



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:43 AM UTC

PDB ID : 4H1Z / pdb_00004h1z
Title : Crystal structure of putative isomerase from *Sinorhizobium meliloti*, open loop conformation (target EFI-502104)
Authors : Patskovsky, Y.; Toro, R.; Bhosle, R.; Hillerich, B.; Seidel, R.D.; Washington, E.; Scott Glenn, A.; Chowdhury, S.; Evans, B.; Hammonds, J.; Zencheck, W.D.; Imker, H.J.; Gerlt, J.A.; Almo, S.C.; Enzyme Function Initiative (EFI)
Deposited on : 2012-09-11
Resolution : 2.01 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

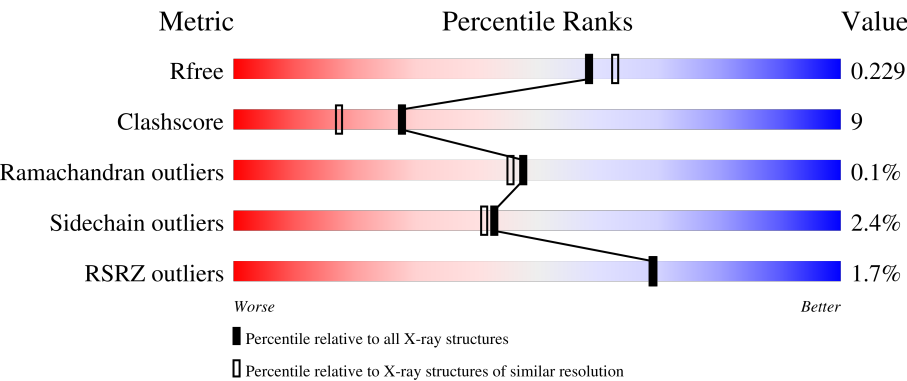
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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.



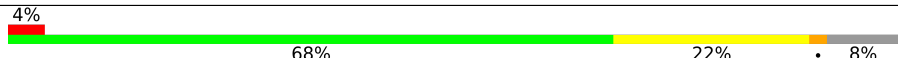
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	412	2% 72% 19% 9%
1	B	412	2% 77% 16% 6%
1	C	412	% 73% 17% 8%
1	D	412	% 76% 15% 8%

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Mol	Chain	Length	Quality of chain
1	E	412	
1	F	412	
1	G	412	
1	H	412	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	B	404	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 24833 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 Enolase Q92Zs5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	0
			2880	1824	517	526	13			
1	B	388	Total	C	N	O	S	0	1	0
			2982	1892	535	542	13			
1	C	378	Total	C	N	O	S	0	1	0
			2909	1843	525	528	13			
1	D	379	Total	C	N	O	S	0	1	0
			2913	1845	526	529	13			
1	E	381	Total	C	N	O	S	0	1	0
			2926	1854	528	531	13			
1	F	374	Total	C	N	O	S	0	4	0
			2880	1829	515	523	13			
1	G	380	Total	C	N	O	S	0	1	0
			2916	1849	524	530	13			
1	H	378	Total	C	N	O	S	0	0	0
			2891	1833	519	526	13			

There are 184 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP Q92ZS5
A	-21	HIS	-	expression tag	UNP Q92ZS5
A	-20	HIS	-	expression tag	UNP Q92ZS5
A	-19	HIS	-	expression tag	UNP Q92ZS5
A	-18	HIS	-	expression tag	UNP Q92ZS5
A	-17	HIS	-	expression tag	UNP Q92ZS5
A	-16	HIS	-	expression tag	UNP Q92ZS5
A	-15	SER	-	expression tag	UNP Q92ZS5
A	-14	SER	-	expression tag	UNP Q92ZS5
A	-13	GLY	-	expression tag	UNP Q92ZS5
A	-12	VAL	-	expression tag	UNP Q92ZS5
A	-11	ASP	-	expression tag	UNP Q92ZS5
A	-10	LEU	-	expression tag	UNP Q92ZS5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	GLY	-	expression tag	UNP Q92ZS5
A	-8	THR	-	expression tag	UNP Q92ZS5
A	-7	GLU	-	expression tag	UNP Q92ZS5
A	-6	ASN	-	expression tag	UNP Q92ZS5
A	-5	LEU	-	expression tag	UNP Q92ZS5
A	-4	TYR	-	expression tag	UNP Q92ZS5
A	-3	PHE	-	expression tag	UNP Q92ZS5
A	-2	GLN	-	expression tag	UNP Q92ZS5
A	-1	SER	-	expression tag	UNP Q92ZS5
A	0	MET	-	expression tag	UNP Q92ZS5
B	-22	MET	-	expression tag	UNP Q92ZS5
B	-21	HIS	-	expression tag	UNP Q92ZS5
B	-20	HIS	-	expression tag	UNP Q92ZS5
B	-19	HIS	-	expression tag	UNP Q92ZS5
B	-18	HIS	-	expression tag	UNP Q92ZS5
B	-17	HIS	-	expression tag	UNP Q92ZS5
B	-16	HIS	-	expression tag	UNP Q92ZS5
B	-15	SER	-	expression tag	UNP Q92ZS5
B	-14	SER	-	expression tag	UNP Q92ZS5
B	-13	GLY	-	expression tag	UNP Q92ZS5
B	-12	VAL	-	expression tag	UNP Q92ZS5
B	-11	ASP	-	expression tag	UNP Q92ZS5
B	-10	LEU	-	expression tag	UNP Q92ZS5
B	-9	GLY	-	expression tag	UNP Q92ZS5
B	-8	THR	-	expression tag	UNP Q92ZS5
B	-7	GLU	-	expression tag	UNP Q92ZS5
B	-6	ASN	-	expression tag	UNP Q92ZS5
B	-5	LEU	-	expression tag	UNP Q92ZS5
B	-4	TYR	-	expression tag	UNP Q92ZS5
B	-3	PHE	-	expression tag	UNP Q92ZS5
B	-2	GLN	-	expression tag	UNP Q92ZS5
B	-1	SER	-	expression tag	UNP Q92ZS5
B	0	MET	-	expression tag	UNP Q92ZS5
C	-22	MET	-	expression tag	UNP Q92ZS5
C	-21	HIS	-	expression tag	UNP Q92ZS5
C	-20	HIS	-	expression tag	UNP Q92ZS5
C	-19	HIS	-	expression tag	UNP Q92ZS5
C	-18	HIS	-	expression tag	UNP Q92ZS5
C	-17	HIS	-	expression tag	UNP Q92ZS5
C	-16	HIS	-	expression tag	UNP Q92ZS5
C	-15	SER	-	expression tag	UNP Q92ZS5
C	-14	SER	-	expression tag	UNP Q92ZS5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-13	GLY	-	expression tag	UNP Q92ZS5
C	-12	VAL	-	expression tag	UNP Q92ZS5
C	-11	ASP	-	expression tag	UNP Q92ZS5
C	-10	LEU	-	expression tag	UNP Q92ZS5
C	-9	GLY	-	expression tag	UNP Q92ZS5
C	-8	THR	-	expression tag	UNP Q92ZS5
C	-7	GLU	-	expression tag	UNP Q92ZS5
C	-6	ASN	-	expression tag	UNP Q92ZS5
C	-5	LEU	-	expression tag	UNP Q92ZS5
C	-4	TYR	-	expression tag	UNP Q92ZS5
C	-3	PHE	-	expression tag	UNP Q92ZS5
C	-2	GLN	-	expression tag	UNP Q92ZS5
C	-1	SER	-	expression tag	UNP Q92ZS5
C	0	MET	-	expression tag	UNP Q92ZS5
D	-22	MET	-	expression tag	UNP Q92ZS5
D	-21	HIS	-	expression tag	UNP Q92ZS5
D	-20	HIS	-	expression tag	UNP Q92ZS5
D	-19	HIS	-	expression tag	UNP Q92ZS5
D	-18	HIS	-	expression tag	UNP Q92ZS5
D	-17	HIS	-	expression tag	UNP Q92ZS5
D	-16	HIS	-	expression tag	UNP Q92ZS5
D	-15	SER	-	expression tag	UNP Q92ZS5
D	-14	SER	-	expression tag	UNP Q92ZS5
D	-13	GLY	-	expression tag	UNP Q92ZS5
D	-12	VAL	-	expression tag	UNP Q92ZS5
D	-11	ASP	-	expression tag	UNP Q92ZS5
D	-10	LEU	-	expression tag	UNP Q92ZS5
D	-9	GLY	-	expression tag	UNP Q92ZS5
D	-8	THR	-	expression tag	UNP Q92ZS5
D	-7	GLU	-	expression tag	UNP Q92ZS5
D	-6	ASN	-	expression tag	UNP Q92ZS5
D	-5	LEU	-	expression tag	UNP Q92ZS5
D	-4	TYR	-	expression tag	UNP Q92ZS5
D	-3	PHE	-	expression tag	UNP Q92ZS5
D	-2	GLN	-	expression tag	UNP Q92ZS5
D	-1	SER	-	expression tag	UNP Q92ZS5
D	0	MET	-	expression tag	UNP Q92ZS5
E	-22	MET	-	expression tag	UNP Q92ZS5
E	-21	HIS	-	expression tag	UNP Q92ZS5
E	-20	HIS	-	expression tag	UNP Q92ZS5
E	-19	HIS	-	expression tag	UNP Q92ZS5
E	-18	HIS	-	expression tag	UNP Q92ZS5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-17	HIS	-	expression tag	UNP Q92ZS5
E	-16	HIS	-	expression tag	UNP Q92ZS5
E	-15	SER	-	expression tag	UNP Q92ZS5
E	-14	SER	-	expression tag	UNP Q92ZS5
E	-13	GLY	-	expression tag	UNP Q92ZS5
E	-12	VAL	-	expression tag	UNP Q92ZS5
E	-11	ASP	-	expression tag	UNP Q92ZS5
E	-10	LEU	-	expression tag	UNP Q92ZS5
E	-9	GLY	-	expression tag	UNP Q92ZS5
E	-8	THR	-	expression tag	UNP Q92ZS5
E	-7	GLU	-	expression tag	UNP Q92ZS5
E	-6	ASN	-	expression tag	UNP Q92ZS5
E	-5	LEU	-	expression tag	UNP Q92ZS5
E	-4	TYR	-	expression tag	UNP Q92ZS5
E	-3	PHE	-	expression tag	UNP Q92ZS5
E	-2	GLN	-	expression tag	UNP Q92ZS5
E	-1	SER	-	expression tag	UNP Q92ZS5
E	0	MET	-	expression tag	UNP Q92ZS5
F	-22	MET	-	expression tag	UNP Q92ZS5
F	-21	HIS	-	expression tag	UNP Q92ZS5
F	-20	HIS	-	expression tag	UNP Q92ZS5
F	-19	HIS	-	expression tag	UNP Q92ZS5
F	-18	HIS	-	expression tag	UNP Q92ZS5
F	-17	HIS	-	expression tag	UNP Q92ZS5
F	-16	HIS	-	expression tag	UNP Q92ZS5
F	-15	SER	-	expression tag	UNP Q92ZS5
F	-14	SER	-	expression tag	UNP Q92ZS5
F	-13	GLY	-	expression tag	UNP Q92ZS5
F	-12	VAL	-	expression tag	UNP Q92ZS5
F	-11	ASP	-	expression tag	UNP Q92ZS5
F	-10	LEU	-	expression tag	UNP Q92ZS5
F	-9	GLY	-	expression tag	UNP Q92ZS5
F	-8	THR	-	expression tag	UNP Q92ZS5
F	-7	GLU	-	expression tag	UNP Q92ZS5
F	-6	ASN	-	expression tag	UNP Q92ZS5
F	-5	LEU	-	expression tag	UNP Q92ZS5
F	-4	TYR	-	expression tag	UNP Q92ZS5
F	-3	PHE	-	expression tag	UNP Q92ZS5
F	-2	GLN	-	expression tag	UNP Q92ZS5
F	-1	SER	-	expression tag	UNP Q92ZS5
F	0	MET	-	expression tag	UNP Q92ZS5
G	-22	MET	-	expression tag	UNP Q92ZS5

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-21	HIS	-	expression tag	UNP Q92ZS5
G	-20	HIS	-	expression tag	UNP Q92ZS5
G	-19	HIS	-	expression tag	UNP Q92ZS5
G	-18	HIS	-	expression tag	UNP Q92ZS5
G	-17	HIS	-	expression tag	UNP Q92ZS5
G	-16	HIS	-	expression tag	UNP Q92ZS5
G	-15	SER	-	expression tag	UNP Q92ZS5
G	-14	SER	-	expression tag	UNP Q92ZS5
G	-13	GLY	-	expression tag	UNP Q92ZS5
G	-12	VAL	-	expression tag	UNP Q92ZS5
G	-11	ASP	-	expression tag	UNP Q92ZS5
G	-10	LEU	-	expression tag	UNP Q92ZS5
G	-9	GLY	-	expression tag	UNP Q92ZS5
G	-8	THR	-	expression tag	UNP Q92ZS5
G	-7	GLU	-	expression tag	UNP Q92ZS5
G	-6	ASN	-	expression tag	UNP Q92ZS5
G	-5	LEU	-	expression tag	UNP Q92ZS5
G	-4	TYR	-	expression tag	UNP Q92ZS5
G	-3	PHE	-	expression tag	UNP Q92ZS5
G	-2	GLN	-	expression tag	UNP Q92ZS5
G	-1	SER	-	expression tag	UNP Q92ZS5
G	0	MET	-	expression tag	UNP Q92ZS5
H	-22	MET	-	expression tag	UNP Q92ZS5
H	-21	HIS	-	expression tag	UNP Q92ZS5
H	-20	HIS	-	expression tag	UNP Q92ZS5
H	-19	HIS	-	expression tag	UNP Q92ZS5
H	-18	HIS	-	expression tag	UNP Q92ZS5
H	-17	HIS	-	expression tag	UNP Q92ZS5
H	-16	HIS	-	expression tag	UNP Q92ZS5
H	-15	SER	-	expression tag	UNP Q92ZS5
H	-14	SER	-	expression tag	UNP Q92ZS5
H	-13	GLY	-	expression tag	UNP Q92ZS5
H	-12	VAL	-	expression tag	UNP Q92ZS5
H	-11	ASP	-	expression tag	UNP Q92ZS5
H	-10	LEU	-	expression tag	UNP Q92ZS5
H	-9	GLY	-	expression tag	UNP Q92ZS5
H	-8	THR	-	expression tag	UNP Q92ZS5
H	-7	GLU	-	expression tag	UNP Q92ZS5
H	-6	ASN	-	expression tag	UNP Q92ZS5
H	-5	LEU	-	expression tag	UNP Q92ZS5
H	-4	TYR	-	expression tag	UNP Q92ZS5
H	-3	PHE	-	expression tag	UNP Q92ZS5

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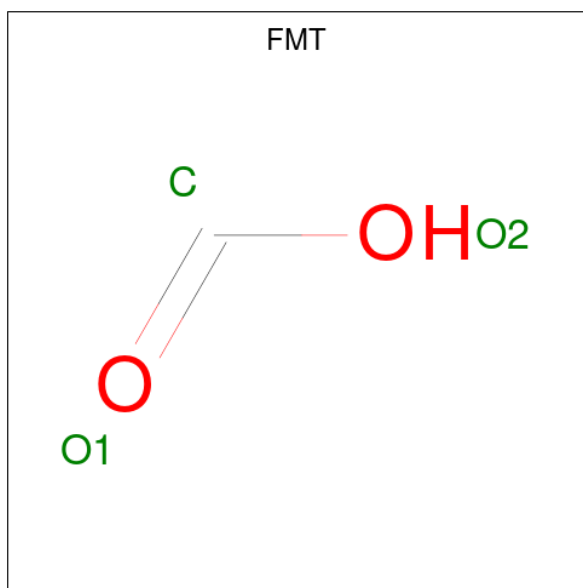
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Chain	Residue	Modelled	Actual	Comment	Reference
H	-2	GLN	-	expression tag	UNP Q92ZS5
H	-1	SER	-	expression tag	UNP Q92ZS5
H	0	MET	-	expression tag	UNP Q92ZS5

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	E	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0
2	G	1	Total Mg 1 1	0	0
2	H	1	Total Mg 1 1	0	0

- Molecule 3 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).

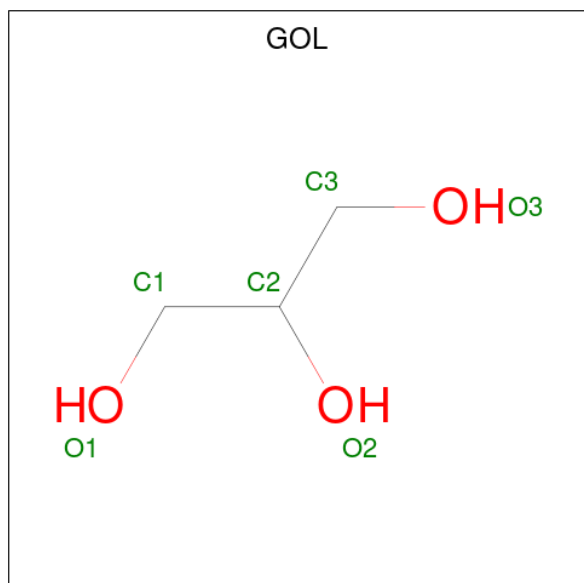


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 3 1 2	0	0
3	B	1	Total C O 3 1 2	0	0
3	C	1	Total C O 3 1 2	0	0
3	E	1	Total C O 3 1 2	0	0
3	F	1	Total C O 3 1 2	0	0
3	G	1	Total C O 3 1 2	0	0
3	H	1	Total C O 3 1 2	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 2 2	0	1
4	C	2	Total Cl 3 3	0	1
4	D	1	Total Cl 2 2	0	1
4	G	1	Total Cl 2 2	0	1

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



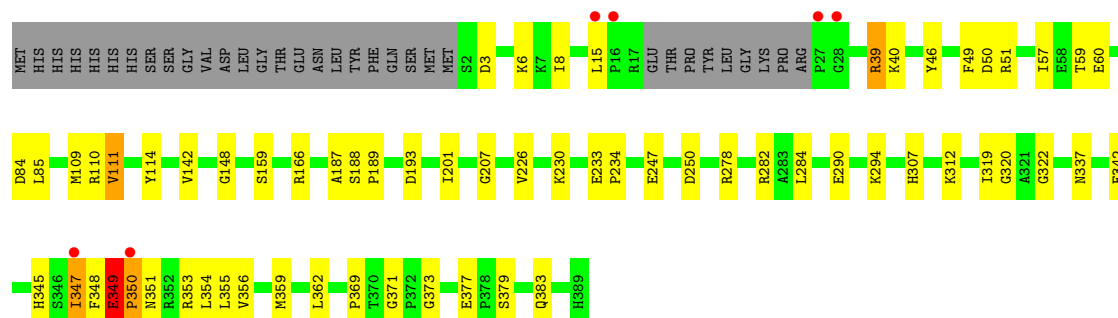
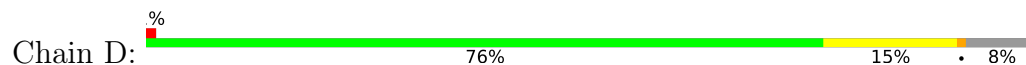
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

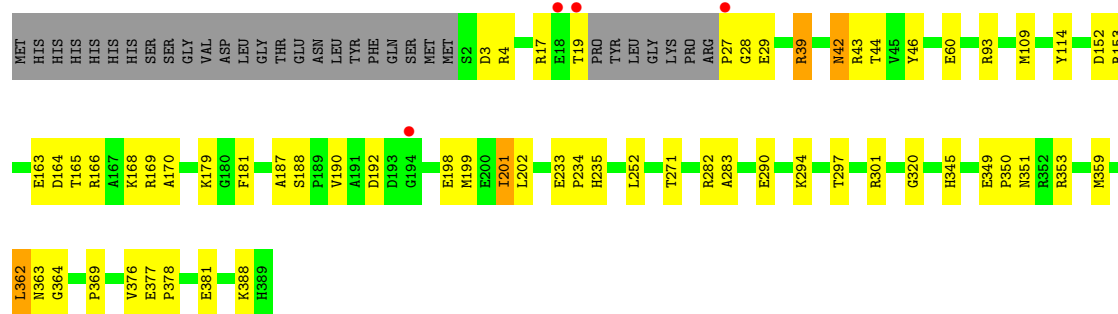
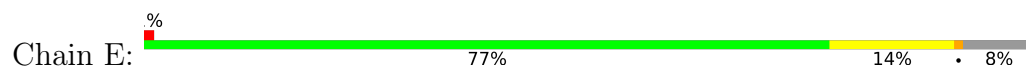
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	135	Total	O	0	0
			135	135		
6	B	237	Total	O	0	0
			237	237		
6	C	161	Total	O	0	0
			161	161		
6	D	168	Total	O	0	0
			168	168		
6	E	213	Total	O	0	0
			213	213		
6	F	201	Total	O	0	0
			201	201		
6	G	233	Total	O	0	0
			233	233		
6	H	143	Total	O	0	1
			144	144		



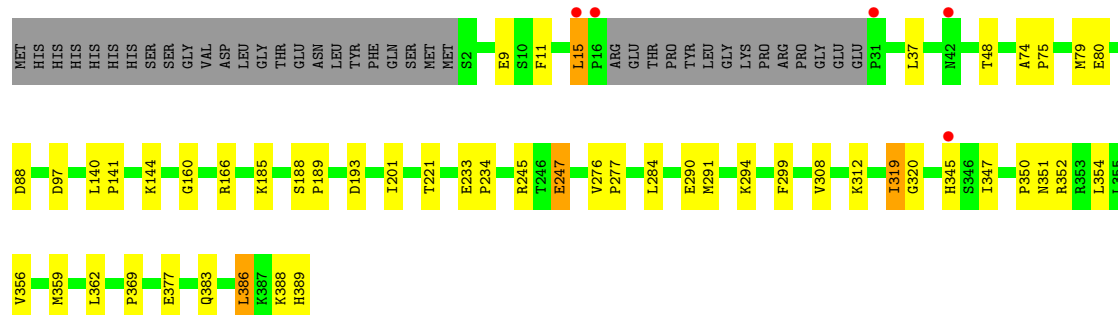
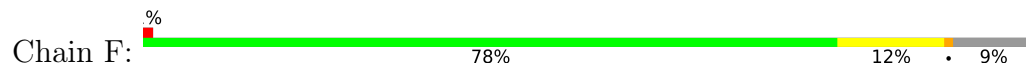
• Molecule 1: Enolase Q92Zs5



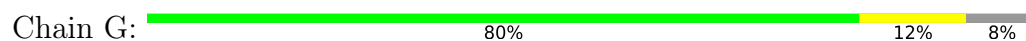
• Molecule 1: Enolase Q92Zs5

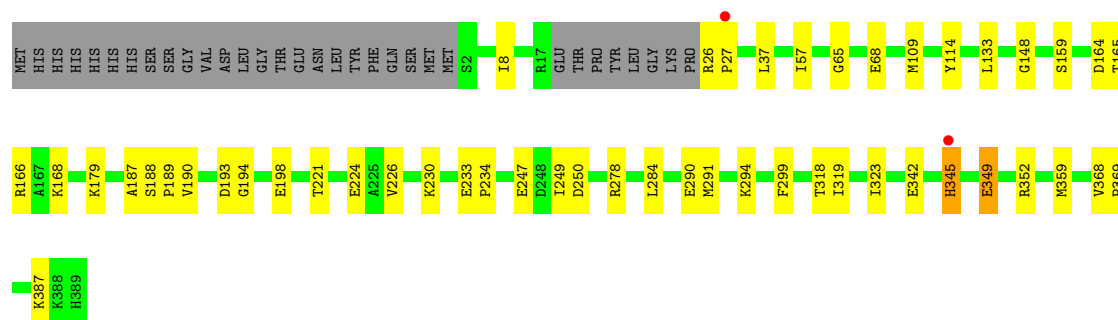


• Molecule 1: Enolase Q92Zs5

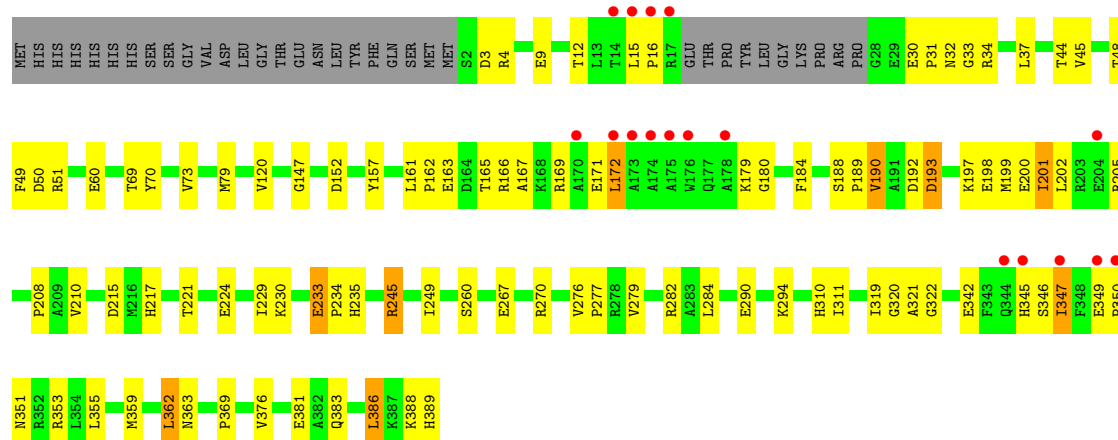


• Molecule 1: Enolase Q92Zs5





● Molecule 1: Enolase Q92Zs5



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	102.72Å 121.70Å 137.69Å 90.00° 99.79° 90.00°	Depositor
Resolution (Å)	45.02 – 2.01 45.02 – 2.01	Depositor EDS
% Data completeness (in resolution range)	98.4 (45.02-2.01) 98.5 (45.02-2.01)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.172 , 0.229 0.173 , 0.229	Depositor DCC
R_{free} test set	6578 reflections (2.98%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.498	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	24833	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.57	2/2942 (0.1%)	0.86	4/3990 (0.1%)
1	B	0.63	4/3053 (0.1%)	0.88	4/4144 (0.1%)
1	C	0.64	2/2975 (0.1%)	0.90	4/4032 (0.1%)
1	D	0.63	3/2980 (0.1%)	0.91	9/4039 (0.2%)
1	E	0.60	1/2993 (0.0%)	0.89	3/4057 (0.1%)
1	F	0.66	2/2955 (0.1%)	0.92	3/4007 (0.1%)
1	G	0.66	3/2983 (0.1%)	0.94	5/4045 (0.1%)
1	H	0.66	4/2954 (0.1%)	0.92	9/4006 (0.2%)
All	All	0.63	21/23835 (0.1%)	0.90	41/32320 (0.1%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	39[A]	ARG	CA-C	9.09	1.64	1.52
1	D	39[B]	ARG	CA-C	9.09	1.64	1.52
1	B	24	LYS	C-N	5.62	1.41	1.33
1	C	349	GLU	C-N	5.57	1.41	1.34
1	G	349	GLU	C-N	5.48	1.40	1.34
1	H	15	LEU	C-N	5.46	1.41	1.33
1	H	233	GLU	C-N	5.44	1.40	1.33
1	H	162	PRO	N-CD	5.39	1.55	1.47
1	D	350	PRO	N-CD	5.38	1.55	1.47
1	B	31	PRO	N-CD	5.37	1.55	1.47
1	A	350	PRO	N-CD	5.36	1.55	1.47
1	C	350	PRO	N-CD	5.33	1.55	1.47
1	G	26	ARG	C-N	5.32	1.40	1.34
1	A	349	GLU	C-N	5.31	1.40	1.33
1	B	30	GLU	C-N	5.28	1.40	1.33
1	F	350	PRO	N-CD	5.21	1.55	1.47
1	F	319	ILE	C-O	-5.17	1.19	1.24
1	B	25	PRO	N-CD	5.16	1.54	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	27	PRO	N-CD	5.13	1.54	1.47
1	E	188	SER	C-N	5.11	1.41	1.34
1	H	16	PRO	N-CD	5.03	1.54	1.47

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	347	ILE	N-CA-C	10.10	120.02	110.53
1	D	39[A]	ARG	CA-C-O	8.65	130.92	120.62
1	D	39[B]	ARG	CA-C-O	8.65	130.92	120.62
1	F	140	LEU	CA-C-N	-7.98	111.79	119.85
1	F	140	LEU	C-N-CA	-7.98	111.79	119.85
1	A	339	ASP	N-CA-C	6.98	118.68	111.14
1	D	371	GLY	CA-C-N	6.76	126.72	120.03
1	D	371	GLY	C-N-CA	6.76	126.72	120.03
1	B	24	LYS	CA-C-N	-6.76	112.80	119.76
1	B	24	LYS	C-N-CA	-6.76	112.80	119.76
1	D	207	GLY	CA-C-N	6.23	125.85	119.56
1	D	207	GLY	C-N-CA	6.23	125.85	119.56
1	C	347	ILE	N-CA-C	6.21	116.36	110.53
1	C	349	GLU	CA-C-N	-6.12	113.38	119.56
1	C	349	GLU	C-N-CA	-6.12	113.38	119.56
1	H	15	LEU	CA-C-N	-6.11	113.48	119.78
1	H	15	LEU	C-N-CA	-6.11	113.48	119.78
1	G	26	ARG	CA-C-N	-6.00	112.73	118.97
1	G	26	ARG	C-N-CA	-6.00	112.73	118.97
1	A	347	ILE	N-CA-C	5.95	116.12	110.53
1	H	179	LYS	N-CA-C	5.80	119.33	112.72
1	B	30	GLU	CA-C-N	-5.74	113.86	119.78
1	B	30	GLU	C-N-CA	-5.74	113.86	119.78
1	F	221	THR	N-CA-C	-5.69	103.19	110.53
1	G	349	GLU	CA-C-N	-5.69	113.81	119.56
1	G	349	GLU	C-N-CA	-5.69	113.81	119.56
1	A	349	GLU	CA-C-N	-5.50	113.23	119.28
1	A	349	GLU	C-N-CA	-5.50	113.23	119.28
1	H	208	PRO	N-CA-C	-5.46	109.00	114.68
1	H	161	LEU	CA-C-N	-5.45	113.03	119.84
1	H	161	LEU	C-N-CA	-5.45	113.03	119.84
1	D	349	GLU	CA-C-N	-5.29	113.17	119.05
1	D	349	GLU	C-N-CA	-5.29	113.17	119.05
1	D	347	ILE	CB-CA-C	-5.28	105.11	112.02
1	E	19	THR	N-CA-CB	-5.24	102.60	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	233	GLU	CA-C-N	-5.13	113.64	119.28
1	H	233	GLU	C-N-CA	-5.13	113.64	119.28
1	G	164	ASP	N-CA-C	5.06	119.45	113.28
1	C	217	HIS	N-CA-C	5.03	118.38	111.54
1	E	188	SER	CA-C-N	-5.01	113.49	119.05
1	E	188	SER	C-N-CA	-5.01	113.49	119.05

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	2880	0	2872	62	0
1	B	2982	0	2977	49	0
1	C	2909	0	2908	63	0
1	D	2913	0	2907	55	0
1	E	2926	0	2920	53	0
1	F	2880	0	2881	38	0
1	G	2916	0	2908	38	0
1	H	2891	0	2880	75	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
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	3	0	1	0	0
3	B	3	0	1	1	0
3	C	3	0	1	0	0
3	E	3	0	1	0	0
3	F	3	0	1	0	0
3	G	3	0	1	0	0
3	H	3	0	1	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	2	0	0	0	0
4	C	3	0	0	0	0
4	D	2	0	0	0	0
4	G	2	0	0	0	0
5	B	6	0	8	5	0
6	A	135	0	0	3	0
6	B	237	0	0	3	0
6	C	161	0	0	6	0
6	D	168	0	0	1	0
6	E	213	0	0	5	0
6	F	201	0	0	4	0
6	G	233	0	0	4	0
6	H	144	0	0	3	0
All	All	24833	0	23268	419	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 9.

All (419) 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:A:15:LEU:HD11	6:A:605:HOH:O	1.49	1.11
1:H:349:GLU:HB3	1:H:350:PRO:HD3	1.47	0.95
1:D:166:ARG:HG2	1:D:201:ILE:HG21	1.49	0.95
1:H:383:GLN:HA	1:H:386:LEU:HD22	1.50	0.93
1:A:163:GLU:HB2	1:A:169:ARG:CG	1.99	0.92
1:B:31:PRO:HG3	1:B:37:LEU:HD23	1.52	0.91
1:E:166:ARG:HG2	1:E:201:ILE:HD12	1.49	0.91
1:A:163:GLU:HB2	1:A:169:ARG:HG2	1.53	0.91
1:B:344:GLN:HB3	1:B:347:ILE:HD12	1.54	0.89
3:B:403:FMT:H	6:B:638:HOH:O	1.73	0.89
1:C:359:MET:HB2	1:C:369:PRO:HG3	1.55	0.87
1:D:319:ILE:HA	1:D:347:ILE:HD12	1.56	0.87
1:F:15:LEU:HD21	1:F:354:LEU:HD11	1.56	0.85
1:G:345:HIS:CD2	1:G:345:HIS:H	1.90	0.85
1:E:349:GLU:HB2	1:E:350:PRO:HD3	1.57	0.84
1:G:188:SER:OG	1:G:189:PRO:HD3	1.79	0.83
1:C:353:ARG:NH1	1:C:381:GLU:HG2	1.95	0.82
1:E:39:ARG:HG3	1:E:46:TYR:CE1	2.16	0.81
1:E:362:LEU:O	1:E:362:LEU:HD23	1.79	0.81
1:F:15:LEU:CD2	1:F:354:LEU:HD11	2.12	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:250:ASP:HB3	1:H:282:ARG:HH11	1.49	0.77
1:G:221:THR:OG1	1:G:224:GLU:HG3	1.84	0.77
1:E:39:ARG:HG3	1:E:46:TYR:HE1	1.49	0.76
1:H:200:GLU:HB2	1:H:235:HIS:CE1	2.21	0.76
1:A:37:LEU:C	1:A:37:LEU:HD12	2.12	0.74
1:D:39[A]:ARG:HG3	1:D:46:TYR:CE1	2.22	0.74
1:H:362:LEU:HD23	1:H:362:LEU:O	1.87	0.74
1:A:37:LEU:HD12	1:A:37:LEU:O	1.89	0.73
1:H:349:GLU:CB	1:H:350:PRO:HD3	2.18	0.73
1:A:163:GLU:HB2	1:A:169:ARG:HG3	1.71	0.73
1:E:359:MET:HB2	1:E:369:PRO:HG3	1.72	0.72
1:E:3:ASP:O	1:E:4:ARG:HD3	1.89	0.72
1:A:335:LEU:HB2	1:A:338:VAL:HG23	1.71	0.72
1:G:250:ASP:HB3	1:H:282:ARG:NH1	2.04	0.72
1:B:186:PHE:HE1	1:B:199:MET:CE	2.02	0.71
1:E:198:GLU:O	1:E:201:ILE:HG22	1.90	0.71
1:C:43:ARG:HG2	1:C:43:ARG:HH11	1.55	0.71
1:G:345:HIS:H	1:G:345:HIS:HD2	1.38	0.71
1:C:319:ILE:HA	1:C:347:ILE:HG13	1.72	0.71
1:B:199:MET:HG2	1:B:235:HIS:HB2	1.73	0.71
1:H:31:PRO:HG3	1:H:48:THR:HG21	1.72	0.70
1:E:163:GLU:HB2	1:E:169:ARG:HG2	1.73	0.69
1:F:15:LEU:HD21	1:F:354:LEU:CD1	2.23	0.68
1:A:291:MET:HE2	1:A:299:PHE:CD2	2.28	0.68
1:H:322:GLY:HA3	1:H:355:LEU:HD11	1.74	0.68
1:G:37:LEU:HD12	1:G:37:LEU:C	2.19	0.68
1:B:21:TYR:O	1:B:22:LEU:HB2	1.91	0.68
1:F:290:GLU:H	1:F:294:LYS:HZ2	1.42	0.67
1:A:387:LYS:HD3	6:A:544:HOH:O	1.94	0.67
1:C:197:LYS:HE2	1:C:200:GLU:OE1	1.94	0.67
1:F:166:ARG:HG2	1:F:201:ILE:HG21	1.76	0.67
1:H:359:MET:HB2	1:H:369:PRO:HG3	1.77	0.66
1:H:163:GLU:HB2	1:H:169:ARG:HG2	1.77	0.66
1:H:229:ILE:O	1:H:233:GLU:HG3	1.95	0.66
1:A:347:ILE:HD12	1:A:347:ILE:C	2.21	0.66
1:G:187:ALA:O	1:G:190:VAL:HG22	1.96	0.66
1:G:345:HIS:CD2	1:G:345:HIS:N	2.64	0.65
1:B:199:MET:CG	1:B:235:HIS:HB2	2.27	0.65
1:H:50:ASP:HB3	1:H:319:ILE:HG21	1.78	0.65
1:H:201:ILE:HG23	1:H:202:LEU:N	2.11	0.65
1:G:290:GLU:H	1:G:294:LYS:HZ2	1.45	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:388:LYS:HG3	1:H:389:HIS:N	2.12	0.64
1:H:349:GLU:HB3	1:H:350:PRO:CD	2.25	0.64
1:D:322:GLY:HA3	1:D:355:LEU:HD11	1.80	0.63
1:A:34:ARG:HD2	1:C:88:ASP:OD2	1.99	0.62
1:F:290:GLU:H	1:F:294:LYS:NZ	1.95	0.62
1:B:199:MET:HE2	1:B:199:MET:HA	1.81	0.62
1:D:49:PHE:CD1	1:D:49:PHE:O	2.53	0.62
1:A:322:GLY:HA3	1:A:355:LEU:HD11	1.80	0.62
1:A:359:MET:HB2	1:A:369:PRO:HG3	1.81	0.61
1:B:362:LEU:O	1:B:362:LEU:HD23	2.01	0.61
1:H:198:GLU:O	1:H:201:ILE:HG22	2.00	0.61
1:D:347:ILE:O	1:D:350:PRO:HD2	2.00	0.61
1:A:163:GLU:HA	1:A:163:GLU:OE1	2.01	0.60
1:E:349:GLU:CB	1:E:350:PRO:HD3	2.31	0.60
1:B:31:PRO:HA	1:B:37:LEU:HB3	1.84	0.60
1:D:319:ILE:HG23	1:D:347:ILE:HD12	1.82	0.60
1:C:43:ARG:HG2	1:C:43:ARG:NH1	2.17	0.60
1:D:250:ASP:HB3	1:E:282[A]:ARG:HE	1.67	0.60
1:A:215:ASP:HA	1:A:241:GLU:HB3	1.84	0.60
1:E:187:ALA:O	1:E:190:VAL:HG22	2.01	0.60
1:G:290:GLU:H	1:G:294:LYS:NZ	2.00	0.60
1:F:88:ASP:HB2	6:F:641:HOH:O	2.00	0.59
1:D:159:SER:HB2	1:D:342:GLU:OE2	2.03	0.59
1:G:359:MET:HB2	1:G:369:PRO:HG3	1.83	0.59
1:C:345:HIS:ND1	1:C:349:GLU:HG3	2.18	0.59
1:H:200:GLU:HA	1:H:235:HIS:ND1	2.18	0.59
1:A:201:ILE:HG23	1:A:202:LEU:N	2.18	0.58
1:C:28:GLY:O	1:C:40:LYS:NZ	2.36	0.58
1:A:335:LEU:HB2	1:A:338:VAL:CG2	2.33	0.58
1:A:188:SER:N	1:A:189:PRO:CD	2.67	0.58
1:B:110:ARG:HH21	1:F:247[A]:GLU:CD	2.12	0.58
1:D:142:VAL:HG23	1:D:373:GLY:C	2.29	0.57
1:G:179:LYS:HE3	6:G:625:HOH:O	2.03	0.57
1:H:345:HIS:O	1:H:349:GLU:HB2	2.04	0.57
1:E:201:ILE:HG23	1:E:202:LEU:N	2.19	0.57
1:B:17:ARG:HD2	1:B:50:ASP:OD1	2.04	0.57
6:E:704:HOH:O	1:G:247[B]:GLU:HG3	2.04	0.57
1:G:249:ILE:HG12	6:G:727:HOH:O	2.04	0.57
1:G:37:LEU:HD12	1:G:37:LEU:O	2.04	0.57
1:B:300:MET:HB2	5:B:404:GOL:H12	1.87	0.57
1:E:297:THR:O	1:E:301:ARG:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:290:GLU:H	1:D:294:LYS:NZ	2.03	0.56
1:H:201:ILE:HG23	1:H:202:LEU:H	1.70	0.56
1:C:176:TRP:HZ2	1:C:345:HIS:CD2	2.23	0.56
1:G:387:LYS:NZ	6:G:676:HOH:O	2.37	0.56
1:H:188:SER:OG	1:H:189:PRO:HD3	2.05	0.56
1:C:377:GLU:HB2	1:C:378:PRO:HD2	1.86	0.56
1:H:319:ILE:HA	1:H:347:ILE:HG21	1.88	0.56
1:B:188:SER:OG	1:B:189:PRO:HD3	2.06	0.56
1:E:164:ASP:HB3	1:E:168:LYS:HZ1	1.70	0.56
1:F:188:SER:N	1:F:189:PRO:CD	2.68	0.56
1:H:152:ASP:HB2	6:H:625:HOH:O	2.04	0.56
1:B:28:GLY:O	1:B:40:LYS:HE3	2.06	0.56
1:D:166:ARG:HG2	1:D:201:ILE:CG2	2.31	0.56
1:A:164:ASP:OD1	1:A:165:THR:HG23	2.05	0.56
1:E:320:GLY:O	1:E:351:ASN:ND2	2.39	0.56
6:C:565:HOH:O	1:D:282:ARG:HD3	2.06	0.55
1:E:42:ASN:ND2	1:E:44:THR:H	2.04	0.55
1:C:350:PRO:HG2	1:C:351:ASN:HD22	1.71	0.55
1:H:163:GLU:HG2	1:H:172:LEU:HD13	1.87	0.55
1:E:39:ARG:CG	1:E:46:TYR:CE1	2.89	0.55
1:F:383:GLN:HA	1:F:386:LEU:HD22	1.88	0.55
1:D:50:ASP:CG	1:D:319:ILE:HG21	2.30	0.55
1:E:17:ARG:HB2	6:E:702:HOH:O	2.06	0.55
1:H:290:GLU:H	1:H:294:LYS:NZ	2.04	0.55
1:H:49:PHE:HE2	1:H:51:ARG:CZ	2.19	0.55
1:C:377:GLU:HB2	1:C:378:PRO:CD	2.36	0.54
1:E:42:ASN:C	1:E:42:ASN:HD22	2.14	0.54
1:H:49:PHE:N	1:H:49:PHE:CD1	2.75	0.54
1:D:6:LYS:HD2	1:D:60:GLU:OE2	2.06	0.54
1:D:319:ILE:CA	1:D:347:ILE:HD12	2.33	0.54
1:D:345:HIS:CE1	1:D:349:GLU:HB2	2.42	0.54
1:F:312:LYS:HE2	6:F:625:HOH:O	2.08	0.54
1:C:31:PRO:HG3	1:C:48:THR:HG21	1.89	0.54
1:D:278:ARG:HB3	1:D:284:LEU:HD22	1.90	0.54
1:D:356:VAL:CG2	1:D:377:GLU:HG3	2.37	0.54
1:F:308:VAL:HG23	6:F:549:HOH:O	2.07	0.54
1:A:37:LEU:C	1:A:37:LEU:CD1	2.80	0.54
1:E:166:ARG:HG2	1:E:201:ILE:CD1	2.32	0.54
1:A:89:PHE:CE1	1:C:38:VAL:HG23	2.43	0.54
1:D:379:SER:O	1:D:383:GLN:HG3	2.09	0.53
1:D:39[A]:ARG:HG3	1:D:46:TYR:HE1	1.70	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:278:ARG:HB3	1:C:284:LEU:CD2	2.39	0.53
1:D:319:ILE:HG12	1:D:347:ILE:HD11	1.90	0.53
1:C:278:ARG:HB3	1:C:284:LEU:HD22	1.90	0.53
1:D:8:ILE:HG12	1:D:57:ILE:HG12	1.89	0.53
1:F:166:ARG:HB2	6:F:701:HOH:O	2.06	0.53
1:H:290:GLU:H	1:H:294:LYS:HZ2	1.57	0.53
1:C:220:HIS:HB2	1:C:243:PRO:O	2.09	0.53
1:A:335:LEU:CB	1:A:338:VAL:HG23	2.39	0.53
1:E:345:HIS:H	1:E:345:HIS:CD2	2.26	0.53
1:G:349:GLU:O	1:G:352:ARG:HG3	2.09	0.53
1:H:188:SER:N	1:H:189:PRO:CD	2.71	0.53
1:A:171:GLU:O	1:A:174:ALA:HB3	2.10	0.52
1:C:247:GLU:CD	1:C:247:GLU:H	2.18	0.52
1:F:9:GLU:OE2	1:F:388:LYS:HE3	2.09	0.52
1:H:37:LEU:C	1:H:37:LEU:HD12	2.34	0.52
1:D:60:GLU:OE2	1:D:60:GLU:HA	2.08	0.52
1:C:198:GLU:O	1:C:201:ILE:HG22	2.10	0.52
1:D:348:PHE:C	1:D:348:PHE:CD1	2.87	0.52
1:E:377:GLU:HB2	1:E:378:PRO:HD2	1.90	0.52
1:A:32:ASN:C	1:A:34:ARG:H	2.18	0.52
1:F:245:ARG:HB3	1:F:247[A]:GLU:OE1	2.09	0.52
1:H:193:ASP:OD1	1:H:193:ASP:N	2.42	0.52
1:B:344:GLN:CB	1:B:347:ILE:HD12	2.34	0.52
1:C:109:MET:HE3	1:C:114:TYR:CD2	2.45	0.52
1:A:74:ALA:N	1:A:75:PRO:CD	2.72	0.51
1:C:197:LYS:CE	1:C:200:GLU:OE1	2.57	0.51
1:D:290:GLU:H	1:D:294:LYS:HZ2	1.58	0.51
1:H:362:LEU:CD2	1:H:362:LEU:N	2.73	0.51
1:C:362:LEU:O	1:C:362:LEU:HG	2.09	0.51
1:G:247[A]:GLU:HG2	1:H:310:HIS:CE1	2.45	0.51
1:C:321:ALA:O	1:C:355:LEU:HD11	2.10	0.51
1:A:43:ARG:NH2	1:C:104:ASP:OD1	2.39	0.51
1:H:205:ARG:HD2	1:H:205:ARG:O	2.10	0.51
1:H:284:LEU:N	1:H:284:LEU:HD23	2.26	0.51
1:B:186:PHE:HE1	1:B:199:MET:HE1	1.73	0.51
1:A:165:THR:OG1	1:A:168:LYS:HG3	2.11	0.51
1:B:180:GLY:HA3	1:B:363:ASN:HD22	1.75	0.51
1:C:176:TRP:HZ2	1:C:345:HIS:CG	2.29	0.51
1:F:188:SER:N	1:F:189:PRO:HD2	2.25	0.51
1:F:356:VAL:CG2	1:F:377:GLU:HG3	2.40	0.51
1:E:201:ILE:CG2	1:E:202:LEU:N	2.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:VAL:HG23	1:C:89:PHE:CE1	2.46	0.51
1:B:359:MET:HB2	1:B:369:PRO:HG3	1.93	0.51
1:H:163:GLU:HB2	1:H:169:ARG:CG	2.39	0.51
1:C:282:ARG:CD	6:C:520:HOH:O	2.60	0.50
1:D:320:GLY:O	1:D:351:ASN:ND2	2.40	0.50
1:F:320:GLY:O	1:F:351:ASN:ND2	2.44	0.50
1:G:8:ILE:HG12	1:G:57:ILE:HG12	1.93	0.50
1:A:290:GLU:H	1:A:294:LYS:HZ2	1.58	0.50
1:B:358:ASP:OD2	1:B:370:THR:OG1	2.29	0.50
1:F:345:HIS:CD2	1:F:345:HIS:H	2.28	0.50
1:H:167:ALA:O	1:H:171:GLU:HG2	2.11	0.50
1:H:120:VAL:HG11	1:H:270:ARG:HB3	1.93	0.50
1:A:111:VAL:HG22	6:C:521:HOH:O	2.11	0.50
1:H:157:TYR:CZ	1:H:342:GLU:HG3	2.46	0.50
1:F:359:MET:HB2	1:F:369:PRO:HG3	1.94	0.50
1:H:171:GLU:O	1:H:172:LEU:C	2.55	0.50
1:F:141:PRO:HG2	1:F:144:LYS:HG2	1.94	0.50
1:B:186:PHE:CE1	1:B:199:MET:CE	2.89	0.49
1:B:300:MET:CB	5:B:404:GOL:H12	2.42	0.49
1:A:13:LEU:HD13	1:A:386:LEU:HD23	1.95	0.49
1:H:276:VAL:HB	1:H:277:PRO:HD3	1.93	0.49
1:C:56:ARG:NH1	6:C:544:HOH:O	2.35	0.49
1:B:11:PHE:N	1:B:11:PHE:CD1	2.80	0.49
1:A:201:ILE:CG2	1:A:202:LEU:N	2.75	0.49
1:A:286:ILE:HG12	1:A:312:LYS:HB2	1.95	0.49
1:C:347:ILE:O	1:C:351:ASN:ND2	2.45	0.49
1:D:188:SER:N	1:D:189:PRO:CD	2.75	0.49
1:E:252:LEU:HG	1:E:283:ALA:HB1	1.95	0.48
1:A:233:GLU:N	1:A:234:PRO:CD	2.77	0.48
1:A:278:ARG:HB3	1:A:284:LEU:HD22	1.95	0.48
1:E:353:ARG:NH1	1:E:381:GLU:OE2	2.46	0.48
1:A:362:LEU:HD23	1:A:362:LEU:N	2.27	0.48
1:H:166:ARG:HG2	1:H:201:ILE:HD12	1.95	0.48
1:C:31:PRO:HA	1:C:37:LEU:HB3	1.95	0.48
1:B:349:GLU:HB2	1:B:350:PRO:HD3	1.95	0.48
1:E:42:ASN:HD22	1:E:43:ARG:N	2.11	0.48
1:E:290:GLU:H	1:E:294:LYS:NZ	2.12	0.48
1:E:109:MET:HE3	1:E:114:TYR:CD2	2.49	0.47
1:C:187:ALA:HB1	1:C:189:PRO:HD2	1.96	0.47
1:D:49:PHE:HE1	1:D:51:ARG:CZ	2.27	0.47
1:E:42:ASN:ND2	1:E:42:ASN:C	2.72	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:163:GLU:HB2	1:E:169:ARG:CG	2.41	0.47
1:D:166:ARG:NE	1:D:193:ASP:OD2	2.47	0.47
1:A:32:ASN:O	1:A:34:ARG:N	2.47	0.47
1:B:353:ARG:NE	1:B:381:GLU:OE1	2.47	0.47
1:C:347:ILE:HD13	1:C:347:ILE:N	2.29	0.47
1:E:164:ASP:HB3	1:E:168:LYS:NZ	2.28	0.47
1:B:181:PHE:CD2	1:B:364:GLY:HA2	2.50	0.47
1:D:111:VAL:HG13	6:D:598:HOH:O	2.14	0.47
1:A:3:ASP:CG	1:A:59:THR:HB	2.40	0.47
1:A:187:ALA:HB1	1:A:189:PRO:HD2	1.97	0.47
1:E:27:PRO:O	1:E:29:GLU:N	2.47	0.47
1:E:60:GLU:OE1	1:E:60:GLU:HA	2.15	0.47
1:A:97:ASP:HA	1:D:148:GLY:HA3	1.96	0.47
1:B:2:SER:HB3	1:B:4:ARG:HE	1.80	0.47
1:C:165:THR:HG23	1:C:168:LYS:HD2	1.96	0.47
1:H:230:LYS:O	1:H:234:PRO:HD3	2.15	0.47
1:A:320:GLY:O	1:A:351:ASN:ND2	2.47	0.47
1:B:23:GLY:O	1:B:39:ARG:NH2	2.48	0.47
1:C:201:ILE:HG23	1:C:202:LEU:N	2.29	0.46
1:D:50:ASP:CG	1:D:319:ILE:CG2	2.89	0.46
1:F:247[A]:GLU:CD	1:F:247[A]:GLU:H	2.23	0.46
1:H:30:GLU:H	1:H:30:GLU:CD	2.23	0.46
1:H:362:LEU:HD23	1:H:362:LEU:C	2.37	0.46
1:A:290:GLU:H	1:A:294:LYS:NZ	2.13	0.46
1:A:379:SER:O	1:A:383:GLN:HG3	2.15	0.46
1:D:233:GLU:N	1:D:234:PRO:CD	2.78	0.46
1:E:233:GLU:N	1:E:234:PRO:CD	2.78	0.46
1:A:32:ASN:C	1:A:32:ASN:OD1	2.58	0.46
1:C:12:THR:HG23	1:C:79:MET:CE	2.45	0.46
1:E:3:ASP:C	1:E:4:ARG:HD3	2.40	0.46
1:G:194:GLY:HA2	6:G:577:HOH:O	2.14	0.46
1:A:151:ARG:NH1	1:A:332:SER:O	2.49	0.46
1:B:188:SER:N	1:B:189:PRO:CD	2.78	0.46
1:C:200:GLU:HG3	1:C:235:HIS:CE1	2.51	0.46
1:H:37:LEU:HD12	1:H:37:LEU:O	2.15	0.46
1:H:201:ILE:CG2	1:H:202:LEU:N	2.78	0.46
1:A:50:ASP:HB3	1:A:319:ILE:CG2	2.46	0.46
1:A:229:ILE:O	1:A:233:GLU:HG3	2.16	0.46
1:A:276:VAL:HB	1:A:277:PRO:HD3	1.98	0.46
1:C:166:ARG:HG2	1:C:201:ILE:HD12	1.98	0.46
1:C:224:GLU:OE1	1:D:312:LYS:HE3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:385:LEU:O	1:C:387:LYS:HE2	2.16	0.46
1:E:271:THR:CG2	6:E:573:HOH:O	2.64	0.46
1:C:226:VAL:O	1:C:230:LYS:HG2	2.16	0.46
1:B:183:SER:C	1:B:184:PHE:CD1	2.94	0.45
1:C:76:ARG:HA	1:C:79:MET:HE3	1.98	0.45
1:F:284:LEU:N	1:F:284:LEU:HD23	2.31	0.45
1:E:199:MET:HG3	1:E:235:HIS:HB2	1.98	0.45
1:A:362:LEU:N	1:A:362:LEU:CD2	2.80	0.45
1:D:187:ALA:HB1	1:D:189:PRO:HD2	1.98	0.45
1:F:74:ALA:N	1:F:75:PRO:CD	2.79	0.45
1:H:180:GLY:HA3	1:H:363:ASN:HA	1.97	0.45
1:C:233:GLU:N	1:C:234:PRO:CD	2.79	0.45
1:H:147:GLY:HA2	6:H:633:HOH:O	2.15	0.45
1:C:166:ARG:HG2	1:C:201:ILE:CD1	2.47	0.45
1:D:351:ASN:HA	1:D:354:LEU:HD12	1.98	0.45
1:E:362:LEU:HD23	1:E:362:LEU:C	2.38	0.45
1:A:187:ALA:C	1:A:189:PRO:HD2	2.42	0.45
1:G:291:MET:HE2	1:G:299:PHE:CD2	2.52	0.45
1:F:356:VAL:HG22	1:F:377:GLU:HG3	1.98	0.45
1:B:76:ARG:HB2	1:F:80:GLU:HG2	1.98	0.45
1:E:93:ARG:NH2	6:E:630:HOH:O	2.38	0.45
1:E:152:ASP:O	1:E:153:ARG:HD3	2.16	0.45
1:F:352:ARG:HG2	1:F:352:ARG:HH11	1.82	0.45
1:D:110:ARG:NH2	1:H:245:ARG:HG3	2.32	0.44
1:D:319:ILE:HG23	1:D:347:ILE:CD1	2.46	0.44
1:A:154:ILE:O	1:A:155:ALA:C	2.60	0.44
1:F:233:GLU:N	1:F:234:PRO:CD	2.80	0.44
1:B:163:GLU:HB2	1:B:169:ARG:HG2	1.99	0.44
1:B:188:SER:N	1:B:189:PRO:HD2	2.32	0.44
1:D:109:MET:HE3	1:D:114:TYR:CD2	2.52	0.44
1:E:377:GLU:HB2	1:E:378:PRO:CD	2.48	0.44
1:B:290:GLU:H	1:B:294:LYS:NZ	2.15	0.44
1:C:179:LYS:O	1:C:363:ASN:HA	2.18	0.44
1:D:3:ASP:OD1	1:D:59:THR:HB	2.18	0.44
1:D:345:HIS:CE1	1:D:349:GLU:OE1	2.70	0.44
1:C:17:ARG:O	1:C:17:ARG:HG3	2.17	0.44
1:E:165:THR:HG23	1:E:168:LYS:HE3	1.98	0.44
1:E:282[A]:ARG:HA	6:E:711:HOH:O	2.17	0.44
1:B:3:ASP:O	1:B:4:ARG:HD3	2.18	0.44
1:H:320:GLY:O	1:H:351:ASN:ND2	2.50	0.44
1:C:203:ARG:HA	1:C:203:ARG:HD3	1.73	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:322:GLY:HA3	1:C:355:LEU:HD11	2.00	0.44
1:D:226:VAL:O	1:D:230:LYS:HG2	2.18	0.44
1:G:278:ARG:HB3	1:G:284:LEU:HD22	2.00	0.44
1:G:318:THR:OG1	1:G:319:ILE:N	2.50	0.44
1:H:49:PHE:CD2	1:H:49:PHE:O	2.70	0.44
1:H:198:GLU:O	1:H:201:ILE:CG2	2.65	0.44
1:A:181:PHE:CE2	1:A:364:GLY:HA2	2.53	0.43
1:B:184:PHE:CD1	1:B:184:PHE:N	2.86	0.43
1:A:347:ILE:HD12	1:A:347:ILE:O	2.17	0.43
1:B:279:VAL:HG22	1:B:311:ILE:CD1	2.48	0.43
1:C:176:TRP:CZ2	1:C:345:HIS:HB2	2.54	0.43
1:G:37:LEU:C	1:G:37:LEU:CD1	2.90	0.43
1:H:44:THR:HG22	1:H:45:VAL:N	2.33	0.43
1:H:50:ASP:HB3	1:H:319:ILE:CG2	2.46	0.43
1:G:109:MET:HE3	1:G:114:TYR:CD2	2.54	0.43
1:G:226:VAL:O	1:G:230:LYS:HG2	2.18	0.43
1:B:21:TYR:CE2	1:B:22:LEU:HG	2.53	0.43
1:B:186:PHE:CD2	1:B:198:GLU:HG2	2.54	0.43
1:D:159:SER:HB3	1:D:342:GLU:HG2	2.01	0.43
1:D:362:LEU:N	1:D:362:LEU:HD23	2.34	0.43
1:D:84:ASP:O	1:H:34:ARG:HD3	2.18	0.43
1:H:193:ASP:O	1:H:197:LYS:HG3	2.19	0.43
1:C:201:ILE:CG2	1:C:202:LEU:N	2.82	0.43
1:A:377:GLU:HB2	1:A:378:PRO:HD2	2.00	0.43
6:A:547:HOH:O	1:C:111:VAL:HG22	2.19	0.43
5:B:404:GOL:C3	6:B:689:HOH:O	2.66	0.43
1:G:188:SER:HG	1:G:189:PRO:HD3	1.81	0.43
1:G:166:ARG:HG2	1:G:198:GLU:OE1	2.19	0.42
1:H:279:VAL:HG22	1:H:311:ILE:CD1	2.48	0.42
1:B:300:MET:HB2	5:B:404:GOL:C1	2.49	0.42
1:E:164:ASP:CB	1:E:168:LYS:HZ1	2.32	0.42
1:F:276:VAL:HB	1:F:277:PRO:CD	2.49	0.42
1:G:165:THR:OG1	1:G:168:LYS:HG3	2.19	0.42
1:H:267:GLU:HG2	6:H:510:HOH:O	2.19	0.42
1:H:284:LEU:HD23	1:H:284:LEU:H	1.83	0.42
1:E:179:LYS:O	1:E:363:ASN:HA	2.19	0.42
1:F:166:ARG:NE	1:F:193:ASP:OD2	2.47	0.42
1:F:362:LEU:HD23	1:F:362:LEU:N	2.35	0.42
1:A:55:VAL:CG2	1:A:82:ILE:HD13	2.49	0.42
1:B:376[B]:VAL:HG12	1:B:377:GLU:N	2.34	0.42
1:F:291:MET:HE2	1:F:299:PHE:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:159:SER:HB2	1:G:342:GLU:OE2	2.19	0.42
1:H:199:MET:HB3	1:H:235:HIS:HB2	2.00	0.42
1:B:379:SER:O	1:B:383:GLN:HG3	2.20	0.42
1:E:164:ASP:CB	1:E:168:LYS:NZ	2.82	0.42
1:F:319:ILE:HA	1:F:347:ILE:HD13	2.01	0.42
1:H:190:VAL:O	1:H:190:VAL:CG1	2.67	0.42
1:A:157:TYR:CZ	1:A:342:GLU:HG3	2.55	0.42
1:C:176:TRP:CZ2	1:C:345:HIS:CG	3.07	0.42
1:F:97:ASP:HA	1:G:148:GLY:HA3	2.01	0.42
1:H:32:ASN:O	1:H:33:GLY:C	2.63	0.42
1:C:268:GLU:OE1	1:C:268:GLU:N	2.42	0.42
1:C:304:ALA:O	1:C:308:VAL:HG23	2.20	0.42
1:F:386:LEU:HD12	1:F:386:LEU:HA	1.92	0.42
1:H:221:THR:OG1	1:H:224:GLU:HG3	2.20	0.42
1:E:165:THR:OG1	1:E:168:LYS:HG3	2.20	0.42
1:F:11:PHE:CB	1:F:386:LEU:HG	2.50	0.42
1:G:188:SER:N	1:G:189:PRO:CD	2.83	0.42
1:H:3:ASP:O	1:H:4:ARG:HD2	2.19	0.42
1:C:32:ASN:OD1	1:C:32:ASN:C	2.63	0.42
1:G:65:GLY:HA3	1:G:133:LEU:HG	2.01	0.42
1:G:368:VAL:HA	1:G:369:PRO:HD3	1.91	0.41
1:A:83:ASP:OD2	1:A:389:HIS:ND1	2.37	0.41
1:B:318:THR:OG1	1:B:319:ILE:N	2.53	0.41
1:A:32:ASN:C	1:A:34:ARG:N	2.78	0.41
1:A:94:ASP:HA	1:A:95:PRO:HD3	1.88	0.41
1:A:362:LEU:HD23	1:A:362:LEU:H	1.85	0.41
1:D:356:VAL:HG22	1:D:377:GLU:HG3	2.01	0.41
1:B:133:LEU:HD23	1:B:133:LEU:HA	1.92	0.41
1:C:110:ARG:NH2	6:C:661:HOH:O	2.47	0.41
1:B:199:MET:HE3	1:B:199:MET:HB2	1.79	0.41
1:C:50:ASP:HB3	1:C:319:ILE:HG21	2.02	0.41
1:C:171:GLU:CD	1:C:205:ARG:HH21	2.28	0.41
1:E:27:PRO:O	1:E:29:GLU:HG3	2.21	0.41
1:H:69:THR:OG1	1:H:70:TYR:N	2.53	0.41
1:H:321:ALA:HA	1:H:351:ASN:HD22	1.85	0.41
1:A:94:ASP:O	1:A:97:ASP:HB2	2.21	0.41
1:A:171:GLU:O	1:A:174:ALA:N	2.53	0.41
1:E:27:PRO:C	1:E:29:GLU:N	2.78	0.41
1:F:160:GLY:HA2	1:F:185:LYS:O	2.20	0.41
1:B:192:ASP:HB2	6:B:733:HOH:O	2.21	0.41
1:G:166:ARG:NH2	1:G:193:ASP:CB	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:353:ARG:NH1	1:H:381:GLU:OE1	2.53	0.41
1:B:185:LYS:HA	1:B:213:ALA:O	2.20	0.41
1:B:297:THR:HA	5:B:404:GOL:H11	2.02	0.41
1:B:322:GLY:HA3	1:B:355:LEU:HD11	2.02	0.41
1:C:222:ALA:O	1:C:226:VAL:HG23	2.20	0.41
1:C:282:ARG:NE	6:C:520:HOH:O	2.53	0.41
1:C:345:HIS:CE1	1:C:349:GLU:HG3	2.55	0.41
1:D:307:HIS:HA	1:D:337:ASN:ND2	2.36	0.41
1:D:359:MET:HB2	1:D:369:PRO:HG3	2.01	0.41
1:E:349:GLU:CB	1:E:350:PRO:CD	2.99	0.41
1:G:233:GLU:N	1:G:234:PRO:CD	2.84	0.41
1:H:184:PHE:HE1	1:H:210:VAL:HG11	1.85	0.41
1:B:94:ASP:HA	1:B:95:PRO:HD3	1.96	0.41
1:D:85:LEU:HD22	1:H:73:VAL:HG11	2.03	0.41
1:D:142:VAL:HG23	1:D:373:GLY:O	2.21	0.41
1:E:170:ALA:HB2	1:E:201:ILE:HD13	2.02	0.41
1:H:12:THR:HG23	1:H:79:MET:CE	2.51	0.41
1:H:321:ALA:HA	1:H:351:ASN:ND2	2.36	0.41
1:C:332:SER:HB3	1:C:338:VAL:HG11	2.02	0.40
1:D:362:LEU:HD23	1:D:362:LEU:H	1.86	0.40
1:D:347:ILE:C	1:D:350:PRO:HD2	2.46	0.40
1:D:353:ARG:HG3	1:D:354:LEU:HG	2.03	0.40
1:G:68:GLU:HB2	1:G:323:ILE:HB	2.04	0.40
1:H:9:GLU:OE2	1:H:388:LYS:NZ	2.47	0.40
1:H:215:ASP:O	1:H:217:HIS:HD2	2.05	0.40
1:C:37:LEU:C	1:C:37:LEU:HD12	2.46	0.40
1:E:181:PHE:CD2	1:E:364:GLY:HA2	2.57	0.40
1:F:79:MET:SD	1:F:389:HIS:HB2	2.61	0.40
1:H:349:GLU:CB	1:H:350:PRO:CD	2.92	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	372/412 (90%)	363 (98%)	8 (2%)	1 (0%)	36	31
1	B	387/412 (94%)	378 (98%)	9 (2%)	0	100	100
1	C	375/412 (91%)	367 (98%)	8 (2%)	0	100	100
1	D	376/412 (91%)	366 (97%)	10 (3%)	0	100	100
1	E	378/412 (92%)	369 (98%)	8 (2%)	1 (0%)	36	31
1	F	374/412 (91%)	366 (98%)	8 (2%)	0	100	100
1	G	377/412 (92%)	367 (97%)	10 (3%)	0	100	100
1	H	374/412 (91%)	362 (97%)	12 (3%)	0	100	100
All	All	3013/3296 (91%)	2938 (98%)	73 (2%)	2 (0%)	48	45

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	33	GLY
1	E	28	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	290/323 (90%)	284 (98%)	6 (2%)	47	47
1	B	301/323 (93%)	293 (97%)	8 (3%)	39	37
1	C	293/323 (91%)	282 (96%)	11 (4%)	29	23
1	D	293/323 (91%)	289 (99%)	4 (1%)	59	61
1	E	294/323 (91%)	287 (98%)	7 (2%)	43	41
1	F	291/323 (90%)	285 (98%)	6 (2%)	47	47
1	G	293/323 (91%)	292 (100%)	1 (0%)	86	86
1	H	290/323 (90%)	276 (95%)	14 (5%)	23	15
All	All	2345/2584 (91%)	2288 (98%)	57 (2%)	43	41

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	37	LEU
1	A	153	ARG
1	A	201	ILE
1	A	345	HIS
1	A	347	ILE
1	B	11	PHE
1	B	60	GLU
1	B	197	LYS
1	B	199	MET
1	B	260	SER
1	B	312	LYS
1	B	339	ASP
1	B	362	LEU
1	C	15	LEU
1	C	40	LYS
1	C	48	THR
1	C	165	THR
1	C	192	ASP
1	C	247	GLU
1	C	254	ARG
1	C	345	HIS
1	C	346	SER
1	C	353	ARG
1	C	362	LEU
1	D	15	LEU
1	D	111	VAL
1	D	247	GLU
1	D	349	GLU
1	E	39	ARG
1	E	42	ASN
1	E	192	ASP
1	E	201	ILE
1	E	362	LEU
1	E	376	VAL
1	E	388	LYS
1	F	15	LEU
1	F	37	LEU
1	F	48	THR
1	F	247[A]	GLU
1	F	247[B]	GLU
1	F	386	LEU
1	G	345	HIS

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Mol	Chain	Res	Type
1	H	60	GLU
1	H	165	THR
1	H	172	LEU
1	H	190	VAL
1	H	192	ASP
1	H	193	ASP
1	H	201	ILE
1	H	245	ARG
1	H	249	ILE
1	H	260	SER
1	H	346	SER
1	H	362	LEU
1	H	376	VAL
1	H	386	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	235	HIS
1	A	273	HIS
1	A	310	HIS
1	A	337	ASN
1	A	363	ASN
1	B	235	HIS
1	B	310	HIS
1	B	363	ASN
1	C	177	GLN
1	C	235	HIS
1	C	273	HIS
1	C	351	ASN
1	D	337	ASN
1	D	345	HIS
1	E	42	ASN
1	E	310	HIS
1	E	345	HIS
1	F	307	HIS
1	F	345	HIS
1	F	351	ASN
1	G	345	HIS
1	H	217	HIS
1	H	351	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 17 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	FMT	H	402	2	2,2,2	0.74	0	1,1,1	0.56	0
3	FMT	F	402	2	2,2,2	0.60	0	1,1,1	0.73	0
3	FMT	C	404	2	2,2,2	0.70	0	1,1,1	0.67	0
3	FMT	G	403	2	2,2,2	0.74	0	1,1,1	0.60	0
3	FMT	E	402	2	2,2,2	0.71	0	1,1,1	0.60	0
5	GOL	B	404	-	5,5,5	0.18	0	5,5,5	0.72	0
3	FMT	A	402	2	2,2,2	0.68	0	1,1,1	0.64	0
3	FMT	B	403	2	2,2,2	0.84	0	1,1,1	0.65	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	B	404	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	404	GOL	O1-C1-C2-C3
5	B	404	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	404	GOL	5	0
3	B	403	FMT	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	376/412 (91%)	0.13	7 (1%) 66 66	26, 44, 78, 119	0
1	B	388/412 (94%)	-0.29	7 (1%) 67 68	22, 34, 67, 100	1 (0%)
1	C	378/412 (91%)	-0.09	4 (1%) 78 78	20, 38, 80, 131	1 (0%)
1	D	379/412 (91%)	-0.09	6 (1%) 70 70	24, 40, 69, 114	1 (0%)
1	E	381/412 (92%)	-0.30	4 (1%) 79 80	20, 34, 66, 113	1 (0%)
1	F	374/412 (90%)	-0.22	5 (1%) 75 75	20, 34, 67, 103	4 (1%)
1	G	380/412 (92%)	-0.41	2 (0%) 87 87	19, 31, 60, 111	1 (0%)
1	H	378/412 (91%)	0.04	17 (4%) 38 37	25, 39, 84, 128	0
All	All	3034/3296 (92%)	-0.16	52 (1%) 69 69	19, 37, 71, 131	9 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	19	THR	6.1
1	H	15	LEU	4.8
1	F	15	LEU	4.6
1	D	15	LEU	4.0
1	E	27	PRO	3.9
1	C	16	PRO	3.7
1	H	175	ALA	3.6
1	A	347	ILE	3.6
1	F	31	PRO	3.5
1	A	15	LEU	3.5
1	A	16	PRO	3.4
1	C	350	PRO	3.3
1	H	174	ALA	3.3
1	E	18	GLU	3.2
1	H	349	GLU	2.9
1	H	345	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	H	178	ALA	2.9
1	G	27	PRO	2.8
1	D	27	PRO	2.7
1	B	15	LEU	2.6
1	G	345	HIS	2.6
1	A	345	HIS	2.6
1	H	344	GLN	2.6
1	F	42	ASN	2.6
1	F	16	PRO	2.6
1	H	16	PRO	2.5
1	H	17	ARG	2.5
1	A	350	PRO	2.4
1	A	339	ASP	2.3
1	B	24	LYS	2.3
1	A	41	ALA	2.3
1	H	14	THR	2.3
1	C	49	PHE	2.2
1	B	27	PRO	2.2
1	D	347	ILE	2.2
1	C	362	LEU	2.2
1	F	345	HIS	2.2
1	H	347	ILE	2.2
1	B	33	GLY	2.2
1	E	194	GLY	2.2
1	H	172	LEU	2.2
1	H	176	TRP	2.2
1	D	28	GLY	2.1
1	B	22	LEU	2.1
1	B	21	TYR	2.1
1	H	204	GLU	2.1
1	H	173	ALA	2.1
1	D	16	PRO	2.1
1	H	350	PRO	2.1
1	H	170	ALA	2.1
1	B	25	PRO	2.0
1	D	350	PRO	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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
5	GOL	B	404	6/6	0.81	0.27	20,20,20,20	0
3	FMT	A	402	3/3	0.91	0.10	59,59,62,63	0
4	CL	C	403	1/1	0.94	0.08	58,58,58,58	0
3	FMT	C	404	3/3	0.96	0.06	37,37,44,48	0
3	FMT	H	402	3/3	0.96	0.09	34,34,48,65	0
3	FMT	B	403	3/3	0.97	0.07	36,36,39,39	0
3	FMT	E	402	3/3	0.97	0.08	27,27,37,37	0
4	CL	D	402[A]	1/1	0.97	0.05	27,27,27,27	1
4	CL	D	402[B]	1/1	0.97	0.05	28,28,28,28	1
3	FMT	F	402	3/3	0.97	0.07	31,31,39,43	0
4	CL	C	401[B]	1/1	0.98	0.07	29,29,29,29	1
3	FMT	G	403	3/3	0.98	0.05	22,22,38,40	0
4	CL	B	402[A]	1/1	0.98	0.10	27,27,27,27	1
4	CL	B	402[B]	1/1	0.98	0.10	26,26,26,26	1
4	CL	C	401[A]	1/1	0.98	0.07	30,30,30,30	1
2	MG	C	402	1/1	0.99	0.02	33,33,33,33	0
2	MG	D	401	1/1	0.99	0.02	29,29,29,29	0
2	MG	E	401	1/1	0.99	0.03	27,27,27,27	0
2	MG	F	401	1/1	0.99	0.03	32,32,32,32	0
2	MG	G	402	1/1	0.99	0.02	23,23,23,23	0
2	MG	H	401	1/1	0.99	0.02	37,37,37,37	0
4	CL	G	401[A]	1/1	0.99	0.03	23,23,23,23	1
4	CL	G	401[B]	1/1	0.99	0.03	23,23,23,23	1
2	MG	A	401	1/1	0.99	0.02	41,41,41,41	0
2	MG	B	401	1/1	1.00	0.03	24,24,24,24	0

6.5 Other polymers ⓘ

There are no such residues in this entry.