



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:35 AM UTC

PDB ID : 4TVU / pdb_00004tvu
Title : Crystal structure of trehalose synthase from *Deinococcus radiodurans* reveals a closed conformation for catalysis of the intramolecular isomerization
Authors : Wang, Y.L.; Chow, S.Y.; Lin, Y.T.; Liaw, S.H.
Deposited on : 2014-06-28
Resolution : 2.70 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:

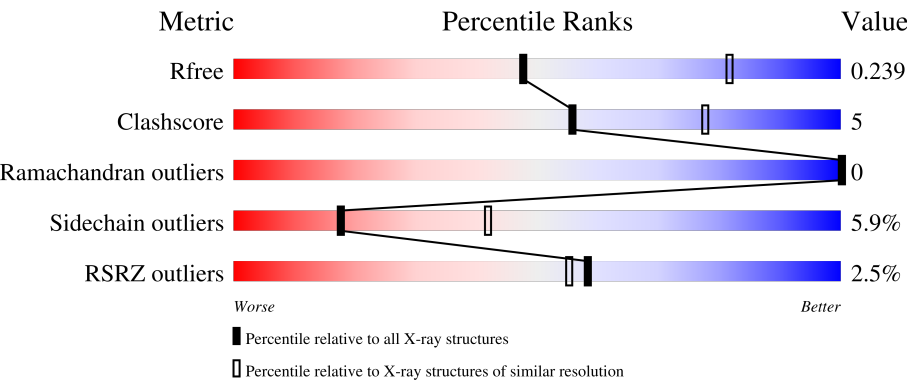
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	571	83%11%...
1	B	571	2% 84%10%..
1	C	571	2% 85%9%..
1	D	571	3% 85%10%..
1	E	571	% 85%9%..

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Mol	Chain	Length	Quality of chain
1	F	571	 2% 85% 9% ..
1	G	571	 % 84% 10% . .
1	H	571	 9% 81% 12% ..

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
4	TRS	B	602	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 36291 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 Trehalose synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			
1	B	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			
1	C	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			
1	D	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			
1	E	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			
1	F	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			
1	G	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			
1	H	548	Total	C	N	O	S	0	0	0
			4404	2818	751	819	16			

There are 184 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	expression tag	UNP I3NX86
A	0	VAL	-	expression tag	UNP I3NX86
A	1	PRO	-	expression tag	UNP I3NX86
A	97	TRP	ARG	engineered mutation	UNP I3NX86
A	313	ILE	THR	engineered mutation	UNP I3NX86
A	380	VAL	ILE	engineered mutation	UNP I3NX86
A	553	SER	-	expression tag	UNP I3NX86
A	554	ARG	-	expression tag	UNP I3NX86
A	555	VAL	-	expression tag	UNP I3NX86
A	556	ASP	-	expression tag	UNP I3NX86
A	557	LYS	-	expression tag	UNP I3NX86
A	558	LEU	-	expression tag	UNP I3NX86
A	559	ALA	-	expression tag	UNP I3NX86

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Chain	Residue	Modelled	Actual	Comment	Reference
A	560	ALA	-	expression tag	UNP I3NX86
A	561	ALA	-	expression tag	UNP I3NX86
A	562	LEU	-	expression tag	UNP I3NX86
A	563	GLU	-	expression tag	UNP I3NX86
A	564	HIS	-	expression tag	UNP I3NX86
A	565	HIS	-	expression tag	UNP I3NX86
A	566	HIS	-	expression tag	UNP I3NX86
A	567	HIS	-	expression tag	UNP I3NX86
A	568	HIS	-	expression tag	UNP I3NX86
A	569	HIS	-	expression tag	UNP I3NX86
B	-1	MET	-	expression tag	UNP I3NX86
B	0	VAL	-	expression tag	UNP I3NX86
B	1	PRO	-	expression tag	UNP I3NX86
B	97	TRP	ARG	engineered mutation	UNP I3NX86
B	313	ILE	THR	engineered mutation	UNP I3NX86
B	380	VAL	ILE	engineered mutation	UNP I3NX86
B	553	SER	-	expression tag	UNP I3NX86
B	554	ARG	-	expression tag	UNP I3NX86
B	555	VAL	-	expression tag	UNP I3NX86
B	556	ASP	-	expression tag	UNP I3NX86
B	557	LYS	-	expression tag	UNP I3NX86
B	558	LEU	-	expression tag	UNP I3NX86
B	559	ALA	-	expression tag	UNP I3NX86
B	560	ALA	-	expression tag	UNP I3NX86
B	561	ALA	-	expression tag	UNP I3NX86
B	562	LEU	-	expression tag	UNP I3NX86
B	563	GLU	-	expression tag	UNP I3NX86
B	564	HIS	-	expression tag	UNP I3NX86
B	565	HIS	-	expression tag	UNP I3NX86
B	566	HIS	-	expression tag	UNP I3NX86
B	567	HIS	-	expression tag	UNP I3NX86
B	568	HIS	-	expression tag	UNP I3NX86
B	569	HIS	-	expression tag	UNP I3NX86
C	-1	MET	-	expression tag	UNP I3NX86
C	0	VAL	-	expression tag	UNP I3NX86
C	1	PRO	-	expression tag	UNP I3NX86
C	97	TRP	ARG	engineered mutation	UNP I3NX86
C	313	ILE	THR	engineered mutation	UNP I3NX86
C	380	VAL	ILE	engineered mutation	UNP I3NX86
C	553	SER	-	expression tag	UNP I3NX86
C	554	ARG	-	expression tag	UNP I3NX86
C	555	VAL	-	expression tag	UNP I3NX86

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Chain	Residue	Modelled	Actual	Comment	Reference
C	556	ASP	-	expression tag	UNP I3NX86
C	557	LYS	-	expression tag	UNP I3NX86
C	558	LEU	-	expression tag	UNP I3NX86
C	559	ALA	-	expression tag	UNP I3NX86
C	560	ALA	-	expression tag	UNP I3NX86
C	561	ALA	-	expression tag	UNP I3NX86
C	562	LEU	-	expression tag	UNP I3NX86
C	563	GLU	-	expression tag	UNP I3NX86
C	564	HIS	-	expression tag	UNP I3NX86
C	565	HIS	-	expression tag	UNP I3NX86
C	566	HIS	-	expression tag	UNP I3NX86
C	567	HIS	-	expression tag	UNP I3NX86
C	568	HIS	-	expression tag	UNP I3NX86
C	569	HIS	-	expression tag	UNP I3NX86
D	-1	MET	-	expression tag	UNP I3NX86
D	0	VAL	-	expression tag	UNP I3NX86
D	1	PRO	-	expression tag	UNP I3NX86
D	97	TRP	ARG	engineered mutation	UNP I3NX86
D	313	ILE	THR	engineered mutation	UNP I3NX86
D	380	VAL	ILE	engineered mutation	UNP I3NX86
D	553	SER	-	expression tag	UNP I3NX86
D	554	ARG	-	expression tag	UNP I3NX86
D	555	VAL	-	expression tag	UNP I3NX86
D	556	ASP	-	expression tag	UNP I3NX86
D	557	LYS	-	expression tag	UNP I3NX86
D	558	LEU	-	expression tag	UNP I3NX86
D	559	ALA	-	expression tag	UNP I3NX86
D	560	ALA	-	expression tag	UNP I3NX86
D	561	ALA	-	expression tag	UNP I3NX86
D	562	LEU	-	expression tag	UNP I3NX86
D	563	GLU	-	expression tag	UNP I3NX86
D	564	HIS	-	expression tag	UNP I3NX86
D	565	HIS	-	expression tag	UNP I3NX86
D	566	HIS	-	expression tag	UNP I3NX86
D	567	HIS	-	expression tag	UNP I3NX86
D	568	HIS	-	expression tag	UNP I3NX86
D	569	HIS	-	expression tag	UNP I3NX86
E	-1	MET	-	expression tag	UNP I3NX86
E	0	VAL	-	expression tag	UNP I3NX86
E	1	PRO	-	expression tag	UNP I3NX86
E	97	TRP	ARG	engineered mutation	UNP I3NX86
E	313	ILE	THR	engineered mutation	UNP I3NX86

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Chain	Residue	Modelled	Actual	Comment	Reference
E	380	VAL	ILE	engineered mutation	UNP I3NX86
E	553	SER	-	expression tag	UNP I3NX86
E	554	ARG	-	expression tag	UNP I3NX86
E	555	VAL	-	expression tag	UNP I3NX86
E	556	ASP	-	expression tag	UNP I3NX86
E	557	LYS	-	expression tag	UNP I3NX86
E	558	LEU	-	expression tag	UNP I3NX86
E	559	ALA	-	expression tag	UNP I3NX86
E	560	ALA	-	expression tag	UNP I3NX86
E	561	ALA	-	expression tag	UNP I3NX86
E	562	LEU	-	expression tag	UNP I3NX86
E	563	GLU	-	expression tag	UNP I3NX86
E	564	HIS	-	expression tag	UNP I3NX86
E	565	HIS	-	expression tag	UNP I3NX86
E	566	HIS	-	expression tag	UNP I3NX86
E	567	HIS	-	expression tag	UNP I3NX86
E	568	HIS	-	expression tag	UNP I3NX86
E	569	HIS	-	expression tag	UNP I3NX86
F	-1	MET	-	expression tag	UNP I3NX86
F	0	VAL	-	expression tag	UNP I3NX86
F	1	PRO	-	expression tag	UNP I3NX86
F	97	TRP	ARG	engineered mutation	UNP I3NX86
F	313	ILE	THR	engineered mutation	UNP I3NX86
F	380	VAL	ILE	engineered mutation	UNP I3NX86
F	553	SER	-	expression tag	UNP I3NX86
F	554	ARG	-	expression tag	UNP I3NX86
F	555	VAL	-	expression tag	UNP I3NX86
F	556	ASP	-	expression tag	UNP I3NX86
F	557	LYS	-	expression tag	UNP I3NX86
F	558	LEU	-	expression tag	UNP I3NX86
F	559	ALA	-	expression tag	UNP I3NX86
F	560	ALA	-	expression tag	UNP I3NX86
F	561	ALA	-	expression tag	UNP I3NX86
F	562	LEU	-	expression tag	UNP I3NX86
F	563	GLU	-	expression tag	UNP I3NX86
F	564	HIS	-	expression tag	UNP I3NX86
F	565	HIS	-	expression tag	UNP I3NX86
F	566	HIS	-	expression tag	UNP I3NX86
F	567	HIS	-	expression tag	UNP I3NX86
F	568	HIS	-	expression tag	UNP I3NX86
F	569	HIS	-	expression tag	UNP I3NX86
G	-1	MET	-	expression tag	UNP I3NX86

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	VAL	-	expression tag	UNP I3NX86
G	1	PRO	-	expression tag	UNP I3NX86
G	97	TRP	ARG	engineered mutation	UNP I3NX86
G	313	ILE	THR	engineered mutation	UNP I3NX86
G	380	VAL	ILE	engineered mutation	UNP I3NX86
G	553	SER	-	expression tag	UNP I3NX86
G	554	ARG	-	expression tag	UNP I3NX86
G	555	VAL	-	expression tag	UNP I3NX86
G	556	ASP	-	expression tag	UNP I3NX86
G	557	LYS	-	expression tag	UNP I3NX86
G	558	LEU	-	expression tag	UNP I3NX86
G	559	ALA	-	expression tag	UNP I3NX86
G	560	ALA	-	expression tag	UNP I3NX86
G	561	ALA	-	expression tag	UNP I3NX86
G	562	LEU	-	expression tag	UNP I3NX86
G	563	GLU	-	expression tag	UNP I3NX86
G	564	HIS	-	expression tag	UNP I3NX86
G	565	HIS	-	expression tag	UNP I3NX86
G	566	HIS	-	expression tag	UNP I3NX86
G	567	HIS	-	expression tag	UNP I3NX86
G	568	HIS	-	expression tag	UNP I3NX86
G	569	HIS	-	expression tag	UNP I3NX86
H	-1	MET	-	expression tag	UNP I3NX86
H	0	VAL	-	expression tag	UNP I3NX86
H	1	PRO	-	expression tag	UNP I3NX86
H	97	TRP	ARG	engineered mutation	UNP I3NX86
H	313	ILE	THR	engineered mutation	UNP I3NX86
H	380	VAL	ILE	engineered mutation	UNP I3NX86
H	553	SER	-	expression tag	UNP I3NX86
H	554	ARG	-	expression tag	UNP I3NX86
H	555	VAL	-	expression tag	UNP I3NX86
H	556	ASP	-	expression tag	UNP I3NX86
H	557	LYS	-	expression tag	UNP I3NX86
H	558	LEU	-	expression tag	UNP I3NX86
H	559	ALA	-	expression tag	UNP I3NX86
H	560	ALA	-	expression tag	UNP I3NX86
H	561	ALA	-	expression tag	UNP I3NX86
H	562	LEU	-	expression tag	UNP I3NX86
H	563	GLU	-	expression tag	UNP I3NX86
H	564	HIS	-	expression tag	UNP I3NX86
H	565	HIS	-	expression tag	UNP I3NX86
H	566	HIS	-	expression tag	UNP I3NX86

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Chain	Residue	Modelled	Actual	Comment	Reference
H	567	HIS	-	expression tag	UNP I3NX86
H	568	HIS	-	expression tag	UNP I3NX86
H	569	HIS	-	expression tag	UNP I3NX86

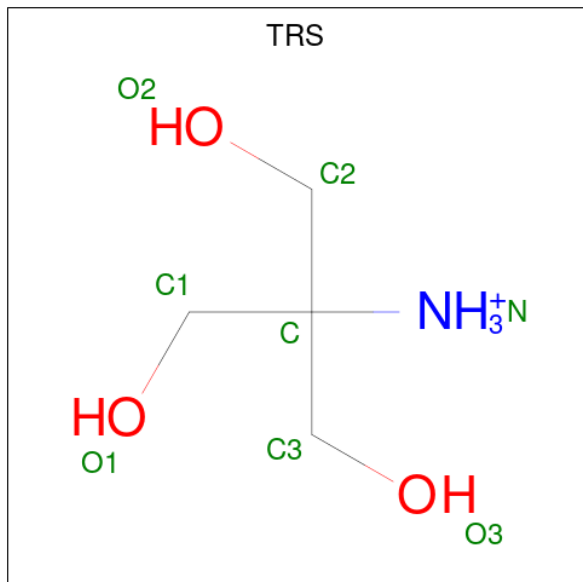
- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		
4	B	1	Total	C	N	O	0	0
			8	4	1	3		
4	C	1	Total	C	N	O	0	0
			8	4	1	3		
4	D	1	Total	C	N	O	0	0
			8	4	1	3		
4	E	1	Total	C	N	O	0	0
			8	4	1	3		
4	F	1	Total	C	N	O	0	0
			8	4	1	3		
4	G	1	Total	C	N	O	0	0
			8	4	1	3		
4	H	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	157	Total	O	0	0
			157	157		
5	B	147	Total	O	0	0
			147	147		

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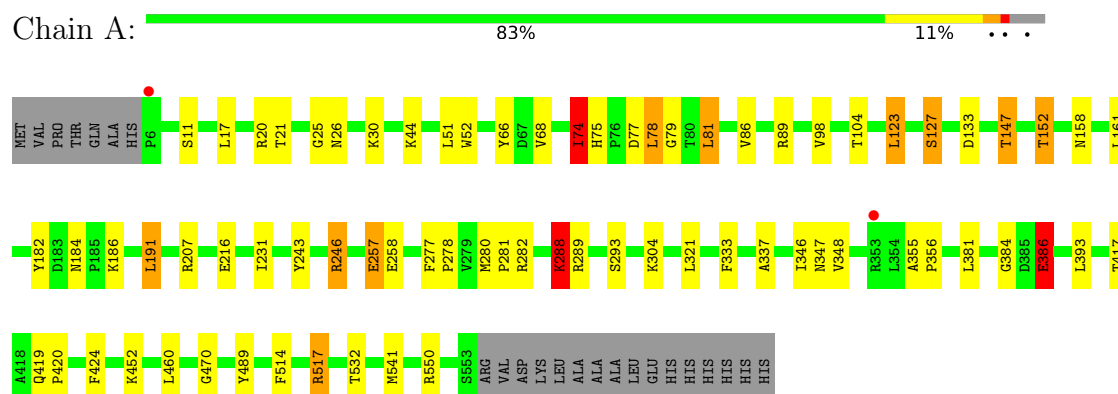
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	107	Total 107	O 107	0	0
5	D	99	Total 99	O 99	0	0
5	E	193	Total 193	O 193	0	0
5	F	85	Total 85	O 85	0	0
5	G	133	Total 133	O 133	0	0
5	H	58	Total 58	O 58	0	0

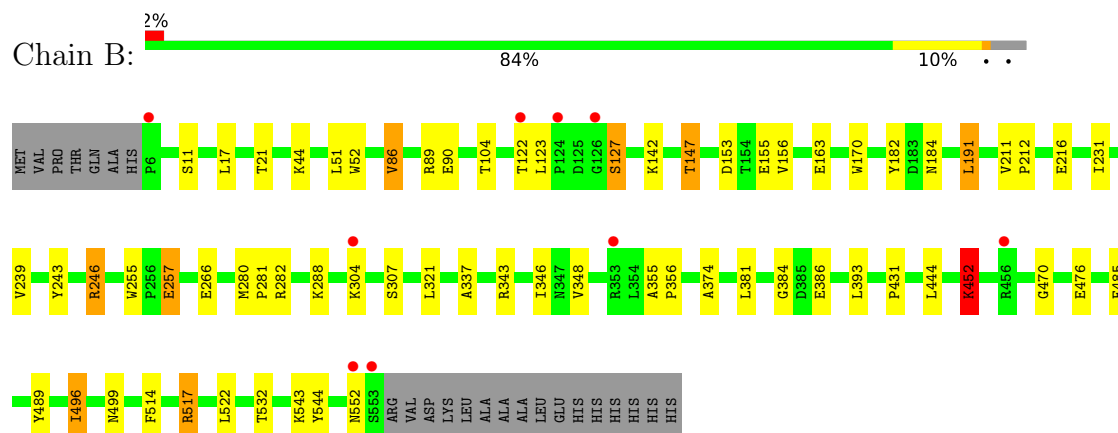
3 Residue-property plots

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.

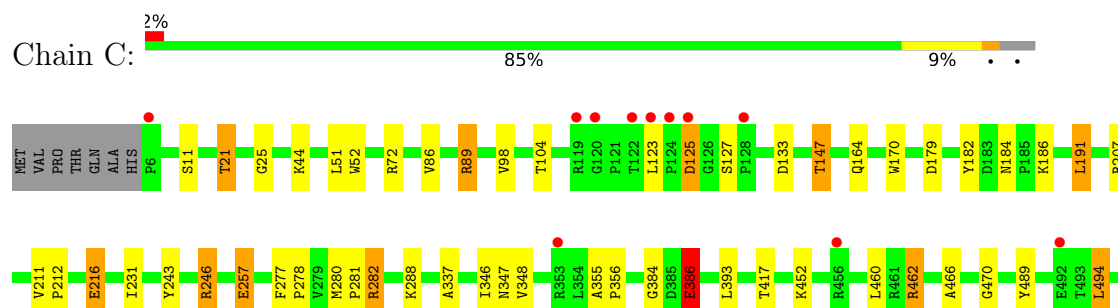
• Molecule 1: Trehalose synthase



• Molecule 1: Trehalose synthase

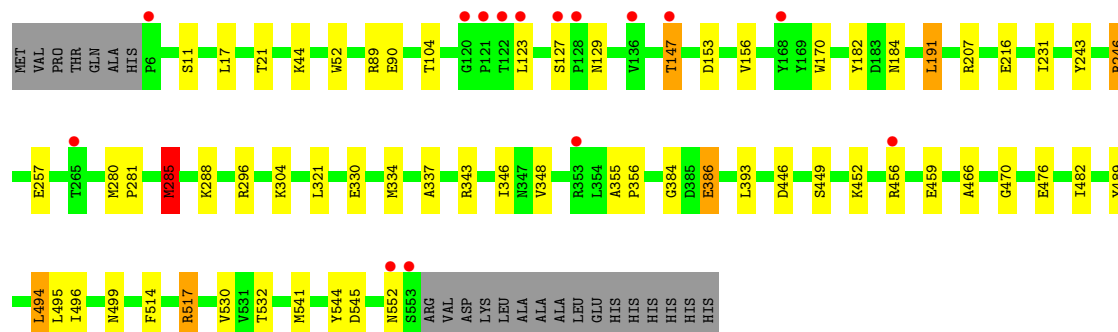
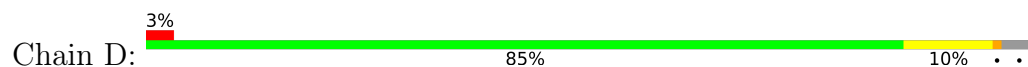


• Molecule 1: Trehalose synthase

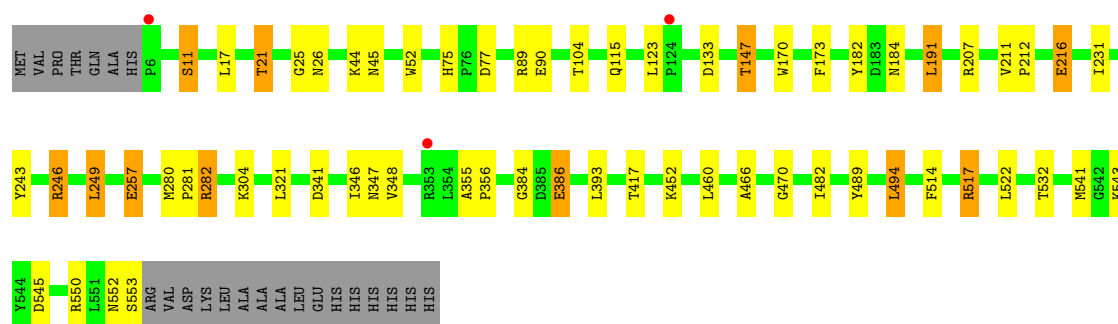
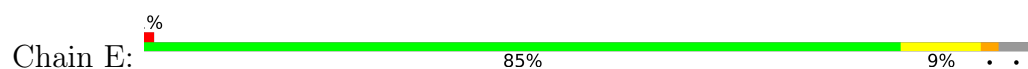




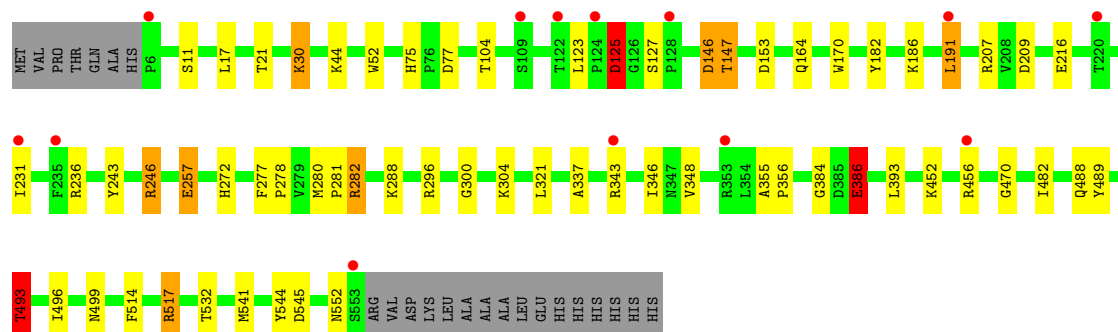
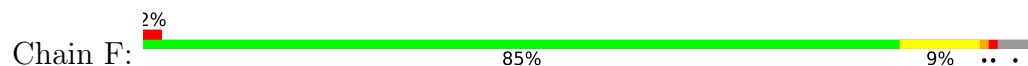
• Molecule 1: Trehalose synthase



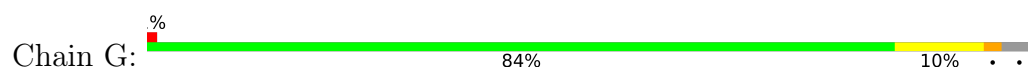
• Molecule 1: Trehalose synthase

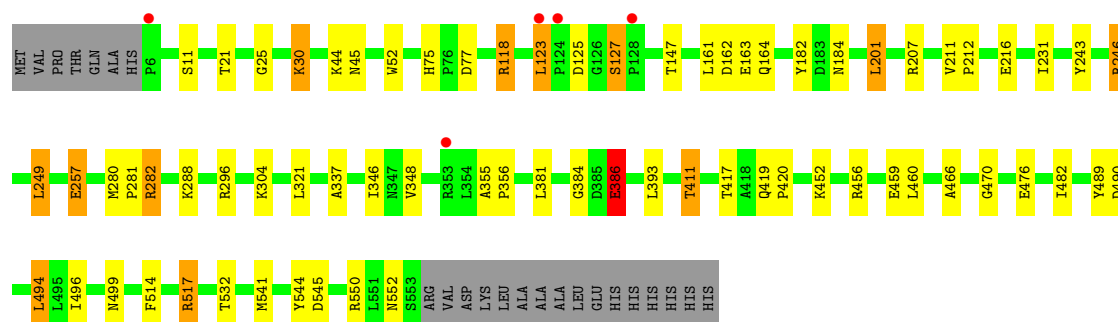


• Molecule 1: Trehalose synthase

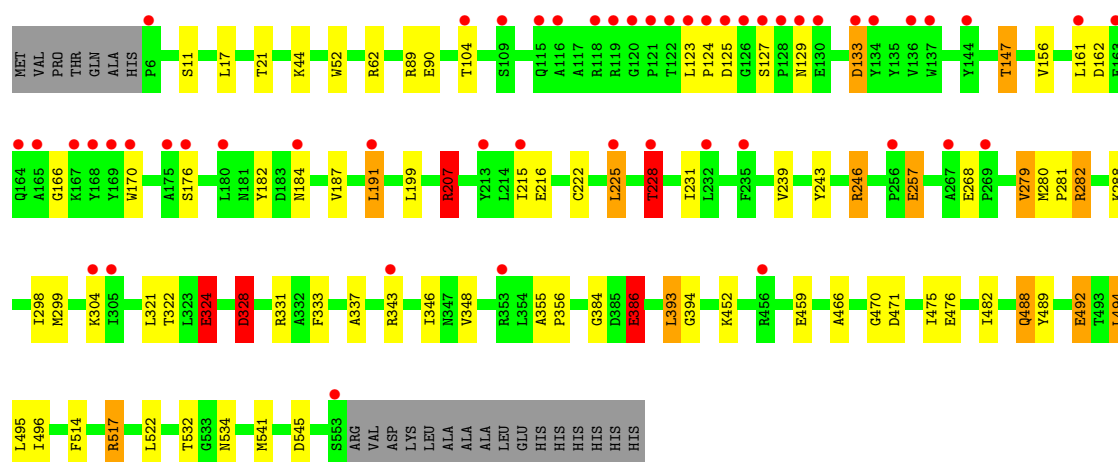
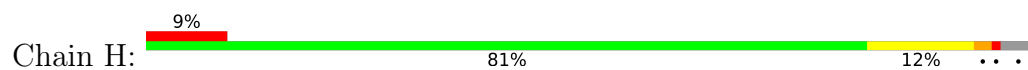


• Molecule 1: Trehalose synthase





• Molecule 1: Trehalose synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	95.61Å 195.90Å 130.99Å 90.00° 92.89° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 30.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.6 (30.00-2.70) 97.6 (30.00-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.04 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.194 , 0.236 0.197 , 0.239	Depositor DCC
R_{free} test set	6455 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.522	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	36291	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.65% 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: TRS, MG, CA

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	0.71	0/4537	0.87	7/6182 (0.1%)
1	B	0.72	0/4537	0.87	8/6182 (0.1%)
1	C	0.68	0/4537	0.86	5/6182 (0.1%)
1	D	0.66	0/4537	0.82	1/6182 (0.0%)
1	E	0.71	0/4537	0.84	2/6182 (0.0%)
1	F	0.67	0/4537	0.84	3/6182 (0.0%)
1	G	0.69	1/4537 (0.0%)	0.85	4/6182 (0.1%)
1	H	0.70	2/4537 (0.0%)	0.90	11/6182 (0.2%)
All	All	0.69	3/36296 (0.0%)	0.86	41/49456 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	324	GLU	CA-CB	5.99	1.63	1.53
1	G	490	ASP	CA-C	5.79	1.60	1.52
1	H	488	GLN	N-CA	5.02	1.52	1.46

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	152	THR	N-CA-CB	-9.38	96.34	110.12
1	H	475	ILE	CG1-CB-CG2	-7.26	88.91	110.70
1	D	285	MET	CG-SD-CE	-7.25	84.94	100.90
1	H	279	VAL	CB-CA-C	-7.09	100.77	112.26
1	G	490	ASP	CB-CA-C	7.09	124.52	110.42
1	B	452	LYS	CD-CE-NZ	7.08	134.56	111.90
1	H	246	ARG	CA-CB-CG	-7.07	99.97	114.10
1	B	239	VAL	N-CA-CB	6.94	119.46	110.57
1	A	74	ILE	CB-CA-C	6.86	120.78	110.63
1	H	328	ASP	N-CA-CB	6.72	120.57	110.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	207	ARG	CG-CD-NE	6.37	126.02	112.00
1	G	45	ASN	CB-CA-C	-6.04	100.75	110.79
1	E	543	LYS	CD-CE-NZ	5.95	130.94	111.90
1	G	411	THR	CB-CA-C	5.88	119.07	109.90
1	B	452	LYS	CG-CD-CE	-5.79	97.98	111.30
1	H	228	THR	N-CA-CB	5.76	118.37	110.01
1	C	543	LYS	CG-CD-CE	-5.72	98.14	111.30
1	H	239	VAL	N-CA-CB	5.69	119.12	110.58
1	A	347	ASN	CB-CA-C	-5.62	103.23	112.06
1	A	30	LYS	CD-CE-NZ	-5.60	93.97	111.90
1	B	255	TRP	CA-C-N	5.50	124.96	119.24
1	B	255	TRP	C-N-CA	5.50	124.96	119.24
1	H	386	GLU	CB-CA-C	-5.48	100.49	110.63
1	F	493	THR	N-CA-CB	5.46	119.28	110.65
1	B	543	LYS	CG-CD-CE	-5.37	98.94	111.30
1	C	386	GLU	CB-CA-C	-5.28	100.86	110.63
1	E	347	ASN	CB-CA-C	-5.26	103.80	112.06
1	B	155	GLU	CB-CG-CD	5.26	121.54	112.60
1	F	125	ASP	CB-CA-C	5.22	117.83	109.27
1	B	86	VAL	N-CA-CB	5.21	118.39	110.58
1	C	347	ASN	CB-CA-C	-5.21	103.88	112.06
1	C	541	MET	CG-SD-CE	5.20	112.35	100.90
1	F	386	GLU	CB-CA-C	-5.17	101.06	110.63
1	A	288	LYS	CA-CB-CG	5.15	124.40	114.10
1	H	488	GLN	CB-CA-C	5.15	119.13	109.71
1	H	207	ARG	CB-CG-CD	5.13	123.10	111.30
1	G	386	GLU	CB-CA-C	-5.13	101.14	110.63
1	C	125	ASP	CB-CA-C	-5.12	99.94	110.38
1	H	488	GLN	N-CA-CB	5.04	119.14	110.68
1	A	541	MET	CG-SD-CE	5.01	111.92	100.90
1	A	386	GLU	CB-CA-C	-5.01	101.37	110.63

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	4404	0	4201	38	0
1	B	4404	0	4201	32	0
1	C	4404	0	4201	40	0
1	D	4404	0	4201	47	0
1	E	4404	0	4201	48	0
1	F	4404	0	4201	44	0
1	G	4404	0	4201	47	0
1	H	4404	0	4201	58	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	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	8	0	12	1	0
4	B	8	0	12	0	0
4	C	8	0	12	0	0
4	D	8	0	12	0	0
4	E	8	0	12	1	0
4	F	8	0	12	0	0
4	G	8	0	12	0	0
4	H	8	0	12	0	0
5	A	157	0	0	1	0
5	B	147	0	0	3	0
5	C	107	0	0	4	0
5	D	99	0	0	3	0
5	E	193	0	0	6	0
5	F	85	0	0	1	0
5	G	133	0	0	1	0
5	H	58	0	0	6	0
All	All	36291	0	33704	325	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 5.

All (325) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:182:TYR:OH	1:D:191:LEU:CD1	2.07	1.02
1:C:182:TYR:OH	1:C:191:LEU:CD1	2.08	1.01
1:E:482:ILE:HG23	1:E:541:MET:HE1	1.47	0.96
1:F:482:ILE:HG23	1:F:541:MET:HE1	1.48	0.96
1:G:482:ILE:HG23	1:G:541:MET:HE1	1.48	0.95
1:H:482:ILE:HG23	1:H:541:MET:HE1	1.48	0.95
1:G:118:ARG:NH1	1:G:162:ASP:OD2	2.01	0.93
1:D:482:ILE:HG23	1:D:541:MET:HE1	1.48	0.92
1:B:123:LEU:HB2	5:B:819:HOH:O	1.71	0.87
1:D:182:TYR:OH	1:D:191:LEU:HD13	1.75	0.86
1:B:127:SER:N	5:B:819:HOH:O	2.08	0.85
1:H:125:ASP:HB2	5:H:718:HOH:O	1.76	0.85
1:D:52:TRP:CZ2	1:D:207:ARG:HD3	2.12	0.85
1:F:52:TRP:CZ2	1:F:207:ARG:HD3	2.10	0.85
1:C:182:TYR:OH	1:C:191:LEU:HD13	1.76	0.84
1:A:52:TRP:CZ2	1:A:207:ARG:HD3	2.11	0.84
1:G:52:TRP:CZ2	1:G:207:ARG:HD3	2.13	0.84
1:C:52:TRP:CZ2	1:C:207:ARG:HD3	2.14	0.83
1:E:52:TRP:CZ2	1:E:207:ARG:HD3	2.14	0.81
1:D:182:TYR:OH	1:D:191:LEU:HD11	1.80	0.81
1:C:182:TYR:OH	1:C:191:LEU:HD11	1.81	0.77
1:H:52:TRP:CZ2	1:H:207:ARG:HD3	2.20	0.76
1:C:179:ASP:HB3	5:C:807:HOH:O	1.87	0.74
1:C:216:GLU:OE1	5:C:807:HOH:O	2.05	0.73
1:H:184:ASN:OD1	1:H:187:VAL:HG23	1.89	0.73
1:B:374:ALA:CB	1:B:496:ILE:HD13	2.18	0.72
1:H:199:LEU:HD22	1:H:246:ARG:HD3	1.72	0.71
1:H:215:ILE:HB	1:H:228:THR:HG22	1.74	0.69
1:B:444:LEU:O	1:B:452:LYS:NZ	2.25	0.69
1:H:492:GLU:HA	1:H:492:GLU:OE1	1.93	0.69
1:E:216:GLU:HG2	5:E:745:HOH:O	1.94	0.68
1:B:485:PHE:CZ	1:B:496:ILE:HG12	2.28	0.68
1:G:296:ARG:NH1	5:G:770:HOH:O	2.27	0.68
1:H:222:CYS:O	5:H:708:HOH:O	2.12	0.68
1:E:482:ILE:CG2	1:E:541:MET:HE1	2.23	0.67
1:F:482:ILE:HG12	1:F:541:MET:HE3	1.77	0.67
1:E:482:ILE:HG12	1:E:541:MET:HE3	1.77	0.67
1:E:75:HIS:ND1	1:E:77:ASP:OD1	2.28	0.66
1:G:482:ILE:HG12	1:G:541:MET:HE3	1.76	0.66
1:E:341:ASP:HB3	5:E:868:HOH:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:482:ILE:CG2	1:F:541:MET:HE1	2.25	0.66
1:D:482:ILE:CG2	1:D:541:MET:HE1	2.25	0.66
1:F:30:LYS:HE3	1:F:77:ASP:CB	2.24	0.66
1:H:482:ILE:HG12	1:H:541:MET:HE3	1.76	0.66
1:D:482:ILE:HG12	1:D:541:MET:HE3	1.77	0.65
1:F:125:ASP:OD1	1:F:125:ASP:O	2.15	0.65
1:G:118:ARG:NH1	1:G:162:ASP:CG	2.54	0.65
1:H:133:ASP:OD1	1:H:133:ASP:N	2.29	0.65
1:H:328:ASP:OD1	1:H:331:ARG:NH1	2.30	0.65
1:A:288:LYS:HG3	1:A:337:ALA:HB1	1.78	0.65
1:G:30:LYS:HE3	1:G:77:ASP:CB	2.27	0.65
1:D:285:MET:HE1	1:D:330:GLU:HB3	1.79	0.64
1:F:75:HIS:ND1	1:F:77:ASP:OD1	2.31	0.64
1:G:207:ARG:HG3	1:G:249:LEU:HD13	1.79	0.64
1:H:243:TYR:HB3	1:H:246:ARG:HD2	1.79	0.64
1:D:285:MET:HE3	1:D:334:MET:HG2	1.80	0.64
1:H:482:ILE:CG2	1:H:541:MET:HE1	2.25	0.64
1:E:21:THR:HG22	5:E:725:HOH:O	1.97	0.63
1:E:207:ARG:HG3	1:E:249:LEU:HD13	1.81	0.63
1:G:75:HIS:ND1	1:G:77:ASP:OD1	2.32	0.62
1:A:74:ILE:HD11	1:A:79:GLY:O	2.00	0.62
1:G:482:ILE:CG2	1:G:541:MET:HE1	2.24	0.61
1:H:333:PHE:CD1	1:H:333:PHE:C	2.78	0.61
1:C:545:ASP:OD2	1:D:545:ASP:OD2	2.18	0.61
1:H:282:ARG:HD3	1:H:282:ARG:N	2.15	0.61
1:F:236:ARG:CZ	1:F:272:HIS:CD2	2.84	0.61
1:G:282:ARG:HD3	1:G:282:ARG:N	2.15	0.60
1:A:75:HIS:ND1	1:A:77:ASP:OD1	2.34	0.59
1:F:30:LYS:HE3	1:F:77:ASP:HB3	1.84	0.58
1:A:25:GLY:HA3	1:E:417:THR:HG21	1.85	0.58
1:A:280:MET:HB3	1:A:281:PRO:HD3	1.85	0.58
1:C:282:ARG:HD3	1:C:282:ARG:N	2.18	0.57
1:D:52:TRP:CE2	1:D:207:ARG:HD3	2.39	0.57
1:F:488:GLN:HG2	1:F:493:THR:HB	1.86	0.57
1:C:104:THR:HG22	1:C:191:LEU:HD22	1.87	0.57
1:H:225:LEU:HB2	1:H:228:THR:HG23	1.85	0.56
1:E:282:ARG:HD3	1:E:282:ARG:N	2.19	0.56
1:C:553:SER:O	1:H:156:VAL:HG22	2.05	0.56
1:F:282:ARG:HD3	1:F:282:ARG:N	2.20	0.56
1:G:52:TRP:CE2	1:G:207:ARG:HD3	2.40	0.56
1:C:417:THR:HG21	1:G:25:GLY:HA3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:52:TRP:CE2	1:E:207:ARG:HD3	2.41	0.55
1:D:104:THR:HG22	1:D:191:LEU:HD22	1.88	0.55
1:B:431:PRO:HD3	1:E:45:ASN:ND2	2.21	0.55
1:H:123:LEU:HB3	1:H:124:PRO:HD2	1.88	0.55
1:C:72:ARG:NH1	5:C:711:HOH:O	2.40	0.55
1:G:30:LYS:HE3	1:G:77:ASP:HB3	1.89	0.55
1:A:74:ILE:HD11	1:A:79:GLY:C	2.32	0.55
1:D:343:ARG:NH2	1:G:459:GLU:OE1	2.30	0.55
1:D:343:ARG:NH2	1:G:459:GLU:HB3	2.22	0.55
1:F:52:TRP:CE2	1:F:207:ARG:HD3	2.40	0.54
1:B:374:ALA:HB1	1:B:496:ILE:HD13	1.88	0.54
1:F:123:LEU:C	1:F:125:ASP:H	2.15	0.54
1:F:280:MET:HB3	1:F:281:PRO:HD3	1.89	0.54
1:D:153:ASP:O	1:G:550:ARG:NH2	2.36	0.54
1:D:280:MET:HB3	1:D:281:PRO:HD3	1.89	0.54
1:C:52:TRP:CE2	1:C:207:ARG:HD3	2.41	0.54
1:D:456:ARG:NH1	5:D:795:HOH:O	2.38	0.54
1:H:534:ASN:HB3	5:H:734:HOH:O	2.07	0.54
1:H:280:MET:HB3	1:H:281:PRO:HD3	1.90	0.53
1:C:280:MET:HB3	1:C:281:PRO:HD3	1.90	0.53
1:B:280:MET:HB3	1:B:281:PRO:HD3	1.90	0.53
1:A:52:TRP:CE2	1:A:207:ARG:HD3	2.42	0.53
1:G:280:MET:HB3	1:G:281:PRO:HD3	1.90	0.53
1:E:280:MET:HB3	1:E:281:PRO:HD3	1.91	0.53
1:H:222:CYS:HB3	5:H:708:HOH:O	2.09	0.53
1:G:470:GLY:HA2	1:G:489:TYR:HB2	1.91	0.52
1:H:222:CYS:HA	1:H:225:LEU:HD21	1.90	0.52
1:C:462:ARG:NH2	1:H:394:GLY:HA3	2.24	0.52
1:A:419:GLN:HG3	1:A:420:PRO:HD2	1.91	0.52
1:E:470:GLY:HA2	1:E:489:TYR:HB2	1.92	0.52
1:D:470:GLY:HA2	1:D:489:TYR:HB2	1.92	0.51
1:E:466:ALA:HB3	1:E:494:LEU:HD22	1.92	0.51
1:D:123:LEU:HD21	1:D:129:ASN:HD22	1.76	0.51
1:G:419:GLN:HG3	1:G:420:PRO:HD2	1.92	0.51
1:H:162:ASP:HB3	5:H:714:HOH:O	2.11	0.50
1:H:470:GLY:HA2	1:H:489:TYR:HB2	1.94	0.50
1:H:482:ILE:HG12	1:H:541:MET:CE	2.41	0.50
1:F:236:ARG:CZ	1:F:272:HIS:HD2	2.25	0.50
1:C:470:GLY:HA2	1:C:489:TYR:HB2	1.93	0.50
1:G:545:ASP:OD2	1:H:545:ASP:OD2	2.30	0.50
1:H:355:ALA:HB3	1:H:356:PRO:HD3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:ASN:C	1:E:26:ASN:HA	2.37	0.50
1:F:470:GLY:HA2	1:F:489:TYR:HB2	1.94	0.50
1:G:482:ILE:HG12	1:G:541:MET:CE	2.42	0.50
1:C:182:TYR:HB3	1:C:231:ILE:CD1	2.42	0.50
1:F:482:ILE:HG12	1:F:541:MET:CE	2.42	0.50
1:B:470:GLY:HA2	1:B:489:TYR:HB2	1.93	0.49
1:E:482:ILE:HG12	1:E:541:MET:CE	2.42	0.49
1:F:30:LYS:HE3	1:F:77:ASP:HB2	1.93	0.49
1:C:466:ALA:HB3	1:C:494:LEU:HD22	1.95	0.49
1:H:279:VAL:CG1	1:H:299:MET:SD	3.01	0.49
1:A:470:GLY:HA2	1:A:489:TYR:HB2	1.94	0.49
1:B:182:TYR:HB3	1:B:231:ILE:CD1	2.43	0.49
1:H:182:TYR:HB3	1:H:231:ILE:CD1	2.42	0.49
1:H:343:ARG:NH2	1:H:393:LEU:O	2.39	0.49
1:H:322:THR:OG1	1:H:324:GLU:CG	2.61	0.49
1:C:460:LEU:HD13	1:H:343:ARG:CD	2.43	0.48
1:D:482:ILE:HG12	1:D:541:MET:CE	2.42	0.48
1:D:89:ARG:NH1	1:D:90:GLU:OE1	2.47	0.48
1:D:296:ARG:HD3	5:D:719:HOH:O	2.12	0.48
1:F:182:TYR:HB3	1:F:231:ILE:CD1	2.43	0.48
1:B:17:LEU:C	1:B:17:LEU:HD12	2.38	0.48
1:E:355:ALA:HB3	1:E:356:PRO:HD3	1.96	0.48
1:F:243:TYR:HB3	1:F:246:ARG:HG2	1.95	0.48
1:H:466:ALA:HB3	1:H:494:LEU:HD22	1.95	0.48
1:B:153:ASP:O	1:E:550:ARG:NH2	2.39	0.48
1:C:460:LEU:HD13	1:H:343:ARG:HD2	1.95	0.48
1:E:182:TYR:HB3	1:E:231:ILE:CD1	2.43	0.48
1:G:466:ALA:HB3	1:G:494:LEU:HD22	1.94	0.48
1:G:30:LYS:HE3	1:G:77:ASP:HB2	1.95	0.48
1:E:243:TYR:HB3	1:E:246:ARG:HG2	1.96	0.48
1:E:173:PHE:CE1	4:E:602:TRS:H12	2.49	0.47
1:A:182:TYR:HB3	1:A:231:ILE:CD1	2.43	0.47
1:B:243:TYR:HB3	1:B:246:ARG:HG2	1.95	0.47
1:H:215:ILE:H	1:H:228:THR:HG22	1.79	0.47
1:F:300:GLY:C	5:F:701:HOH:O	2.57	0.47
1:F:355:ALA:HB3	1:F:356:PRO:HD3	1.97	0.47
1:H:279:VAL:HG11	1:H:299:MET:SD	2.54	0.47
1:A:81:LEU:O	1:A:81:LEU:HD22	2.15	0.47
1:D:182:TYR:HB3	1:D:231:ILE:CD1	2.44	0.47
1:D:243:TYR:HB3	1:D:246:ARG:HG2	1.96	0.47
1:D:466:ALA:HB3	1:D:494:LEU:HD22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:482:ILE:CG2	1:E:541:MET:CE	2.92	0.47
1:G:118:ARG:NH1	1:G:162:ASP:OD1	2.47	0.47
1:G:182:TYR:HB3	1:G:231:ILE:CD1	2.45	0.47
1:G:243:TYR:HB3	1:G:246:ARG:HG2	1.97	0.47
1:G:355:ALA:HB3	1:G:356:PRO:HD3	1.96	0.47
1:H:104:THR:HG22	1:H:191:LEU:CD2	2.45	0.47
1:H:104:THR:HG22	1:H:191:LEU:HD22	1.97	0.47
1:A:104:THR:HG22	1:A:191:LEU:CD2	2.45	0.47
1:C:25:GLY:HA3	1:G:417:THR:HG21	1.97	0.47
1:A:75:HIS:HB3	1:A:78:LEU:HD22	1.97	0.47
1:A:104:THR:HG22	1:A:191:LEU:HD22	1.97	0.47
1:C:417:THR:HG21	1:G:25:GLY:CA	2.45	0.46
1:B:355:ALA:HB3	1:B:356:PRO:HD3	1.95	0.46
1:C:21:THR:HG22	5:C:720:HOH:O	2.14	0.46
1:C:243:TYR:HB3	1:C:246:ARG:HG2	1.96	0.46
1:F:104:THR:HG22	1:F:191:LEU:CD2	2.46	0.46
1:F:125:ASP:OD1	1:F:125:ASP:C	2.59	0.46
1:A:355:ALA:HB3	1:A:356:PRO:HD3	1.98	0.46
1:A:384:GLY:N	1:A:386:GLU:OE2	2.48	0.46
1:C:460:LEU:HD13	1:H:343:ARG:NE	2.31	0.46
1:E:45:ASN:HB3	5:E:832:HOH:O	2.15	0.46
1:A:243:TYR:HB3	1:A:246:ARG:HG2	1.96	0.46
1:B:89:ARG:NH1	1:B:90:GLU:OE1	2.48	0.46
1:A:26:ASN:HA	1:E:26:ASN:C	2.41	0.46
1:A:52:TRP:CD1	1:A:52:TRP:C	2.93	0.46
1:E:384:GLY:N	1:E:386:GLU:OE2	2.49	0.46
1:C:355:ALA:HB3	1:C:356:PRO:HD3	1.98	0.45
1:D:384:GLY:N	1:D:386:GLU:OE2	2.49	0.45
1:E:89:ARG:NH1	1:E:90:GLU:OE1	2.48	0.45
1:H:166:GLY:N	5:H:714:HOH:O	2.49	0.45
1:H:384:GLY:N	1:H:386:GLU:OE2	2.49	0.45
1:A:289:ARG:HG3	1:A:333:PHE:CZ	2.50	0.45
1:B:257:GLU:H	1:B:257:GLU:CD	2.25	0.45
1:F:104:THR:HG22	1:F:191:LEU:HD22	1.98	0.45
1:G:482:ILE:CG2	1:G:541:MET:CE	2.92	0.45
1:H:123:LEU:HD11	1:H:129:ASN:HA	1.97	0.45
1:H:304:LYS:H	1:H:304:LYS:HD2	1.81	0.45
1:D:288:LYS:HG3	1:D:337:ALA:HB1	1.99	0.45
1:E:104:THR:HG22	1:E:191:LEU:CD2	2.47	0.45
1:E:545:ASP:OD2	1:F:545:ASP:OD2	2.34	0.45
1:B:104:THR:HG22	1:B:191:LEU:HD22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:TRP:CD1	1:B:52:TRP:C	2.95	0.45
1:D:156:VAL:HG13	1:G:552:ASN:HB3	1.98	0.45
1:H:89:ARG:NH1	1:H:90:GLU:OE1	2.49	0.45
1:B:170:TRP:NE1	1:B:216:GLU:HG2	2.31	0.45
1:F:482:ILE:CG2	1:F:541:MET:CE	2.94	0.45
1:H:322:THR:OG1	1:H:324:GLU:HG2	2.17	0.45
1:B:266:GLU:OE1	1:B:307:SER:OG	2.25	0.45
1:D:355:ALA:HB3	1:D:356:PRO:HD3	1.99	0.45
1:E:104:THR:HG22	1:E:191:LEU:HD22	1.98	0.45
1:H:52:TRP:CD1	1:H:52:TRP:C	2.95	0.45
1:H:482:ILE:CG2	1:H:541:MET:CE	2.93	0.45
1:B:104:THR:HG22	1:B:191:LEU:CD2	2.46	0.44
1:E:321:LEU:HD23	1:E:321:LEU:C	2.42	0.44
1:H:62:ARG:HE	1:H:176:SER:CB	2.30	0.44
1:B:156:VAL:HG13	1:E:552:ASN:HB3	1.99	0.44
1:F:52:TRP:CD1	1:F:52:TRP:C	2.94	0.44
1:E:115:GLN:HG3	5:E:801:HOH:O	2.16	0.44
1:B:343:ARG:HD3	1:E:460:LEU:HD13	1.99	0.44
1:F:384:GLY:N	1:F:386:GLU:OE2	2.49	0.44
1:D:482:ILE:CG2	1:D:541:MET:CE	2.93	0.44
1:A:25:GLY:CA	1:E:417:THR:HG21	2.47	0.44
1:A:460:LEU:HD13	1:F:343:ARG:HD3	2.00	0.44
1:D:499:ASN:O	1:D:544:TYR:HA	2.18	0.44
1:F:246:ARG:N	1:F:246:ARG:HD3	2.33	0.43
1:G:384:GLY:N	1:G:386:GLU:OE2	2.50	0.43
1:A:66:TYR:CD1	4:A:602:TRS:H31	2.53	0.43
1:B:288:LYS:HG3	1:B:337:ALA:HB1	2.00	0.43
1:E:514:PHE:HB3	1:E:517:ARG:HG3	1.99	0.43
1:E:184:ASN:OD1	1:E:184:ASN:C	2.62	0.43
1:A:417:THR:HG21	1:E:25:GLY:HA3	1.99	0.43
1:D:17:LEU:C	1:D:17:LEU:HD12	2.44	0.43
1:D:343:ARG:NH1	1:G:456:ARG:HG3	2.33	0.43
1:D:530:VAL:HG23	5:D:706:HOH:O	2.18	0.43
1:E:257:GLU:H	1:E:257:GLU:CD	2.26	0.43
1:H:17:LEU:HD12	1:H:17:LEU:C	2.43	0.43
1:H:321:LEU:C	1:H:321:LEU:HD23	2.44	0.43
1:A:257:GLU:CD	1:A:257:GLU:H	2.27	0.43
1:D:246:ARG:HD3	1:D:246:ARG:N	2.34	0.43
1:F:517:ARG:HB3	1:F:552:ASN:O	2.19	0.43
1:H:147:THR:HG21	1:H:170:TRP:HH2	1.84	0.43
1:A:321:LEU:C	1:A:321:LEU:HD23	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:343:ARG:HD2	1:G:460:LEU:HD13	2.00	0.43
1:F:321:LEU:C	1:F:321:LEU:HD23	2.44	0.43
1:H:495:LEU:C	1:H:496:ILE:HD13	2.43	0.43
1:C:288:LYS:HG3	1:C:337:ALA:HB1	2.01	0.42
1:F:17:LEU:HD12	1:F:17:LEU:C	2.43	0.42
1:A:147:THR:HG23	1:A:158:ASN:HD22	1.84	0.42
1:C:52:TRP:CD1	1:C:52:TRP:C	2.96	0.42
1:E:147:THR:HG21	1:E:170:TRP:HH2	1.84	0.42
1:F:257:GLU:CD	1:F:257:GLU:H	2.27	0.42
1:B:321:LEU:C	1:B:321:LEU:HD23	2.45	0.42
1:B:517:ARG:HB3	1:B:552:ASN:O	2.19	0.42
1:D:285:MET:HE3	1:D:334:MET:CG	2.49	0.42
1:H:222:CYS:HA	1:H:225:LEU:CD2	2.49	0.42
1:B:184:ASN:OD1	1:B:184:ASN:C	2.63	0.42
1:C:257:GLU:CD	1:C:257:GLU:H	2.27	0.42
1:D:321:LEU:C	1:D:321:LEU:HD23	2.43	0.42
1:H:288:LYS:HG3	1:H:337:ALA:HB1	2.02	0.42
1:A:246:ARG:N	1:A:246:ARG:HD3	2.34	0.42
1:D:52:TRP:CD1	1:D:52:TRP:C	2.97	0.42
1:E:482:ILE:CG1	1:E:541:MET:HE3	2.48	0.42
1:H:514:PHE:HB3	1:H:517:ARG:HG3	2.00	0.42
1:B:499:ASN:O	1:B:544:TYR:HA	2.20	0.42
1:C:384:GLY:N	1:C:386:GLU:OE2	2.49	0.42
1:E:17:LEU:HD12	1:E:17:LEU:C	2.44	0.42
1:G:482:ILE:CG1	1:G:541:MET:HE3	2.48	0.42
1:C:514:PHE:HB3	1:C:517:ARG:HG3	2.01	0.42
1:B:147:THR:HG21	1:B:170:TRP:HH2	1.84	0.41
1:G:123:LEU:HD23	1:G:127:SER:HB2	2.01	0.41
1:G:257:GLU:H	1:G:257:GLU:CD	2.28	0.41
1:H:257:GLU:CD	1:H:257:GLU:H	2.28	0.41
1:A:20:ARG:HD2	1:A:424:PHE:CE2	2.55	0.41
1:D:482:ILE:CG1	1:D:541:MET:HE3	2.49	0.41
1:C:184:ASN:OD1	1:C:184:ASN:C	2.63	0.41
1:C:211:VAL:N	1:C:212:PRO:CD	2.83	0.41
1:E:11:SER:OG	5:E:865:HOH:O	2.11	0.41
1:C:553:SER:O	1:H:156:VAL:CG2	2.69	0.41
1:F:277:PHE:N	1:F:278:PRO:CD	2.83	0.41
1:G:499:ASN:O	1:G:544:TYR:HA	2.20	0.41
1:D:123:LEU:HD21	1:D:129:ASN:ND2	2.36	0.41
1:G:288:LYS:HG3	1:G:337:ALA:HB1	2.02	0.41
1:A:550:ARG:NH2	1:F:153:ASP:O	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:147:THR:HG21	1:C:170:TRP:HH2	1.86	0.41
1:D:517:ARG:HB3	1:D:552:ASN:O	2.20	0.41
1:C:517:ARG:HB3	1:C:552:ASN:O	2.21	0.41
1:D:147:THR:HG21	1:D:170:TRP:HH2	1.85	0.41
1:G:201:LEU:HA	1:G:201:LEU:HD12	1.81	0.41
1:A:277:PHE:N	1:A:278:PRO:CD	2.83	0.41
1:D:184:ASN:OD1	1:D:184:ASN:C	2.64	0.41
1:A:288:LYS:CG	1:A:337:ALA:HB1	2.49	0.41
1:F:123:LEU:C	1:F:125:ASP:N	2.79	0.41
1:F:147:THR:HG21	1:F:170:TRP:HH2	1.86	0.41
1:F:207:ARG:NH2	1:F:209:ASP:OD2	2.54	0.41
1:F:288:LYS:HG3	1:F:337:ALA:HB1	2.02	0.41
1:G:125:ASP:OD1	1:G:127:SER:OG	2.39	0.41
1:G:514:PHE:HB3	1:G:517:ARG:HG3	2.02	0.41
1:A:123:LEU:HD23	1:A:127:SER:HB2	2.01	0.41
1:C:246:ARG:N	1:C:246:ARG:HD3	2.35	0.41
1:C:277:PHE:N	1:C:278:PRO:CD	2.84	0.41
1:E:211:VAL:N	1:E:212:PRO:CD	2.84	0.41
1:F:499:ASN:O	1:F:544:TYR:HA	2.20	0.41
1:G:52:TRP:CD1	1:G:52:TRP:C	2.99	0.41
1:G:211:VAL:N	1:G:212:PRO:CD	2.84	0.41
1:A:514:PHE:HB3	1:A:517:ARG:HG3	2.02	0.40
1:E:517:ARG:HB3	1:E:552:ASN:O	2.21	0.40
1:A:184:ASN:C	1:A:184:ASN:OD1	2.64	0.40
1:A:293:SER:HB3	5:A:820:HOH:O	2.22	0.40
5:B:709:HOH:O	1:E:553:SER:C	2.63	0.40
1:C:499:ASN:O	1:C:544:TYR:HA	2.21	0.40
1:F:146:ASP:OD1	1:F:146:ASP:N	2.53	0.40
1:C:89:ARG:NE	1:C:89:ARG:HA	2.37	0.40
1:D:495:LEU:C	1:D:496:ILE:HD13	2.46	0.40
1:D:514:PHE:HB3	1:D:517:ARG:HG3	2.02	0.40
1:F:514:PHE:HB3	1:F:517:ARG:HG3	2.03	0.40
1:B:384:GLY:N	1:B:386:GLU:OE2	2.52	0.40
1:D:446:ASP:HB3	1:D:449:SER:HB3	2.04	0.40
1:E:246:ARG:N	1:E:246:ARG:HD3	2.35	0.40
1:G:321:LEU:HD23	1:G:321:LEU:C	2.46	0.40
1:B:211:VAL:N	1:B:212:PRO:CD	2.84	0.40
1:B:514:PHE:HB3	1:B:517:ARG:HG3	2.03	0.40
1:G:184:ASN:OD1	1:G:184:ASN:C	2.64	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	546/571 (96%)	532 (97%)	14 (3%)	0	100	100
1	B	546/571 (96%)	532 (97%)	14 (3%)	0	100	100
1	C	546/571 (96%)	531 (97%)	15 (3%)	0	100	100
1	D	546/571 (96%)	532 (97%)	14 (3%)	0	100	100
1	E	546/571 (96%)	531 (97%)	15 (3%)	0	100	100
1	F	546/571 (96%)	530 (97%)	16 (3%)	0	100	100
1	G	546/571 (96%)	533 (98%)	13 (2%)	0	100	100
1	H	546/571 (96%)	533 (98%)	13 (2%)	0	100	100
All	All	4368/4568 (96%)	4254 (97%)	114 (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	465/484 (96%)	430 (92%)	35 (8%)	12	31
1	B	465/484 (96%)	440 (95%)	25 (5%)	20	45
1	C	465/484 (96%)	436 (94%)	29 (6%)	16	39
1	D	465/484 (96%)	444 (96%)	21 (4%)	24	52
1	E	465/484 (96%)	443 (95%)	22 (5%)	23	51
1	F	465/484 (96%)	438 (94%)	27 (6%)	18	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	465/484 (96%)	435 (94%)	30 (6%)	15	37
1	H	465/484 (96%)	433 (93%)	32 (7%)	14	34
All	All	3720/3872 (96%)	3499 (94%)	221 (6%)	18	42

All (221) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	17	LEU
1	A	21	THR
1	A	44	LYS
1	A	51	LEU
1	A	68	VAL
1	A	74	ILE
1	A	78	LEU
1	A	81	LEU
1	A	86	VAL
1	A	89	ARG
1	A	98	VAL
1	A	123	LEU
1	A	127	SER
1	A	133	ASP
1	A	147	THR
1	A	152	THR
1	A	161	LEU
1	A	186	LYS
1	A	191	LEU
1	A	216	GLU
1	A	246	ARG
1	A	257	GLU
1	A	258	GLU
1	A	282	ARG
1	A	288	LYS
1	A	304	LYS
1	A	346	ILE
1	A	348	VAL
1	A	381	LEU
1	A	386	GLU
1	A	393	LEU
1	A	452	LYS
1	A	517	ARG

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Mol	Chain	Res	Type
1	A	532	THR
1	B	11	SER
1	B	21	THR
1	B	44	LYS
1	B	51	LEU
1	B	86	VAL
1	B	122	THR
1	B	127	SER
1	B	142	LYS
1	B	147	THR
1	B	163	GLU
1	B	191	LEU
1	B	246	ARG
1	B	257	GLU
1	B	282	ARG
1	B	304	LYS
1	B	346	ILE
1	B	348	VAL
1	B	381	LEU
1	B	393	LEU
1	B	452	LYS
1	B	476	GLU
1	B	496	ILE
1	B	517	ARG
1	B	522	LEU
1	B	532	THR
1	C	11	SER
1	C	21	THR
1	C	44	LYS
1	C	51	LEU
1	C	86	VAL
1	C	89	ARG
1	C	98	VAL
1	C	123	LEU
1	C	125	ASP
1	C	127	SER
1	C	133	ASP
1	C	147	THR
1	C	164	GLN
1	C	186	LYS
1	C	191	LEU
1	C	216	GLU

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Mol	Chain	Res	Type
1	C	246	ARG
1	C	257	GLU
1	C	282	ARG
1	C	346	ILE
1	C	348	VAL
1	C	386	GLU
1	C	393	LEU
1	C	452	LYS
1	C	462	ARG
1	C	494	LEU
1	C	517	ARG
1	C	522	LEU
1	C	532	THR
1	D	11	SER
1	D	21	THR
1	D	44	LYS
1	D	127	SER
1	D	147	THR
1	D	191	LEU
1	D	216	GLU
1	D	246	ARG
1	D	257	GLU
1	D	285	MET
1	D	304	LYS
1	D	346	ILE
1	D	348	VAL
1	D	386	GLU
1	D	393	LEU
1	D	452	LYS
1	D	459	GLU
1	D	476	GLU
1	D	494	LEU
1	D	517	ARG
1	D	532	THR
1	E	11	SER
1	E	21	THR
1	E	44	LYS
1	E	123	LEU
1	E	133	ASP
1	E	147	THR
1	E	191	LEU
1	E	216	GLU

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Mol	Chain	Res	Type
1	E	246	ARG
1	E	249	LEU
1	E	257	GLU
1	E	282	ARG
1	E	304	LYS
1	E	346	ILE
1	E	348	VAL
1	E	386	GLU
1	E	393	LEU
1	E	452	LYS
1	E	494	LEU
1	E	517	ARG
1	E	522	LEU
1	E	532	THR
1	F	11	SER
1	F	21	THR
1	F	30	LYS
1	F	44	LYS
1	F	125	ASP
1	F	127	SER
1	F	146	ASP
1	F	147	THR
1	F	164	GLN
1	F	186	LYS
1	F	191	LEU
1	F	216	GLU
1	F	246	ARG
1	F	257	GLU
1	F	282	ARG
1	F	296	ARG
1	F	304	LYS
1	F	346	ILE
1	F	348	VAL
1	F	386	GLU
1	F	393	LEU
1	F	452	LYS
1	F	456	ARG
1	F	493	THR
1	F	496	ILE
1	F	517	ARG
1	F	532	THR
1	G	11	SER

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Mol	Chain	Res	Type
1	G	21	THR
1	G	30	LYS
1	G	44	LYS
1	G	118	ARG
1	G	123	LEU
1	G	127	SER
1	G	147	THR
1	G	161	LEU
1	G	163	GLU
1	G	164	GLN
1	G	201	LEU
1	G	216	GLU
1	G	246	ARG
1	G	249	LEU
1	G	257	GLU
1	G	282	ARG
1	G	304	LYS
1	G	346	ILE
1	G	348	VAL
1	G	381	LEU
1	G	386	GLU
1	G	393	LEU
1	G	411	THR
1	G	452	LYS
1	G	476	GLU
1	G	494	LEU
1	G	496	ILE
1	G	517	ARG
1	G	532	THR
1	H	11	SER
1	H	21	THR
1	H	44	LYS
1	H	127	SER
1	H	133	ASP
1	H	147	THR
1	H	161	LEU
1	H	191	LEU
1	H	207	ARG
1	H	216	GLU
1	H	225	LEU
1	H	228	THR
1	H	257	GLU

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Mol	Chain	Res	Type
1	H	268	GLU
1	H	282	ARG
1	H	298	ILE
1	H	324	GLU
1	H	328	ASP
1	H	346	ILE
1	H	348	VAL
1	H	386	GLU
1	H	393	LEU
1	H	452	LYS
1	H	459	GLU
1	H	471	ASP
1	H	476	GLU
1	H	488	GLN
1	H	492	GLU
1	H	494	LEU
1	H	517	ARG
1	H	522	LEU
1	H	532	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	HIS
1	E	45	ASN
1	E	192	HIS
1	E	488	GLN
1	F	369	ASN
1	G	45	ASN
1	G	192	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 16 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	$\# Z > 2$
4	TRS	D	602	-	7,7,7	0.36	0	9,9,9	0.66	0
4	TRS	F	602	-	7,7,7	0.48	0	9,9,9	0.94	0
4	TRS	A	602	-	7,7,7	0.58	0	9,9,9	0.84	0
4	TRS	E	602	-	7,7,7	0.56	0	9,9,9	0.79	0
4	TRS	B	602	-	7,7,7	0.53	0	9,9,9	1.33	3 (33%)
4	TRS	G	602	-	7,7,7	0.41	0	9,9,9	0.98	0
4	TRS	C	602	-	7,7,7	0.64	0	9,9,9	0.88	0
4	TRS	H	602	-	7,7,7	0.35	0	9,9,9	0.74	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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	TRS	D	602	-	-	8/9/9/9	-
4	TRS	F	602	-	-	7/9/9/9	-
4	TRS	A	602	-	-	7/9/9/9	-
4	TRS	E	602	-	-	7/9/9/9	-
4	TRS	B	602	-	-	9/9/9/9	-
4	TRS	G	602	-	-	1/9/9/9	-
4	TRS	C	602	-	-	6/9/9/9	-
4	TRS	H	602	-	-	6/9/9/9	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	602	TRS	C2-C-N	2.13	113.59	108.17
4	B	602	TRS	O1-C1-C	2.12	116.79	110.88
4	B	602	TRS	C3-C-C2	-2.03	105.25	110.66

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	602	TRS	C2-C-C1-O1
4	A	602	TRS	C3-C-C1-O1
4	B	602	TRS	C2-C-C1-O1
4	B	602	TRS	C3-C-C1-O1
4	B	602	TRS	C1-C-C2-O2
4	B	602	TRS	C3-C-C2-O2
4	C	602	TRS	C2-C-C1-O1
4	C	602	TRS	C3-C-C1-O1
4	C	602	TRS	N-C-C1-O1
4	D	602	TRS	C3-C-C1-O1
4	F	602	TRS	C3-C-C1-O1
4	F	602	TRS	C1-C-C3-O3
4	F	602	TRS	C2-C-C3-O3
4	F	602	TRS	N-C-C3-O3
4	H	602	TRS	N-C-C2-O2
4	H	602	TRS	N-C-C3-O3
4	A	602	TRS	C1-C-C3-O3
4	D	602	TRS	C2-C-C1-O1
4	F	602	TRS	C2-C-C1-O1
4	H	602	TRS	C3-C-C2-O2
4	H	602	TRS	C1-C-C3-O3
4	A	602	TRS	N-C-C1-O1
4	A	602	TRS	N-C-C3-O3
4	B	602	TRS	N-C-C2-O2
4	B	602	TRS	N-C-C3-O3
4	C	602	TRS	C2-C-C3-O3
4	C	602	TRS	N-C-C3-O3
4	D	602	TRS	N-C-C2-O2
4	E	602	TRS	N-C-C1-O1
4	E	602	TRS	C1-C-C2-O2
4	A	602	TRS	C2-C-C3-O3
4	B	602	TRS	C1-C-C3-O3
4	B	602	TRS	C2-C-C3-O3

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Mol	Chain	Res	Type	Atoms
4	C	602	TRS	C1-C-C3-O3
4	D	602	TRS	C1-C-C2-O2
4	E	602	TRS	C2-C-C1-O1
4	E	602	TRS	C3-C-C1-O1
4	E	602	TRS	C3-C-C2-O2
4	H	602	TRS	C1-C-C2-O2
4	B	602	TRS	N-C-C1-O1
4	D	602	TRS	N-C-C1-O1
4	D	602	TRS	C3-C-C2-O2
4	E	602	TRS	N-C-C2-O2
4	F	602	TRS	N-C-C1-O1
4	A	602	TRS	C1-C-C2-O2
4	D	602	TRS	C2-C-C3-O3
4	E	602	TRS	C2-C-C3-O3
4	F	602	TRS	C1-C-C2-O2
4	H	602	TRS	C2-C-C3-O3
4	D	602	TRS	C1-C-C3-O3
4	G	602	TRS	N-C-C3-O3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	602	TRS	1	0
4	E	602	TRS	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	548/571 (95%)	-0.32	2 (0%) 88 87	7, 28, 47, 80	3 (0%)
1	B	548/571 (95%)	-0.18	9 (1%) 70 68	8, 30, 54, 85	3 (0%)
1	C	548/571 (95%)	0.24	11 (2%) 65 62	9, 40, 64, 85	3 (0%)
1	D	548/571 (95%)	0.19	15 (2%) 56 53	11, 41, 71, 99	3 (0%)
1	E	548/571 (95%)	-0.39	3 (0%) 87 86	7, 25, 42, 84	3 (0%)
1	F	548/571 (95%)	0.17	13 (2%) 59 56	9, 38, 66, 110	3 (0%)
1	G	548/571 (95%)	-0.09	5 (0%) 81 80	11, 34, 57, 115	3 (0%)
1	H	548/571 (95%)	0.66	51 (9%) 14 12	14, 50, 84, 151	3 (0%)
All	All	4384/4568 (95%)	0.04	109 (2%) 58 55	7, 35, 68, 151	24 (0%)

All (109) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	128	PRO	4.8
1	B	353	ARG	4.5
1	H	124	PRO	4.4
1	F	6	PRO	4.2
1	A	353	ARG	3.9
1	G	353	ARG	3.9
1	H	353	ARG	3.9
1	H	175	ALA	3.9
1	H	136	VAL	3.9
1	D	353	ARG	3.9
1	C	6	PRO	3.8
1	F	353	ARG	3.8
1	H	119	ARG	3.8
1	H	456	ARG	3.8
1	H	168	TYR	3.7
1	H	121	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	H	169	TYR	3.6
1	H	235	PHE	3.6
1	H	161	LEU	3.5
1	D	552	ASN	3.5
1	D	456	ARG	3.4
1	G	124	PRO	3.4
1	A	6	PRO	3.4
1	B	304	LYS	3.3
1	C	353	ARG	3.3
1	H	127	SER	3.3
1	H	6	PRO	3.2
1	H	129	ASN	3.2
1	D	123	LEU	3.2
1	G	123	LEU	3.2
1	B	124	PRO	3.2
1	E	353	ARG	3.1
1	F	124	PRO	3.1
1	G	6	PRO	3.1
1	F	191	LEU	3.1
1	H	133	ASP	3.1
1	F	456	ARG	3.1
1	D	553	SER	3.0
1	E	124	PRO	3.0
1	D	265	THR	2.9
1	H	184	ASN	2.8
1	F	343	ARG	2.8
1	H	137	TRP	2.8
1	H	122	THR	2.8
1	C	123	LEU	2.8
1	H	126	GLY	2.8
1	C	124	PRO	2.8
1	H	553	SER	2.8
1	H	232	LEU	2.7
1	B	122	THR	2.6
1	H	123	LEU	2.6
1	H	120	GLY	2.6
1	H	118	ARG	2.6
1	D	121	PRO	2.6
1	B	553	SER	2.6
1	H	130	GLU	2.6
1	D	120	GLY	2.6
1	H	267	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	128	PRO	2.5
1	F	553	SER	2.5
1	C	119	ARG	2.5
1	H	225	LEU	2.5
1	H	343	ARG	2.5
1	F	128	PRO	2.5
1	H	116	ALA	2.5
1	H	144	TYR	2.4
1	H	180	LEU	2.4
1	H	215	ILE	2.4
1	H	170	TRP	2.4
1	D	147	THR	2.4
1	H	104	THR	2.4
1	C	492	GLU	2.4
1	C	120	GLY	2.4
1	H	164	GLN	2.4
1	H	176	SER	2.4
1	H	304	LYS	2.4
1	C	128	PRO	2.4
1	H	115	GLN	2.4
1	D	122	THR	2.3
1	B	552	ASN	2.3
1	B	126	GLY	2.3
1	B	456	ARG	2.3
1	F	235	PHE	2.3
1	D	127	SER	2.3
1	B	6	PRO	2.3
1	H	191	LEU	2.3
1	E	6	PRO	2.3
1	H	125	ASP	2.3
1	C	456	ARG	2.3
1	H	109	SER	2.3
1	F	122	THR	2.2
1	H	163	GLU	2.2
1	D	6	PRO	2.2
1	H	256	PRO	2.2
1	H	305	ILE	2.2
1	C	125	ASP	2.2
1	C	122	THR	2.2
1	H	134	TYR	2.2
1	H	165	ALA	2.1
1	H	213	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	136	VAL	2.1
1	G	128	PRO	2.1
1	H	269	PRO	2.1
1	F	231	ILE	2.1
1	H	228	THR	2.1
1	F	109	SER	2.0
1	H	167	LYS	2.0
1	F	220	THR	2.0
1	D	168	TYR	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
4	TRS	H	602	8/8	0.83	0.12	48,52,52,54	0
4	TRS	A	602	8/8	0.91	0.11	29,31,34,41	0
4	TRS	B	602	8/8	0.92	0.08	24,26,27,27	0
2	CA	F	600	1/1	0.92	0.06	39,39,39,39	0
4	TRS	F	602	8/8	0.93	0.07	29,32,33,33	0
4	TRS	G	602	8/8	0.93	0.08	28,31,32,33	0
4	TRS	E	602	8/8	0.93	0.08	21,21,22,22	0
4	TRS	C	602	8/8	0.95	0.07	27,29,30,32	0
4	TRS	D	602	8/8	0.95	0.07	34,34,35,35	0
2	CA	H	600	1/1	0.95	0.05	67,67,67,67	0
3	MG	B	601	1/1	0.96	0.07	8,8,8,8	0
2	CA	B	600	1/1	0.96	0.05	36,36,36,36	0
2	CA	D	600	1/1	0.96	0.04	44,44,44,44	0
3	MG	A	601	1/1	0.97	0.10	14,14,14,14	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	A	600	1/1	0.97	0.05	32,32,32,32	0
3	MG	F	601	1/1	0.98	0.04	20,20,20,20	0
3	MG	H	601	1/1	0.98	0.04	28,28,28,28	0
3	MG	G	601	1/1	0.99	0.06	16,16,16,16	0
2	CA	C	600	1/1	0.99	0.02	38,38,38,38	0
2	CA	G	600	1/1	0.99	0.04	30,30,30,30	0
3	MG	E	601	1/1	0.99	0.06	7,7,7,7	0
2	CA	E	600	1/1	0.99	0.02	25,25,25,25	0
3	MG	C	601	1/1	1.00	0.02	12,12,12,12	0
3	MG	D	601	1/1	1.00	0.04	10,10,10,10	0

6.5 Other polymers [i](#)

There are no such residues in this entry.