



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:27 AM UTC

PDB ID : 6DHJ / pdb_00006dhj
Title : Crystal structures of cyanuric acid hydrolase from Moorella thermoacetica
Authors : Shi, K.; Aihara, H.
Deposited on : 2018-05-20
Resolution : 3.20 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

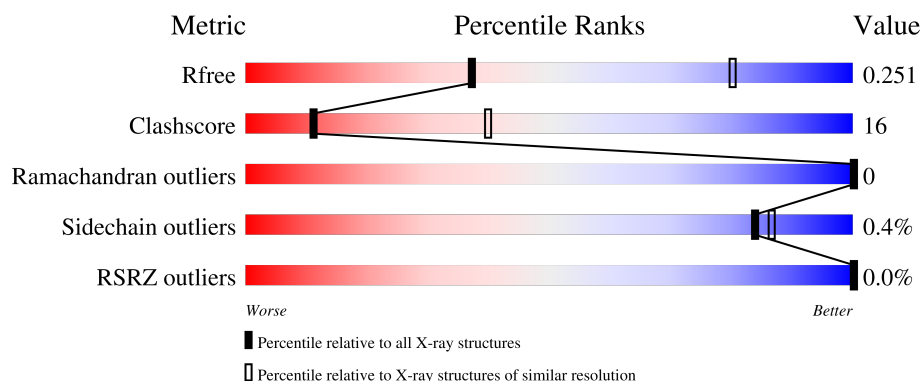
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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

Mol	Chain	Length	Quality of chain
1	A	370	 74% 25% ..
1	B	370	 75% 23% .
1	C	370	 72% 25% .
1	D	370	 75% 23% ..
1	E	370	 72% 26% .

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Mol	Chain	Length	Quality of chain
1	F	370	 71%26%•
1	G	370	 68%29%•
1	H	370	 77%21%•

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
2	CIT	H	401	-	X	-	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 21566 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 Cyanuric acid amidohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	365	Total	C	N	O	S	0	0	0
			2707	1677	484	533	13			
1	B	363	Total	C	N	O	S	0	0	0
			2688	1666	480	529	13			
1	C	362	Total	C	N	O	S	0	0	0
			2680	1660	479	528	13			
1	D	364	Total	C	N	O	S	0	0	0
			2697	1671	481	532	13			
1	E	364	Total	C	N	O	S	0	0	0
			2697	1671	481	532	13			
1	F	357	Total	C	N	O	S	0	0	0
			2647	1640	473	522	12			
1	G	359	Total	C	N	O	S	0	0	0
			2658	1645	476	525	12			
1	H	363	Total	C	N	O	S	0	0	0
			2688	1666	480	529	13			

There are 224 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	46	SER	CYS	engineered mutation	UNP Q2RGM7
A	53	CYS	GLY	engineered mutation	UNP Q2RGM7
A	103	ALA	GLN	engineered mutation	UNP Q2RGM7
A	104	ALA	GLU	engineered mutation	UNP Q2RGM7
A	107	ALA	LYS	engineered mutation	UNP Q2RGM7
A	162	ALA	CYS	engineered mutation	UNP Q2RGM7
A	218	VAL	CYS	engineered mutation	UNP Q2RGM7
A	235	CYS	GLU	engineered mutation	UNP Q2RGM7
A	279	ILE	LEU	engineered mutation	UNP Q2RGM7
A	280	ARG	LYS	engineered mutation	UNP Q2RGM7
A	281	SER	PHE	engineered mutation	UNP Q2RGM7
A	?	-	CYS	deletion	UNP Q2RGM7
A	?	-	CYS	deletion	UNP Q2RGM7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	PRO	deletion	UNP Q2RGM7
A	?	-	PRO	deletion	UNP Q2RGM7
A	?	-	ALA	deletion	UNP Q2RGM7
A	283	ASP	GLU	engineered mutation	UNP Q2RGM7
A	290	MET	LEU	engineered mutation	UNP Q2RGM7
A	291	ASP	ALA	engineered mutation	UNP Q2RGM7
A	292	ARG	LYS	engineered mutation	UNP Q2RGM7
A	368	LEU	-	expression tag	UNP Q2RGM7
A	369	GLU	-	expression tag	UNP Q2RGM7
A	370	HIS	-	expression tag	UNP Q2RGM7
A	371	HIS	-	expression tag	UNP Q2RGM7
A	372	HIS	-	expression tag	UNP Q2RGM7
A	373	HIS	-	expression tag	UNP Q2RGM7
A	374	HIS	-	expression tag	UNP Q2RGM7
A	375	HIS	-	expression tag	UNP Q2RGM7
B	46	SER	CYS	engineered mutation	UNP Q2RGM7
B	53	CYS	GLY	engineered mutation	UNP Q2RGM7
B	103	ALA	GLN	engineered mutation	UNP Q2RGM7
B	104	ALA	GLU	engineered mutation	UNP Q2RGM7
B	107	ALA	LYS	engineered mutation	UNP Q2RGM7
B	162	ALA	CYS	engineered mutation	UNP Q2RGM7
B	218	VAL	CYS	engineered mutation	UNP Q2RGM7
B	235	CYS	GLU	engineered mutation	UNP Q2RGM7
B	279	ILE	LEU	engineered mutation	UNP Q2RGM7
B	280	ARG	LYS	engineered mutation	UNP Q2RGM7
B	281	SER	PHE	engineered mutation	UNP Q2RGM7
B	?	-	CYS	deletion	UNP Q2RGM7
B	?	-	CYS	deletion	UNP Q2RGM7
B	?	-	PRO	deletion	UNP Q2RGM7
B	?	-	PRO	deletion	UNP Q2RGM7
B	?	-	ALA	deletion	UNP Q2RGM7
B	283	ASP	GLU	engineered mutation	UNP Q2RGM7
B	290	MET	LEU	engineered mutation	UNP Q2RGM7
B	291	ASP	ALA	engineered mutation	UNP Q2RGM7
B	292	ARG	LYS	engineered mutation	UNP Q2RGM7
B	368	LEU	-	expression tag	UNP Q2RGM7
B	369	GLU	-	expression tag	UNP Q2RGM7
B	370	HIS	-	expression tag	UNP Q2RGM7
B	371	HIS	-	expression tag	UNP Q2RGM7
B	372	HIS	-	expression tag	UNP Q2RGM7
B	373	HIS	-	expression tag	UNP Q2RGM7
B	374	HIS	-	expression tag	UNP Q2RGM7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	375	HIS	-	expression tag	UNP Q2RGM7
C	46	SER	CYS	engineered mutation	UNP Q2RGM7
C	53	CYS	GLY	engineered mutation	UNP Q2RGM7
C	103	ALA	GLN	engineered mutation	UNP Q2RGM7
C	104	ALA	GLU	engineered mutation	UNP Q2RGM7
C	107	ALA	LYS	engineered mutation	UNP Q2RGM7
C	162	ALA	CYS	engineered mutation	UNP Q2RGM7
C	218	VAL	CYS	engineered mutation	UNP Q2RGM7
C	235	CYS	GLU	engineered mutation	UNP Q2RGM7
C	279	ILE	LEU	engineered mutation	UNP Q2RGM7
C	280	ARG	LYS	engineered mutation	UNP Q2RGM7
C	281	SER	PHE	engineered mutation	UNP Q2RGM7
C	?	-	CYS	deletion	UNP Q2RGM7
C	?	-	CYS	deletion	UNP Q2RGM7
C	?	-	PRO	deletion	UNP Q2RGM7
C	?	-	PRO	deletion	UNP Q2RGM7
C	?	-	ALA	deletion	UNP Q2RGM7
C	283	ASP	GLU	engineered mutation	UNP Q2RGM7
C	290	MET	LEU	engineered mutation	UNP Q2RGM7
C	291	ASP	ALA	engineered mutation	UNP Q2RGM7
C	292	ARG	LYS	engineered mutation	UNP Q2RGM7
C	368	LEU	-	expression tag	UNP Q2RGM7
C	369	GLU	-	expression tag	UNP Q2RGM7
C	370	HIS	-	expression tag	UNP Q2RGM7
C	371	HIS	-	expression tag	UNP Q2RGM7
C	372	HIS	-	expression tag	UNP Q2RGM7
C	373	HIS	-	expression tag	UNP Q2RGM7
C	374	HIS	-	expression tag	UNP Q2RGM7
C	375	HIS	-	expression tag	UNP Q2RGM7
D	46	SER	CYS	engineered mutation	UNP Q2RGM7
D	53	CYS	GLY	engineered mutation	UNP Q2RGM7
D	103	ALA	GLN	engineered mutation	UNP Q2RGM7
D	104	ALA	GLU	engineered mutation	UNP Q2RGM7
D	107	ALA	LYS	engineered mutation	UNP Q2RGM7
D	162	ALA	CYS	engineered mutation	UNP Q2RGM7
D	218	VAL	CYS	engineered mutation	UNP Q2RGM7
D	235	CYS	GLU	engineered mutation	UNP Q2RGM7
D	279	ILE	LEU	engineered mutation	UNP Q2RGM7
D	280	ARG	LYS	engineered mutation	UNP Q2RGM7
D	281	SER	PHE	engineered mutation	UNP Q2RGM7
D	?	-	CYS	deletion	UNP Q2RGM7
D	?	-	CYS	deletion	UNP Q2RGM7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	PRO	deletion	UNP Q2RGM7
D	?	-	PRO	deletion	UNP Q2RGM7
D	?	-	ALA	deletion	UNP Q2RGM7
D	283	ASP	GLU	engineered mutation	UNP Q2RGM7
D	290	MET	LEU	engineered mutation	UNP Q2RGM7
D	291	ASP	ALA	engineered mutation	UNP Q2RGM7
D	292	ARG	LYS	engineered mutation	UNP Q2RGM7
D	368	LEU	-	expression tag	UNP Q2RGM7
D	369	GLU	-	expression tag	UNP Q2RGM7
D	370	HIS	-	expression tag	UNP Q2RGM7
D	371	HIS	-	expression tag	UNP Q2RGM7
D	372	HIS	-	expression tag	UNP Q2RGM7
D	373	HIS	-	expression tag	UNP Q2RGM7
D	374	HIS	-	expression tag	UNP Q2RGM7
D	375	HIS	-	expression tag	UNP Q2RGM7
E	46	SER	CYS	engineered mutation	UNP Q2RGM7
E	53	CYS	GLY	engineered mutation	UNP Q2RGM7
E	103	ALA	GLN	engineered mutation	UNP Q2RGM7
E	104	ALA	GLU	engineered mutation	UNP Q2RGM7
E	107	ALA	LYS	engineered mutation	UNP Q2RGM7
E	162	ALA	CYS	engineered mutation	UNP Q2RGM7
E	218	VAL	CYS	engineered mutation	UNP Q2RGM7
E	235	CYS	GLU	engineered mutation	UNP Q2RGM7
E	279	ILE	LEU	engineered mutation	UNP Q2RGM7
E	280	ARG	LYS	engineered mutation	UNP Q2RGM7
E	281	SER	PHE	engineered mutation	UNP Q2RGM7
E	?	-	CYS	deletion	UNP Q2RGM7
E	?	-	CYS	deletion	UNP Q2RGM7
E	?	-	PRO	deletion	UNP Q2RGM7
E	?	-	PRO	deletion	UNP Q2RGM7
E	?	-	ALA	deletion	UNP Q2RGM7
E	283	ASP	GLU	engineered mutation	UNP Q2RGM7
E	290	MET	LEU	engineered mutation	UNP Q2RGM7
E	291	ASP	ALA	engineered mutation	UNP Q2RGM7
E	292	ARG	LYS	engineered mutation	UNP Q2RGM7
E	368	LEU	-	expression tag	UNP Q2RGM7
E	369	GLU	-	expression tag	UNP Q2RGM7
E	370	HIS	-	expression tag	UNP Q2RGM7
E	371	HIS	-	expression tag	UNP Q2RGM7
E	372	HIS	-	expression tag	UNP Q2RGM7
E	373	HIS	-	expression tag	UNP Q2RGM7
E	374	HIS	-	expression tag	UNP Q2RGM7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	375	HIS	-	expression tag	UNP Q2RGM7
F	46	SER	CYS	engineered mutation	UNP Q2RGM7
F	53	CYS	GLY	engineered mutation	UNP Q2RGM7
F	103	ALA	GLN	engineered mutation	UNP Q2RGM7
F	104	ALA	GLU	engineered mutation	UNP Q2RGM7
F	107	ALA	LYS	engineered mutation	UNP Q2RGM7
F	162	ALA	CYS	engineered mutation	UNP Q2RGM7
F	218	VAL	CYS	engineered mutation	UNP Q2RGM7
F	235	CYS	GLU	engineered mutation	UNP Q2RGM7
F	279	ILE	LEU	engineered mutation	UNP Q2RGM7
F	280	ARG	LYS	engineered mutation	UNP Q2RGM7
F	281	SER	PHE	engineered mutation	UNP Q2RGM7
F	?	-	CYS	deletion	UNP Q2RGM7
F	?	-	CYS	deletion	UNP Q2RGM7
F	?	-	PRO	deletion	UNP Q2RGM7
F	?	-	PRO	deletion	UNP Q2RGM7
F	?	-	ALA	deletion	UNP Q2RGM7
F	283	ASP	GLU	engineered mutation	UNP Q2RGM7
F	290	MET	LEU	engineered mutation	UNP Q2RGM7
F	291	ASP	ALA	engineered mutation	UNP Q2RGM7
F	292	ARG	LYS	engineered mutation	UNP Q2RGM7
F	368	LEU	-	expression tag	UNP Q2RGM7
F	369	GLU	-	expression tag	UNP Q2RGM7
F	370	HIS	-	expression tag	UNP Q2RGM7
F	371	HIS	-	expression tag	UNP Q2RGM7
F	372	HIS	-	expression tag	UNP Q2RGM7
F	373	HIS	-	expression tag	UNP Q2RGM7
F	374	HIS	-	expression tag	UNP Q2RGM7
F	375	HIS	-	expression tag	UNP Q2RGM7
G	46	SER	CYS	engineered mutation	UNP Q2RGM7
G	53	CYS	GLY	engineered mutation	UNP Q2RGM7
G	103	ALA	GLN	engineered mutation	UNP Q2RGM7
G	104	ALA	GLU	engineered mutation	UNP Q2RGM7
G	107	ALA	LYS	engineered mutation	UNP Q2RGM7
G	162	ALA	CYS	engineered mutation	UNP Q2RGM7
G	218	VAL	CYS	engineered mutation	UNP Q2RGM7
G	235	CYS	GLU	engineered mutation	UNP Q2RGM7
G	279	ILE	LEU	engineered mutation	UNP Q2RGM7
G	280	ARG	LYS	engineered mutation	UNP Q2RGM7
G	281	SER	PHE	engineered mutation	UNP Q2RGM7
G	?	-	CYS	deletion	UNP Q2RGM7
G	?	-	CYS	deletion	UNP Q2RGM7

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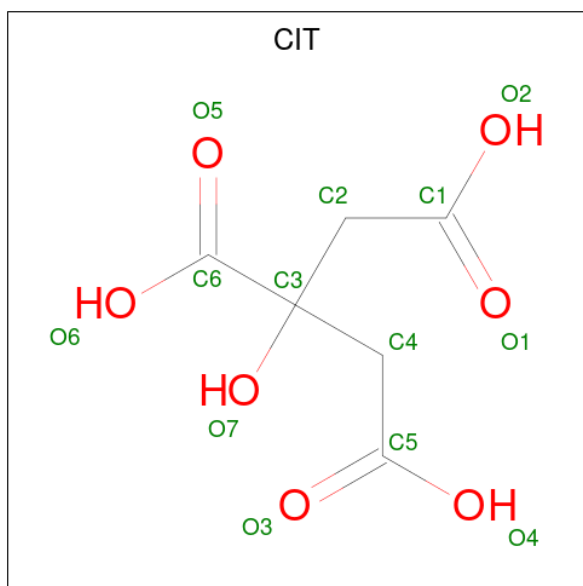
Chain	Residue	Modelled	Actual	Comment	Reference
G	?	-	PRO	deletion	UNP Q2RGM7
G	?	-	PRO	deletion	UNP Q2RGM7
G	?	-	ALA	deletion	UNP Q2RGM7
G	283	ASP	GLU	engineered mutation	UNP Q2RGM7
G	290	MET	LEU	engineered mutation	UNP Q2RGM7
G	291	ASP	ALA	engineered mutation	UNP Q2RGM7
G	292	ARG	LYS	engineered mutation	UNP Q2RGM7
G	368	LEU	-	expression tag	UNP Q2RGM7
G	369	GLU	-	expression tag	UNP Q2RGM7
G	370	HIS	-	expression tag	UNP Q2RGM7
G	371	HIS	-	expression tag	UNP Q2RGM7
G	372	HIS	-	expression tag	UNP Q2RGM7
G	373	HIS	-	expression tag	UNP Q2RGM7
G	374	HIS	-	expression tag	UNP Q2RGM7
G	375	HIS	-	expression tag	UNP Q2RGM7
H	46	SER	CYS	engineered mutation	UNP Q2RGM7
H	53	CYS	GLY	engineered mutation	UNP Q2RGM7
H	103	ALA	GLN	engineered mutation	UNP Q2RGM7
H	104	ALA	GLU	engineered mutation	UNP Q2RGM7
H	107	ALA	LYS	engineered mutation	UNP Q2RGM7
H	162	ALA	CYS	engineered mutation	UNP Q2RGM7
H	218	VAL	CYS	engineered mutation	UNP Q2RGM7
H	235	CYS	GLU	engineered mutation	UNP Q2RGM7
H	279	ILE	LEU	engineered mutation	UNP Q2RGM7
H	280	ARG	LYS	engineered mutation	UNP Q2RGM7
H	281	SER	PHE	engineered mutation	UNP Q2RGM7
H	?	-	CYS	deletion	UNP Q2RGM7
H	?	-	CYS	deletion	UNP Q2RGM7
H	?	-	PRO	deletion	UNP Q2RGM7
H	?	-	PRO	deletion	UNP Q2RGM7
H	?	-	ALA	deletion	UNP Q2RGM7
H	283	ASP	GLU	engineered mutation	UNP Q2RGM7
H	290	MET	LEU	engineered mutation	UNP Q2RGM7
H	291	ASP	ALA	engineered mutation	UNP Q2RGM7
H	292	ARG	LYS	engineered mutation	UNP Q2RGM7
H	368	LEU	-	expression tag	UNP Q2RGM7
H	369	GLU	-	expression tag	UNP Q2RGM7
H	370	HIS	-	expression tag	UNP Q2RGM7
H	371	HIS	-	expression tag	UNP Q2RGM7
H	372	HIS	-	expression tag	UNP Q2RGM7
H	373	HIS	-	expression tag	UNP Q2RGM7
H	374	HIS	-	expression tag	UNP Q2RGM7

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Chain	Residue	Modelled	Actual	Comment	Reference
H	375	HIS	-	expression tag	UNP Q2RGM7

- Molecule 2 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	C	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		
2	E	1	Total	C	O	0	0
			13	6	7		
2	F	1	Total	C	O	0	0
			13	6	7		
2	G	1	Total	C	O	0	0
			13	6	7		
2	H	1	Total	C	O	0	0
			13	6	7		

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.

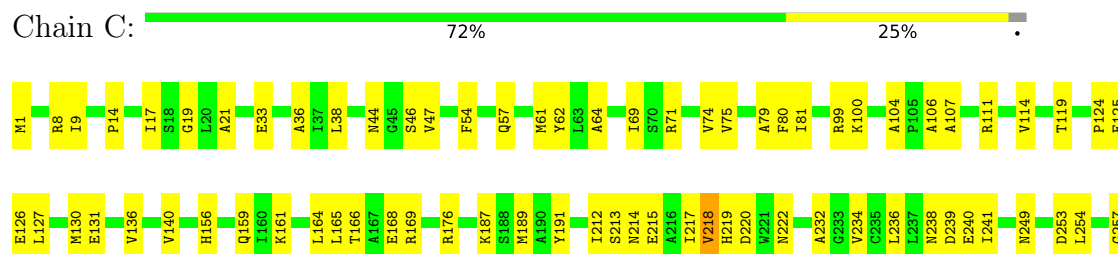
• Molecule 1: Cyanuric acid amidohydrolase



• Molecule 1: Cyanuric acid amidohydrolase



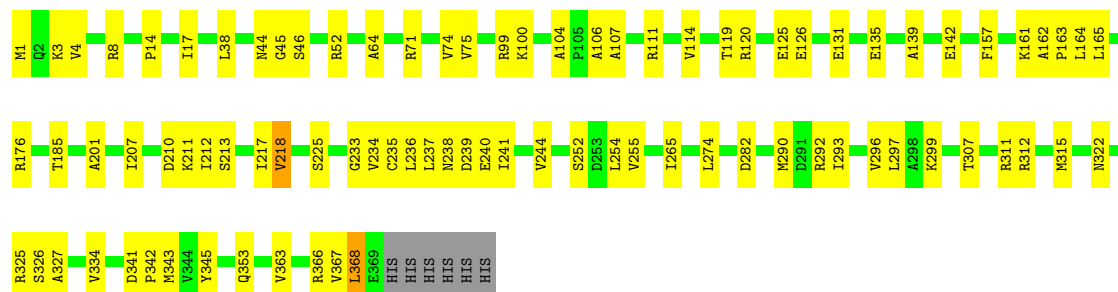
• Molecule 1: Cyanuric acid amidohydrolase





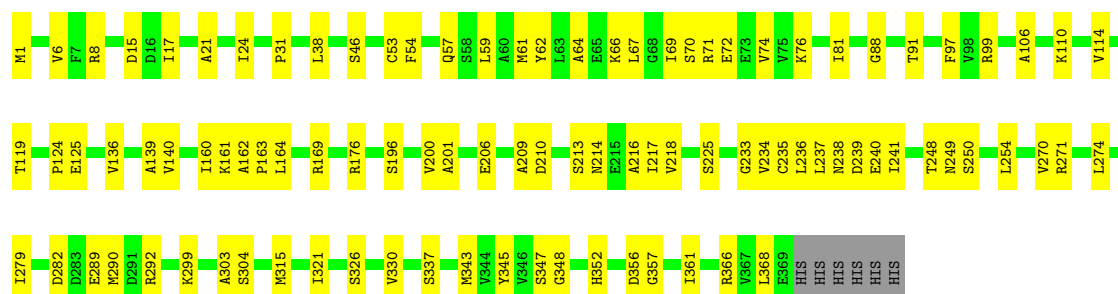
• Molecule 1: Cyanuric acid amidohydrolase

Chain D: 75% 23% ..



• Molecule 1: Cyanuric acid amidohydrolase

Chain E: 72% 26% .



• Molecule 1: Cyanuric acid amidohydrolase

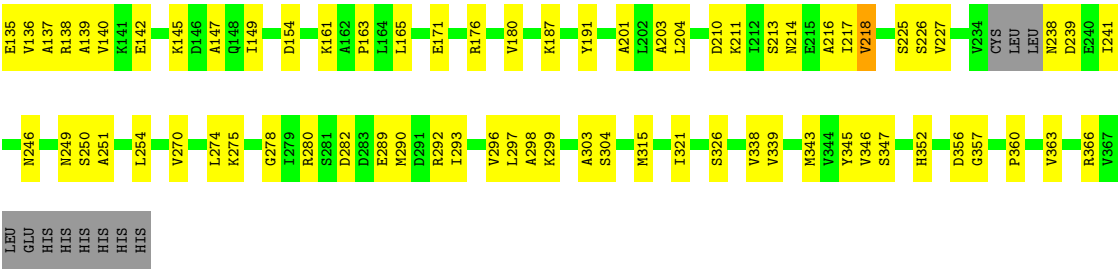
Chain F: 71% 26% .



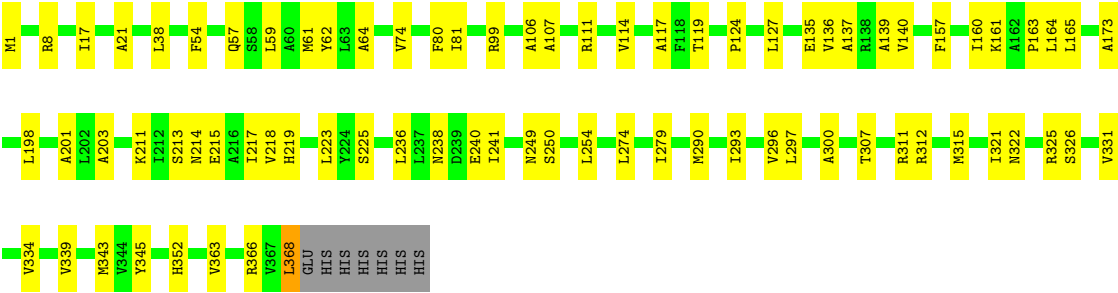
• Molecule 1: Cyanuric acid amidohydrolase

Chain G: 68% 29% .





● Molecule 1: Cyanuric acid amidohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.91Å 164.56Å 202.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	127.80 – 3.20 127.80 – 3.20	Depositor EDS
% Data completeness (in resolution range)	95.6 (127.80-3.20) 95.7 (127.80-3.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 3.19Å)	Xtriage
Refinement program	PHENIX (1.15.2_3472: ???)	Depositor
R, R_{free}	0.208 , 0.250 0.213 , 0.251	Depositor DCC
R_{free} test set	2807 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	118.7	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 107.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21566	wwPDB-VP
Average B, all atoms (Å ²)	140.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CIT

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	0.25	0/2740	0.57	0/3710
1	B	0.28	0/2720	0.61	1/3683 (0.0%)
1	C	0.25	0/2712	0.59	0/3672
1	D	0.29	0/2729	0.65	0/3695
1	E	0.24	0/2729	0.57	1/3695 (0.0%)
1	F	0.21	0/2678	0.52	0/3624
1	G	0.20	0/2689	0.51	0/3639
1	H	0.28	0/2720	0.62	1/3683 (0.0%)
All	All	0.25	0/21717	0.58	3/29401 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	240	GLU	N-CA-C	7.70	121.99	109.59
1	B	240	GLU	N-CA-C	5.70	118.77	109.59
1	E	209	ALA	N-CA-C	5.54	117.31	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2707	0	2733	95	0
1	B	2688	0	2718	94	1
1	C	2680	0	2707	90	0
1	D	2697	0	2722	87	0
1	E	2697	0	2725	95	1
1	F	2647	0	2675	97	0
1	G	2658	0	2681	95	0
1	H	2688	0	2712	89	0
2	A	13	0	5	2	0
2	B	13	0	5	3	0
2	C	13	0	5	0	0
2	D	13	0	5	2	0
2	E	13	0	5	3	0
2	F	13	0	5	2	0
2	G	13	0	5	2	0
2	H	13	0	5	0	0
All	All	21566	0	21713	704	1

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 16.

All (704) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:LYS:HG2	1:C:238:ASN:HB2	1.29	1.08
1:H:160:ILE:HG12	1:H:241:ILE:HD13	1.37	1.06
1:H:80:PHE:O	1:H:161:LYS:NZ	1.88	1.05
1:E:161:LYS:HE3	1:E:233:GLY:HA2	1.38	1.05
1:G:315:MET:HE2	1:G:326:SER:HB3	1.44	0.97
1:A:57:GLN:HE21	1:E:57:GLN:HE21	1.10	0.97
1:A:111:ARG:HH22	1:A:370:HIS:HB2	1.30	0.96
1:F:119:THR:HB	1:F:135:GLU:HG3	1.47	0.96
1:C:161:LYS:CG	1:C:238:ASN:HB2	1.95	0.95
1:F:107:ALA:HB1	1:F:108:PRO:HD2	1.51	0.92
1:G:217:ILE:O	1:G:352:HIS:NE2	2.02	0.92
1:H:1:MET:HE2	1:H:106:ALA:HA	1.53	0.90
1:E:110:LYS:HE3	1:E:248:THR:HG22	1.55	0.89
1:H:274:LEU:HD13	1:H:290:MET:HE1	1.52	0.89
1:G:201:ALA:HB2	1:G:225:SER:HB2	1.55	0.87
1:G:296:VAL:HG22	1:G:363:VAL:HG22	1.57	0.86
1:A:93:HIS:HB3	1:A:343:MET:HE3	1.57	0.85
1:E:1:MET:HE2	1:E:106:ALA:HA	1.58	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:315:MET:HE2	1:H:326:SER:HB3	1.58	0.84
1:D:239:ASP:O	1:D:240:GLU:HG3	1.76	0.84
1:B:161:LYS:HG3	1:B:238:ASN:HB3	1.61	0.83
1:A:111:ARG:NH2	1:A:370:HIS:HB2	1.93	0.82
1:F:164:LEU:O	1:F:169:ARG:NH1	2.12	0.82
1:H:163:PRO:CB	1:H:238:ASN:OD1	2.24	0.82
1:F:210:ASP:OD1	1:F:211:LYS:HE3	1.79	0.82
1:H:1:MET:CE	1:H:106:ALA:HA	2.10	0.82
1:C:274:LEU:HD13	1:C:290:MET:HE1	1.62	0.81
1:F:315:MET:HE2	1:F:326:SER:HB3	1.62	0.81
1:C:1:MET:CE	1:C:106:ALA:HB2	2.09	0.81
1:E:162:ALA:O	1:E:233:GLY:N	2.12	0.81
1:G:119:THR:HB	1:G:135:GLU:HG3	1.62	0.81
1:F:139:ALA:HB3	1:F:241:ILE:HD12	1.63	0.80
1:G:210:ASP:OD1	1:G:211:LYS:HE3	1.81	0.80
1:A:210:ASP:OD1	1:A:211:LYS:HE3	1.80	0.79
1:G:163:PRO:HD3	1:G:239:ASP:OD1	1.82	0.79
1:F:8:ARG:HG2	1:F:343:MET:HE3	1.65	0.79
1:C:75:VAL:HG11	1:C:236:LEU:HD12	1.64	0.78
1:G:211:LYS:HA	1:G:211:LYS:HE2	1.63	0.78
1:D:210:ASP:OD1	1:D:211:LYS:HE3	1.83	0.78
1:B:129:ARG:NH2	1:B:180:VAL:HA	1.98	0.78
1:E:119:THR:HG22	1:E:139:ALA:HB2	1.66	0.78
1:A:211:LYS:HE2	1:A:211:LYS:HA	1.65	0.77
1:C:161:LYS:HG2	1:C:238:ASN:CB	2.12	0.77
1:F:1:MET:CE	1:F:106:ALA:HB2	2.14	0.77
1:E:1:MET:CE	1:E:106:ALA:HA	2.14	0.77
1:G:99:ARG:HD3	1:G:114:VAL:HG13	1.66	0.77
1:D:233:GLY:H	1:D:238:ASN:HD22	1.31	0.76
1:B:300:ALA:HB1	1:B:331:VAL:HG21	1.67	0.76
1:C:161:LYS:CD	1:C:238:ASN:HB2	2.15	0.76
1:E:161:LYS:O	1:E:239:ASP:HA	1.86	0.75
1:A:93:HIS:CB	1:A:343:MET:HE3	2.16	0.75
1:B:1:MET:HE2	1:B:106:ALA:HA	1.68	0.75
1:D:162:ALA:O	1:D:238:ASN:ND2	2.20	0.74
1:G:99:ARG:HD3	1:G:114:VAL:CG1	2.17	0.74
1:A:368:LEU:HD22	1:A:369:GLU:H	1.50	0.74
1:F:119:THR:HB	1:F:135:GLU:CG	2.17	0.74
1:D:211:LYS:HA	1:D:211:LYS:HE2	1.69	0.74
1:G:315:MET:HE2	1:G:326:SER:CB	2.18	0.74
1:D:322:ASN:HB2	1:D:325:ARG:HG3	1.68	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:LEU:HD22	1:A:369:GLU:N	2.03	0.73
1:D:163:PRO:HA	1:D:238:ASN:HD21	1.54	0.73
1:F:139:ALA:CB	1:F:241:ILE:HD12	2.19	0.72
1:E:201:ALA:HB2	1:E:225:SER:HB2	1.71	0.72
1:E:315:MET:HE2	1:E:326:SER:HB3	1.72	0.72
1:B:61:MET:HE3	1:C:71:ARG:CZ	2.20	0.72
1:F:1:MET:HE1	1:F:106:ALA:HB2	1.71	0.72
1:H:99:ARG:HD3	1:H:114:VAL:HG13	1.71	0.72
1:E:213:SER:O	1:E:217:ILE:HG13	1.90	0.71
1:B:214:ASN:HA	1:B:217:ILE:HD12	1.72	0.71
1:E:110:LYS:HD2	1:E:110:LYS:N	2.06	0.71
1:H:214:ASN:HA	1:H:217:ILE:HD12	1.72	0.71
1:A:1:MET:HE3	1:A:106:ALA:HB2	1.71	0.71
1:F:211:LYS:HE2	1:F:211:LYS:HA	1.73	0.71
1:H:307:THR:HG22	1:H:312:ARG:HA	1.73	0.71
1:C:61:MET:HG2	1:C:71:ARG:HD3	1.72	0.69
1:A:119:THR:HB	1:A:135:GLU:HG3	1.73	0.69
1:G:282:ASP:O	1:G:290:MET:HG2	1.91	0.69
1:B:1:MET:HE1	1:B:106:ALA:CB	2.21	0.69
1:F:81:ILE:HD11	1:F:161:LYS:HB2	1.74	0.69
1:B:214:ASN:HA	1:B:217:ILE:CD1	2.23	0.69
1:A:1:MET:CE	1:A:106:ALA:HB2	2.23	0.69
1:A:107:ALA:HB3	1:A:111:ARG:HG3	1.75	0.69
1:H:8:ARG:HG2	1:H:343:MET:HE3	1.73	0.69
1:E:254:LEU:CD2	1:E:366:ARG:HG2	2.23	0.69
1:B:8:ARG:HG2	1:B:343:MET:HE3	1.76	0.68
1:B:83:SER:OG	2:B:401:CIT:O7	2.11	0.68
1:C:119:THR:OG1	1:C:239:ASP:OD1	2.11	0.68
1:E:53:CYS:SG	1:E:234:VAL:HG21	2.33	0.68
1:A:282:ASP:O	1:A:290:MET:HG2	1.93	0.68
1:C:79:ALA:HA	1:C:240:GLU:OE1	1.94	0.68
1:G:213:SER:H	1:G:216:ALA:HB3	1.59	0.68
1:D:161:LYS:HG2	1:D:238:ASN:HB2	1.76	0.68
1:G:119:THR:HB	1:G:135:GLU:CG	2.23	0.68
1:H:163:PRO:HB3	1:H:238:ASN:OD1	1.92	0.67
1:E:234:VAL:HG13	1:E:235:CYS:SG	2.34	0.67
1:D:119:THR:OG1	1:D:239:ASP:OD2	2.11	0.67
1:B:75:VAL:HG11	1:B:236:LEU:HD12	1.76	0.67
1:E:279:ILE:HD11	1:E:292:ARG:HB2	1.77	0.67
1:B:299:LYS:HE2	1:B:353:GLN:OE1	1.96	0.66
1:A:129:ARG:HG2	1:A:214:ASN:ND2	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:38:LEU:HD11	1:D:345:TYR:CZ	2.30	0.66
1:A:166:THR:OG1	1:A:168:GLU:OE2	2.12	0.66
1:C:75:VAL:CG1	1:C:236:LEU:HD12	2.25	0.66
1:C:124:PRO:HB3	1:C:169:ARG:HB3	1.78	0.66
1:C:315:MET:HE2	1:C:326:SER:HB3	1.78	0.66
1:H:296:VAL:HG22	1:H:363:VAL:HG22	1.77	0.66
1:A:328:ARG:NH1	2:A:401:CIT:O6	2.30	0.65
1:C:213:SER:O	1:C:217:ILE:HG13	1.96	0.65
1:D:163:PRO:HA	1:D:238:ASN:ND2	2.10	0.65
1:B:61:MET:HE3	1:C:71:ARG:NH2	2.11	0.65
1:E:161:LYS:CE	1:E:233:GLY:HA2	2.22	0.65
1:F:139:ALA:HB3	1:F:241:ILE:CD1	2.27	0.65
1:H:211:LYS:HE2	1:H:223:LEU:HD21	1.78	0.65
1:A:105:PRO:HG3	1:D:341:ASP:HB2	1.78	0.65
1:B:1:MET:CE	1:B:106:ALA:HA	2.26	0.65
1:E:282:ASP:O	1:E:290:MET:HG2	1.96	0.65
1:G:1:MET:HE1	1:G:106:ALA:HB2	1.79	0.65
1:B:64:ALA:HB1	1:B:69:ILE:O	1.97	0.65
1:H:99:ARG:HD3	1:H:114:VAL:CG1	2.26	0.65
1:B:166:THR:OG1	1:B:168:GLU:OE2	2.15	0.65
1:B:271:ARG:O	1:B:275:LYS:HG3	1.97	0.65
1:H:81:ILE:CD1	1:H:161:LYS:HD2	2.27	0.65
1:A:64:ALA:CA	1:A:74:VAL:HG21	2.27	0.65
1:B:1:MET:HE1	1:B:106:ALA:HB2	1.79	0.65
1:B:75:VAL:CG1	1:B:236:LEU:HD12	2.27	0.65
1:E:214:ASN:HA	1:E:217:ILE:HD12	1.79	0.65
1:E:304:SER:OG	1:E:356:ASP:OD2	2.14	0.65
1:A:279:ILE:HG23	1:A:289:GLU:HB2	1.79	0.64
1:B:300:ALA:CB	1:B:331:VAL:HG21	2.26	0.64
1:F:315:MET:HE2	1:F:326:SER:CB	2.27	0.64
1:B:129:ARG:HG2	1:B:214:ASN:ND2	2.13	0.64
1:A:145:LYS:NZ	1:G:171:GLU:HG2	2.13	0.64
1:B:279:ILE:HD11	1:B:292:ARG:HB2	1.78	0.64
1:D:233:GLY:H	1:D:238:ASN:ND2	1.95	0.64
1:A:315:MET:HE2	1:A:326:SER:HB3	1.79	0.64
1:E:164:LEU:O	1:E:169:ARG:NH1	2.31	0.64
1:C:131:GLU:OE1	1:C:131:GLU:N	2.31	0.64
1:D:1:MET:CE	1:D:106:ALA:HB2	2.28	0.64
1:H:201:ALA:HB2	1:H:225:SER:HB2	1.78	0.63
1:B:21:ALA:HA	1:B:62:TYR:OH	1.98	0.63
1:E:17:ILE:HG22	1:E:59:LEU:HD23	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:161:LYS:NZ	1:D:236:LEU:HD21	2.12	0.63
2:E:401:CIT:O4	2:E:401:CIT:O7	2.16	0.63
1:H:81:ILE:HD13	1:H:161:LYS:HD2	1.79	0.63
1:B:215:GLU:OE2	1:B:215:GLU:N	2.32	0.63
1:G:139:ALA:HB3	1:G:241:ILE:HD12	1.80	0.63
1:H:163:PRO:CG	1:H:238:ASN:OD1	2.47	0.62
1:F:156:HIS:CE1	1:F:257:GLY:HA2	2.34	0.62
1:H:161:LYS:CE	1:H:236:LEU:HD21	2.29	0.62
1:A:119:THR:HB	1:A:135:GLU:CG	2.29	0.62
1:C:161:LYS:HD3	1:C:238:ASN:HB2	1.82	0.62
1:G:303:ALA:HB2	1:G:357:GLY:HA3	1.82	0.62
1:H:213:SER:O	1:H:217:ILE:HG13	1.99	0.62
1:B:161:LYS:CG	1:B:238:ASN:HB3	2.29	0.62
1:E:88:GLY:C	1:H:315:MET:HE3	2.24	0.62
1:D:201:ALA:HB2	1:D:225:SER:HB2	1.80	0.62
1:H:160:ILE:HG12	1:H:241:ILE:CD1	2.24	0.62
1:H:236:LEU:H	1:H:236:LEU:HD23	1.64	0.62
1:D:99:ARG:HD3	1:D:114:VAL:HG13	1.82	0.61
2:F:401:CIT:O3	2:F:401:CIT:O7	2.17	0.61
1:H:161:LYS:HE2	1:H:236:LEU:HD21	1.82	0.61
1:D:218:VAL:HG13	1:D:218:VAL:O	1.98	0.61
1:E:1:MET:HE3	1:E:250:SER:N	2.14	0.61
1:F:107:ALA:HB3	1:F:111:ARG:CZ	2.31	0.61
1:G:122:PHE:CE1	1:G:132:GLN:HA	2.35	0.61
1:C:61:MET:CG	1:C:71:ARG:HD3	2.31	0.61
1:A:10:PRO:CA	1:A:343:MET:HE1	2.31	0.61
1:C:127:LEU:HD12	1:C:165:LEU:HG	1.81	0.61
1:D:292:ARG:HD3	1:D:368:LEU:HD22	1.82	0.61
1:C:166:THR:OG1	1:C:168:GLU:OE2	2.15	0.60
1:C:214:ASN:HA	1:C:217:ILE:HD12	1.82	0.60
1:A:163:PRO:HA	1:A:236:LEU:CD2	2.31	0.60
1:C:127:LEU:HD12	1:C:165:LEU:CD2	2.31	0.60
1:G:138:ARG:NH2	1:G:142:GLU:HG3	2.16	0.60
1:C:1:MET:HE1	1:C:106:ALA:HB2	1.83	0.60
1:E:254:LEU:HD23	1:E:366:ARG:HA	1.84	0.60
1:F:299:LYS:HE2	1:F:353:GLN:OE1	2.01	0.60
1:C:8:ARG:HG2	1:C:343:MET:HE3	1.84	0.60
1:H:1:MET:HE1	1:H:106:ALA:CB	2.31	0.60
1:H:163:PRO:HG3	1:H:238:ASN:OD1	2.01	0.60
1:H:315:MET:HE2	1:H:326:SER:CB	2.29	0.60
1:C:1:MET:HE3	1:C:106:ALA:HB2	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:MET:HG3	1:C:104:ALA:O	2.00	0.60
1:A:265:ILE:HA	1:A:334:VAL:HG21	1.83	0.60
1:B:161:LYS:HG3	1:B:238:ASN:CB	2.31	0.60
1:E:254:LEU:HD21	1:E:366:ARG:HG2	1.83	0.60
1:H:274:LEU:CD1	1:H:290:MET:HE1	2.30	0.60
1:B:315:MET:HE2	1:B:326:SER:HB3	1.85	0.59
1:F:156:HIS:ND1	1:F:257:GLY:HA2	2.18	0.59
1:G:274:LEU:HD11	1:G:339:VAL:HG12	1.84	0.59
1:E:218:VAL:HG12	1:E:218:VAL:O	2.02	0.59
1:F:130:MET:HG2	1:F:214:ASN:OD1	2.02	0.59
1:G:125:GLU:CD	1:G:176:ARG:HH22	2.10	0.59
1:H:161:LYS:HZ3	1:H:236:LEU:HD11	1.68	0.59
1:G:161:LYS:O	1:G:239:ASP:HA	2.03	0.59
1:D:296:VAL:HG22	1:D:363:VAL:HG22	1.84	0.58
1:A:125:GLU:CD	1:A:176:ARG:HH22	2.11	0.58
1:G:125:GLU:O	1:G:129:ARG:HD2	2.04	0.58
1:E:17:ILE:HD12	1:E:17:ILE:C	2.28	0.58
1:F:57:GLN:CD	1:H:57:GLN:HE21	2.12	0.58
1:D:161:LYS:HZ3	1:D:236:LEU:HD21	1.69	0.58
1:E:1:MET:HE1	1:E:106:ALA:CB	2.34	0.58
1:B:52:ARG:HD3	2:B:401:CIT:O4	2.04	0.57
1:F:296:VAL:HG22	1:F:363:VAL:HG22	1.85	0.57
1:A:319:SER:HB3	1:H:322:ASN:HD21	1.68	0.57
1:D:282:ASP:O	1:D:290:MET:HG2	2.04	0.57
1:F:127:LEU:HD12	1:F:165:LEU:CD2	2.34	0.57
1:H:127:LEU:HD12	1:H:165:LEU:CD2	2.34	0.57
1:H:214:ASN:HA	1:H:217:ILE:CD1	2.34	0.57
1:B:57:GLN:OE1	1:C:54:PHE:HE1	1.87	0.57
1:G:137:ALA:HA	1:G:203:ALA:HB2	1.86	0.57
1:A:57:GLN:NE2	1:E:57:GLN:HE21	1.92	0.57
1:B:201:ALA:HB2	1:B:225:SER:HB2	1.85	0.57
1:F:119:THR:OG1	1:F:239:ASP:HB2	2.04	0.57
1:H:218:VAL:HG13	1:H:219:HIS:ND1	2.18	0.57
1:D:235:CYS:HB2	1:D:236:LEU:HD22	1.86	0.57
1:E:21:ALA:HA	1:E:62:TYR:OH	2.04	0.57
1:F:69:ILE:HD11	1:F:74:VAL:HG22	1.87	0.57
1:G:304:SER:OG	1:G:356:ASP:OD2	2.15	0.57
1:E:8:ARG:HG2	1:E:343:MET:HE3	1.86	0.57
1:A:61:MET:HE3	1:E:71:ARG:CZ	2.35	0.57
1:A:164:LEU:N	1:A:236:LEU:HD22	2.20	0.57
1:H:161:LYS:NZ	1:H:236:LEU:HD21	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:218:VAL:HG23	1:F:219:HIS:CE1	2.39	0.57
1:A:64:ALA:HA	1:A:74:VAL:HG21	1.87	0.56
1:F:133:VAL:HG13	1:F:199:GLY:HA2	1.85	0.56
1:G:84:GLY:N	2:G:401:CIT:O4	2.36	0.56
1:E:214:ASN:HA	1:E:217:ILE:CD1	2.35	0.56
1:E:234:VAL:HG13	1:E:235:CYS:N	2.20	0.56
1:B:50:PHE:CE2	1:C:14:PRO:HD2	2.40	0.56
1:G:299:LYS:HD2	1:G:347:SER:OG	2.05	0.56
1:H:136:VAL:O	1:H:140:VAL:HG23	2.06	0.56
1:A:1:MET:CE	1:B:310:GLY:HA3	2.36	0.56
1:G:8:ARG:HG2	1:G:343:MET:HE3	1.86	0.56
1:E:110:LYS:HG3	1:E:248:THR:HG22	1.88	0.56
1:B:71:ARG:CZ	1:C:61:MET:HE3	2.36	0.56
1:C:64:ALA:CA	1:C:74:VAL:HG21	2.35	0.56
1:A:315:MET:HE2	1:A:326:SER:CB	2.36	0.55
1:B:127:LEU:HD12	1:B:165:LEU:CD2	2.36	0.55
1:C:299:LYS:HD2	1:C:347:SER:OG	2.06	0.55
1:D:1:MET:HE1	1:D:106:ALA:HB2	1.88	0.55
1:D:315:MET:HE1	1:D:327:ALA:HA	1.87	0.55
1:F:119:THR:HG22	1:F:139:ALA:HB2	1.89	0.55
1:E:124:PRO:HG3	1:E:169:ARG:HB3	1.89	0.55
1:F:254:LEU:HD23	1:F:366:ARG:HA	1.89	0.55
1:A:99:ARG:HD3	1:A:114:VAL:HG13	1.88	0.55
1:C:130:MET:SD	1:C:212:ILE:HG22	2.47	0.55
1:C:300:ALA:CB	1:C:331:VAL:HG21	2.36	0.55
1:E:213:SER:H	1:E:216:ALA:HB3	1.72	0.55
1:E:279:ILE:HD12	1:E:289:GLU:O	2.07	0.55
1:G:297:LEU:HD23	1:G:345:TYR:HB3	1.89	0.55
1:B:99:ARG:HD3	1:B:114:VAL:CG1	2.36	0.55
1:F:1:MET:HE2	1:F:106:ALA:HB2	1.87	0.55
1:G:254:LEU:CD2	1:G:366:ARG:HG2	2.36	0.55
1:H:38:LEU:HD11	1:H:345:TYR:CZ	2.41	0.55
1:G:125:GLU:OE1	1:G:176:ARG:NH2	2.40	0.55
1:H:218:VAL:HG13	1:H:219:HIS:CE1	2.42	0.55
1:B:69:ILE:HB	1:B:73:GLU:OE1	2.07	0.55
1:B:235:CYS:HB2	1:B:236:LEU:HD22	1.88	0.54
1:C:38:LEU:HD21	1:C:159:GLN:HG3	1.88	0.54
1:D:236:LEU:O	1:D:237:LEU:C	2.49	0.54
1:D:315:MET:HE2	1:D:326:SER:CB	2.36	0.54
1:G:1:MET:HE3	1:G:249:ASN:C	2.33	0.54
1:D:14:PRO:HD2	1:G:50:PHE:CE2	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:119:THR:HB	1:D:135:GLU:HG3	1.88	0.54
1:F:163:PRO:HD3	1:F:239:ASP:OD1	2.07	0.54
1:G:138:ARG:HD3	1:G:138:ARG:C	2.33	0.54
1:H:127:LEU:HD12	1:H:165:LEU:HD23	1.89	0.54
1:H:21:ALA:HA	1:H:62:TYR:OH	2.07	0.54
1:D:4:VAL:H	1:D:252:SER:HB3	1.71	0.54
1:F:17:ILE:C	1:F:17:ILE:HD12	2.32	0.54
1:C:75:VAL:HG11	1:C:236:LEU:CD1	2.37	0.54
1:D:64:ALA:CA	1:D:74:VAL:HG21	2.36	0.54
1:B:99:ARG:HD3	1:B:114:VAL:HG13	1.90	0.54
1:D:52:ARG:NH2	1:D:234:VAL:HG11	2.23	0.54
1:D:161:LYS:HD3	1:D:240:GLU:OE2	2.08	0.54
1:H:17:ILE:HG22	1:H:59:LEU:HD23	1.90	0.54
1:B:119:THR:HG23	1:B:241:ILE:CD1	2.38	0.53
1:E:61:MET:CG	1:E:71:ARG:HD3	2.37	0.53
1:G:211:LYS:HE2	1:G:211:LYS:CA	2.37	0.53
1:A:137:ALA:HA	1:A:203:ALA:HB2	1.89	0.53
1:A:299:LYS:HE2	1:A:353:GLN:OE1	2.08	0.53
1:B:213:SER:O	1:B:217:ILE:HG13	2.07	0.53
1:F:24:ILE:HD12	1:F:62:TYR:HE2	1.72	0.53
1:G:138:ARG:HD3	1:G:139:ALA:N	2.23	0.53
1:A:17:ILE:HD12	1:A:17:ILE:C	2.33	0.53
1:D:265:ILE:HA	1:D:334:VAL:HG21	1.89	0.53
1:A:99:ARG:HD3	1:A:114:VAL:CG1	2.38	0.53
1:G:138:ARG:CZ	1:G:142:GLU:HG3	2.39	0.53
1:A:81:ILE:CD1	1:A:161:LYS:HD3	2.39	0.53
1:F:214:ASN:HA	1:F:217:ILE:HD12	1.90	0.53
1:G:213:SER:O	1:G:217:ILE:HG13	2.09	0.53
1:C:107:ALA:HB3	1:C:111:ARG:HG3	1.91	0.53
1:D:99:ARG:HD3	1:D:114:VAL:CG1	2.38	0.53
1:E:46:SER:HB2	1:F:319:SER:OG	2.08	0.53
1:G:21:ALA:HA	1:G:62:TYR:OH	2.08	0.53
1:C:319:SER:HB3	1:D:322:ASN:HD21	1.73	0.53
1:G:127:LEU:HD12	1:G:165:LEU:CD2	2.39	0.53
1:D:1:MET:HE3	1:D:106:ALA:HB2	1.92	0.52
1:D:239:ASP:C	1:D:240:GLU:HG3	2.35	0.52
1:F:99:ARG:HD3	1:F:114:VAL:HG13	1.90	0.52
1:A:164:LEU:HG	1:A:236:LEU:CD2	2.39	0.52
1:A:270:VAL:HG22	1:A:361:ILE:CD1	2.39	0.52
1:A:275:LYS:HA	1:A:279:ILE:O	2.08	0.52
1:C:99:ARG:HD3	1:C:114:VAL:HG13	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:321:ILE:HG12	1:H:321:ILE:HG12	1.90	0.52
1:B:125:GLU:CD	1:B:176:ARG:HH22	2.17	0.52
1:F:279:ILE:HD13	1:F:293:ILE:HG13	1.91	0.52
1:A:10:PRO:N	1:A:343:MET:HE1	2.24	0.52
1:A:319:SER:HB3	1:H:322:ASN:ND2	2.24	0.52
1:D:368:LEU:HD12	1:D:368:LEU:O	2.09	0.52
1:G:4:VAL:CG2	1:G:114:VAL:HG12	2.39	0.52
1:G:191:TYR:CE1	1:G:218:VAL:HB	2.44	0.52
1:B:191:TYR:CE1	1:B:218:VAL:HG21	2.44	0.52
1:D:1:MET:HG3	1:D:104:ALA:O	2.08	0.52
1:D:254:LEU:HD23	1:D:366:ARG:HA	1.91	0.52
1:F:64:ALA:HB1	1:F:69:ILE:O	2.10	0.52
1:H:107:ALA:CB	1:H:111:ARG:HG3	2.39	0.52
1:A:1:MET:HE3	1:B:310:GLY:HA3	1.89	0.52
1:B:129:ARG:HG2	1:B:214:ASN:HD22	1.75	0.52
1:B:279:ILE:HD12	1:B:289:GLU:O	2.08	0.52
1:D:8:ARG:HG2	1:D:343:MET:HE3	1.90	0.52
1:F:83:SER:OG	2:F:401:CIT:O5	2.06	0.52
1:F:137:ALA:HA	1:F:203:ALA:HB2	1.91	0.52
1:A:163:PRO:HA	1:A:236:LEU:HD22	1.92	0.52
1:D:75:VAL:HG11	1:D:236:LEU:HD12	1.91	0.52
1:F:107:ALA:HB1	1:F:108:PRO:CD	2.33	0.52
1:G:38:LEU:HD11	1:G:345:TYR:CE2	2.45	0.52
1:D:17:ILE:C	1:D:17:ILE:HD12	2.35	0.52
1:H:119:THR:HB	1:H:135:GLU:HG3	1.90	0.52
1:E:348:GLY:H	2:E:401:CIT:C6	2.23	0.52
1:B:300:ALA:HB1	1:B:331:VAL:CG2	2.37	0.51
1:C:46:SER:OG	1:C:47:VAL:N	2.42	0.51
1:D:71:ARG:CZ	1:G:61:MET:HE3	2.41	0.51
1:G:214:ASN:HA	1:G:217:ILE:CD1	2.41	0.51
1:D:164:LEU:HD12	1:D:164:LEU:C	2.36	0.51
1:A:145:LYS:HZ1	1:G:171:GLU:HG2	1.73	0.51
1:C:215:GLU:O	1:C:219:HIS:HB2	2.11	0.51
1:F:253:ASP:O	1:F:254:LEU:HD23	2.11	0.51
1:A:125:GLU:OE1	1:A:176:ARG:NH2	2.40	0.51
1:A:217:ILE:O	1:A:217:ILE:HG22	2.11	0.51
1:E:125:GLU:OE2	1:E:176:ARG:NH2	2.28	0.51
1:H:124:PRO:HG2	1:H:173:ALA:HB2	1.93	0.51
1:H:236:LEU:HD23	1:H:236:LEU:N	2.26	0.51
1:A:64:ALA:N	1:A:74:VAL:HG21	2.26	0.51
1:E:274:LEU:HD13	1:E:290:MET:HE1	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:215:GLU:O	1:H:219:HIS:HB2	2.10	0.51
1:A:255:VAL:HG13	1:A:367:VAL:CG1	2.41	0.51
1:C:282:ASP:O	1:C:290:MET:HG2	2.11	0.51
1:F:99:ARG:HD3	1:F:114:VAL:CG1	2.41	0.51
1:A:163:PRO:HA	1:A:236:LEU:HD23	1.93	0.50
1:G:1:MET:HE3	1:G:250:SER:N	2.26	0.50
1:D:3:LYS:HB3	1:D:100:LYS:HG3	1.92	0.50
1:D:75:VAL:CG1	1:D:236:LEU:HD12	2.41	0.50
1:G:113:ALA:HB3	1:G:149:ILE:HD11	1.93	0.50
1:H:160:ILE:CG1	1:H:241:ILE:HD13	2.26	0.50
1:B:274:LEU:HD21	1:B:293:ILE:HD11	1.93	0.50
1:B:300:ALA:HB2	1:B:331:VAL:HG11	1.93	0.50
1:H:322:ASN:HB2	1:H:325:ARG:HG3	1.93	0.50
1:B:299:LYS:HD2	1:B:347:SER:O	2.12	0.50
1:E:61:MET:HG2	1:E:71:ARG:HD3	1.94	0.50
1:A:164:LEU:HG	1:A:236:LEU:CD1	2.42	0.50
1:F:71:ARG:NH2	1:H:61:MET:HE3	2.27	0.50
1:B:61:MET:HG3	1:B:71:ARG:HD3	1.94	0.50
1:E:217:ILE:O	1:E:352:HIS:NE2	2.40	0.50
1:A:321:ILE:HG12	1:F:321:ILE:HG12	1.92	0.50
1:C:187:LYS:HG2	1:C:191:TYR:HE2	1.75	0.50
1:D:52:ARG:CZ	1:D:234:VAL:HG11	2.41	0.50
1:H:1:MET:HE3	1:H:250:SER:CA	2.42	0.50
1:C:126:GLU:OE2	1:C:131:GLU:HB3	2.11	0.50
1:C:254:LEU:CD2	1:C:366:ARG:HG2	2.41	0.50
1:F:95:THR:OG1	1:F:295:ASN:ND2	2.43	0.50
1:G:142:GLU:OE1	1:G:142:GLU:HA	2.12	0.50
1:D:52:ARG:NH1	2:D:401:CIT:O4	2.45	0.50
1:D:142:GLU:HA	1:D:142:GLU:OE1	2.12	0.50
1:H:119:THR:HG22	1:H:139:ALA:HB2	1.93	0.49
1:D:44:ASN:OD1	1:D:46:SER:OG	2.24	0.49
1:E:303:ALA:HB2	1:E:357:GLY:HA3	1.93	0.49
1:F:81:ILE:CD1	1:F:161:LYS:HB2	2.41	0.49
1:F:274:LEU:HD13	1:F:290:MET:HE1	1.93	0.49
1:A:130:MET:HG2	1:A:214:ASN:OD1	2.12	0.49
1:E:1:MET:HE3	1:E:249:ASN:C	2.37	0.49
1:E:91:THR:OG1	1:H:311:ARG:HB3	2.12	0.49
1:H:17:ILE:HD12	1:H:17:ILE:C	2.38	0.49
1:C:17:ILE:HD12	1:C:17:ILE:C	2.38	0.49
1:E:237:LEU:HD12	1:E:237:LEU:N	2.28	0.49
1:F:136:VAL:O	1:F:140:VAL:HG23	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:160:ILE:HD11	1:E:200:VAL:CG2	2.42	0.49
1:H:163:PRO:CG	1:H:238:ASN:CG	2.85	0.49
1:B:81:ILE:HD12	1:B:159:GLN:HG2	1.95	0.49
1:C:21:ALA:HA	1:C:62:TYR:OH	2.13	0.49
1:C:125:GLU:CD	1:C:176:ARG:HH22	2.21	0.49
1:C:127:LEU:HD12	1:C:165:LEU:CG	2.41	0.49
1:F:8:ARG:CG	1:F:343:MET:HE3	2.41	0.49
1:A:191:TYR:CE1	1:A:218:VAL:HG21	2.48	0.49
1:A:299:LYS:HB2	1:A:360:PRO:HD2	1.94	0.49
1:F:274:LEU:HD22	1:F:293:ILE:HD11	1.95	0.49
1:G:17:ILE:C	1:G:17:ILE:HD12	2.38	0.49
1:A:10:PRO:HA	1:A:343:MET:CE	2.43	0.48
1:C:64:ALA:N	1:C:74:VAL:HG21	2.28	0.48
1:F:57:GLN:NE2	1:H:54:PHE:HE1	2.11	0.48
1:G:107:ALA:HB1	1:G:111:ARG:CZ	2.42	0.48
1:B:83:SER:HG	2:B:401:CIT:HO7	1.50	0.48
1:G:204:LEU:CD1	1:G:227:VAL:HG21	2.43	0.48
1:H:107:ALA:HB3	1:H:111:ARG:HG3	1.95	0.48
1:C:99:ARG:HD3	1:C:114:VAL:CG1	2.42	0.48
1:C:136:VAL:O	1:C:140:VAL:HG23	2.12	0.48
1:E:53:CYS:SG	1:E:234:VAL:CG2	3.00	0.48
1:F:87:GLU:HG2	1:F:325:ARG:HB3	1.95	0.48
1:H:218:VAL:HG22	1:H:218:VAL:O	2.14	0.48
1:B:38:LEU:HD11	1:B:345:TYR:CZ	2.47	0.48
1:B:61:MET:CG	1:B:71:ARG:HD3	2.43	0.48
1:B:119:THR:HG23	1:B:241:ILE:HD11	1.96	0.48
1:C:80:PHE:O	1:C:161:LYS:NZ	2.43	0.48
1:A:136:VAL:O	1:A:140:VAL:HG23	2.14	0.48
1:E:15:ASP:HB3	1:E:54:PHE:CZ	2.48	0.48
1:F:156:HIS:CD2	1:F:256:ILE:HG23	2.49	0.48
1:D:120:ARG:HG2	1:D:135:GLU:CD	2.37	0.48
1:F:156:HIS:CG	1:F:256:ILE:HG23	2.49	0.48
1:D:255:VAL:HG13	1:D:367:VAL:CG1	2.43	0.48
1:D:299:LYS:HE2	1:D:353:GLN:OE1	2.14	0.48
1:F:282:ASP:O	1:F:290:MET:HG2	2.13	0.48
1:E:124:PRO:CG	1:E:169:ARG:HB3	2.44	0.48
1:B:81:ILE:HD12	1:B:159:GLN:CG	2.43	0.47
1:D:119:THR:HG22	1:D:139:ALA:HB2	1.96	0.47
1:E:299:LYS:HD2	1:E:347:SER:OG	2.14	0.47
1:A:64:ALA:HA	1:A:74:VAL:CG2	2.43	0.47
1:B:17:ILE:C	1:B:17:ILE:HD12	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:36:ALA:CB	1:C:79:ALA:HB3	2.44	0.47
1:F:213:SER:O	1:F:217:ILE:HG13	2.14	0.47
1:H:164:LEU:C	1:H:164:LEU:HD12	2.39	0.47
1:A:119:THR:CG2	1:A:239:ASP:HB2	2.45	0.47
1:B:119:THR:CG2	1:B:241:ILE:HD11	2.44	0.47
1:E:161:LYS:HG3	1:E:236:LEU:HD21	1.96	0.47
1:B:81:ILE:CD1	1:B:159:GLN:HG3	2.44	0.47
1:E:206:GLU:OE1	1:E:225:SER:OG	2.29	0.47
1:F:191:TYR:CE1	1:F:218:VAL:HG21	2.49	0.47
1:D:126:GLU:OE2	1:D:131:GLU:HB3	2.14	0.47
1:E:62:TYR:CZ	1:E:66:LYS:HG3	2.49	0.47
1:F:214:ASN:HA	1:F:217:ILE:CD1	2.45	0.47
1:H:163:PRO:HG3	1:H:238:ASN:O	2.13	0.47
1:C:33:GLU:OE1	1:C:100:LYS:HG2	2.14	0.47
1:C:303:ALA:HB2	1:C:357:GLY:HA3	1.97	0.47
1:A:240:GLU:C	1:A:241:ILE:HD12	2.39	0.47
1:C:38:LEU:HG	1:C:81:ILE:HB	1.96	0.47
1:C:321:ILE:CD1	1:G:321:ILE:HG12	2.45	0.47
1:E:69:ILE:HG12	1:E:70:SER:H	1.80	0.47
1:G:81:ILE:HD11	1:G:161:LYS:HB2	1.95	0.47
1:G:123:LEU:HB2	1:G:126:GLU:HG3	1.96	0.47
1:G:274:LEU:CD2	1:G:293:ILE:HD11	2.45	0.47
1:G:278:GLY:O	1:G:292:ARG:NH2	2.47	0.47
1:C:218:VAL:HG13	1:C:218:VAL:O	2.14	0.47
1:E:236:LEU:HD12	1:E:238:ASN:OD1	2.15	0.47
1:F:122:PHE:CE1	1:F:132:GLN:HA	2.49	0.47
1:H:1:MET:HE3	1:H:250:SER:N	2.30	0.47
1:A:69:ILE:HB	1:A:73:GLU:OE1	2.15	0.47
1:B:1:MET:HG3	1:B:2:GLN:N	2.30	0.47
1:C:119:THR:O	1:C:239:ASP:OD2	2.32	0.47
1:E:24:ILE:HD12	1:E:62:TYR:HE2	1.80	0.47
1:B:129:ARG:HH22	1:B:180:VAL:HA	1.75	0.47
1:F:9:ILE:CG2	1:F:19:GLY:HA3	2.45	0.47
1:G:125:GLU:HA	1:G:180:VAL:HG12	1.97	0.47
1:G:163:PRO:HG3	1:G:238:ASN:OD1	2.15	0.47
1:A:129:ARG:HG2	1:A:214:ASN:HD22	1.80	0.46
1:A:163:PRO:HD3	1:A:239:ASP:OD1	2.14	0.46
1:B:130:MET:HG2	1:B:214:ASN:OD1	2.14	0.46
1:E:234:VAL:CG1	1:E:235:CYS:N	2.78	0.46
1:G:38:LEU:HD11	1:G:345:TYR:CZ	2.50	0.46
1:B:236:LEU:HD23	1:B:236:LEU:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:9:ILE:HD11	1:F:23:LEU:HD12	1.98	0.46
1:G:297:LEU:CD2	1:G:345:TYR:HB3	2.45	0.46
1:A:211:LYS:HE2	1:A:211:LYS:CA	2.39	0.46
1:C:161:LYS:HG2	1:C:238:ASN:HD22	1.80	0.46
1:E:64:ALA:HB1	1:E:69:ILE:O	2.16	0.46
1:F:107:ALA:HB3	1:F:111:ARG:NH1	2.31	0.46
1:A:207:ILE:HG13	1:A:212:ILE:HD11	1.97	0.46
1:G:6:VAL:HG22	1:G:97:PHE:CD2	2.51	0.46
1:G:136:VAL:O	1:G:140:VAL:HG23	2.16	0.46
1:G:139:ALA:CB	1:G:241:ILE:HD12	2.44	0.46
1:G:275:LYS:O	1:G:280:ARG:NH1	2.49	0.46
1:A:234:VAL:O	1:A:236:LEU:HG	2.16	0.46
1:B:81:ILE:CD1	1:B:159:GLN:CG	2.94	0.46
1:B:315:MET:HE2	1:B:326:SER:CB	2.44	0.46
1:E:24:ILE:HG21	1:E:66:LYS:HE2	1.96	0.46
1:E:218:VAL:O	1:E:218:VAL:CG1	2.63	0.46
1:F:279:ILE:CD1	1:F:293:ILE:HG13	2.45	0.46
1:H:198:LEU:HD11	1:H:352:HIS:CE1	2.50	0.46
1:C:38:LEU:HA	1:C:81:ILE:O	2.16	0.46
1:D:274:LEU:CD2	1:D:293:ILE:HD11	2.45	0.46
1:F:21:ALA:HA	1:F:62:TYR:OH	2.16	0.46
1:G:187:LYS:HG2	1:G:191:TYR:HE2	1.81	0.46
1:B:125:GLU:O	1:B:129:ARG:CZ	2.64	0.46
1:C:161:LYS:HG2	1:C:238:ASN:ND2	2.30	0.46
1:E:271:ARG:HA	1:E:274:LEU:HD12	1.98	0.46
1:F:87:GLU:HB3	1:F:329:ALA:HB2	1.97	0.46
1:F:254:LEU:CD2	1:F:366:ARG:HG2	2.45	0.46
1:A:52:ARG:NH1	2:A:401:CIT:O3	2.49	0.46
1:B:81:ILE:HD13	1:B:159:GLN:NE2	2.31	0.46
1:D:236:LEU:HD23	1:D:236:LEU:N	2.31	0.46
1:G:120:ARG:HG2	1:G:135:GLU:HG2	1.98	0.46
1:B:274:LEU:CD2	1:B:293:ILE:HD11	2.46	0.45
1:C:214:ASN:HA	1:C:217:ILE:CD1	2.46	0.45
1:C:279:ILE:CD1	1:C:293:ILE:HG13	2.46	0.45
1:F:69:ILE:HB	1:F:73:GLU:OE1	2.16	0.45
1:A:234:VAL:HG23	1:A:235:CYS:N	2.31	0.45
1:B:337:SER:HB3	1:D:334:VAL:HG22	1.98	0.45
1:A:322:ASN:HB3	1:A:325:ARG:HG3	1.98	0.45
1:C:64:ALA:HB1	1:C:69:ILE:O	2.16	0.45
1:E:240:GLU:C	1:E:241:ILE:HD12	2.41	0.45
1:F:126:GLU:OE2	1:F:131:GLU:HG2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:126:GLU:OE2	1:G:131:GLU:HG2	2.16	0.45
1:C:156:HIS:CE1	1:C:257:GLY:HA2	2.51	0.45
1:G:107:ALA:HB1	1:G:108:PRO:HD2	1.98	0.45
1:H:119:THR:HB	1:H:135:GLU:CG	2.46	0.45
1:B:303:ALA:HB2	1:B:357:GLY:HA3	1.98	0.45
1:G:81:ILE:HD13	1:G:161:LYS:HD3	1.98	0.45
1:B:36:ALA:CB	1:B:79:ALA:HB3	2.47	0.45
1:B:165:LEU:O	1:B:185:THR:HG22	2.17	0.45
1:F:299:LYS:HD3	1:F:347:SER:OG	2.17	0.45
1:H:279:ILE:CD1	1:H:293:ILE:HG13	2.47	0.45
1:A:270:VAL:HG22	1:A:361:ILE:HD13	1.97	0.45
1:C:240:GLU:C	1:C:241:ILE:HD12	2.41	0.45
1:C:298:ALA:O	1:C:346:VAL:HA	2.17	0.45
1:D:161:LYS:O	1:D:238:ASN:O	2.35	0.45
1:E:270:VAL:HG22	1:E:361:ILE:CD1	2.46	0.45
1:G:290:MET:HE3	1:G:338:VAL:O	2.17	0.45
1:G:303:ALA:HA	1:G:356:ASP:OD1	2.17	0.45
1:F:8:ARG:HG2	1:F:343:MET:CE	2.40	0.45
1:A:7:PHE:CD1	1:A:29:ILE:HD11	2.51	0.45
1:A:201:ALA:HB2	1:A:225:SER:HB2	1.99	0.45
1:D:45:GLY:HA2	2:D:401:CIT:O3	2.17	0.45
1:F:1:MET:HB3	1:F:102:VAL:O	2.17	0.45
1:B:136:VAL:O	1:B:140:VAL:HG23	2.16	0.44
1:C:321:ILE:HD11	1:G:321:ILE:HG12	1.98	0.44
1:E:17:ILE:HG22	1:E:59:LEU:CD2	2.46	0.44
1:G:140:VAL:HG23	1:G:241:ILE:HD13	1.98	0.44
1:A:72:GLU:O	1:A:76:LYS:HG3	2.17	0.44
1:B:164:LEU:HD12	1:B:164:LEU:C	2.43	0.44
1:B:253:ASP:O	1:B:367:VAL:HG22	2.17	0.44
1:F:160:ILE:HD11	1:F:200:VAL:CG2	2.47	0.44
1:H:137:ALA:HA	1:H:203:ALA:HB2	1.99	0.44
1:C:236:LEU:N	1:C:236:LEU:HD23	2.32	0.44
1:D:38:LEU:HD11	1:D:345:TYR:OH	2.17	0.44
1:G:191:TYR:OH	1:G:218:VAL:HG11	2.18	0.44
1:H:64:ALA:CA	1:H:74:VAL:HG21	2.46	0.44
1:A:282:ASP:HA	1:A:290:MET:HE2	1.98	0.44
1:C:294:VAL:HB	1:C:364:ILE:HG22	1.99	0.44
1:C:300:ALA:HB1	1:C:331:VAL:HG21	2.00	0.44
1:D:292:ARG:NH1	1:D:368:LEU:HD23	2.32	0.44
1:E:99:ARG:NH2	1:E:114:VAL:O	2.46	0.44
1:E:163:PRO:HD3	1:E:239:ASP:OD1	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:VAL:CG1	1:B:236:LEU:CD1	2.95	0.44
1:D:254:LEU:CD2	1:D:366:ARG:HG2	2.47	0.44
1:F:217:ILE:O	1:F:217:ILE:HG22	2.17	0.44
1:B:161:LYS:HE3	1:B:233:GLY:HA3	1.99	0.44
1:E:38:LEU:HA	1:E:81:ILE:O	2.18	0.44
1:F:211:LYS:HE2	1:F:211:LYS:CA	2.46	0.44
1:G:1:MET:HE3	1:G:250:SER:CA	2.48	0.44
1:G:254:LEU:HD23	1:G:366:ARG:HG2	1.99	0.44
1:B:1:MET:HE3	1:B:250:SER:CA	2.48	0.44
1:C:75:VAL:CG1	1:C:236:LEU:CD1	2.94	0.44
1:E:31:PRO:HB2	1:E:67:LEU:HD21	2.00	0.44
1:E:38:LEU:HD11	1:E:345:TYR:CZ	2.53	0.44
1:E:69:ILE:HG23	1:E:74:VAL:HG23	1.99	0.44
1:H:274:LEU:HD13	1:H:290:MET:CE	2.37	0.44
1:A:61:MET:SD	1:E:61:MET:SD	3.16	0.44
1:A:93:HIS:HB2	1:A:343:MET:HE3	1.97	0.44
1:A:105:PRO:HD3	1:D:341:ASP:OD2	2.17	0.44
1:B:20:LEU:HD21	1:B:63:LEU:HD21	2.00	0.44
1:E:110:LYS:HE3	1:E:248:THR:CG2	2.35	0.44
1:E:164:LEU:HD12	1:E:164:LEU:C	2.43	0.44
1:H:368:LEU:C	1:H:368:LEU:HD23	2.41	0.44
1:D:64:ALA:HA	1:D:74:VAL:HG21	1.99	0.43
1:D:125:GLU:CD	1:D:176:ARG:HH22	2.25	0.43
1:H:1:MET:HE3	1:H:249:ASN:C	2.42	0.43
1:B:81:ILE:HD13	1:B:159:GLN:HE21	1.81	0.43
1:D:315:MET:HE2	1:D:326:SER:HB2	1.99	0.43
1:G:145:LYS:HE2	1:G:145:LYS:HB3	1.87	0.43
1:H:161:LYS:HZ3	1:H:236:LEU:CD1	2.28	0.43
1:F:71:ARG:CZ	1:H:61:MET:HE3	2.47	0.43
1:A:119:THR:O	1:A:238:ASN:HB2	2.19	0.43
1:D:213:SER:O	1:D:217:ILE:HG13	2.18	0.43
1:D:292:ARG:HD3	1:D:368:LEU:CD2	2.45	0.43
1:E:110:LYS:CE	1:E:248:THR:HG22	2.39	0.43
1:F:87:GLU:CB	1:F:329:ALA:HB2	2.48	0.43
1:F:119:THR:CB	1:F:135:GLU:HG3	2.34	0.43
1:A:148:GLN:OE1	1:A:249:ASN:ND2	2.52	0.43
1:B:57:GLN:NE2	1:C:57:GLN:HG2	2.34	0.43
1:C:189:MET:HE2	1:C:232:ALA:HB3	2.00	0.43
1:D:157:PHE:HB3	1:D:244:VAL:HB	2.00	0.43
1:H:217:ILE:HG22	1:H:217:ILE:O	2.18	0.43
1:B:311:ARG:HG3	1:D:342:PRO:HD2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:274:LEU:HD22	1:D:293:ILE:HD11	2.01	0.43
1:F:54:PHE:HE1	1:H:57:GLN:OE1	2.00	0.43
1:H:38:LEU:HD11	1:H:345:TYR:CE2	2.53	0.43
1:D:307:THR:HG22	1:D:312:ARG:HA	2.00	0.43
1:E:6:VAL:HG22	1:E:97:PHE:CD2	2.54	0.43
1:A:164:LEU:HG	1:A:236:LEU:HD22	2.00	0.43
1:B:368:LEU:N	1:B:368:LEU:HD12	2.33	0.43
1:E:38:LEU:HD11	1:E:345:TYR:CE2	2.54	0.43
1:E:140:VAL:HG21	1:E:200:VAL:HA	2.01	0.43
1:E:236:LEU:C	1:E:237:LEU:HD12	2.44	0.43
1:F:111:ARG:HB3	1:F:252:SER:O	2.19	0.43
1:F:3:LYS:HB3	1:F:100:LYS:HG3	2.01	0.42
1:H:1:MET:HE1	1:H:106:ALA:HA	1.95	0.42
1:C:164:LEU:C	1:C:164:LEU:HD12	2.44	0.42
1:C:234:VAL:O	1:C:234:VAL:HG22	2.19	0.42
1:E:136:VAL:HG11	1:E:196:SER:O	2.19	0.42
1:F:120:ARG:HG2	1:F:135:GLU:CD	2.44	0.42
1:B:145:LYS:HE2	1:B:145:LYS:HB3	1.95	0.42
1:C:164:LEU:HD21	1:C:234:VAL:HB	2.01	0.42
1:G:3:LYS:HG3	1:G:251:ALA:O	2.20	0.42
1:G:9:ILE:HD11	1:G:23:LEU:HD12	2.01	0.42
1:A:93:HIS:CB	1:A:343:MET:CE	2.94	0.42
1:B:161:LYS:HG3	1:B:238:ASN:CG	2.44	0.42
1:C:9:ILE:CG2	1:C:19:GLY:HA3	2.49	0.42
1:D:165:LEU:O	1:D:185:THR:HG22	2.18	0.42
1:G:9:ILE:HG21	1:G:19:GLY:HA3	2.02	0.42
1:H:17:ILE:HG22	1:H:59:LEU:CD2	2.48	0.42
1:B:161:LYS:CE	1:B:233:GLY:HA3	2.48	0.42
1:C:119:THR:N	1:C:239:ASP:OD2	2.53	0.42
1:D:315:MET:HE2	1:D:326:SER:C	2.44	0.42
1:E:125:GLU:CD	1:E:176:ARG:HH22	2.20	0.42
1:E:315:MET:HE1	1:E:330:VAL:HG21	2.01	0.42
1:F:201:ALA:HB2	1:F:225:SER:HB2	2.00	0.42
1:H:300:ALA:CB	1:H:331:VAL:HG21	2.49	0.42
1:B:125:GLU:HA	1:B:129:ARG:NH2	2.34	0.42
1:D:368:LEU:HA	1:D:368:LEU:HD13	1.78	0.42
1:H:117:ALA:O	1:H:241:ILE:HB	2.19	0.42
1:D:233:GLY:HA3	1:D:238:ASN:HB3	2.01	0.42
1:E:136:VAL:O	1:E:140:VAL:HG23	2.19	0.42
1:G:4:VAL:HG22	1:G:114:VAL:HG12	2.01	0.42
1:G:52:ARG:HH11	2:G:401:CIT:C5	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:LEU:HD12	1:B:165:LEU:HD23	2.01	0.42
1:A:10:PRO:HA	1:A:343:MET:HE3	2.02	0.42
1:C:44:ASN:OD1	1:C:325:ARG:NH2	2.53	0.42
1:E:72:GLU:O	1:E:76:LYS:HG3	2.19	0.42
1:E:348:GLY:N	2:E:401:CIT:O6	2.51	0.42
1:F:133:VAL:HG13	1:F:199:GLY:CA	2.48	0.42
1:A:52:ARG:HD2	1:A:82:MET:O	2.20	0.42
1:F:113:ALA:HB3	1:F:149:ILE:HD11	2.02	0.42
1:G:270:VAL:HG12	1:G:338:VAL:HG11	2.02	0.42
1:G:298:ALA:O	1:G:346:VAL:HA	2.20	0.42
1:A:342:PRO:HD2	1:F:311:ARG:HG3	2.02	0.41
1:B:91:THR:OG1	1:D:311:ARG:HB3	2.19	0.41
1:F:274:LEU:CD2	1:F:293:ILE:HD11	2.50	0.41
1:A:254:LEU:HD23	1:A:366:ARG:HA	2.02	0.41
1:C:1:MET:HE1	1:C:249:ASN:ND2	2.35	0.41
1:F:57:GLN:HE22	1:H:54:PHE:HE1	1.68	0.41
1:G:81:ILE:CD1	1:G:161:LYS:HB2	2.49	0.41
1:G:289:GLU:CD	1:G:292:ARG:HH12	2.28	0.41
1:B:217:ILE:O	1:B:217:ILE:HG22	2.20	0.41
1:D:207:ILE:CD1	1:D:212:ILE:HD11	2.50	0.41
1:F:9:ILE:CD1	1:F:23:LEU:HD12	2.49	0.41
1:F:274:LEU:HD23	1:F:363:VAL:HG21	2.03	0.41
1:H:64:ALA:HB2	1:H:74:VAL:HG21	2.03	0.41
1:C:64:ALA:HA	1:C:74:VAL:HG21	2.02	0.41
1:D:297:LEU:HD23	1:D:345:TYR:HB3	2.01	0.41
1:F:127:LEU:HD12	1:F:165:LEU:HD21	2.02	0.41
1:G:115:GLY:N	1:G:147:ALA:HB2	2.36	0.41
1:H:238:ASN:CG	1:H:238:ASN:O	2.63	0.41
1:C:80:PHE:O	1:C:161:LYS:HE3	2.21	0.41
1:F:119:THR:HG22	1:F:139:ALA:CB	2.51	0.41
1:F:368:LEU:C	1:F:368:LEU:HD23	2.45	0.41
1:A:234:VAL:HG23	1:A:235:CYS:H	1.85	0.41
1:B:120:ARG:HG2	1:B:135:GLU:CD	2.44	0.41
1:C:217:ILE:HG22	1:C:217:ILE:O	2.20	0.41
1:C:274:LEU:CD2	1:C:293:ILE:HD11	2.50	0.41
1:D:107:ALA:HB3	1:D:111:ARG:HG3	2.02	0.41
1:E:61:MET:HG3	1:E:71:ARG:HD3	2.02	0.41
1:G:1:MET:CE	1:G:106:ALA:HB2	2.50	0.41
1:G:4:VAL:HG21	1:G:114:VAL:HG12	2.02	0.41
1:H:1:MET:HE1	1:H:106:ALA:HB2	2.01	0.41
1:H:157:PHE:CZ	1:H:297:LEU:HD13	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:ARG:HD2	1:B:82:MET:O	2.21	0.41
1:A:132:GLN:O	1:A:136:VAL:HG23	2.21	0.41
1:F:1:MET:HE3	1:F:250:SER:N	2.36	0.41
1:A:119:THR:HG21	1:A:239:ASP:HB2	2.02	0.41
1:C:253:ASP:O	1:C:254:LEU:HD23	2.21	0.41
1:D:3:LYS:HG3	1:D:252:SER:HA	2.03	0.41
1:D:240:GLU:C	1:D:241:ILE:HD12	2.46	0.41
1:F:1:MET:HE3	1:F:250:SER:CA	2.50	0.41
1:F:57:GLN:NE2	1:H:54:PHE:CE1	2.88	0.41
1:A:95:THR:OG1	1:A:295:ASN:ND2	2.54	0.41
1:B:1:MET:CE	1:B:106:ALA:CA	2.98	0.41
1:B:368:LEU:HD12	1:B:368:LEU:H	1.85	0.41
1:E:110:LYS:HG3	1:E:248:THR:CG2	2.50	0.41
1:G:154:ASP:HB3	1:G:246:ASN:O	2.21	0.41
1:C:119:THR:OG1	1:C:239:ASP:CG	2.64	0.40
1:D:211:LYS:HE2	1:D:211:LYS:CA	2.44	0.40
1:A:164:LEU:HD12	1:A:164:LEU:C	2.47	0.40
1:G:9:ILE:CG2	1:G:19:GLY:HA3	2.51	0.40
1:G:282:ASP:OD1	1:G:290:MET:HE2	2.20	0.40
1:H:254:LEU:HD23	1:H:366:ARG:HA	2.03	0.40
1:A:274:LEU:HD11	1:A:339:VAL:HG12	2.03	0.40
1:C:220:ASP:OD1	1:C:222:ASN:HB2	2.21	0.40
1:E:31:PRO:HG2	1:E:67:LEU:HD21	2.02	0.40
1:A:170:ILE:HG12	1:A:180:VAL:HG21	2.02	0.40
1:E:361:ILE:O	1:E:361:ILE:HG23	2.21	0.40
1:F:335:ILE:O	1:F:339:VAL:HG22	2.21	0.40
1:G:226:SER:O	1:G:360:PRO:HB3	2.21	0.40
1:H:274:LEU:HD11	1:H:339:VAL:HG12	2.03	0.40
1:D:217:ILE:O	1:D:217:ILE:HG22	2.21	0.40
1:E:337:SER:HB3	1:H:334:VAL:HG22	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:ARG:NH1	1:E:210:ASP:OD2[4_455]	1.97	0.23

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	363/370 (98%)	351 (97%)	12 (3%)	0	100	100
1	B	361/370 (98%)	349 (97%)	12 (3%)	0	100	100
1	C	360/370 (97%)	345 (96%)	15 (4%)	0	100	100
1	D	362/370 (98%)	347 (96%)	15 (4%)	0	100	100
1	E	362/370 (98%)	351 (97%)	11 (3%)	0	100	100
1	F	353/370 (95%)	344 (98%)	9 (2%)	0	100	100
1	G	355/370 (96%)	347 (98%)	8 (2%)	0	100	100
1	H	361/370 (98%)	349 (97%)	12 (3%)	0	100	100
All	All	2877/2960 (97%)	2783 (97%)	94 (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	288/293 (98%)	286 (99%)	2 (1%)	76	83
1	B	286/293 (98%)	286 (100%)	0	100	100
1	C	285/293 (97%)	284 (100%)	1 (0%)	84	86
1	D	287/293 (98%)	285 (99%)	2 (1%)	76	83
1	E	287/293 (98%)	286 (100%)	1 (0%)	86	88
1	F	281/293 (96%)	280 (100%)	1 (0%)	84	86

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	282/293 (96%)	281 (100%)	1 (0%)	84	86
1	H	286/293 (98%)	285 (100%)	1 (0%)	86	88
All	All	2282/2344 (97%)	2273 (100%)	9 (0%)	84	86

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	368	LEU
1	A	370	HIS
1	C	218	VAL
1	D	218	VAL
1	D	368	LEU
1	E	368	LEU
1	F	218	VAL
1	G	218	VAL
1	H	368	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	GLN
1	A	148	GLN
1	A	183	ASN
1	A	214	ASN
1	A	249	ASN
1	B	2	GLN
1	B	57	GLN
1	B	159	GLN
1	B	214	ASN
1	C	57	GLN
1	C	322	ASN
1	D	238	ASN
1	D	322	ASN
1	E	148	GLN
1	F	2	GLN
1	F	57	GLN
1	F	134	ASN
1	F	249	ASN
1	F	322	ASN
1	G	2	GLN

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Mol	Chain	Res	Type
1	H	57	GLN
1	H	322	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

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	$\# Z > 2$
2	CIT	H	401	-	12,12,12	1.28	2 (16%)	17,17,17	1.74	6 (35%)
2	CIT	E	401	-	12,12,12	1.12	0	17,17,17	1.37	1 (5%)
2	CIT	D	401	-	12,12,12	1.23	1 (8%)	17,17,17	1.96	6 (35%)
2	CIT	F	401	-	12,12,12	1.00	0	17,17,17	1.62	2 (11%)
2	CIT	B	401	-	12,12,12	1.07	0	17,17,17	1.87	5 (29%)
2	CIT	G	401	-	12,12,12	1.09	0	17,17,17	1.59	2 (11%)
2	CIT	C	401	-	12,12,12	1.06	0	17,17,17	1.92	3 (17%)
2	CIT	A	401	-	12,12,12	1.19	0	17,17,17	1.53	5 (29%)

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
2	CIT	H	401	-	-	10/16/16/16	-
2	CIT	E	401	-	-	9/16/16/16	-
2	CIT	D	401	-	-	7/16/16/16	-
2	CIT	F	401	-	-	9/16/16/16	-
2	CIT	B	401	-	-	7/16/16/16	-
2	CIT	G	401	-	-	11/16/16/16	-
2	CIT	C	401	-	-	2/16/16/16	-
2	CIT	A	401	-	-	9/16/16/16	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	CIT	C3-C6	2.15	1.55	1.53
2	H	401	CIT	C3-C6	2.09	1.55	1.53
2	H	401	CIT	O5-C6	2.07	1.28	1.22

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	CIT	O6-C6-C3	5.22	123.15	113.14
2	D	401	CIT	O6-C6-C3	4.99	122.71	113.14
2	G	401	CIT	O6-C6-C3	4.12	121.05	113.14
2	F	401	CIT	O6-C6-C3	4.08	120.96	113.14
2	B	401	CIT	O6-C6-C3	3.95	120.71	113.14
2	E	401	CIT	O6-C6-C3	3.68	120.21	113.14
2	H	401	CIT	O6-C6-C3	3.67	120.17	113.14
2	B	401	CIT	O7-C3-C2	-2.83	102.92	109.38
2	B	401	CIT	O4-C5-C4	2.71	122.92	114.35
2	H	401	CIT	O2-C1-O1	-2.59	116.68	123.33
2	D	401	CIT	O2-C1-O1	-2.53	116.84	123.33
2	A	401	CIT	O2-C1-C2	2.47	122.18	114.35
2	D	401	CIT	C3-C4-C5	-2.36	107.46	113.92
2	D	401	CIT	O4-C5-C4	2.35	121.78	114.35
2	A	401	CIT	O2-C1-O1	-2.31	117.39	123.33
2	C	401	CIT	C2-C3-C6	2.29	115.11	110.03
2	D	401	CIT	O2-C1-C2	2.27	121.55	114.35
2	A	401	CIT	O5-C6-C3	2.25	126.45	122.09
2	F	401	CIT	O2-C1-C2	2.22	121.38	114.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	401	CIT	O4-C5-C4	2.22	121.37	114.35
2	H	401	CIT	O2-C1-C2	2.16	121.18	114.35
2	H	401	CIT	O4-C5-O3	-2.15	117.80	123.33
2	B	401	CIT	O2-C1-C2	2.15	121.16	114.35
2	B	401	CIT	C2-C3-C6	2.07	114.61	110.03
2	H	401	CIT	O7-C3-C6	-2.05	106.06	108.96
2	A	401	CIT	O6-C6-O5	-2.04	117.34	123.86
2	D	401	CIT	O6-C6-O5	-2.02	117.39	123.86
2	C	401	CIT	O2-C1-O1	-2.01	118.16	123.33
2	G	401	CIT	O2-C1-C2	2.01	120.70	114.35
2	A	401	CIT	O4-C5-C4	2.00	120.69	114.35

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	CIT	O7-C3-C4-C5
2	A	401	CIT	C6-C3-C4-C5
2	A	401	CIT	C2-C3-C6-O5
2	A	401	CIT	C2-C3-C6-O6
2	A	401	CIT	O7-C3-C6-O5
2	A	401	CIT	O7-C3-C6-O6
2	B	401	CIT	O7-C3-C6-O5
2	B	401	CIT	O7-C3-C6-O6
2	B	401	CIT	C4-C3-C6-O5
2	B	401	CIT	C4-C3-C6-O6
2	D	401	CIT	C2-C3-C4-C5
2	D	401	CIT	O7-C3-C4-C5
2	D	401	CIT	C6-C3-C4-C5
2	D	401	CIT	O7-C3-C6-O5
2	D	401	CIT	O7-C3-C6-O6
2	D	401	CIT	C4-C3-C6-O5
2	D	401	CIT	C4-C3-C6-O6
2	E	401	CIT	C2-C3-C4-C5
2	E	401	CIT	O7-C3-C4-C5
2	E	401	CIT	C6-C3-C4-C5
2	E	401	CIT	O7-C3-C6-O5
2	E	401	CIT	O7-C3-C6-O6
2	E	401	CIT	C4-C3-C6-O5
2	E	401	CIT	C4-C3-C6-O6
2	F	401	CIT	O7-C3-C6-O5
2	F	401	CIT	O7-C3-C6-O6

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Mol	Chain	Res	Type	Atoms
2	F	401	CIT	C4-C3-C6-O5
2	F	401	CIT	C4-C3-C6-O6
2	G	401	CIT	C2-C3-C6-O5
2	G	401	CIT	C2-C3-C6-O6
2	G	401	CIT	O7-C3-C6-O5
2	G	401	CIT	O7-C3-C6-O6
2	H	401	CIT	C1-C2-C3-O7
2	H	401	CIT	C1-C2-C3-C4
2	H	401	CIT	C1-C2-C3-C6
2	H	401	CIT	C2-C3-C6-O5
2	H	401	CIT	C2-C3-C6-O6
2	H	401	CIT	O7-C3-C6-O5
2	H	401	CIT	O7-C3-C6-O6
2	A	401	CIT	C2-C3-C4-C5
2	F	401	CIT	C1-C2-C3-C4
2	H	401	CIT	C4-C3-C6-O6
2	F	401	CIT	C1-C2-C3-C6
2	F	401	CIT	C1-C2-C3-O7
2	E	401	CIT	C3-C4-C5-O3
2	E	401	CIT	C3-C4-C5-O4
2	F	401	CIT	C2-C3-C6-O5
2	F	401	CIT	C2-C3-C6-O6
2	H	401	CIT	C4-C3-C6-O5
2	G	401	CIT	C1-C2-C3-C6
2	C	401	CIT	C1-C2-C3-O7
2	C	401	CIT	C1-C2-C3-C4
2	G	401	CIT	C4-C3-C6-O6
2	G	401	CIT	O7-C3-C4-C5
2	G	401	CIT	C1-C2-C3-C4
2	G	401	CIT	C6-C3-C4-C5
2	H	401	CIT	C6-C3-C4-C5
2	B	401	CIT	C1-C2-C3-O7
2	A	401	CIT	O2-C1-C2-C3
2	A	401	CIT	O1-C1-C2-C3
2	G	401	CIT	O1-C1-C2-C3
2	G	401	CIT	O2-C1-C2-C3
2	B	401	CIT	O1-C1-C2-C3
2	B	401	CIT	O2-C1-C2-C3

There are no ring outliers.

6 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	401	CIT	3	0
2	D	401	CIT	2	0
2	F	401	CIT	2	0
2	B	401	CIT	3	0
2	G	401	CIT	2	0
2	A	401	CIT	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	365/370 (98%)	-0.88	0 100 100	62, 112, 196, 251	0
1	B	363/370 (98%)	-0.77	0 100 100	68, 114, 179, 254	0
1	C	362/370 (97%)	-0.70	0 100 100	86, 133, 196, 242	0
1	D	364/370 (98%)	-0.73	0 100 100	75, 114, 166, 230	0
1	E	364/370 (98%)	-0.81	0 100 100	84, 137, 210, 259	0
1	F	357/370 (96%)	-0.70	1 (0%) 90 81	91, 165, 230, 306	0
1	G	359/370 (97%)	-0.74	0 100 100	90, 166, 236, 297	0
1	H	363/370 (98%)	-0.68	0 100 100	77, 126, 179, 254	0
All	All	2897/2960 (97%)	-0.75	1 (0%) 100 100	62, 131, 212, 306	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	232	ALA	2.6

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	CIT	D	401	13/13	0.70	0.10	93,100,109,110	0
2	CIT	E	401	13/13	0.72	0.09	96,109,113,115	0
2	CIT	B	401	13/13	0.85	0.07	73,85,90,91	0
2	CIT	G	401	13/13	0.85	0.06	126,140,153,154	0
2	CIT	H	401	13/13	0.85	0.08	87,111,122,123	0
2	CIT	C	401	13/13	0.86	0.07	87,105,119,119	0
2	CIT	F	401	13/13	0.89	0.06	101,139,147,148	0
2	CIT	A	401	13/13	0.91	0.06	84,104,116,116	0

6.5 Other polymers ⓘ

There are no such residues in this entry.