621:PHE:CG	1:E:622:PRO:HA	2.54	0.43
1:G:339:GLN:H	1:G:339:GLN:NE2	2.17	0.43
1:G:621:PHE:CG	1:G:622:PRO:HA	2.53	0.43
1:H:549:ASP:OD2	1:H:604:HIS:NE2	2.42	0.43
1:B:164:ARG:HA	1:B:165:PRO:HD3	1.82	0.43
1:H:318:ASP:OD1	1:H:318:ASP:N	2.51	0.43
1:B:621:PHE:CG	1:B:622:PRO:HA	2.54	0.43
1:D:621:PHE:CG	1:D:622:PRO:HA	2.54	0.43
1:C:527:VAL:HA	1:C:528:PRO:HD3	1.88	0.42
1:F:232:VAL:HB	1:F:344:GLY:HA2	1.99	0.42
1:G:566:LEU:HA	1:G:567:PRO:HD3	1.92	0.42
1:A:164:ARG:HA	1:A:165:PRO:HD3	1.86	0.42
1:G:449:LEU:HG	1:G:457:LEU:HD22	2.02	0.42
1:H:480:PHE:CE1	1:H:622:PRO:HD2	2.55	0.42
1:A:331:LEU:HD23	1:A:331:LEU:HA	1.94	0.42
1:C:583:LYS:HA	1:C:584:PRO:HD2	1.94	0.42
1:G:365:MET:HB2	1:G:420:ILE:HG23	2.01	0.42
1:B:228:GLU:O	1:B:229:SER:C	2.63	0.42
1:B:461:HIS:HA	1:B:462:PRO:HD3	1.85	0.42
1:B:532:LEU:HA	1:B:658:THR:HG23	2.01	0.42
1:D:339:GLN:NE2	1:D:370:HIS:H	2.17	0.42
1:C:164:ARG:HA	1:C:165:PRO:HD3	1.86	0.42
1:F:478:VAL:HA	1:F:479:PRO:HD3	1.95	0.42
1:A:153:LYS:HA	1:A:153:LYS:HD2	1.80	0.42
1:A:339:GLN:HE22	1:A:369:ARG:HB2	1.84	0.42
1:A:480:PHE:CE1	1:A:622:PRO:HD2	2.54	0.42
1:E:92:THR:HG23	1:E:148:GLN:HB2	2.01	0.42
1:F:447:THR:HA	1:F:448:PRO:HD3	1.86	0.42
1:C:228:GLU:O	1:C:229:SER:C	2.63	0.41
1:G:205:SER:HB3	1:G:206:ALA:H	1.63	0.41
1:A:479:PRO:HA	1:A:621:PHE:O	2.20	0.41
1:B:339:GLN:NE2	1:B:370:HIS:H	2.18	0.41
1:F:449:LEU:HG	1:F:457:LEU:HD22	2.02	0.41
1:H:339:GLN:H	1:H:339:GLN:NE2	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:SER:HA	1:B:228:GLU:O	2.21	0.41
1:H:369:ARG:H	1:H:369:ARG:HG3	1.61	0.41
1:C:478:VAL:HA	1:C:479:PRO:HD3	1.93	0.41
1:H:235:TRP:HA	1:H:241:PHE:HB2	2.02	0.41
1:A:228:GLU:O	1:A:229:SER:C	2.63	0.41
1:E:232:VAL:HB	1:E:344:GLY:HA2	2.01	0.41
1:B:527:VAL:HA	1:B:528:PRO:HD3	1.90	0.41
1:E:339:GLN:H	1:E:339:GLN:NE2	2.19	0.41
1:B:201:MET:HB2	1:B:222:LEU:HD11	2.02	0.41
1:G:461:HIS:HA	1:G:462:PRO:HD3	1.73	0.41
1:H:204:SER:HA	1:H:228:GLU:O	2.21	0.41
1:C:544:VAL:HB	1:C:617:GLN:HG2	2.02	0.41
1:F:365:MET:HB2	1:F:420:ILE:HG23	2.02	0.41
1:B:96:ILE:HA	1:B:97:PRO:HD3	1.89	0.41
1:C:566:LEU:HA	1:C:567:PRO:HD3	1.90	0.41
1:E:274:ASP:HA	1:E:277:ASN:HB2	2.02	0.41
1:E:365:MET:HB2	1:E:420:ILE:HG23	2.02	0.41
1:F:228:GLU:O	1:F:229:SER:C	2.64	0.41
1:F:566:LEU:HA	1:F:567:PRO:HD3	1.93	0.41
1:G:96:ILE:HA	1:G:97:PRO:HD3	1.95	0.41
1:G:339:GLN:NE2	1:G:370:HIS:H	2.18	0.41
1:H:110:THR:HA	1:H:111:PRO:HD3	1.98	0.41
1:H:447:THR:HA	1:H:448:PRO:HD3	1.87	0.41
1:C:480:PHE:CE1	1:C:622:PRO:HD2	2.56	0.41
1:H:662:LEU:HA	1:H:663:PRO:HD3	1.99	0.41
1:H:339:GLN:NE2	1:H:370:HIS:H	2.18	0.40
1:D:104:PRO:HB3	1:D:406:LEU:HB3	2.02	0.40
1:C:201:MET:HB2	1:C:222:LEU:HD11	2.02	0.40
1:E:425:ASP:OD2	1:E:523:ARG:NH1	2.48	0.40
1:E:527:VAL:HA	1:E:528:PRO:HD3	1.91	0.40
1:H:201:MET:HB2	1:H:222:LEU:HD11	2.02	0.40
1:H:365:MET:HB2	1:H:420:ILE:HG23	2.02	0.40
1:H:512:THR:HG23	1:H:615:GLN:HG3	2.03	0.40
1:E:479:PRO:HA	1:E:621:PHE:O	2.20	0.40
1:F:104:PRO:HB3	1:F:406:LEU:HB3	2.03	0.40

There are no symmetry-related clashes.

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	608/652 (93%)	571 (94%)	36 (6%)	1 (0%)	43	72
1	B	608/652 (93%)	574 (94%)	33 (5%)	1 (0%)	43	72
1	C	608/652 (93%)	580 (95%)	27 (4%)	1 (0%)	43	72
1	D	608/652 (93%)	576 (95%)	31 (5%)	1 (0%)	43	72
1	E	608/652 (93%)	573 (94%)	33 (5%)	2 (0%)	36	66
1	F	608/652 (93%)	578 (95%)	29 (5%)	1 (0%)	43	72
1	G	608/652 (93%)	574 (94%)	33 (5%)	1 (0%)	43	72
1	H	608/652 (93%)	571 (94%)	36 (6%)	1 (0%)	43	72
All	All	4864/5216 (93%)	4597 (94%)	258 (5%)	9 (0%)	43	72

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	344	GLY
1	B	344	GLY
1	C	344	GLY
1	D	344	GLY
1	E	344	GLY
1	F	344	GLY
1	G	344	GLY
1	H	344	GLY
1	E	237	GLY

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	511/543 (94%)	475 (93%)	36 (7%)	14	40
1	B	511/543 (94%)	473 (93%)	38 (7%)	13	37
1	C	511/543 (94%)	475 (93%)	36 (7%)	14	40
1	D	511/543 (94%)	477 (93%)	34 (7%)	15	42
1	E	511/543 (94%)	472 (92%)	39 (8%)	12	36
1	F	511/543 (94%)	468 (92%)	43 (8%)	10	32
1	G	511/543 (94%)	473 (93%)	38 (7%)	13	37
1	H	511/543 (94%)	474 (93%)	37 (7%)	13	39
All	All	4088/4344 (94%)	3787 (93%)	301 (7%)	13	37

All (301) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	SER
1	A	71	GLN
1	A	72	GLN
1	A	93	VAL
1	A	101	ARG
1	A	107	LEU
1	A	115	LYS
1	A	120	ARG
1	A	130	VAL
1	A	140	GLU
1	A	229	SER
1	A	242	HIS
1	A	256	SER
1	A	266	ASP
1	A	293	LEU
1	A	319	LYS
1	A	323	GLN
1	A	331	LEU
1	A	339	GLN
1	A	353	LYS
1	A	369	ARG
1	A	384	GLU
1	A	412	SER
1	A	432	ARG
1	A	442	CYS

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Mol	Chain	Res	Type
1	A	453	ASP
1	A	496	TRP
1	A	541	SER
1	A	583	LYS
1	A	585	GLU
1	A	588	GLN
1	A	619	SER
1	A	620	TRP
1	A	629	GLN
1	A	644	THR
1	A	666	LYS
1	B	54	SER
1	B	65	VAL
1	B	71	GLN
1	B	98	LYS
1	B	107	LEU
1	B	110	THR
1	B	115	LYS
1	B	120	ARG
1	B	130	VAL
1	B	201	MET
1	B	205	SER
1	B	229	SER
1	B	242	HIS
1	B	256	SER
1	B	266	ASP
1	B	293	LEU
1	B	319	LYS
1	B	323	GLN
1	B	331	LEU
1	B	339	GLN
1	B	346	ILE
1	B	412	SER
1	B	432	ARG
1	B	442	CYS
1	B	453	ASP
1	B	469	SER
1	B	496	TRP
1	B	541	SER
1	B	579	LYS
1	B	583	LYS
1	B	588	GLN

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Mol	Chain	Res	Type
1	B	619	SER
1	B	620	TRP
1	B	629	GLN
1	B	639	LYS
1	B	644	THR
1	B	655	LYS
1	B	666	LYS
1	C	65	VAL
1	C	71	GLN
1	C	98	LYS
1	C	101	ARG
1	C	102	ASN
1	C	115	LYS
1	C	120	ARG
1	C	130	VAL
1	C	195	SER
1	C	201	MET
1	C	205	SER
1	C	229	SER
1	C	242	HIS
1	C	256	SER
1	C	286	SER
1	C	293	LEU
1	C	319	LYS
1	C	323	GLN
1	C	331	LEU
1	C	339	GLN
1	C	353	LYS
1	C	384	GLU
1	C	412	SER
1	C	432	ARG
1	C	442	CYS
1	C	453	ASP
1	C	496	TRP
1	C	541	SER
1	C	583	LYS
1	C	619	SER
1	C	620	TRP
1	C	629	GLN
1	C	639	LYS
1	C	644	THR
1	C	655	LYS

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Mol	Chain	Res	Type
1	C	666	LYS
1	D	65	VAL
1	D	71	GLN
1	D	98	LYS
1	D	115	LYS
1	D	120	ARG
1	D	130	VAL
1	D	201	MET
1	D	205	SER
1	D	229	SER
1	D	242	HIS
1	D	256	SER
1	D	266	ASP
1	D	286	SER
1	D	293	LEU
1	D	319	LYS
1	D	323	GLN
1	D	331	LEU
1	D	339	GLN
1	D	346	ILE
1	D	353	LYS
1	D	412	SER
1	D	432	ARG
1	D	442	CYS
1	D	453	ASP
1	D	496	TRP
1	D	541	SER
1	D	583	LYS
1	D	588	GLN
1	D	619	SER
1	D	620	TRP
1	D	629	GLN
1	D	644	THR
1	D	655	LYS
1	D	666	LYS
1	E	54	SER
1	E	65	VAL
1	E	71	GLN
1	E	107	LEU
1	E	115	LYS
1	E	120	ARG
1	E	130	VAL

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Mol	Chain	Res	Type
1	E	201	MET
1	E	204	SER
1	E	205	SER
1	E	229	SER
1	E	242	HIS
1	E	256	SER
1	E	286	SER
1	E	293	LEU
1	E	319	LYS
1	E	323	GLN
1	E	331	LEU
1	E	339	GLN
1	E	346	ILE
1	E	353	LYS
1	E	384	GLU
1	E	412	SER
1	E	420	ILE
1	E	432	ARG
1	E	442	CYS
1	E	453	ASP
1	E	460	THR
1	E	496	TRP
1	E	541	SER
1	E	583	LYS
1	E	588	GLN
1	E	619	SER
1	E	620	TRP
1	E	629	GLN
1	E	639	LYS
1	E	644	THR
1	E	655	LYS
1	E	666	LYS
1	F	54	SER
1	F	65	VAL
1	F	71	GLN
1	F	101	ARG
1	F	107	LEU
1	F	115	LYS
1	F	120	ARG
1	F	130	VAL
1	F	140	GLU
1	F	195	SER

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Mol	Chain	Res	Type
1	F	201	MET
1	F	204	SER
1	F	229	SER
1	F	242	HIS
1	F	256	SER
1	F	266	ASP
1	F	293	LEU
1	F	319	LYS
1	F	331	LEU
1	F	339	GLN
1	F	346	ILE
1	F	353	LYS
1	F	384	GLU
1	F	412	SER
1	F	417	ASP
1	F	420	ILE
1	F	432	ARG
1	F	442	CYS
1	F	453	ASP
1	F	458	SER
1	F	465	ASP
1	F	490	SER
1	F	496	TRP
1	F	541	SER
1	F	583	LYS
1	F	588	GLN
1	F	619	SER
1	F	620	TRP
1	F	629	GLN
1	F	639	LYS
1	F	644	THR
1	F	655	LYS
1	F	666	LYS
1	G	54	SER
1	G	65	VAL
1	G	71	GLN
1	G	93	VAL
1	G	107	LEU
1	G	115	LYS
1	G	120	ARG
1	G	130	VAL
1	G	195	SER

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Mol	Chain	Res	Type
1	G	201	MET
1	G	205	SER
1	G	229	SER
1	G	242	HIS
1	G	256	SER
1	G	266	ASP
1	G	293	LEU
1	G	319	LYS
1	G	323	GLN
1	G	331	LEU
1	G	339	GLN
1	G	346	ILE
1	G	369	ARG
1	G	412	SER
1	G	427	LYS
1	G	432	ARG
1	G	442	CYS
1	G	453	ASP
1	G	496	TRP
1	G	541	SER
1	G	583	LYS
1	G	588	GLN
1	G	619	SER
1	G	620	TRP
1	G	629	GLN
1	G	639	LYS
1	G	644	THR
1	G	655	LYS
1	G	666	LYS
1	H	54	SER
1	H	65	VAL
1	H	71	GLN
1	H	98	LYS
1	H	107	LEU
1	H	115	LYS
1	H	120	ARG
1	H	130	VAL
1	H	201	MET
1	H	229	SER
1	H	242	HIS
1	H	256	SER
1	H	266	ASP

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Mol	Chain	Res	Type
1	H	293	LEU
1	H	319	LYS
1	H	323	GLN
1	H	331	LEU
1	H	339	GLN
1	H	346	ILE
1	H	353	LYS
1	H	412	SER
1	H	432	ARG
1	H	438	CYS
1	H	442	CYS
1	H	453	ASP
1	H	469	SER
1	H	496	TRP
1	H	541	SER
1	H	583	LYS
1	H	588	GLN
1	H	619	SER
1	H	620	TRP
1	H	629	GLN
1	H	639	LYS
1	H	644	THR
1	H	655	LYS
1	H	666	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (57) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	133	GLN
1	A	191	ASN
1	A	339	GLN
1	A	441	ASN
1	A	499	GLN
1	A	652	HIS
1	B	113	ASN
1	B	133	GLN
1	B	191	ASN
1	B	248	GLN
1	B	295	GLN
1	B	339	GLN
1	B	391	HIS

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Mol	Chain	Res	Type
1	B	499	GLN
1	B	648	GLN
1	B	652	HIS
1	C	72	GLN
1	C	113	ASN
1	C	133	GLN
1	C	191	ASN
1	C	248	GLN
1	C	339	GLN
1	C	648	GLN
1	C	652	HIS
1	D	133	GLN
1	D	191	ASN
1	D	248	GLN
1	D	339	GLN
1	D	414	HIS
1	D	648	GLN
1	D	652	HIS
1	E	133	GLN
1	E	191	ASN
1	E	248	GLN
1	E	339	GLN
1	E	391	HIS
1	F	99	ASN
1	F	113	ASN
1	F	133	GLN
1	F	191	ASN
1	F	248	GLN
1	F	339	GLN
1	F	391	HIS
1	F	648	GLN
1	G	133	GLN
1	G	191	ASN
1	G	339	GLN
1	G	391	HIS
1	G	499	GLN
1	G	648	GLN
1	G	652	HIS
1	H	133	GLN
1	H	191	ASN
1	H	339	GLN
1	H	391	HIS

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Mol	Chain	Res	Type
1	H	648	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	612/652 (93%)	-0.31	3 (0%) 87 82	20, 20, 20, 20	0
1	B	612/652 (93%)	-0.33	6 (0%) 79 72	20, 20, 20, 20	0
1	C	612/652 (93%)	-0.24	9 (1%) 72 63	20, 20, 20, 20	0
1	D	612/652 (93%)	-0.16	9 (1%) 72 63	20, 20, 20, 20	0
1	E	612/652 (93%)	-0.05	3 (0%) 87 82	20, 20, 20, 20	0
1	F	612/652 (93%)	0.21	13 (2%) 63 54	20, 20, 20, 20	0
1	G	612/652 (93%)	-0.10	4 (0%) 84 77	20, 20, 20, 20	0
1	H	612/652 (93%)	-0.10	5 (0%) 82 75	20, 20, 20, 20	0
All	All	4896/5216 (93%)	-0.14	52 (1%) 78 70	20, 20, 20, 20	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	441	ASN	4.8
1	H	50	HIS	4.0
1	F	71	GLN	3.6
1	D	50	HIS	3.6
1	G	71	GLN	3.6
1	C	441	ASN	3.5
1	F	50	HIS	3.4
1	B	441	ASN	3.3
1	A	71	GLN	3.2
1	F	65	VAL	3.2
1	C	50	HIS	3.2
1	B	71	GLN	3.1
1	B	50	HIS	3.0
1	F	441	ASN	3.0
1	D	441	ASN	2.9
1	F	102	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	72	GLN	2.7
1	D	440	SER	2.6
1	C	65	VAL	2.6
1	D	265	ASN	2.6
1	E	266	ASP	2.5
1	E	71	GLN	2.5
1	D	71	GLN	2.4
1	H	51	ASP	2.4
1	C	58	GLY	2.3
1	F	654	GLY	2.3
1	G	50	HIS	2.3
1	D	57	THR	2.3
1	G	432	ARG	2.3
1	F	64	SER	2.2
1	F	563	GLY	2.2
1	F	101	ARG	2.2
1	H	441	ASN	2.2
1	E	50	HIS	2.2
1	F	453	ASP	2.2
1	D	414	HIS	2.1
1	C	440	SER	2.1
1	H	651	HIS	2.1
1	C	59	SER	2.1
1	F	619	SER	2.1
1	C	71	GLN	2.1
1	D	58	GLY	2.1
1	C	72	GLN	2.1
1	F	99	ASN	2.1
1	D	432	ARG	2.1
1	F	432	ARG	2.1
1	B	65	VAL	2.0
1	G	65	VAL	2.0
1	A	265	ASN	2.0
1	H	266	ASP	2.0
1	B	64	SER	2.0
1	C	295	GLN	2.0

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

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.