



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:38 AM UTC

PDB ID : 4UB9 / pdb_00004ub9
Title : Structural and catalytic characterization of molinate hydrolase
Authors : Leite, J.P.; Duarte, M.; Paiva, A.; Ferreira-da-Silva, F.; Matias, P.M.; Nunes, O.; Gales, L.
Deposited on : 2014-08-12
Resolution : 2.27 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

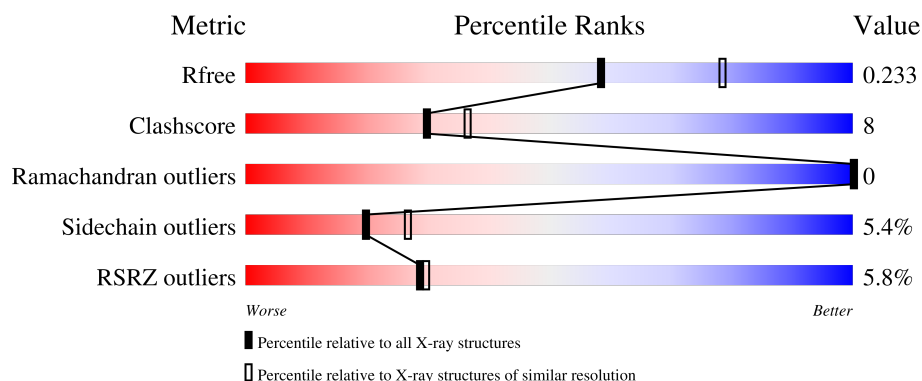
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	496	80% 12% • 6%
1	B	496	81% 11% • 6%
1	C	496	80% 12% • 6%
1	D	496	81% 11% • 6%
1	E	496	79% 12% • 6%

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Mol	Chain	Length	Quality of chain
1	F	496	4% 79% 12% • 6%
1	G	496	3% 80% 11% •• 6%
1	H	496	31% 75% 15% • 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 29391 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 Molinate hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	0
			3597	2276	619	680	22			
1	B	467	Total	C	N	O	S	0	1	0
			3602	2279	619	682	22			
1	C	467	Total	C	N	O	S	0	0	0
			3597	2276	619	680	22			
1	D	467	Total	C	N	O	S	0	0	0
			3597	2276	619	680	22			
1	E	467	Total	C	N	O	S	0	0	0
			3597	2276	619	680	22			
1	F	467	Total	C	N	O	S	0	0	0
			3597	2276	619	680	22			
1	G	467	Total	C	N	O	S	0	1	0
			3597	2276	619	680	22			
1	H	467	Total	C	N	O	S	0	0	0
			3597	2276	619	680	22			

There are 256 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP G2XLB0
A	2	ALA	-	expression tag	UNP G2XLB0
A	3	SER	-	expression tag	UNP G2XLB0
A	4	TRP	-	expression tag	UNP G2XLB0
A	5	SER	-	expression tag	UNP G2XLB0
A	6	HIS	-	expression tag	UNP G2XLB0
A	7	PRO	-	expression tag	UNP G2XLB0
A	8	GLN	-	expression tag	UNP G2XLB0
A	9	PHE	-	expression tag	UNP G2XLB0
A	10	GLU	-	expression tag	UNP G2XLB0
A	11	LYS	-	expression tag	UNP G2XLB0
A	12	ILE	-	expression tag	UNP G2XLB0
A	13	GLU	-	expression tag	UNP G2XLB0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	14	GLY	-	expression tag	UNP G2XLB0
A	15	ARG	-	expression tag	UNP G2XLB0
A	16	ARG	-	expression tag	UNP G2XLB0
A	17	ASP	-	expression tag	UNP G2XLB0
A	18	ARG	-	expression tag	UNP G2XLB0
A	19	GLY	-	expression tag	UNP G2XLB0
A	20	PRO	-	expression tag	UNP G2XLB0
A	21	GLU	-	expression tag	UNP G2XLB0
A	22	PHE	-	expression tag	UNP G2XLB0
A	23	GLU	-	expression tag	UNP G2XLB0
A	24	LEU	-	expression tag	UNP G2XLB0
A	25	GLY	-	expression tag	UNP G2XLB0
A	26	THR	-	expression tag	UNP G2XLB0
A	27	GLU	-	expression tag	UNP G2XLB0
A	28	ASN	-	expression tag	UNP G2XLB0
A	29	LEU	-	expression tag	UNP G2XLB0
A	30	PHE	-	expression tag	UNP G2XLB0
A	31	PHE	-	expression tag	UNP G2XLB0
A	32	GLU	-	expression tag	UNP G2XLB0
B	1	MET	-	initiating methionine	UNP G2XLB0
B	2	ALA	-	expression tag	UNP G2XLB0
B	3	SER	-	expression tag	UNP G2XLB0
B	4	TRP	-	expression tag	UNP G2XLB0
B	5	SER	-	expression tag	UNP G2XLB0
B	6	HIS	-	expression tag	UNP G2XLB0
B	7	PRO	-	expression tag	UNP G2XLB0
B	8	GLN	-	expression tag	UNP G2XLB0
B	9	PHE	-	expression tag	UNP G2XLB0
B	10	GLU	-	expression tag	UNP G2XLB0
B	11	LYS	-	expression tag	UNP G2XLB0
B	12	ILE	-	expression tag	UNP G2XLB0
B	13	GLU	-	expression tag	UNP G2XLB0
B	14	GLY	-	expression tag	UNP G2XLB0
B	15	ARG	-	expression tag	UNP G2XLB0
B	16	ARG	-	expression tag	UNP G2XLB0
B	17	ASP	-	expression tag	UNP G2XLB0
B	18	ARG	-	expression tag	UNP G2XLB0
B	19	GLY	-	expression tag	UNP G2XLB0
B	20	PRO	-	expression tag	UNP G2XLB0
B	21	GLU	-	expression tag	UNP G2XLB0
B	22	PHE	-	expression tag	UNP G2XLB0
B	23	GLU	-	expression tag	UNP G2XLB0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	24	LEU	-	expression tag	UNP G2XLB0
B	25	GLY	-	expression tag	UNP G2XLB0
B	26	THR	-	expression tag	UNP G2XLB0
B	27	GLU	-	expression tag	UNP G2XLB0
B	28	ASN	-	expression tag	UNP G2XLB0
B	29	LEU	-	expression tag	UNP G2XLB0
B	30	PHE	-	expression tag	UNP G2XLB0
B	31	PHE	-	expression tag	UNP G2XLB0
B	32	GLU	-	expression tag	UNP G2XLB0
C	1	MET	-	initiating methionine	UNP G2XLB0
C	2	ALA	-	expression tag	UNP G2XLB0
C	3	SER	-	expression tag	UNP G2XLB0
C	4	TRP	-	expression tag	UNP G2XLB0
C	5	SER	-	expression tag	UNP G2XLB0
C	6	HIS	-	expression tag	UNP G2XLB0
C	7	PRO	-	expression tag	UNP G2XLB0
C	8	GLN	-	expression tag	UNP G2XLB0
C	9	PHE	-	expression tag	UNP G2XLB0
C	10	GLU	-	expression tag	UNP G2XLB0
C	11	LYS	-	expression tag	UNP G2XLB0
C	12	ILE	-	expression tag	UNP G2XLB0
C	13	GLU	-	expression tag	UNP G2XLB0
C	14	GLY	-	expression tag	UNP G2XLB0
C	15	ARG	-	expression tag	UNP G2XLB0
C	16	ARG	-	expression tag	UNP G2XLB0
C	17	ASP	-	expression tag	UNP G2XLB0
C	18	ARG	-	expression tag	UNP G2XLB0
C	19	GLY	-	expression tag	UNP G2XLB0
C	20	PRO	-	expression tag	UNP G2XLB0
C	21	GLU	-	expression tag	UNP G2XLB0
C	22	PHE	-	expression tag	UNP G2XLB0
C	23	GLU	-	expression tag	UNP G2XLB0
C	24	LEU	-	expression tag	UNP G2XLB0
C	25	GLY	-	expression tag	UNP G2XLB0
C	26	THR	-	expression tag	UNP G2XLB0
C	27	GLU	-	expression tag	UNP G2XLB0
C	28	ASN	-	expression tag	UNP G2XLB0
C	29	LEU	-	expression tag	UNP G2XLB0
C	30	PHE	-	expression tag	UNP G2XLB0
C	31	PHE	-	expression tag	UNP G2XLB0
C	32	GLU	-	expression tag	UNP G2XLB0
D	1	MET	-	initiating methionine	UNP G2XLB0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	2	ALA	-	expression tag	UNP G2XLB0
D	3	SER	-	expression tag	UNP G2XLB0
D	4	TRP	-	expression tag	UNP G2XLB0
D	5	SER	-	expression tag	UNP G2XLB0
D	6	HIS	-	expression tag	UNP G2XLB0
D	7	PRO	-	expression tag	UNP G2XLB0
D	8	GLN	-	expression tag	UNP G2XLB0
D	9	PHE	-	expression tag	UNP G2XLB0
D	10	GLU	-	expression tag	UNP G2XLB0
D	11	LYS	-	expression tag	UNP G2XLB0
D	12	ILE	-	expression tag	UNP G2XLB0
D	13	GLU	-	expression tag	UNP G2XLB0
D	14	GLY	-	expression tag	UNP G2XLB0
D	15	ARG	-	expression tag	UNP G2XLB0
D	16	ARG	-	expression tag	UNP G2XLB0
D	17	ASP	-	expression tag	UNP G2XLB0
D	18	ARG	-	expression tag	UNP G2XLB0
D	19	GLY	-	expression tag	UNP G2XLB0
D	20	PRO	-	expression tag	UNP G2XLB0
D	21	GLU	-	expression tag	UNP G2XLB0
D	22	PHE	-	expression tag	UNP G2XLB0
D	23	GLU	-	expression tag	UNP G2XLB0
D	24	LEU	-	expression tag	UNP G2XLB0
D	25	GLY	-	expression tag	UNP G2XLB0
D	26	THR	-	expression tag	UNP G2XLB0
D	27	GLU	-	expression tag	UNP G2XLB0
D	28	ASN	-	expression tag	UNP G2XLB0
D	29	LEU	-	expression tag	UNP G2XLB0
D	30	PHE	-	expression tag	UNP G2XLB0
D	31	PHE	-	expression tag	UNP G2XLB0
D	32	GLU	-	expression tag	UNP G2XLB0
E	1	MET	-	initiating methionine	UNP G2XLB0
E	2	ALA	-	expression tag	UNP G2XLB0
E	3	SER	-	expression tag	UNP G2XLB0
E	4	TRP	-	expression tag	UNP G2XLB0
E	5	SER	-	expression tag	UNP G2XLB0
E	6	HIS	-	expression tag	UNP G2XLB0
E	7	PRO	-	expression tag	UNP G2XLB0
E	8	GLN	-	expression tag	UNP G2XLB0
E	9	PHE	-	expression tag	UNP G2XLB0
E	10	GLU	-	expression tag	UNP G2XLB0
E	11	LYS	-	expression tag	UNP G2XLB0

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Chain	Residue	Modelled	Actual	Comment	Reference
E	12	ILE	-	expression tag	UNP G2XLB0
E	13	GLU	-	expression tag	UNP G2XLB0
E	14	GLY	-	expression tag	UNP G2XLB0
E	15	ARG	-	expression tag	UNP G2XLB0
E	16	ARG	-	expression tag	UNP G2XLB0
E	17	ASP	-	expression tag	UNP G2XLB0
E	18	ARG	-	expression tag	UNP G2XLB0
E	19	GLY	-	expression tag	UNP G2XLB0
E	20	PRO	-	expression tag	UNP G2XLB0
E	21	GLU	-	expression tag	UNP G2XLB0
E	22	PHE	-	expression tag	UNP G2XLB0
E	23	GLU	-	expression tag	UNP G2XLB0
E	24	LEU	-	expression tag	UNP G2XLB0
E	25	GLY	-	expression tag	UNP G2XLB0
E	26	THR	-	expression tag	UNP G2XLB0
E	27	GLU	-	expression tag	UNP G2XLB0
E	28	ASN	-	expression tag	UNP G2XLB0
E	29	LEU	-	expression tag	UNP G2XLB0
E	30	PHE	-	expression tag	UNP G2XLB0
E	31	PHE	-	expression tag	UNP G2XLB0
E	32	GLU	-	expression tag	UNP G2XLB0
F	1	MET	-	initiating methionine	UNP G2XLB0
F	2	ALA	-	expression tag	UNP G2XLB0
F	3	SER	-	expression tag	UNP G2XLB0
F	4	TRP	-	expression tag	UNP G2XLB0
F	5	SER	-	expression tag	UNP G2XLB0
F	6	HIS	-	expression tag	UNP G2XLB0
F	7	PRO	-	expression tag	UNP G2XLB0
F	8	GLN	-	expression tag	UNP G2XLB0
F	9	PHE	-	expression tag	UNP G2XLB0
F	10	GLU	-	expression tag	UNP G2XLB0
F	11	LYS	-	expression tag	UNP G2XLB0
F	12	ILE	-	expression tag	UNP G2XLB0
F	13	GLU	-	expression tag	UNP G2XLB0
F	14	GLY	-	expression tag	UNP G2XLB0
F	15	ARG	-	expression tag	UNP G2XLB0
F	16	ARG	-	expression tag	UNP G2XLB0
F	17	ASP	-	expression tag	UNP G2XLB0
F	18	ARG	-	expression tag	UNP G2XLB0
F	19	GLY	-	expression tag	UNP G2XLB0
F	20	PRO	-	expression tag	UNP G2XLB0
F	21	GLU	-	expression tag	UNP G2XLB0

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Chain	Residue	Modelled	Actual	Comment	Reference
F	22	PHE	-	expression tag	UNP G2XLB0
F	23	GLU	-	expression tag	UNP G2XLB0
F	24	LEU	-	expression tag	UNP G2XLB0
F	25	GLY	-	expression tag	UNP G2XLB0
F	26	THR	-	expression tag	UNP G2XLB0
F	27	GLU	-	expression tag	UNP G2XLB0
F	28	ASN	-	expression tag	UNP G2XLB0
F	29	LEU	-	expression tag	UNP G2XLB0
F	30	PHE	-	expression tag	UNP G2XLB0
F	31	PHE	-	expression tag	UNP G2XLB0
F	32	GLU	-	expression tag	UNP G2XLB0
G	1	MET	-	initiating methionine	UNP G2XLB0
G	2	ALA	-	expression tag	UNP G2XLB0
G	3	SER	-	expression tag	UNP G2XLB0
G	4	TRP	-	expression tag	UNP G2XLB0
G	5	SER	-	expression tag	UNP G2XLB0
G	6	HIS	-	expression tag	UNP G2XLB0
G	7	PRO	-	expression tag	UNP G2XLB0
G	8	GLN	-	expression tag	UNP G2XLB0
G	9	PHE	-	expression tag	UNP G2XLB0
G	10	GLU	-	expression tag	UNP G2XLB0
G	11	LYS	-	expression tag	UNP G2XLB0
G	12	ILE	-	expression tag	UNP G2XLB0
G	13	GLU	-	expression tag	UNP G2XLB0
G	14	GLY	-	expression tag	UNP G2XLB0
G	15	ARG	-	expression tag	UNP G2XLB0
G	16	ARG	-	expression tag	UNP G2XLB0
G	17	ASP	-	expression tag	UNP G2XLB0
G	18	ARG	-	expression tag	UNP G2XLB0
G	19	GLY	-	expression tag	UNP G2XLB0
G	20	PRO	-	expression tag	UNP G2XLB0
G	21	GLU	-	expression tag	UNP G2XLB0
G	22	PHE	-	expression tag	UNP G2XLB0
G	23	GLU	-	expression tag	UNP G2XLB0
G	24	LEU	-	expression tag	UNP G2XLB0
G	25	GLY	-	expression tag	UNP G2XLB0
G	26	THR	-	expression tag	UNP G2XLB0
G	27	GLU	-	expression tag	UNP G2XLB0
G	28	ASN	-	expression tag	UNP G2XLB0
G	29	LEU	-	expression tag	UNP G2XLB0
G	30	PHE	-	expression tag	UNP G2XLB0
G	31	PHE	-	expression tag	UNP G2XLB0

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Chain	Residue	Modelled	Actual	Comment	Reference
G	32	GLU	-	expression tag	UNP G2XLB0
H	1	MET	-	initiating methionine	UNP G2XLB0
H	2	ALA	-	expression tag	UNP G2XLB0
H	3	SER	-	expression tag	UNP G2XLB0
H	4	TRP	-	expression tag	UNP G2XLB0
H	5	SER	-	expression tag	UNP G2XLB0
H	6	HIS	-	expression tag	UNP G2XLB0
H	7	PRO	-	expression tag	UNP G2XLB0
H	8	GLN	-	expression tag	UNP G2XLB0
H	9	PHE	-	expression tag	UNP G2XLB0
H	10	GLU	-	expression tag	UNP G2XLB0
H	11	LYS	-	expression tag	UNP G2XLB0
H	12	ILE	-	expression tag	UNP G2XLB0
H	13	GLU	-	expression tag	UNP G2XLB0
H	14	GLY	-	expression tag	UNP G2XLB0
H	15	ARG	-	expression tag	UNP G2XLB0
H	16	ARG	-	expression tag	UNP G2XLB0
H	17	ASP	-	expression tag	UNP G2XLB0
H	18	ARG	-	expression tag	UNP G2XLB0
H	19	GLY	-	expression tag	UNP G2XLB0
H	20	PRO	-	expression tag	UNP G2XLB0
H	21	GLU	-	expression tag	UNP G2XLB0
H	22	PHE	-	expression tag	UNP G2XLB0
H	23	GLU	-	expression tag	UNP G2XLB0
H	24	LEU	-	expression tag	UNP G2XLB0
H	25	GLY	-	expression tag	UNP G2XLB0
H	26	THR	-	expression tag	UNP G2XLB0
H	27	GLU	-	expression tag	UNP G2XLB0
H	28	ASN	-	expression tag	UNP G2XLB0
H	29	LEU	-	expression tag	UNP G2XLB0
H	30	PHE	-	expression tag	UNP G2XLB0
H	31	PHE	-	expression tag	UNP G2XLB0
H	32	GLU	-	expression tag	UNP G2XLB0

- 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	1	Total Zn 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total 1	Zn 1	0	0
2	E	1	Total 1	Zn 1	0	0
2	F	1	Total 1	Zn 1	0	0
2	G	1	Total 1	Zn 1	0	0
2	H	1	Total 1	Zn 1	0	0

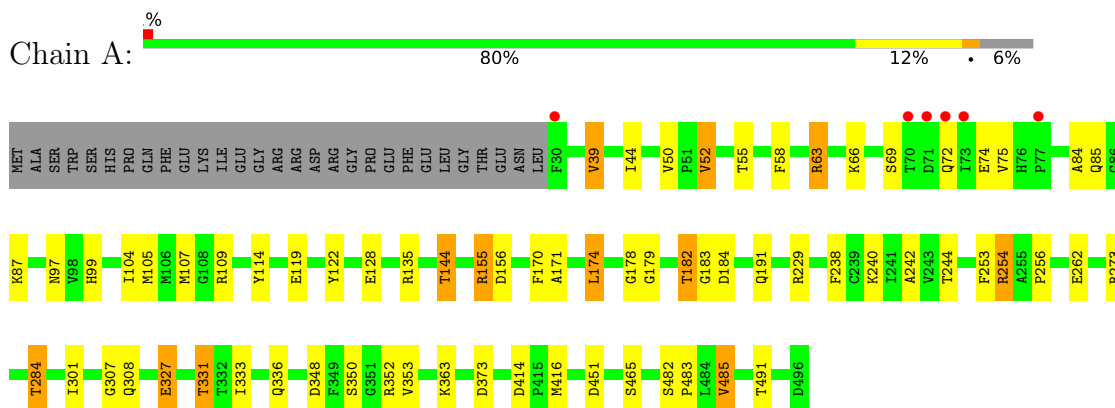
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	116	Total 116	O 116	0	0
3	B	112	Total 112	O 112	0	0
3	C	67	Total 67	O 67	0	0
3	D	83	Total 83	O 83	0	0
3	E	75	Total 75	O 75	0	0
3	F	50	Total 50	O 50	0	0
3	G	54	Total 54	O 54	0	0
3	H	45	Total 45	O 45	0	0

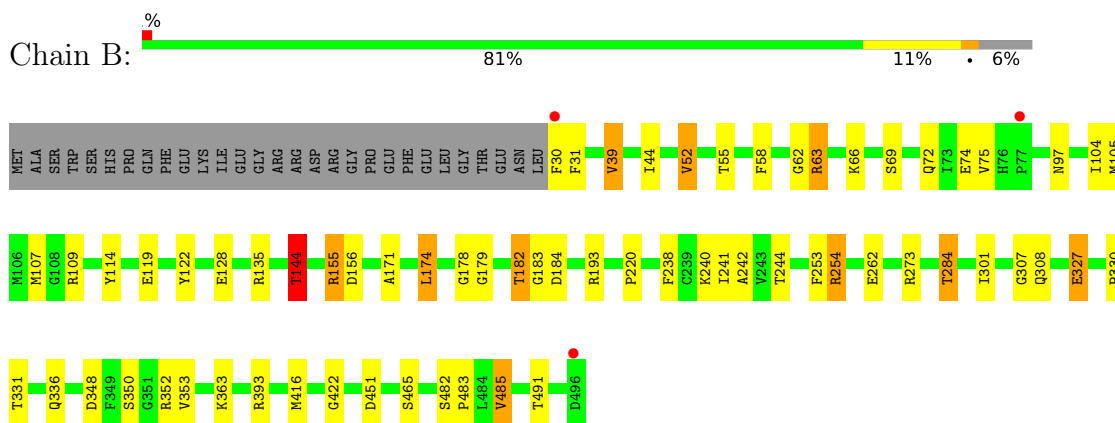
3 Residue-property plots [i](#)

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.

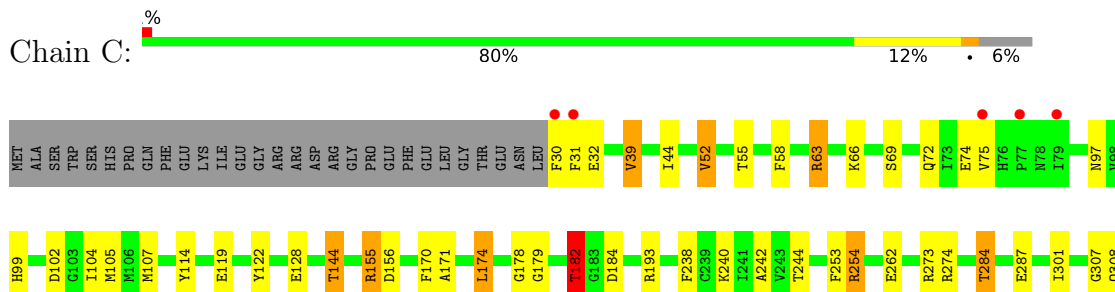
• Molecule 1: Molinate hydrolase



• Molecule 1: Molinate hydrolase

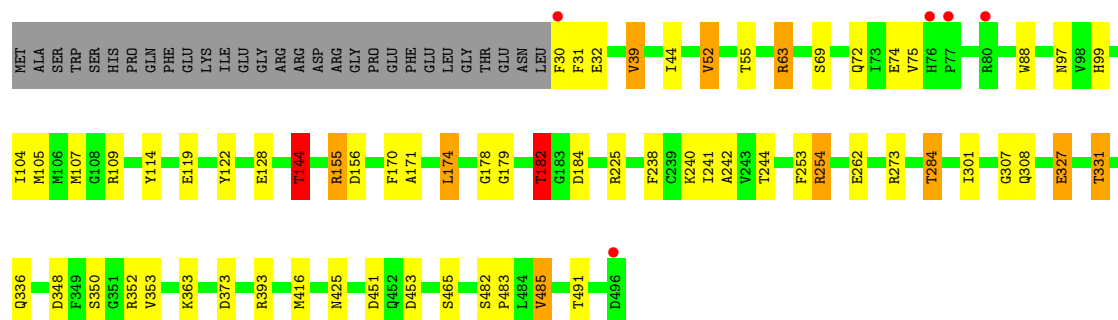
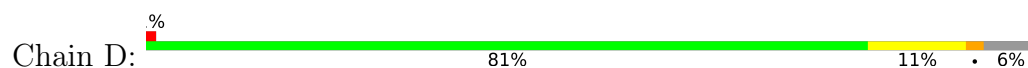


• Molecule 1: Molinate hydrolase

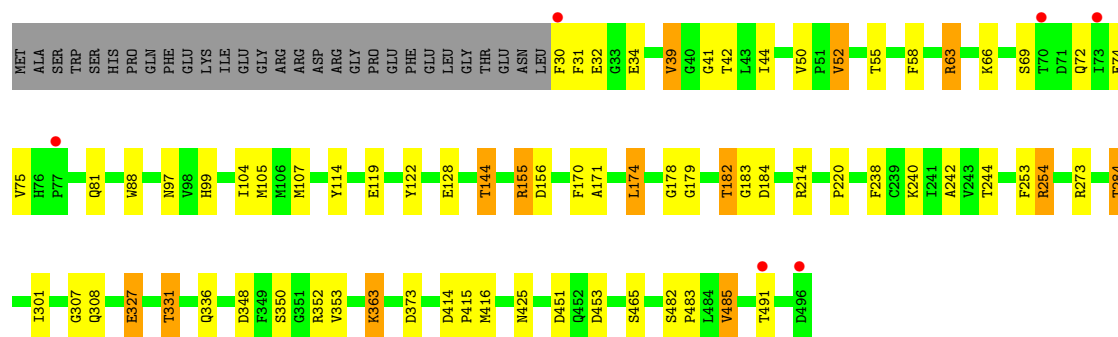
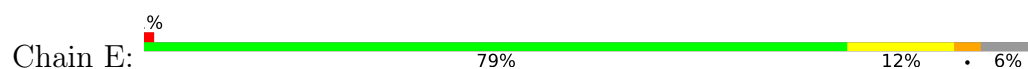




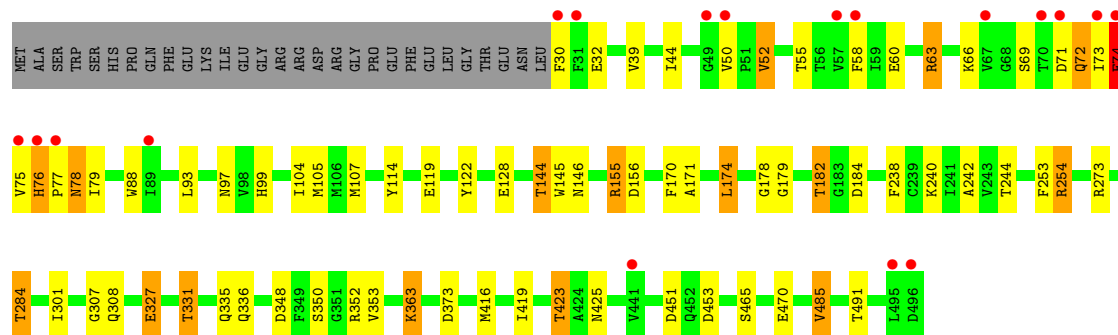
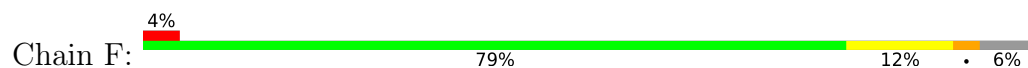
• Molecule 1: Molinate hydrolase



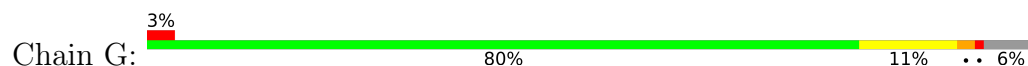
• Molecule 1: Molinate hydrolase

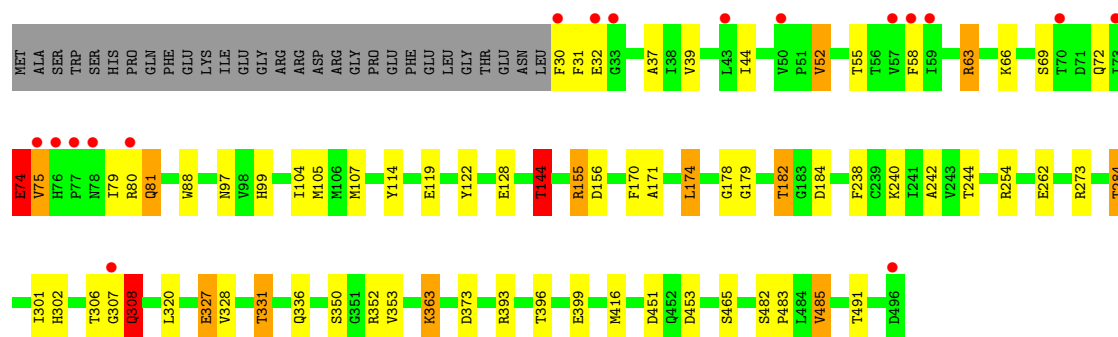


• Molecule 1: Molinate hydrolase

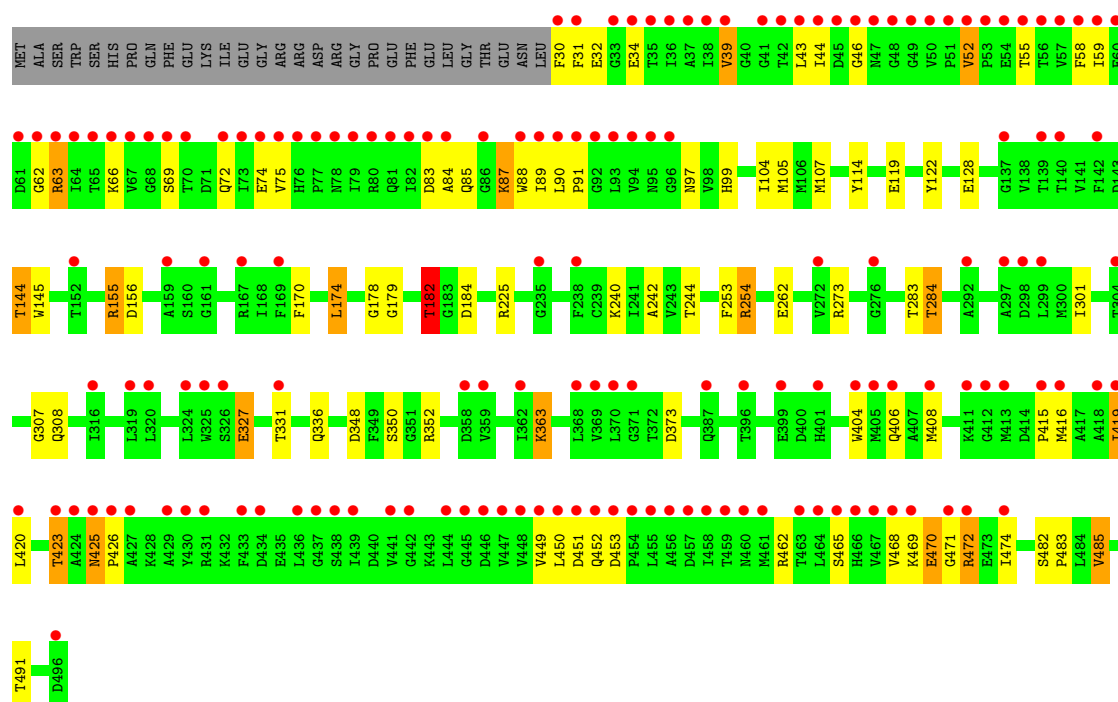
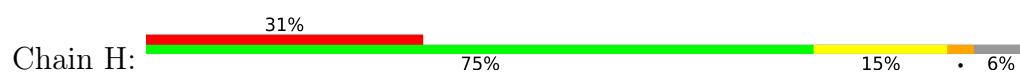


• Molecule 1: Molinate hydrolase





● Molecule 1: Molinate hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	367.70Å 99.08Å 131.34Å 90.00° 109.61° 90.00°	Depositor
Resolution (Å)	123.72 – 2.27 123.72 – 2.27	Depositor EDS
% Data completeness (in resolution range)	94.5 (123.72-2.27) 94.5 (123.72-2.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.27Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.201 , 0.229 0.207 , 0.233	Depositor DCC
R_{free} test set	9639 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 23.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.018 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	29391	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: ZN

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	1.01	0/3673	0.98	3/4989 (0.1%)
1	B	1.00	3/3681 (0.1%)	0.99	6/5000 (0.1%)
1	C	0.99	1/3673 (0.0%)	0.98	6/4989 (0.1%)
1	D	0.97	2/3673 (0.1%)	0.98	6/4989 (0.1%)
1	E	0.95	2/3673 (0.1%)	0.96	5/4989 (0.1%)
1	F	0.91	0/3673	0.98	5/4989 (0.1%)
1	G	0.95	4/3673 (0.1%)	1.01	11/4989 (0.2%)
1	H	0.92	1/3673 (0.0%)	0.97	5/4989 (0.1%)
All	All	0.96	13/29392 (0.0%)	0.98	47/39923 (0.1%)

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	F	0	1
1	G	0	1
All	All	0	2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	31	PHE	N-CA	6.57	1.54	1.46
1	G	307	GLY	C-O	-6.20	1.15	1.23
1	D	31	PHE	N-CA	5.83	1.53	1.45
1	B	330	PRO	N-CA	-5.80	1.39	1.47
1	B	422	GLY	C-O	-5.60	1.18	1.24
1	G	307	GLY	CA-C	-5.54	1.44	1.51
1	C	31	PHE	N-CA	5.47	1.52	1.45
1	G	31	PHE	N-CA	5.45	1.52	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	425	ASN	C-O	-5.40	1.19	1.24
1	E	31	PHE	N-CA	5.37	1.52	1.45
1	H	31	PHE	N-CA	5.29	1.52	1.45
1	E	425	ASN	C-O	-5.11	1.19	1.24
1	G	74	GLU	CA-C	5.11	1.58	1.52

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	307	GLY	CA-C-N	8.30	136.64	121.70
1	G	307	GLY	C-N-CA	8.30	136.64	121.70
1	G	308	GLN	N-CA-CB	-7.95	96.98	110.50
1	G	307	GLY	N-CA-C	7.47	130.87	113.18
1	F	72	GLN	N-CA-C	6.69	118.65	111.36
1	H	32	GLU	N-CA-C	6.66	118.88	108.96
1	G	32	GLU	N-CA-C	6.54	118.70	108.96
1	C	317	GLU	CA-CB-CG	-6.51	101.08	114.10
1	C	32	GLU	N-CA-C	6.38	118.46	108.96
1	F	32	GLU	N-CA-C	6.31	118.36	108.96
1	D	32	GLU	N-CA-C	6.29	118.33	108.96
1	E	32	GLU	N-CA-C	6.28	118.31	108.96
1	C	39	VAL	N-CA-C	6.24	116.85	108.11
1	F	74	GLU	N-CA-C	6.13	119.23	108.75
1	B	39	VAL	N-CA-C	5.98	116.48	108.11
1	A	72	GLN	CA-CB-CG	5.91	125.92	114.10
1	G	74	GLU	CB-CA-C	5.88	120.59	109.37
1	C	356	GLU	CB-CA-C	-5.85	101.08	110.79
1	C	170	PHE	N-CA-C	5.66	117.78	109.24
1	G	306	THR	CB-CA-C	-5.59	102.83	112.44
1	G	170	PHE	N-CA-C	5.57	117.65	109.24
1	E	220	PRO	CA-C-N	5.47	125.56	119.32
1	E	220	PRO	C-N-CA	5.47	125.56	119.32
1	D	170	PHE	N-CA-C	5.41	117.41	109.24
1	H	170	PHE	N-CA-C	5.37	117.84	109.52
1	H	39	VAL	N-CA-C	5.36	115.67	108.17
1	E	170	PHE	N-CA-C	5.35	117.31	109.24
1	A	39	VAL	N-CA-C	5.33	115.63	108.17
1	A	170	PHE	N-CA-C	5.30	117.24	109.24
1	F	170	PHE	N-CA-C	5.29	117.23	109.24
1	B	144	THR	N-CA-CB	-5.28	102.67	110.53
1	E	39	VAL	N-CA-C	5.25	115.47	108.11
1	D	39	VAL	N-CA-C	5.24	115.51	108.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	39	VAL	N-CA-C	5.22	115.42	108.11
1	D	241	ILE	CB-CA-C	-5.21	103.07	110.83
1	C	182	THR	N-CA-CB	-5.20	103.24	111.05
1	G	144	THR	N-CA-CB	-5.20	102.78	110.53
1	B	241	ILE	CB-CA-C	-5.19	103.09	110.83
1	H	182	THR	N-CA-CB	-5.13	103.12	110.87
1	D	144	THR	N-CA-CB	-5.11	102.92	110.53
1	H	470	GLU	N-CA-C	-5.07	104.65	111.54
1	D	182	THR	N-CA-CB	-5.05	103.48	111.05
1	F	39	VAL	N-CA-C	5.03	115.22	108.17
1	B	109	ARG	CG-CD-NE	-5.02	100.95	112.00
1	G	81	GLN	CA-CB-CG	-5.01	104.08	114.10
1	B	220	PRO	CA-C-N	5.00	125.02	119.32
1	B	220	PRO	C-N-CA	5.00	125.02	119.32

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	77	PRO	Peptide
1	G	74	GLU	Peptide

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	3597	0	3555	57	0
1	B	3602	0	3559	47	0
1	C	3597	0	3555	54	0
1	D	3597	0	3555	52	0
1	E	3597	0	3555	54	0
1	F	3597	0	3555	69	1
1	G	3597	0	3545	66	1
1	H	3597	0	3555	97	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	116	0	0	12	0
3	B	112	0	0	9	0
3	C	67	0	0	11	0
3	D	83	0	0	12	0
3	E	75	0	0	11	2
3	F	50	0	0	13	1
3	G	54	0	0	16	1
3	H	45	0	0	21	0
All	All	29391	0	28434	485	4

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 8.

All (485) 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:H:59:ILE:HG22	3:H:641:HOH:O	1.26	1.30
1:H:88:TRP:CZ2	1:H:453:ASP:HB2	1.67	1.30
1:F:331:THR:HG22	3:F:646:HOH:O	1.34	1.27
1:F:58:PHE:CE1	1:F:75:VAL:HG11	1.79	1.16
1:G:331:THR:HG21	1:G:336:GLN:HE21	1.12	1.15
1:H:415:PRO:O	3:H:637:HOH:O	1.66	1.12
1:E:331:THR:HG21	1:E:336:GLN:HE21	1.12	1.12
1:H:331:THR:HG21	1:H:336:GLN:HE21	1.11	1.12
1:H:34:GLU:HG3	1:H:471:GLY:O	1.46	1.11
1:H:59:ILE:CG2	3:H:641:HOH:O	1.83	1.11
1:E:34:GLU:HG2	3:E:662:HOH:O	1.50	1.10
1:C:331:THR:HG21	1:C:336:GLN:HE21	1.13	1.09
1:D:331:THR:HG21	1:D:336:GLN:HE21	1.11	1.09
1:D:105:MET:HG2	3:D:662:HOH:O	1.53	1.09
1:G:37:ALA:HB3	1:G:81:GLN:OE1	1.52	1.08
1:F:331:THR:HG21	1:F:336:GLN:HE21	1.15	1.07
1:B:331:THR:HG21	1:B:336:GLN:HE21	1.10	1.07
1:F:71:ASP:OD1	1:F:72:GLN:NE2	1.88	1.07
1:A:331:THR:HG21	1:A:336:GLN:HE21	1.13	1.04
1:H:34:GLU:CD	1:H:472:ARG:HA	1.81	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:52:VAL:HG21	3:D:671:HOH:O	1.58	1.03
1:G:302:HIS:HA	3:G:654:HOH:O	1.57	1.02
1:D:52:VAL:CG2	3:D:671:HOH:O	2.08	1.02
1:H:34:GLU:OE2	1:H:472:ARG:HA	1.60	0.99
1:H:472:ARG:NH2	1:H:474:ILE:CD1	2.27	0.98
1:A:284:THR:HG21	3:A:660:HOH:O	1.64	0.97
1:F:331:THR:CG2	3:F:646:HOH:O	2.00	0.96
1:F:58:PHE:HE1	1:F:75:VAL:HG11	1.25	0.94
1:H:88:TRP:CE2	1:H:453:ASP:HB2	2.05	0.92
1:G:331:THR:HG22	3:G:644:HOH:O	1.71	0.90
1:H:34:GLU:CG	1:H:471:GLY:O	2.19	0.90
1:H:89:ILE:C	1:H:90:LEU:HD12	1.98	0.89
1:F:145:TRP:O	3:F:644:HOH:O	1.89	0.88
1:B:135:ARG:HD2	3:B:712:HOH:O	1.72	0.88
1:D:331:THR:HG21	1:D:336:GLN:NE2	1.91	0.86
1:B:331:THR:HG21	1:B:336:GLN:NE2	1.90	0.85
1:E:331:THR:HG21	1:E:336:GLN:NE2	1.92	0.85
1:H:83:ASP:OD1	3:H:622:HOH:O	1.95	0.84
1:H:89:ILE:O	1:H:90:LEU:HD12	1.76	0.84
1:C:274:ARG:CD	3:C:602:HOH:O	2.26	0.84
1:B:416:MET:CE	3:B:629:HOH:O	2.24	0.83
1:G:331:THR:HG21	1:G:336:GLN:NE2	1.92	0.83
1:A:331:THR:HG21	1:A:336:GLN:NE2	1.93	0.83
1:A:416:MET:CE	3:A:629:HOH:O	2.27	0.82
1:F:58:PHE:CE1	1:F:75:VAL:CG1	2.62	0.82
1:C:331:THR:HG21	1:C:336:GLN:NE2	1.93	0.82
1:G:75:VAL:HG21	1:G:81:GLN:NE2	1.94	0.82
1:F:331:THR:HG21	1:F:336:GLN:NE2	1.94	0.82
1:H:419:ILE:O	1:H:423:THR:OG1	1.98	0.81
1:H:331:THR:HG21	1:H:336:GLN:NE2	1.92	0.81
1:C:274:ARG:HD3	3:C:602:HOH:O	1.80	0.80
1:H:88:TRP:CZ2	1:H:453:ASP:CB	2.61	0.80
1:D:105:MET:HE3	3:D:662:HOH:O	1.81	0.79
1:E:179:GLY:O	1:E:182:THR:HB	1.81	0.79
1:G:178:GLY:HA3	1:G:182:THR:HG21	1.64	0.79
1:A:179:GLY:O	1:A:182:THR:HB	1.82	0.79
1:H:284:THR:HG21	3:H:645:HOH:O	1.83	0.79
1:C:179:GLY:O	1:C:182:THR:HB	1.83	0.79
1:A:50:VAL:HG12	1:E:414:ASP:OD1	1.82	0.79
1:D:179:GLY:O	1:D:182:THR:HB	1.82	0.78
1:D:178:GLY:HA3	1:D:182:THR:HG21	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:179:GLY:O	1:F:182:THR:HB	1.83	0.78
1:H:449:VAL:HG12	3:H:602:HOH:O	1.82	0.78
1:B:179:GLY:O	1:B:182:THR:HB	1.83	0.78
1:H:179:GLY:O	1:H:182:THR:HB	1.83	0.78
1:H:88:TRP:CE2	1:H:453:ASP:CB	2.66	0.78
1:G:179:GLY:O	1:G:182:THR:HB	1.84	0.78
1:H:178:GLY:HA3	1:H:182:THR:HG21	1.66	0.77
1:H:468:VAL:HA	1:H:472:ARG:O	1.85	0.77
1:H:472:ARG:NH2	1:H:474:ILE:HD12	1.98	0.77
1:A:178:GLY:HA3	1:A:182:THR:HG21	1.66	0.77
1:C:105:MET:SD	3:C:657:HOH:O	2.41	0.77
1:G:399:GLU:HB2	3:G:635:HOH:O	1.83	0.76
1:F:76:HIS:CD2	1:F:78:ASN:HD21	2.03	0.76
1:E:178:GLY:HA3	1:E:182:THR:HG21	1.67	0.76
1:A:262:GLU:HG3	1:B:262:GLU:HG3	1.65	0.76
1:H:34:GLU:CD	1:H:472:ARG:CA	2.58	0.76
1:F:284:THR:HG21	3:F:635:HOH:O	1.86	0.75
1:F:71:ASP:OD1	1:F:72:GLN:N	2.20	0.75
1:C:178:GLY:HA3	1:C:182:THR:HG21	1.67	0.75
1:F:178:GLY:HA3	1:F:182:THR:HG21	1.68	0.75
1:H:415:PRO:O	1:H:419:ILE:HG12	1.82	0.75
1:B:178:GLY:HA3	1:B:182:THR:HG21	1.68	0.74
1:B:393:ARG:HD2	3:B:711:HOH:O	1.86	0.74
1:D:416:MET:CE	3:D:625:HOH:O	2.35	0.74
1:E:416:MET:CE	3:E:625:HOH:O	2.35	0.74
1:G:416:MET:CE	3:G:612:HOH:O	2.35	0.74
1:C:262:GLU:HG3	1:D:262:GLU:HG3	1.69	0.73
1:F:419:ILE:O	1:F:423:THR:OG1	2.07	0.72
1:H:283:THR:OG1	3:H:615:HOH:O	2.06	0.72
1:H:89:ILE:HG13	3:H:636:HOH:O	1.90	0.71
1:D:105:MET:CE	3:D:662:HOH:O	2.37	0.71
1:D:225:ARG:HD2	1:G:320:LEU:HD13	1.74	0.70
1:E:284:THR:CG2	3:E:661:HOH:O	2.40	0.70
1:B:104:ILE:HD13	1:B:107:MET:HE2	1.72	0.70
1:F:244:THR:OG1	1:F:284:THR:HB	1.92	0.70
1:E:104:ILE:HD13	1:E:107:MET:HE2	1.74	0.69
1:F:146:ASN:HA	3:F:644:HOH:O	1.91	0.69
1:A:333:ILE:O	3:A:706:HOH:O	2.10	0.69
1:B:416:MET:HE1	3:B:629:HOH:O	1.91	0.69
1:H:87:LYS:HD3	1:H:451:ASP:HA	1.74	0.69
1:F:104:ILE:HD13	1:F:107:MET:HE2	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:244:THR:OG1	1:D:284:THR:HB	1.92	0.69
1:G:75:VAL:HG21	1:G:81:GLN:HE22	1.54	0.69
1:G:244:THR:OG1	1:G:284:THR:HB	1.93	0.69
1:H:104:ILE:HD13	1:H:107:MET:HE2	1.75	0.69
1:H:174:LEU:HD23	1:H:242:ALA:HB2	1.74	0.69
1:H:244:THR:OG1	1:H:284:THR:HB	1.93	0.69
1:D:104:ILE:HD13	1:D:107:MET:HE2	1.75	0.69
1:G:393:ARG:HD2	3:G:646:HOH:O	1.92	0.68
1:D:174:LEU:HD23	1:D:242:ALA:HB2	1.76	0.68
1:C:104:ILE:HD13	1:C:107:MET:HE2	1.75	0.68
1:H:450:LEU:C	3:H:602:HOH:O	2.36	0.68
1:F:174:LEU:HD23	1:F:242:ALA:HB2	1.75	0.68
1:A:104:ILE:HD13	1:A:107:MET:HE2	1.73	0.68
1:F:76:HIS:O	1:F:79:ILE:HG12	1.94	0.68
1:F:416:MET:CE	3:F:616:HOH:O	2.41	0.68
1:H:99:HIS:HE1	1:H:373:ASP:OD1	1.77	0.68
1:E:416:MET:HE1	3:E:625:HOH:O	1.94	0.68
1:E:244:THR:OG1	1:E:284:THR:HB	1.94	0.67
1:B:244:THR:OG1	1:B:284:THR:HB	1.94	0.67
1:C:244:THR:OG1	1:C:284:THR:HB	1.94	0.67
1:G:104:ILE:HD13	1:G:107:MET:HE2	1.76	0.67
1:F:63:ARG:HH11	1:F:63:ARG:HG3	1.59	0.67
1:C:451:ASP:OD1	1:C:465:SER:OG	2.13	0.67
1:D:52:VAL:HG22	3:D:671:HOH:O	1.85	0.67
1:E:284:THR:HG21	3:E:661:HOH:O	1.95	0.67
1:G:75:VAL:CG2	1:G:81:GLN:NE2	2.57	0.67
1:B:174:LEU:HD23	1:B:242:ALA:HB2	1.75	0.67
1:G:284:THR:HG21	3:G:621:HOH:O	1.95	0.67
1:G:262:GLU:HG3	1:H:262:GLU:HG3	1.77	0.67
1:A:174:LEU:HD23	1:A:242:ALA:HB2	1.76	0.67
1:H:451:ASP:OD1	1:H:465:SER:OG	2.13	0.67
1:A:451:ASP:OD1	1:A:465:SER:OG	2.13	0.66
1:D:174:LEU:HD11	1:D:240:LYS:HE3	1.78	0.66
1:E:451:ASP:OD1	1:E:465:SER:OG	2.13	0.66
1:E:174:LEU:HD23	1:E:242:ALA:HB2	1.76	0.66
1:G:174:LEU:HD23	1:G:242:ALA:HB2	1.77	0.66
1:F:451:ASP:OD1	1:F:465:SER:OG	2.13	0.66
1:G:174:LEU:HD11	1:G:240:LYS:HE3	1.77	0.66
1:F:78:ASN:HD22	1:F:78:ASN:C	2.03	0.66
1:B:451:ASP:OD1	1:B:465:SER:OG	2.14	0.66
1:D:99:HIS:HE1	1:D:373:ASP:OD1	1.79	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:451:ASP:OD1	1:D:465:SER:OG	2.14	0.65
1:E:174:LEU:HD11	1:E:240:LYS:HE3	1.78	0.65
1:A:99:HIS:HE1	1:A:373:ASP:OD1	1.78	0.65
1:B:174:LEU:HD11	1:B:240:LYS:HE3	1.78	0.65
1:F:174:LEU:HD11	1:F:240:LYS:HE3	1.78	0.65
1:G:451:ASP:OD1	1:G:465:SER:OG	2.13	0.65
1:H:85:GLN:HB2	3:H:622:HOH:O	1.96	0.65
1:F:99:HIS:HE1	1:F:373:ASP:OD1	1.80	0.65
1:G:63:ARG:HG3	1:G:63:ARG:HH11	1.62	0.65
1:B:63:ARG:HG3	1:B:63:ARG:HH11	1.62	0.65
1:H:87:LYS:HB3	1:H:450:LEU:O	1.96	0.65
1:A:63:ARG:HH11	1:A:63:ARG:HG3	1.60	0.65
1:E:63:ARG:HG3	1:E:63:ARG:HH11	1.62	0.65
1:G:99:HIS:HE1	1:G:373:ASP:OD1	1.80	0.65
1:C:63:ARG:HG3	1:C:63:ARG:HH11	1.61	0.64
1:C:99:HIS:HE1	1:C:373:ASP:OD1	1.80	0.64
1:H:34:GLU:O	3:H:641:HOH:O	2.15	0.64
1:H:63:ARG:HG3	1:H:63:ARG:HH11	1.61	0.64
1:B:135:ARG:HB3	3:B:712:HOH:O	1.96	0.64
1:D:63:ARG:HG3	1:D:63:ARG:HH11	1.62	0.64
1:C:174:LEU:HD11	1:C:240:LYS:HE3	1.80	0.64
1:C:174:LEU:HD23	1:C:242:ALA:HB2	1.77	0.64
1:H:406:GLN:OE1	3:H:601:HOH:O	2.15	0.64
1:E:99:HIS:HE1	1:E:373:ASP:OD1	1.81	0.64
1:H:425:ASN:N	1:H:425:ASN:HD22	1.95	0.64
1:C:274:ARG:HD2	3:C:602:HOH:O	1.91	0.64
1:H:88:TRP:NE1	1:H:453:ASP:CG	2.56	0.63
1:A:174:LEU:HD11	1:A:240:LYS:HE3	1.79	0.63
1:H:174:LEU:HD11	1:H:240:LYS:HE3	1.80	0.63
1:C:320:LEU:HD13	1:H:225:ARG:CD	2.28	0.63
1:H:472:ARG:NH2	1:H:474:ILE:HD13	2.13	0.63
1:F:335:GLN:NE2	3:F:645:HOH:O	2.32	0.62
1:H:91:PRO:HG2	1:H:423:THR:HG21	1.79	0.62
1:H:182:THR:HG22	1:H:184:ASP:H	1.65	0.62
1:G:37:ALA:HB3	1:G:81:GLN:CD	2.24	0.62
1:H:145:TRP:O	3:H:643:HOH:O	2.16	0.62
1:B:193:ARG:O	1:D:109:ARG:NH1	2.34	0.61
1:F:470:GLU:HB2	3:F:642:HOH:O	1.99	0.61
1:G:182:THR:HG22	1:G:184:ASP:H	1.65	0.61
1:F:58:PHE:CE2	1:F:60:GLU:HG2	2.36	0.61
1:C:102:ASP:HB3	3:C:657:HOH:O	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:THR:HG23	3:A:678:HOH:O	2.00	0.60
1:G:80:ARG:C	1:G:81:GLN:HG2	2.27	0.60
1:G:396:THR:CG2	3:G:635:HOH:O	2.50	0.60
1:F:75:VAL:HG22	1:F:76:HIS:N	2.17	0.60
1:C:416:MET:CE	3:C:618:HOH:O	2.48	0.59
1:F:146:ASN:CA	3:F:644:HOH:O	2.49	0.59
1:A:301:ILE:HA	1:A:327:GLU:HG3	1.85	0.59
1:A:109:ARG:NH1	1:C:193:ARG:O	2.34	0.58
1:E:182:THR:HG22	1:E:184:ASP:H	1.68	0.58
1:H:284:THR:CG2	3:H:645:HOH:O	2.46	0.58
1:B:182:THR:HG22	1:B:184:ASP:H	1.67	0.58
1:A:244:THR:OG1	1:A:284:THR:HB	2.03	0.58
1:E:81:GLN:OE1	3:E:663:HOH:O	2.17	0.57
1:H:301:ILE:HA	1:H:327:GLU:HG3	1.85	0.57
1:A:182:THR:HG22	1:A:184:ASP:H	1.69	0.57
1:C:182:THR:HG22	1:C:184:ASP:H	1.70	0.57
1:F:301:ILE:HA	1:F:327:GLU:HG3	1.86	0.57
1:H:88:TRP:CZ3	1:H:452:GLN:HA	2.39	0.57
1:H:88:TRP:NE1	1:H:453:ASP:OD1	2.37	0.57
1:B:301:ILE:HA	1:B:327:GLU:HG3	1.87	0.57
1:C:393:ARG:HG3	3:C:650:HOH:O	2.05	0.57
1:E:128:GLU:OE1	1:E:485:VAL:HG13	2.05	0.57
1:B:331:THR:HG23	1:B:336:GLN:HB2	1.86	0.57
1:G:301:ILE:HA	1:G:327:GLU:HG3	1.87	0.57
1:H:43:LEU:CA	3:H:623:HOH:O	2.53	0.56
1:H:331:THR:HG23	1:H:336:GLN:HB2	1.87	0.56
1:E:331:THR:HG23	1:E:336:GLN:HB2	1.88	0.56
1:F:76:HIS:CD2	1:F:78:ASN:ND2	2.73	0.56
1:H:46:GLY:O	1:H:420:LEU:CD2	2.54	0.56
1:C:331:THR:HG23	1:C:336:GLN:HB2	1.87	0.56
1:A:331:THR:HG23	1:A:336:GLN:HB2	1.88	0.56
1:D:182:THR:HG22	1:D:184:ASP:H	1.70	0.56
1:E:301:ILE:HA	1:E:327:GLU:HG3	1.88	0.56
1:H:144:THR:HG23	1:H:240:LYS:HE2	1.87	0.55
1:H:472:ARG:HH22	1:H:474:ILE:HD12	1.69	0.55
1:D:144:THR:HG23	1:D:240:LYS:HE2	1.87	0.55
1:D:225:ARG:CD	1:G:320:LEU:HD13	2.36	0.55
1:G:331:THR:CG2	3:G:644:HOH:O	2.41	0.55
1:A:144:THR:HG23	1:A:240:LYS:HE2	1.87	0.55
1:H:52:VAL:HG22	1:H:55:THR:OG1	2.07	0.55
1:G:128:GLU:OE1	1:G:485:VAL:HG13	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:43:LEU:N	3:H:623:HOH:O	2.38	0.55
1:D:105:MET:CG	3:D:662:HOH:O	2.28	0.55
1:F:331:THR:HG23	1:F:336:GLN:HB2	1.89	0.55
1:A:128:GLU:OE1	1:A:485:VAL:HG13	2.07	0.55
1:F:416:MET:HE1	3:F:616:HOH:O	2.03	0.55
1:C:144:THR:HG23	1:C:240:LYS:HE2	1.88	0.55
1:E:144:THR:HG23	1:E:240:LYS:HE2	1.90	0.55
1:B:144:THR:HG23	1:B:240:LYS:HE2	1.89	0.54
1:F:128:GLU:OE1	1:F:485:VAL:HG13	2.07	0.54
1:C:301:ILE:HA	1:C:327:GLU:HG3	1.90	0.54
1:H:128:GLU:OE1	1:H:485:VAL:HG13	2.07	0.54
1:D:128:GLU:OE1	1:D:485:VAL:HG13	2.07	0.54
1:D:174:LEU:CD1	1:D:240:LYS:HE3	2.37	0.54
1:G:174:LEU:CD1	1:G:240:LYS:HE3	2.37	0.54
1:G:396:THR:HG22	3:G:635:HOH:O	2.06	0.54
1:A:416:MET:HE3	3:A:629:HOH:O	2.02	0.54
1:E:105:MET:HE1	1:E:114:TYR:CG	2.43	0.54
1:F:93:LEU:HB3	3:F:643:HOH:O	2.06	0.54
1:E:174:LEU:CD1	1:E:240:LYS:HE3	2.38	0.54
1:H:105:MET:HE1	1:H:114:TYR:CG	2.43	0.54
1:C:52:VAL:HG22	1:C:55:THR:OG1	2.08	0.54
1:F:52:VAL:HG22	1:F:55:THR:OG1	2.07	0.54
1:H:144:THR:HG23	1:H:240:LYS:CE	2.38	0.54
1:D:52:VAL:HG22	1:D:55:THR:OG1	2.07	0.54
1:F:182:THR:HG22	1:F:184:ASP:H	1.71	0.54
1:H:174:LEU:CD1	1:H:240:LYS:HE3	2.38	0.54
1:A:52:VAL:HG22	1:A:55:THR:OG1	2.08	0.54
1:B:97:ASN:OD1	1:B:144:THR:HB	2.09	0.53
1:C:144:THR:HG23	1:C:240:LYS:CE	2.39	0.53
1:G:52:VAL:HG22	1:G:55:THR:OG1	2.08	0.53
1:B:62:GLY:HA3	3:B:688:HOH:O	2.07	0.53
1:C:287:GLU:HB2	3:D:666:HOH:O	2.08	0.53
1:D:331:THR:HG23	1:D:336:GLN:HB2	1.90	0.53
1:F:174:LEU:CD1	1:F:240:LYS:HE3	2.38	0.53
1:B:128:GLU:OE1	1:B:485:VAL:HG13	2.08	0.53
1:C:174:LEU:CD1	1:C:240:LYS:HE3	2.39	0.53
1:D:301:ILE:HA	1:D:327:GLU:HG3	1.90	0.53
1:G:331:THR:HG23	1:G:336:GLN:HB2	1.89	0.53
1:E:34:GLU:CG	3:E:662:HOH:O	2.29	0.53
1:A:174:LEU:CD1	1:A:240:LYS:HE3	2.38	0.53
1:B:144:THR:HG23	1:B:240:LYS:CE	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:GLU:OE1	1:C:485:VAL:HG13	2.09	0.53
1:D:144:THR:HG23	1:D:240:LYS:CE	2.39	0.53
1:F:97:ASN:OD1	1:F:144:THR:HB	2.08	0.53
1:A:63:ARG:HH11	1:A:63:ARG:CG	2.22	0.52
1:B:63:ARG:HH11	1:B:63:ARG:CG	2.23	0.52
1:F:144:THR:HG23	1:F:240:LYS:HE2	1.90	0.52
1:H:63:ARG:HH11	1:H:63:ARG:CG	2.22	0.52
1:B:174:LEU:CD1	1:B:240:LYS:HE3	2.39	0.52
1:F:58:PHE:CZ	1:F:75:VAL:HG11	2.37	0.52
1:A:144:THR:HG23	1:A:240:LYS:CE	2.39	0.52
1:F:63:ARG:HH11	1:F:63:ARG:CG	2.21	0.52
1:E:52:VAL:HG22	1:E:55:THR:OG1	2.09	0.52
1:E:171:ALA:HB2	1:E:238:PHE:CE1	2.45	0.52
1:F:155:ARG:HG3	1:F:156:ASP:N	2.25	0.52
1:C:63:ARG:HH11	1:C:63:ARG:CG	2.23	0.52
1:B:52:VAL:HG22	1:B:55:THR:OG1	2.08	0.52
1:G:75:VAL:CG2	1:G:81:GLN:HE22	2.19	0.52
1:B:155:ARG:HG3	1:B:156:ASP:N	2.25	0.51
1:C:401:HIS:CE1	3:C:658:HOH:O	2.63	0.51
1:G:144:THR:HG23	1:G:240:LYS:HE2	1.91	0.51
1:D:155:ARG:HG3	1:D:156:ASP:N	2.25	0.51
1:A:135:ARG:HD2	3:A:693:HOH:O	2.10	0.51
1:A:256:PRO:HB3	1:A:284:THR:HG22	1.92	0.51
1:D:63:ARG:HH11	1:D:63:ARG:CG	2.24	0.51
1:E:144:THR:HG23	1:E:240:LYS:CE	2.41	0.51
1:G:63:ARG:HH11	1:G:63:ARG:CG	2.23	0.51
1:A:135:ARG:HB3	3:A:693:HOH:O	2.09	0.51
1:A:184:ASP:CG	3:A:696:HOH:O	2.53	0.51
1:D:416:MET:HE3	3:D:625:HOH:O	2.05	0.51
1:H:469:LYS:O	1:H:472:ARG:HB2	2.11	0.51
1:D:393:ARG:HG3	3:D:646:HOH:O	2.10	0.50
1:E:63:ARG:HH11	1:E:63:ARG:CG	2.24	0.50
1:H:44:ILE:HG23	1:H:416:MET:HE1	1.93	0.50
1:H:88:TRP:CD2	1:H:453:ASP:HA	2.46	0.50
1:G:144:THR:HG23	1:G:240:LYS:CE	2.40	0.50
1:F:144:THR:HG23	1:F:240:LYS:CE	2.41	0.50
1:F:146:ASN:CB	3:F:644:HOH:O	2.60	0.50
1:H:84:ALA:HB1	3:H:636:HOH:O	2.11	0.50
1:C:105:MET:HE1	1:C:114:TYR:CG	2.47	0.50
1:G:97:ASN:OD1	1:G:144:THR:HB	2.11	0.50
1:C:97:ASN:OD1	1:C:144:THR:HB	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:105:MET:HE1	1:D:114:TYR:CG	2.46	0.50
1:F:119:GLU:HA	1:F:122:TYR:CZ	2.47	0.50
1:A:171:ALA:HB2	1:A:238:PHE:CE1	2.47	0.50
1:F:308:GLN:HE22	1:F:353:VAL:H	1.58	0.50
1:H:43:LEU:HB2	3:H:623:HOH:O	2.11	0.50
1:D:416:MET:HE1	3:D:625:HOH:O	2.08	0.50
1:G:105:MET:HE1	1:G:114:TYR:CG	2.47	0.50
1:G:174:LEU:HD11	1:G:240:LYS:CE	2.42	0.50
1:A:105:MET:HE1	1:A:114:TYR:CG	2.46	0.49
1:E:42:THR:N	3:E:664:HOH:O	2.45	0.49
1:F:105:MET:HE1	1:F:114:TYR:CG	2.47	0.49
1:C:102:ASP:CB	3:C:657:HOH:O	2.59	0.49
1:H:331:THR:CG2	1:H:336:GLN:HB2	2.43	0.49
1:A:99:HIS:CE1	1:A:373:ASP:OD1	2.64	0.49
1:B:105:MET:HE1	1:B:114:TYR:CG	2.48	0.49
1:E:214:ARG:HD3	3:E:602:HOH:O	2.12	0.49
1:H:155:ARG:HG3	1:H:156:ASP:N	2.28	0.49
1:B:240:LYS:NZ	3:B:656:HOH:O	2.23	0.49
1:C:425:ASN:N	1:C:425:ASN:HD22	2.08	0.49
1:E:97:ASN:OD1	1:E:144:THR:HB	2.11	0.49
1:H:97:ASN:OD1	1:H:144:THR:HB	2.13	0.49
1:A:97:ASN:OD1	1:A:144:THR:HB	2.12	0.49
1:E:331:THR:CG2	1:E:336:GLN:HB2	2.43	0.48
1:G:155:ARG:HG3	1:G:156:ASP:N	2.28	0.48
1:G:302:HIS:CB	3:G:654:HOH:O	2.61	0.48
1:G:416:MET:HE1	3:G:612:HOH:O	2.07	0.48
1:G:119:GLU:HA	1:G:122:TYR:CZ	2.49	0.48
1:A:191:GLN:NE2	3:A:697:HOH:O	2.28	0.48
1:C:331:THR:CG2	1:C:336:GLN:HB2	2.44	0.48
1:E:174:LEU:HD11	1:E:240:LYS:CE	2.43	0.48
1:D:308:GLN:HE22	1:D:353:VAL:H	1.62	0.48
1:A:416:MET:HE1	3:A:629:HOH:O	2.02	0.48
1:B:284:THR:HG21	3:B:653:HOH:O	2.14	0.48
1:G:331:THR:CG2	1:G:336:GLN:HB2	2.44	0.48
1:H:425:ASN:N	1:H:425:ASN:ND2	2.60	0.47
1:F:254:ARG:NH1	1:F:348:ASP:OD2	2.44	0.47
1:D:99:HIS:CE1	1:D:373:ASP:OD1	2.65	0.47
1:G:79:ILE:HG22	1:G:81:GLN:NE2	2.30	0.47
1:E:99:HIS:CE1	1:E:373:ASP:OD1	2.66	0.47
1:C:352:ARG:HD2	1:C:352:ARG:N	2.28	0.47
1:E:253:PHE:O	1:E:307:GLY:HA2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:174:LEU:HD11	1:F:240:LYS:CE	2.43	0.47
1:G:308:GLN:HE22	1:G:353:VAL:H	1.63	0.47
1:H:416:MET:HA	3:H:637:HOH:O	2.14	0.47
1:A:44:ILE:HG23	1:A:416:MET:HE1	1.95	0.47
1:C:99:HIS:CE1	1:C:373:ASP:OD1	2.64	0.47
1:C:308:GLN:HE22	1:C:353:VAL:H	1.63	0.47
1:E:44:ILE:HG23	1:E:416:MET:HE1	1.97	0.47
1:C:254:ARG:NH1	1:C:348:ASP:OD2	2.46	0.46
1:D:97:ASN:OD1	1:D:144:THR:HB	2.14	0.46
1:G:284:THR:CG2	3:G:621:HOH:O	2.60	0.46
1:B:331:THR:CG2	1:B:336:GLN:HB2	2.44	0.46
1:C:119:GLU:HA	1:C:122:TYR:CZ	2.51	0.46
1:D:174:LEU:HD11	1:D:240:LYS:CE	2.44	0.46
1:F:78:ASN:C	1:F:78:ASN:ND2	2.73	0.46
1:H:119:GLU:HA	1:H:122:TYR:CZ	2.51	0.46
1:A:174:LEU:HD11	1:A:240:LYS:CE	2.45	0.46
1:A:331:THR:CG2	1:A:336:GLN:HB2	2.45	0.46
1:B:308:GLN:HE22	1:B:353:VAL:H	1.63	0.46
1:C:155:ARG:HG3	1:C:156:ASP:N	2.30	0.46
1:E:41:GLY:C	3:E:664:HOH:O	2.59	0.46
1:G:328:VAL:HA	3:G:654:HOH:O	2.15	0.45
1:H:88:TRP:CE2	1:H:453:ASP:CA	2.99	0.45
1:A:119:GLU:HA	1:A:122:TYR:CZ	2.52	0.45
1:H:419:ILE:HG12	1:H:419:ILE:H	1.58	0.45
1:A:229:ARG:NH1	3:A:656:HOH:O	2.45	0.45
1:D:44:ILE:HG23	1:D:416:MET:HE1	1.99	0.45
1:D:119:GLU:HA	1:D:122:TYR:CZ	2.52	0.45
1:G:99:HIS:CE1	1:G:373:ASP:OD1	2.65	0.45
1:B:174:LEU:HD11	1:B:240:LYS:CE	2.45	0.45
1:E:308:GLN:HE22	1:E:353:VAL:H	1.65	0.45
1:H:46:GLY:O	1:H:420:LEU:HD21	2.17	0.45
1:D:331:THR:CG2	1:D:336:GLN:HB2	2.46	0.45
1:F:253:PHE:O	1:F:307:GLY:HA2	2.17	0.45
1:H:174:LEU:HD11	1:H:240:LYS:CE	2.45	0.45
1:H:62:GLY:N	3:H:641:HOH:O	2.49	0.45
1:B:119:GLU:HA	1:B:122:TYR:CZ	2.52	0.45
1:B:352:ARG:HD2	1:B:352:ARG:N	2.32	0.45
1:H:34:GLU:CG	1:H:471:GLY:C	2.88	0.45
1:H:99:HIS:CE1	1:H:373:ASP:OD1	2.63	0.45
1:F:75:VAL:HG22	1:F:76:HIS:H	1.81	0.44
1:F:331:THR:HB	1:F:350:SER:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:254:ARG:NH1	1:E:348:ASP:OD2	2.47	0.44
1:E:284:THR:HG23	3:E:661:HOH:O	2.11	0.44
1:F:75:VAL:CG2	1:F:76:HIS:N	2.80	0.44
1:C:174:LEU:HD11	1:C:240:LYS:CE	2.46	0.44
1:B:44:ILE:HG23	1:B:416:MET:HE1	2.00	0.44
1:C:44:ILE:HG23	1:C:416:MET:HE1	1.99	0.44
1:F:331:THR:CG2	1:F:336:GLN:HB2	2.46	0.44
1:A:308:GLN:HE22	1:A:353:VAL:H	1.65	0.44
1:A:50:VAL:CG1	1:E:415:PRO:HD2	2.47	0.44
1:B:393:ARG:HG3	3:B:611:HOH:O	2.17	0.44
1:F:425:ASN:N	1:F:425:ASN:HD22	2.14	0.44
1:D:253:PHE:O	1:D:307:GLY:HA2	2.18	0.44
1:D:331:THR:HB	1:D:350:SER:HB2	1.99	0.44
1:C:171:ALA:HB2	1:C:238:PHE:CE2	2.53	0.44
1:E:58:PHE:HB3	1:E:66:LYS:HB2	2.00	0.44
1:H:253:PHE:O	1:H:307:GLY:HA2	2.17	0.44
1:H:331:THR:HB	1:H:350:SER:HB2	1.99	0.44
1:H:468:VAL:HG13	1:H:472:ARG:N	2.32	0.44
1:E:119:GLU:HA	1:E:122:TYR:CZ	2.52	0.43
1:C:253:PHE:O	1:C:307:GLY:HA2	2.17	0.43
1:E:331:THR:HB	1:E:350:SER:HB2	2.00	0.43
1:F:44:ILE:HG23	1:F:416:MET:HE1	1.99	0.43
1:G:182:THR:CG2	1:G:184:ASP:H	2.30	0.43
1:B:58:PHE:HB3	1:B:66:LYS:HB2	2.01	0.43
1:B:331:THR:HB	1:B:350:SER:HB2	2.00	0.43
1:H:416:MET:CA	3:H:637:HOH:O	2.66	0.43
1:A:155:ARG:HG3	1:A:156:ASP:N	2.32	0.43
1:A:254:ARG:NH1	1:A:348:ASP:OD2	2.46	0.43
1:F:171:ALA:HB2	1:F:238:PHE:CE2	2.54	0.43
1:A:331:THR:HB	1:A:350:SER:HB2	2.00	0.43
1:C:331:THR:HB	1:C:350:SER:HB2	2.01	0.43
1:F:76:HIS:HD2	1:F:78:ASN:ND2	2.17	0.43
1:E:155:ARG:HG3	1:E:156:ASP:N	2.33	0.43
1:G:393:ARG:HG3	3:G:634:HOH:O	2.18	0.43
1:D:254:ARG:NH1	1:D:348:ASP:OD2	2.46	0.43
1:B:182:THR:CG2	1:B:184:ASP:H	2.32	0.42
1:F:352:ARG:N	1:F:352:ARG:HD2	2.34	0.42
1:C:377:THR:OG1	1:C:393:ARG:NH2	2.53	0.42
1:A:482:SER:O	1:A:483:PRO:C	2.60	0.42
1:C:58:PHE:HB3	1:C:66:LYS:HB2	2.01	0.42
1:C:240:LYS:NZ	3:C:645:HOH:O	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:482:SER:O	1:E:483:PRO:C	2.62	0.42
1:H:88:TRP:HE1	1:H:453:ASP:CG	2.25	0.42
1:A:253:PHE:O	1:A:307:GLY:HA2	2.19	0.42
1:B:253:PHE:O	1:B:307:GLY:HA2	2.20	0.42
1:G:88:TRP:CZ2	1:G:453:ASP:HB2	2.54	0.42
1:B:482:SER:O	1:B:483:PRO:C	2.58	0.42
1:D:88:TRP:CZ2	1:D:453:ASP:HB2	2.55	0.42
1:G:331:THR:HB	1:G:350:SER:HB2	2.02	0.42
1:D:171:ALA:HB2	1:D:238:PHE:CE2	2.55	0.42
1:F:99:HIS:CE1	1:F:373:ASP:OD1	2.66	0.42
1:G:302:HIS:CA	3:G:654:HOH:O	2.36	0.42
1:H:182:THR:CG2	1:H:184:ASP:H	2.30	0.42
1:B:254:ARG:NH1	1:B:348:ASP:OD2	2.46	0.41
1:E:182:THR:CG2	1:E:183:GLY:N	2.81	0.41
1:F:73:ILE:HG22	1:F:74:GLU:N	2.35	0.41
1:G:171:ALA:HB2	1:G:238:PHE:CE2	2.55	0.41
1:G:416:MET:HE3	3:G:612:HOH:O	2.11	0.41
1:H:308:GLN:HE22	1:H:352:ARG:HB2	1.84	0.41
1:B:182:THR:CG2	1:B:183:GLY:N	2.83	0.41
1:E:88:TRP:CZ2	1:E:453:ASP:HB2	2.56	0.41
1:A:414:ASP:OD1	1:E:50:VAL:HG12	2.20	0.41
1:H:88:TRP:CE2	1:H:453:ASP:HA	2.55	0.41
1:D:352:ARG:N	1:D:352:ARG:HD2	2.34	0.41
1:A:182:THR:CG2	1:A:183:GLY:N	2.84	0.41
1:E:308:GLN:HE22	1:E:352:ARG:HB2	1.85	0.41
1:G:182:THR:HG22	1:G:184:ASP:N	2.34	0.41
1:A:84:ALA:O	1:A:87:LYS:HB2	2.21	0.41
1:B:171:ALA:HB2	1:B:238:PHE:CE2	2.55	0.41
1:D:182:THR:CG2	1:D:184:ASP:H	2.34	0.41
1:F:146:ASN:CG	3:F:644:HOH:O	2.63	0.41
1:H:363:LYS:HA	1:H:363:LYS:HD2	1.93	0.41
1:C:182:THR:CG2	1:C:184:ASP:H	2.34	0.41
1:D:482:SER:O	1:D:483:PRO:C	2.63	0.41
1:F:58:PHE:HB3	1:F:66:LYS:HB2	2.02	0.41
1:G:363:LYS:HA	1:G:363:LYS:HD2	1.94	0.41
1:H:254:ARG:NH1	1:H:348:ASP:OD2	2.46	0.41
1:A:182:THR:CG2	1:A:184:ASP:H	2.32	0.41
1:A:352:ARG:HD2	1:A:352:ARG:N	2.36	0.41
1:G:79:ILE:O	1:G:81:GLN:HG2	2.21	0.41
1:H:482:SER:O	1:H:483:PRO:C	2.64	0.41
1:E:182:THR:HG23	1:E:183:GLY:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:363:LYS:HA	1:F:363:LYS:HD2	1.93	0.41
1:G:58:PHE:HB3	1:G:66:LYS:HB2	2.02	0.41
1:C:425:ASN:N	1:C:425:ASN:ND2	2.69	0.40
1:G:44:ILE:HG23	1:G:416:MET:HE1	2.02	0.40
1:H:58:PHE:HB3	1:H:66:LYS:HB2	2.03	0.40
1:H:404:TRP:CH2	1:H:408:MET:HG3	2.56	0.40
1:E:363:LYS:HA	1:E:363:LYS:HD2	1.91	0.40
1:G:482:SER:O	1:G:483:PRO:C	2.61	0.40
1:A:155:ARG:HD2	3:A:695:HOH:O	2.20	0.40
1:F:88:TRP:CZ2	1:F:453:ASP:HB2	2.56	0.40
1:G:308:GLN:HE22	1:G:352:ARG:HB2	1.86	0.40
1:A:58:PHE:HB3	1:A:66:LYS:HB2	2.03	0.40
1:C:416:MET:HE1	3:C:618:HOH:O	2.17	0.40
1:H:425:ASN:HB2	1:H:426:PRO:HD3	2.03	0.40

All (4) 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:E:611:HOH:O	3:G:601:HOH:O[3_455]	1.49	0.71
3:E:602:HOH:O	3:F:603:HOH:O[4_558]	1.53	0.67
1:H:452:GLN:NE2	1:H:462:ARG:O[2_657]	1.87	0.33
1:F:74:GLU:O	1:G:74:GLU:OE2[1_565]	2.09	0.11

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	465/496 (94%)	454 (98%)	11 (2%)	0	100	100
1	B	466/496 (94%)	453 (97%)	13 (3%)	0	100	100
1	C	465/496 (94%)	452 (97%)	13 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	465/496 (94%)	453 (97%)	12 (3%)	0	100	100
1	E	465/496 (94%)	451 (97%)	14 (3%)	0	100	100
1	F	465/496 (94%)	451 (97%)	14 (3%)	0	100	100
1	G	465/496 (94%)	449 (97%)	16 (3%)	0	100	100
1	H	465/496 (94%)	450 (97%)	15 (3%)	0	100	100
All	All	3721/3968 (94%)	3613 (97%)	108 (3%)	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	382/407 (94%)	363 (95%)	19 (5%)	22	30
1	B	383/407 (94%)	364 (95%)	19 (5%)	22	30
1	C	382/407 (94%)	362 (95%)	20 (5%)	21	28
1	D	382/407 (94%)	362 (95%)	20 (5%)	21	28
1	E	382/407 (94%)	362 (95%)	20 (5%)	21	28
1	F	382/407 (94%)	361 (94%)	21 (6%)	19	26
1	G	382/407 (94%)	362 (95%)	20 (5%)	21	28
1	H	382/407 (94%)	357 (94%)	25 (6%)	15	20
All	All	3057/3256 (94%)	2893 (95%)	164 (5%)	20	27

All (164) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	VAL
1	A	52	VAL
1	A	63	ARG
1	A	69	SER
1	A	74	GLU

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Mol	Chain	Res	Type
1	A	75	VAL
1	A	85	GLN
1	A	144	THR
1	A	155	ARG
1	A	174	LEU
1	A	182	THR
1	A	254	ARG
1	A	273	ARG
1	A	284	THR
1	A	327	GLU
1	A	331	THR
1	A	363	LYS
1	A	485	VAL
1	A	491	THR
1	B	30	PHE
1	B	39	VAL
1	B	52	VAL
1	B	63	ARG
1	B	69	SER
1	B	72	GLN
1	B	74	GLU
1	B	75	VAL
1	B	144	THR
1	B	155	ARG
1	B	174	LEU
1	B	182	THR
1	B	254	ARG
1	B	273	ARG
1	B	284	THR
1	B	327	GLU
1	B	363	LYS
1	B	485	VAL
1	B	491	THR
1	C	30	PHE
1	C	39	VAL
1	C	52	VAL
1	C	63	ARG
1	C	69	SER
1	C	72	GLN
1	C	74	GLU
1	C	75	VAL
1	C	144	THR

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Mol	Chain	Res	Type
1	C	155	ARG
1	C	174	LEU
1	C	182	THR
1	C	254	ARG
1	C	273	ARG
1	C	284	THR
1	C	327	GLU
1	C	331	THR
1	C	363	LYS
1	C	485	VAL
1	C	491	THR
1	D	30	PHE
1	D	39	VAL
1	D	52	VAL
1	D	63	ARG
1	D	69	SER
1	D	72	GLN
1	D	74	GLU
1	D	75	VAL
1	D	144	THR
1	D	155	ARG
1	D	174	LEU
1	D	182	THR
1	D	254	ARG
1	D	273	ARG
1	D	284	THR
1	D	327	GLU
1	D	331	THR
1	D	363	LYS
1	D	485	VAL
1	D	491	THR
1	E	30	PHE
1	E	39	VAL
1	E	52	VAL
1	E	63	ARG
1	E	69	SER
1	E	72	GLN
1	E	74	GLU
1	E	75	VAL
1	E	144	THR
1	E	155	ARG
1	E	174	LEU

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Mol	Chain	Res	Type
1	E	182	THR
1	E	254	ARG
1	E	273	ARG
1	E	284	THR
1	E	327	GLU
1	E	331	THR
1	E	363	LYS
1	E	485	VAL
1	E	491	THR
1	F	30	PHE
1	F	50	VAL
1	F	52	VAL
1	F	63	ARG
1	F	69	SER
1	F	74	GLU
1	F	76	HIS
1	F	78	ASN
1	F	144	THR
1	F	155	ARG
1	F	174	LEU
1	F	182	THR
1	F	254	ARG
1	F	273	ARG
1	F	284	THR
1	F	327	GLU
1	F	331	THR
1	F	363	LYS
1	F	423	THR
1	F	485	VAL
1	F	491	THR
1	G	30	PHE
1	G	52	VAL
1	G	63	ARG
1	G	69	SER
1	G	72	GLN
1	G	74	GLU
1	G	75	VAL
1	G	144	THR
1	G	155	ARG
1	G	174	LEU
1	G	182	THR
1	G	254	ARG

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Mol	Chain	Res	Type
1	G	273	ARG
1	G	284	THR
1	G	308	GLN
1	G	327	GLU
1	G	331	THR
1	G	363	LYS
1	G	485	VAL
1	G	491	THR
1	H	30	PHE
1	H	39	VAL
1	H	52	VAL
1	H	63	ARG
1	H	69	SER
1	H	72	GLN
1	H	74	GLU
1	H	75	VAL
1	H	87	LYS
1	H	144	THR
1	H	155	ARG
1	H	174	LEU
1	H	182	THR
1	H	254	ARG
1	H	273	ARG
1	H	284	THR
1	H	327	GLU
1	H	363	LYS
1	H	419	ILE
1	H	423	THR
1	H	425	ASN
1	H	470	GLU
1	H	472	ARG
1	H	485	VAL
1	H	491	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (75) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	78	ASN
1	A	99	HIS
1	A	191	GLN
1	A	308	GLN

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Mol	Chain	Res	Type
1	A	329	GLN
1	A	335	GLN
1	A	336	GLN
1	A	391	HIS
1	A	425	ASN
1	B	47	ASN
1	B	78	ASN
1	B	191	GLN
1	B	308	GLN
1	B	329	GLN
1	B	335	GLN
1	B	336	GLN
1	B	425	ASN
1	C	47	ASN
1	C	78	ASN
1	C	99	HIS
1	C	191	GLN
1	C	308	GLN
1	C	329	GLN
1	C	335	GLN
1	C	336	GLN
1	C	425	ASN
1	D	47	ASN
1	D	78	ASN
1	D	99	HIS
1	D	191	GLN
1	D	308	GLN
1	D	329	GLN
1	D	335	GLN
1	D	336	GLN
1	D	391	HIS
1	D	425	ASN
1	E	47	ASN
1	E	78	ASN
1	E	99	HIS
1	E	191	GLN
1	E	308	GLN
1	E	329	GLN
1	E	335	GLN
1	E	336	GLN
1	E	425	ASN
1	F	47	ASN

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Mol	Chain	Res	Type
1	F	76	HIS
1	F	78	ASN
1	F	99	HIS
1	F	191	GLN
1	F	308	GLN
1	F	329	GLN
1	F	335	GLN
1	F	336	GLN
1	F	425	ASN
1	G	47	ASN
1	G	78	ASN
1	G	99	HIS
1	G	191	GLN
1	G	308	GLN
1	G	329	GLN
1	G	336	GLN
1	G	425	ASN
1	H	47	ASN
1	H	78	ASN
1	H	99	HIS
1	H	191	GLN
1	H	302	HIS
1	H	308	GLN
1	H	329	GLN
1	H	335	GLN
1	H	336	GLN
1	H	406	GLN
1	H	425	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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

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	467/496 (94%)	-0.24	6 (1%) 75 76	27, 37, 62, 122	0
1	B	467/496 (94%)	-0.19	3 (0%) 85 86	27, 39, 60, 105	1 (0%)
1	C	467/496 (94%)	-0.08	6 (1%) 75 76	28, 42, 70, 104	0
1	D	467/496 (94%)	-0.04	5 (1%) 78 79	27, 43, 65, 98	0
1	E	467/496 (94%)	-0.10	6 (1%) 75 76	32, 42, 67, 105	0
1	F	467/496 (94%)	0.33	18 (3%) 43 45	32, 53, 80, 126	0
1	G	467/496 (94%)	0.28	17 (3%) 46 48	34, 49, 81, 133	0
1	H	467/496 (94%)	1.48	155 (33%) 1 1	32, 60, 108, 153	0
All	All	3736/3968 (94%)	0.18	216 (5%) 29 30	27, 45, 82, 153	1 (0%)

All (216) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	439	ILE	7.6
1	F	75	VAL	6.7
1	H	44	ILE	6.1
1	H	46	GLY	6.1
1	G	30	PHE	5.9
1	A	70	THR	5.8
1	H	57	VAL	5.6
1	H	89	ILE	5.5
1	H	43	LEU	5.4
1	H	429	ALA	5.4
1	H	73	ILE	5.3
1	H	67	VAL	5.3
1	H	30	PHE	5.3
1	H	68	GLY	5.1
1	H	36	ILE	5.1
1	H	464	LEU	5.0

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Mol	Chain	Res	Type	RSRZ
1	H	450	LEU	5.0
1	H	50	VAL	4.9
1	F	73	ILE	4.9
1	F	30	PHE	4.9
1	H	38	ILE	4.9
1	H	441	VAL	4.8
1	H	420	LEU	4.6
1	H	445	GLY	4.6
1	H	88	TRP	4.5
1	H	52	VAL	4.5
1	H	449	VAL	4.5
1	H	49	GLY	4.5
1	H	93	LEU	4.4
1	H	75	VAL	4.3
1	E	30	PHE	4.2
1	H	42	THR	4.2
1	H	434	ASP	4.2
1	H	94	VAL	4.2
1	H	325	TRP	4.2
1	D	30	PHE	4.2
1	H	404	TRP	4.1
1	H	362	ILE	4.1
1	H	412	GLY	4.0
1	H	64	ILE	4.0
1	H	471	GLY	4.0
1	H	96	GLY	3.9
1	H	58	PHE	3.9
1	H	79	ILE	3.9
1	H	459	THR	3.9
1	H	458	ILE	3.9
1	H	92	GLY	3.9
1	H	84	ALA	3.9
1	H	463	THR	3.8
1	H	39	VAL	3.8
1	H	370	LEU	3.8
1	H	65	THR	3.8
1	H	159	ALA	3.8
1	H	34	GLU	3.8
1	H	31	PHE	3.8
1	H	447	VAL	3.8
1	H	62	GLY	3.7
1	H	455	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
1	H	452	GLN	3.7
1	H	433	PHE	3.7
1	G	78	ASN	3.7
1	C	30	PHE	3.6
1	H	424	ALA	3.6
1	B	30	PHE	3.6
1	H	461	MET	3.6
1	H	56	THR	3.6
1	H	316	ILE	3.5
1	H	448	VAL	3.5
1	H	33	GLY	3.5
1	H	468	VAL	3.4
1	G	73	ILE	3.4
1	H	82	ILE	3.4
1	H	140	THR	3.4
1	F	67	VAL	3.4
1	H	454	PRO	3.3
1	H	51	PRO	3.3
1	H	80	ARG	3.3
1	H	45	ASP	3.3
1	H	139	THR	3.3
1	C	75	VAL	3.2
1	H	137	GLY	3.2
1	H	69	SER	3.2
1	H	235	GLY	3.2
1	H	411	LYS	3.2
1	H	76	HIS	3.2
1	B	496	ASP	3.1
1	A	77	PRO	3.1
1	H	90	LEU	3.1
1	H	35	THR	3.1
1	A	71	ASP	3.1
1	E	496	ASP	3.1
1	H	371	GLY	3.1
1	H	474	ILE	3.1
1	H	54	GLU	3.0
1	H	438	SER	3.0
1	G	496	ASP	3.0
1	H	457	ASP	3.0
1	H	53	PRO	3.0
1	H	169	PHE	3.0
1	F	76	HIS	3.0

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Mol	Chain	Res	Type	RSRZ
1	H	472	ARG	3.0
1	G	307	GLY	3.0
1	A	30	PHE	2.9
1	H	63	ARG	2.9
1	H	37	ALA	2.9
1	H	442	GLY	2.9
1	H	324	LEU	2.9
1	H	406	GLN	2.9
1	H	358	ASP	2.9
1	H	70	THR	2.9
1	H	467	VAL	2.8
1	G	76	HIS	2.8
1	D	77	PRO	2.8
1	H	320	LEU	2.8
1	D	496	ASP	2.7
1	H	453	ASP	2.7
1	H	77	PRO	2.7
1	H	431	ARG	2.7
1	H	444	LEU	2.7
1	H	427	ALA	2.7
1	F	71	ASP	2.7
1	A	73	ILE	2.7
1	H	61	ASP	2.7
1	H	359	VAL	2.7
1	H	59	ILE	2.7
1	F	50	VAL	2.6
1	H	369	VAL	2.6
1	H	466	HIS	2.6
1	H	292	ALA	2.6
1	E	77	PRO	2.6
1	H	297	ALA	2.6
1	F	70	THR	2.6
1	D	80	ARG	2.5
1	F	496	ASP	2.5
1	H	396	THR	2.5
1	H	423	THR	2.5
1	G	75	VAL	2.5
1	H	74	GLU	2.5
1	F	495	LEU	2.5
1	H	496	ASP	2.5
1	H	430	TYR	2.5
1	H	91	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	413	MET	2.5
1	H	66	LYS	2.5
1	C	496	ASP	2.5
1	H	161	GLY	2.5
1	H	304	THR	2.4
1	H	465	SER	2.4
1	H	299	LEU	2.4
1	H	41	GLY	2.4
1	E	70	THR	2.4
1	H	405	MET	2.4
1	H	469	LYS	2.4
1	H	72	GLN	2.4
1	F	49	GLY	2.4
1	H	272	VAL	2.4
1	H	446	ASP	2.3
1	H	331	THR	2.3
1	H	238	PHE	2.3
1	H	436	LEU	2.3
1	F	77	PRO	2.3
1	H	415	PRO	2.3
1	H	48	GLY	2.3
1	H	437	GLY	2.3
1	H	456	ALA	2.3
1	F	74	GLU	2.3
1	H	418	ALA	2.3
1	G	59	ILE	2.3
1	H	298	ASP	2.3
1	H	276	GLY	2.3
1	H	152	THR	2.3
1	G	32	GLU	2.3
1	H	387	GLN	2.3
1	H	399	GLU	2.3
1	H	419	ILE	2.3
1	F	441	VAL	2.3
1	B	77	PRO	2.2
1	D	76	HIS	2.2
1	A	72	GLN	2.2
1	E	73	ILE	2.2
1	H	60	GLU	2.2
1	H	55	THR	2.2
1	G	33	GLY	2.2
1	H	326	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	47	ASN	2.2
1	H	83	ASP	2.2
1	G	77	PRO	2.2
1	H	86	GLY	2.2
1	H	408	MET	2.1
1	H	81	GLN	2.1
1	H	401	HIS	2.1
1	F	89	ILE	2.1
1	F	57	VAL	2.1
1	G	50	VAL	2.1
1	G	57	VAL	2.1
1	F	31	PHE	2.1
1	G	80	ARG	2.1
1	H	167	ARG	2.1
1	H	416	MET	2.1
1	C	79	ILE	2.1
1	H	368	LEU	2.1
1	H	95	ASN	2.1
1	H	460	ASN	2.1
1	C	77	PRO	2.1
1	G	58	PHE	2.1
1	G	43	LEU	2.1
1	H	78	ASN	2.1
1	H	426	PRO	2.1
1	H	142	PHE	2.1
1	H	319	LEU	2.0
1	H	451	ASP	2.0
1	C	31	PHE	2.0
1	F	58	PHE	2.0
1	E	491	THR	2.0
1	G	70	THR	2.0
1	H	425	ASN	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	ZN	D	501	1/1	0.87	0.10	56,56,56,56	1
2	ZN	C	501	1/1	0.89	0.09	65,65,65,65	1
2	ZN	H	501	1/1	0.89	0.09	67,67,67,67	1
2	ZN	G	501	1/1	0.91	0.11	55,55,55,55	1
2	ZN	E	501	1/1	0.92	0.09	60,60,60,60	1
2	ZN	F	501	1/1	0.93	0.08	63,63,63,63	1
2	ZN	A	501	1/1	0.95	0.05	48,48,48,48	1
2	ZN	B	501	1/1	0.97	0.07	55,55,55,55	1

6.5 Other polymers [i](#)

There are no such residues in this entry.