



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:36 AM UTC

PDB ID : 5ING / pdb\_00005ing  
Title : A crotonyl-CoA reductase-carboxylase independent pathway for assembly of unusual alkylmalonyl-CoA polyketide synthase extender unit  
Authors : Valentic, T.R.; Ray, L.; Miyazawa, T.; Song, L.; Withall, D.M.; Milligan, J.C.; Takahashi, S.; Osada, H.; Tsai, S.C.; Challis, G.L.  
Deposited on : 2016-03-07  
Resolution : 2.45 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](mailto: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>.

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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

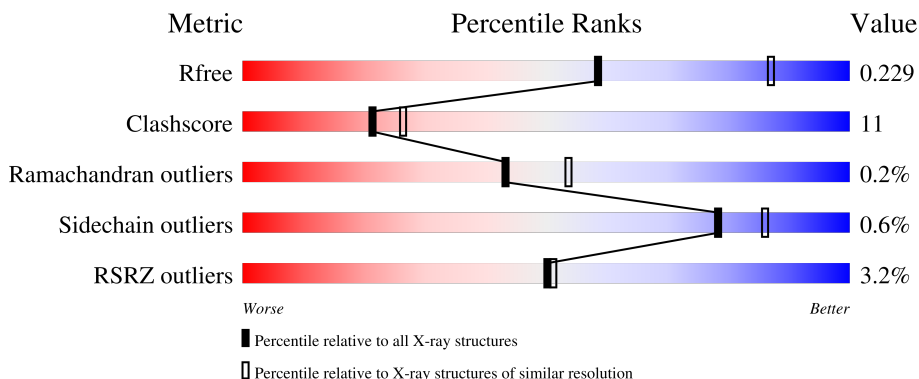
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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.



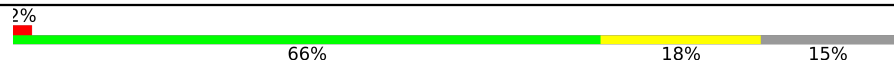
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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  $\geq 3$ , 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	570	 3% 71% 15% • 13%
1	B	570	 3% 72% 15% • 13%
1	C	570	 4% 71% 16% • 12%
1	D	570	 % 71% 15% 14%
1	E	570	 3% 68% 18% 14%

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Mol	Chain	Length	Quality of chain										
1	F	570	 <p>A horizontal bar chart showing the quality distribution of the chain. The bar is divided into four segments: a small red segment (2%), a large green segment (66%), a yellow segment (18%), and a grey segment (15%).</p> <table border="1"><thead><tr><th>Category</th><th>Percentage</th></tr></thead><tbody><tr><td>Red</td><td>2%</td></tr><tr><td>Green</td><td>66%</td></tr><tr><td>Yellow</td><td>18%</td></tr><tr><td>Grey</td><td>15%</td></tr></tbody></table>	Category	Percentage	Red	2%	Green	66%	Yellow	18%	Grey	15%
Category	Percentage												
Red	2%												
Green	66%												
Yellow	18%												
Grey	15%												

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 23360 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 Putative carboxyl transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	496	3760	2366	661	714	6	13	0	0	0
1	B	497	3763	2368	662	714	6	13	0	0	0
1	C	503	3806	2394	670	723	6	13	0	0	0
1	D	488	3700	2329	653	699	6	13	0	0	0
1	E	493	3735	2349	659	709	6	12	0	0	0
1	F	483	3660	2303	648	690	6	13	0	0	0

There are 234 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-37	MSE	-	initiating methionine	UNP A0ACI9
A	-36	GLY	-	expression tag	UNP A0ACI9
A	-35	SER	-	expression tag	UNP A0ACI9
A	-34	SER	-	expression tag	UNP A0ACI9
A	-33	HIS	-	expression tag	UNP A0ACI9
A	-32	HIS	-	expression tag	UNP A0ACI9
A	-31	HIS	-	expression tag	UNP A0ACI9
A	-30	HIS	-	expression tag	UNP A0ACI9
A	-29	HIS	-	expression tag	UNP A0ACI9
A	-28	HIS	-	expression tag	UNP A0ACI9
A	-27	SER	-	expression tag	UNP A0ACI9
A	-26	SER	-	expression tag	UNP A0ACI9
A	-25	GLY	-	expression tag	UNP A0ACI9
A	-24	LEU	-	expression tag	UNP A0ACI9
A	-23	VAL	-	expression tag	UNP A0ACI9
A	-22	PRO	-	expression tag	UNP A0ACI9
A	-21	ARG	-	expression tag	UNP A0ACI9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	GLY	-	expression tag	UNP A0ACI9
A	-19	SER	-	expression tag	UNP A0ACI9
A	-18	HIS	-	expression tag	UNP A0ACI9
A	-17	MSE	-	expression tag	UNP A0ACI9
A	-16	ALA	-	expression tag	UNP A0ACI9
A	-15	SER	-	expression tag	UNP A0ACI9
A	-14	ASP	-	expression tag	UNP A0ACI9
A	-13	SER	-	expression tag	UNP A0ACI9
A	-12	THR	-	expression tag	UNP A0ACI9
A	-11	GLU	-	expression tag	UNP A0ACI9
A	-10	ASN	-	expression tag	UNP A0ACI9
A	-9	LEU	-	expression tag	UNP A0ACI9
A	-8	TYR	-	expression tag	UNP A0ACI9
A	-7	PHE	-	expression tag	UNP A0ACI9
A	-6	GLN	-	expression tag	UNP A0ACI9
A	-5	GLY	-	expression tag	UNP A0ACI9
A	-4	ILE	-	expression tag	UNP A0ACI9
A	-3	ASP	-	expression tag	UNP A0ACI9
A	-2	PRO	-	expression tag	UNP A0ACI9
A	-1	PHE	-	expression tag	UNP A0ACI9
A	0	THR	-	expression tag	UNP A0ACI9
A	1	MSE	-	expression tag	UNP A0ACI9
B	-37	MSE	-	initiating methionine	UNP A0ACI9
B	-36	GLY	-	expression tag	UNP A0ACI9
B	-35	SER	-	expression tag	UNP A0ACI9
B	-34	SER	-	expression tag	UNP A0ACI9
B	-33	HIS	-	expression tag	UNP A0ACI9
B	-32	HIS	-	expression tag	UNP A0ACI9
B	-31	HIS	-	expression tag	UNP A0ACI9
B	-30	HIS	-	expression tag	UNP A0ACI9
B	-29	HIS	-	expression tag	UNP A0ACI9
B	-28	HIS	-	expression tag	UNP A0ACI9
B	-27	SER	-	expression tag	UNP A0ACI9
B	-26	SER	-	expression tag	UNP A0ACI9
B	-25	GLY	-	expression tag	UNP A0ACI9
B	-24	LEU	-	expression tag	UNP A0ACI9
B	-23	VAL	-	expression tag	UNP A0ACI9
B	-22	PRO	-	expression tag	UNP A0ACI9
B	-21	ARG	-	expression tag	UNP A0ACI9
B	-20	GLY	-	expression tag	UNP A0ACI9
B	-19	SER	-	expression tag	UNP A0ACI9
B	-18	HIS	-	expression tag	UNP A0ACI9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	MSE	-	expression tag	UNP A0ACI9
B	-16	ALA	-	expression tag	UNP A0ACI9
B	-15	SER	-	expression tag	UNP A0ACI9
B	-14	ASP	-	expression tag	UNP A0ACI9
B	-13	SER	-	expression tag	UNP A0ACI9
B	-12	THR	-	expression tag	UNP A0ACI9
B	-11	GLU	-	expression tag	UNP A0ACI9
B	-10	ASN	-	expression tag	UNP A0ACI9
B	-9	LEU	-	expression tag	UNP A0ACI9
B	-8	TYR	-	expression tag	UNP A0ACI9
B	-7	PHE	-	expression tag	UNP A0ACI9
B	-6	GLN	-	expression tag	UNP A0ACI9
B	-5	GLY	-	expression tag	UNP A0ACI9
B	-4	ILE	-	expression tag	UNP A0ACI9
B	-3	ASP	-	expression tag	UNP A0ACI9
B	-2	PRO	-	expression tag	UNP A0ACI9
B	-1	PHE	-	expression tag	UNP A0ACI9
B	0	THR	-	expression tag	UNP A0ACI9
B	1	MSE	-	expression tag	UNP A0ACI9
C	-37	MSE	-	initiating methionine	UNP A0ACI9
C	-36	GLY	-	expression tag	UNP A0ACI9
C	-35	SER	-	expression tag	UNP A0ACI9
C	-34	SER	-	expression tag	UNP A0ACI9
C	-33	HIS	-	expression tag	UNP A0ACI9
C	-32	HIS	-	expression tag	UNP A0ACI9
C	-31	HIS	-	expression tag	UNP A0ACI9
C	-30	HIS	-	expression tag	UNP A0ACI9
C	-29	HIS	-	expression tag	UNP A0ACI9
C	-28	HIS	-	expression tag	UNP A0ACI9
C	-27	SER	-	expression tag	UNP A0ACI9
C	-26	SER	-	expression tag	UNP A0ACI9
C	-25	GLY	-	expression tag	UNP A0ACI9
C	-24	LEU	-	expression tag	UNP A0ACI9
C	-23	VAL	-	expression tag	UNP A0ACI9
C	-22	PRO	-	expression tag	UNP A0ACI9
C	-21	ARG	-	expression tag	UNP A0ACI9
C	-20	GLY	-	expression tag	UNP A0ACI9
C	-19	SER	-	expression tag	UNP A0ACI9
C	-18	HIS	-	expression tag	UNP A0ACI9
C	-17	MSE	-	expression tag	UNP A0ACI9
C	-16	ALA	-	expression tag	UNP A0ACI9
C	-15	SER	-	expression tag	UNP A0ACI9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-14	ASP	-	expression tag	UNP A0ACI9
C	-13	SER	-	expression tag	UNP A0ACI9
C	-12	THR	-	expression tag	UNP A0ACI9
C	-11	GLU	-	expression tag	UNP A0ACI9
C	-10	ASN	-	expression tag	UNP A0ACI9
C	-9	LEU	-	expression tag	UNP A0ACI9
C	-8	TYR	-	expression tag	UNP A0ACI9
C	-7	PHE	-	expression tag	UNP A0ACI9
C	-6	GLN	-	expression tag	UNP A0ACI9
C	-5	GLY	-	expression tag	UNP A0ACI9
C	-4	ILE	-	expression tag	UNP A0ACI9
C	-3	ASP	-	expression tag	UNP A0ACI9
C	-2	PRO	-	expression tag	UNP A0ACI9
C	-1	PHE	-	expression tag	UNP A0ACI9
C	0	THR	-	expression tag	UNP A0ACI9
C	1	MSE	-	expression tag	UNP A0ACI9
D	-37	MSE	-	initiating methionine	UNP A0ACI9
D	-36	GLY	-	expression tag	UNP A0ACI9
D	-35	SER	-	expression tag	UNP A0ACI9
D	-34	SER	-	expression tag	UNP A0ACI9
D	-33	HIS	-	expression tag	UNP A0ACI9
D	-32	HIS	-	expression tag	UNP A0ACI9
D	-31	HIS	-	expression tag	UNP A0ACI9
D	-30	HIS	-	expression tag	UNP A0ACI9
D	-29	HIS	-	expression tag	UNP A0ACI9
D	-28	HIS	-	expression tag	UNP A0ACI9
D	-27	SER	-	expression tag	UNP A0ACI9
D	-26	SER	-	expression tag	UNP A0ACI9
D	-25	GLY	-	expression tag	UNP A0ACI9
D	-24	LEU	-	expression tag	UNP A0ACI9
D	-23	VAL	-	expression tag	UNP A0ACI9
D	-22	PRO	-	expression tag	UNP A0ACI9
D	-21	ARG	-	expression tag	UNP A0ACI9
D	-20	GLY	-	expression tag	UNP A0ACI9
D	-19	SER	-	expression tag	UNP A0ACI9
D	-18	HIS	-	expression tag	UNP A0ACI9
D	-17	MSE	-	expression tag	UNP A0ACI9
D	-16	ALA	-	expression tag	UNP A0ACI9
D	-15	SER	-	expression tag	UNP A0ACI9
D	-14	ASP	-	expression tag	UNP A0ACI9
D	-13	SER	-	expression tag	UNP A0ACI9
D	-12	THR	-	expression tag	UNP A0ACI9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-11	GLU	-	expression tag	UNP A0ACI9
D	-10	ASN	-	expression tag	UNP A0ACI9
D	-9	LEU	-	expression tag	UNP A0ACI9
D	-8	TYR	-	expression tag	UNP A0ACI9
D	-7	PHE	-	expression tag	UNP A0ACI9
D	-6	GLN	-	expression tag	UNP A0ACI9
D	-5	GLY	-	expression tag	UNP A0ACI9
D	-4	ILE	-	expression tag	UNP A0ACI9
D	-3	ASP	-	expression tag	UNP A0ACI9
D	-2	PRO	-	expression tag	UNP A0ACI9
D	-1	PHE	-	expression tag	UNP A0ACI9
D	0	THR	-	expression tag	UNP A0ACI9
D	1	MSE	-	expression tag	UNP A0ACI9
E	-37	MSE	-	initiating methionine	UNP A0ACI9
E	-36	GLY	-	expression tag	UNP A0ACI9
E	-35	SER	-	expression tag	UNP A0ACI9
E	-34	SER	-	expression tag	UNP A0ACI9
E	-33	HIS	-	expression tag	UNP A0ACI9
E	-32	HIS	-	expression tag	UNP A0ACI9
E	-31	HIS	-	expression tag	UNP A0ACI9
E	-30	HIS	-	expression tag	UNP A0ACI9
E	-29	HIS	-	expression tag	UNP A0ACI9
E	-28	HIS	-	expression tag	UNP A0ACI9
E	-27	SER	-	expression tag	UNP A0ACI9
E	-26	SER	-	expression tag	UNP A0ACI9
E	-25	GLY	-	expression tag	UNP A0ACI9
E	-24	LEU	-	expression tag	UNP A0ACI9
E	-23	VAL	-	expression tag	UNP A0ACI9
E	-22	PRO	-	expression tag	UNP A0ACI9
E	-21	ARG	-	expression tag	UNP A0ACI9
E	-20	GLY	-	expression tag	UNP A0ACI9
E	-19	SER	-	expression tag	UNP A0ACI9
E	-18	HIS	-	expression tag	UNP A0ACI9
E	-17	MSE	-	expression tag	UNP A0ACI9
E	-16	ALA	-	expression tag	UNP A0ACI9
E	-15	SER	-	expression tag	UNP A0ACI9
E	-14	ASP	-	expression tag	UNP A0ACI9
E	-13	SER	-	expression tag	UNP A0ACI9
E	-12	THR	-	expression tag	UNP A0ACI9
E	-11	GLU	-	expression tag	UNP A0ACI9
E	-10	ASN	-	expression tag	UNP A0ACI9
E	-9	LEU	-	expression tag	UNP A0ACI9

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-8	TYR	-	expression tag	UNP A0ACI9
E	-7	PHE	-	expression tag	UNP A0ACI9
E	-6	GLN	-	expression tag	UNP A0ACI9
E	-5	GLY	-	expression tag	UNP A0ACI9
E	-4	ILE	-	expression tag	UNP A0ACI9
E	-3	ASP	-	expression tag	UNP A0ACI9
E	-2	PRO	-	expression tag	UNP A0ACI9
E	-1	PHE	-	expression tag	UNP A0ACI9
E	0	THR	-	expression tag	UNP A0ACI9
E	1	MSE	-	expression tag	UNP A0ACI9
F	-37	MSE	-	initiating methionine	UNP A0ACI9
F	-36	GLY	-	expression tag	UNP A0ACI9
F	-35	SER	-	expression tag	UNP A0ACI9
F	-34	SER	-	expression tag	UNP A0ACI9
F	-33	HIS	-	expression tag	UNP A0ACI9
F	-32	HIS	-	expression tag	UNP A0ACI9
F	-31	HIS	-	expression tag	UNP A0ACI9
F	-30	HIS	-	expression tag	UNP A0ACI9
F	-29	HIS	-	expression tag	UNP A0ACI9
F	-28	HIS	-	expression tag	UNP A0ACI9
F	-27	SER	-	expression tag	UNP A0ACI9
F	-26	SER	-	expression tag	UNP A0ACI9
F	-25	GLY	-	expression tag	UNP A0ACI9
F	-24	LEU	-	expression tag	UNP A0ACI9
F	-23	VAL	-	expression tag	UNP A0ACI9
F	-22	PRO	-	expression tag	UNP A0ACI9
F	-21	ARG	-	expression tag	UNP A0ACI9
F	-20	GLY	-	expression tag	UNP A0ACI9
F	-19	SER	-	expression tag	UNP A0ACI9
F	-18	HIS	-	expression tag	UNP A0ACI9
F	-17	MSE	-	expression tag	UNP A0ACI9
F	-16	ALA	-	expression tag	UNP A0ACI9
F	-15	SER	-	expression tag	UNP A0ACI9
F	-14	ASP	-	expression tag	UNP A0ACI9
F	-13	SER	-	expression tag	UNP A0ACI9
F	-12	THR	-	expression tag	UNP A0ACI9
F	-11	GLU	-	expression tag	UNP A0ACI9
F	-10	ASN	-	expression tag	UNP A0ACI9
F	-9	LEU	-	expression tag	UNP A0ACI9
F	-8	TYR	-	expression tag	UNP A0ACI9
F	-7	PHE	-	expression tag	UNP A0ACI9
F	-6	GLN	-	expression tag	UNP A0ACI9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-5	GLY	-	expression tag	UNP A0ACI9
F	-4	ILE	-	expression tag	UNP A0ACI9
F	-3	ASP	-	expression tag	UNP A0ACI9
F	-2	PRO	-	expression tag	UNP A0ACI9
F	-1	PHE	-	expression tag	UNP A0ACI9
F	0	THR	-	expression tag	UNP A0ACI9
F	1	MSE	-	expression tag	UNP A0ACI9

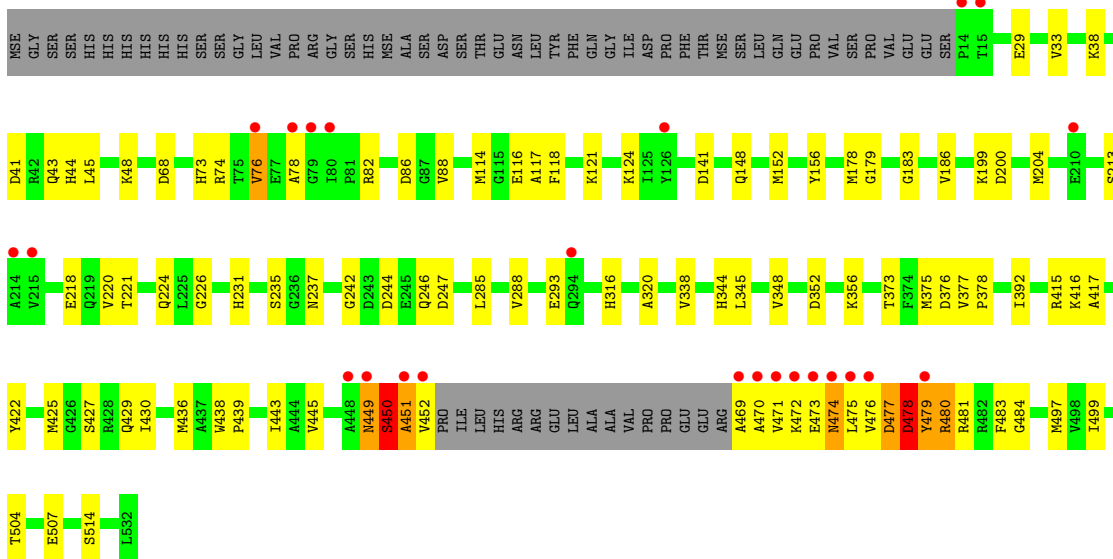
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	166	Total O 166 166	0	0
2	B	214	Total O 214 214	0	0
2	C	156	Total O 156 156	0	0
2	D	152	Total O 152 152	0	0
2	E	136	Total O 136 136	0	0
2	F	112	Total O 112 112	0	0

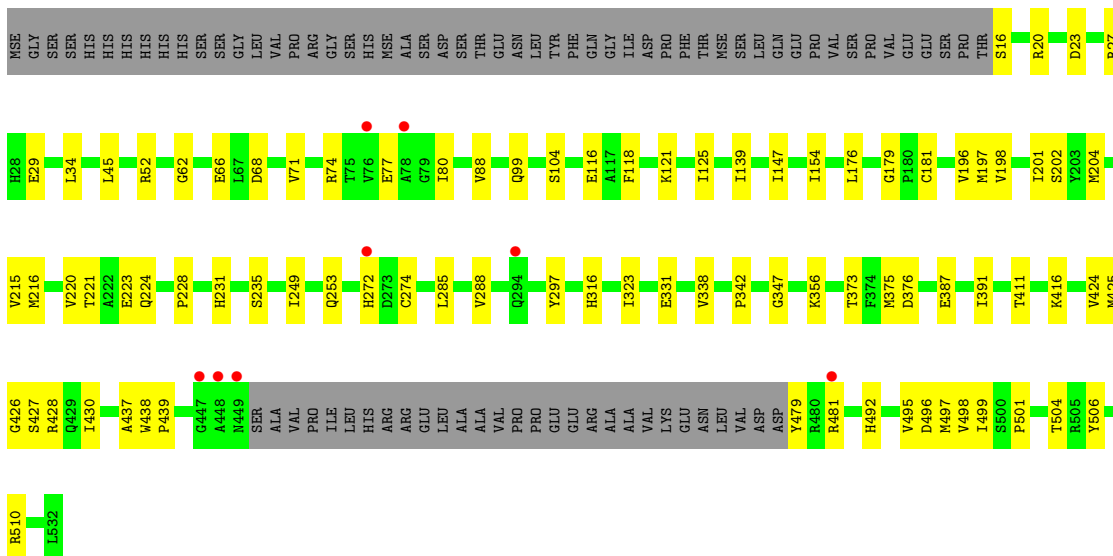




- Molecule 1: Putative carboxyl transferase

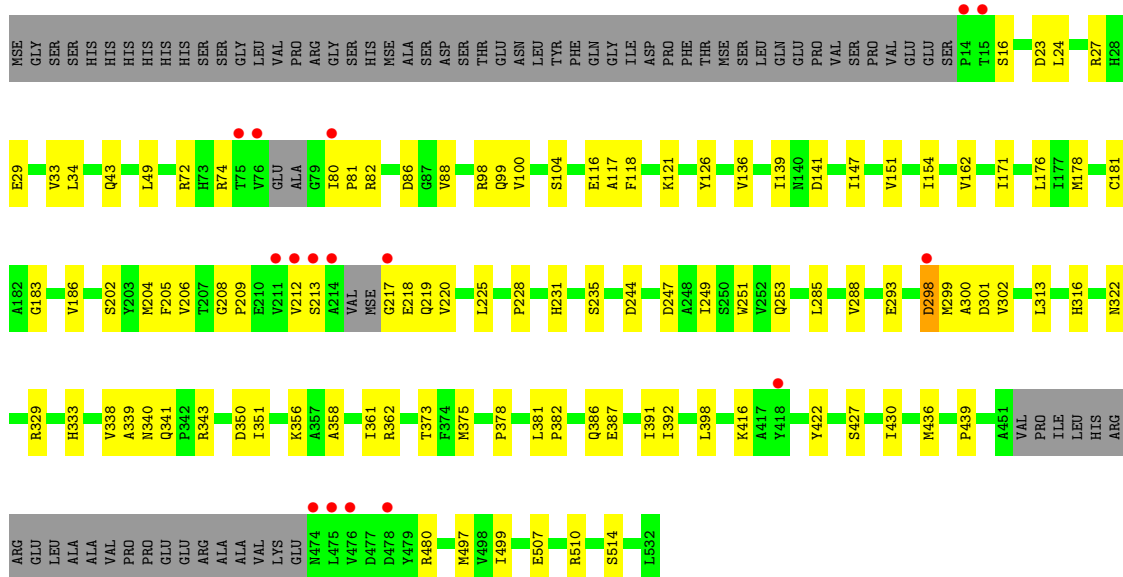


- Molecule 1: Putative carboxyl transferase

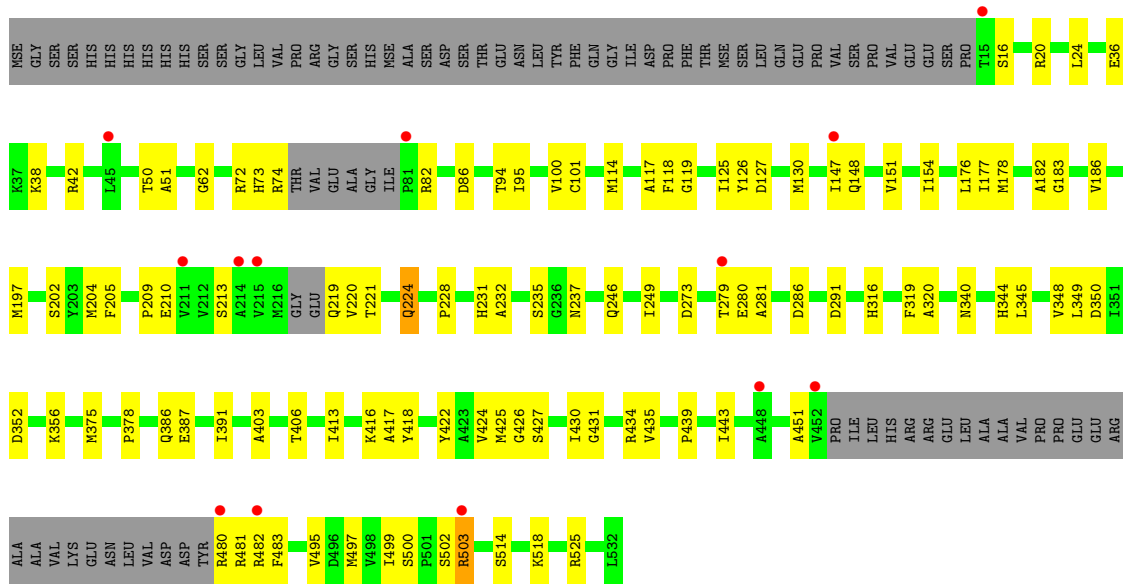


- Molecule 1: Putative carboxyl transferase





● Molecule 1: Putative carboxyl transferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.56Å 163.44Å 186.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	81.72 – 2.45 81.72 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.7 (81.72-2.45) 90.2 (81.72-2.45)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 2.45Å)	Xtrriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, $R_{free}$	0.191 , 0.230 0.191 , 0.229	Depositor DCC
$R_{free}$ test set	2000 reflections (1.48%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.2	Xtrriage
Anisotropy	0.228	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 46.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	23360	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality i

### 5.1 Standard geometry i

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.30	0/3826	0.62	4/5179 (0.1%)
1	B	0.30	0/3829	0.57	0/5184
1	C	0.36	0/3872	0.79	14/5242 (0.3%)
1	D	0.29	0/3765	0.57	0/5095
1	E	0.38	0/3800	0.64	0/5143
1	F	0.37	1/3722 (0.0%)	0.67	2/5033 (0.0%)
All	All	0.34	1/22814 (0.0%)	0.65	20/30876 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	C	0	1
1	E	0	2
All	All	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	503	ARG	CG-CD	5.34	1.68	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	450	SER	CB-CA-C	-18.90	84.87	111.14
1	C	78	ALA	N-CA-C	11.42	125.07	111.82
1	C	450	SER	N-CA-C	-10.86	97.49	112.25
1	F	503	ARG	CG-CD-NE	7.77	129.10	112.00
1	C	78	ALA	CB-CA-C	-7.10	97.04	110.67
1	A	34	LEU	O-C-N	6.96	130.08	122.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	383	GLY	N-CA-C	6.64	128.92	113.18
1	C	449	ASN	CB-CA-C	-6.38	98.99	110.35
1	C	76	VAL	CA-C-N	-6.02	111.17	122.73
1	C	76	VAL	C-N-CA	-6.02	111.17	122.73
1	F	147	ILE	CG1-CB-CG2	5.91	128.44	110.70
1	C	451	ALA	N-CA-C	-5.78	98.06	108.58
1	C	480	ARG	NE-CZ-NH1	-5.64	115.86	121.50
1	C	478	ASP	CA-C-N	5.55	131.52	121.92
1	C	478	ASP	C-N-CA	5.55	131.52	121.92
1	C	479	TYR	CA-CB-CG	5.38	123.58	113.90
1	C	450	SER	CA-C-N	5.32	130.99	122.32
1	C	450	SER	C-N-CA	5.32	130.99	122.32
1	A	33	VAL	CA-C-N	5.00	127.39	120.29
1	A	33	VAL	C-N-CA	5.00	127.39	120.29

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	198	VAL	Peptide
1	A	452	VAL	Peptide
1	C	450	SER	Peptide
1	E	217	GLY	Peptide
1	E	298	ASP	Peptide

## 5.2 Too-close contacts [\(i\)](#)

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	3760	0	3697	64	0
1	B	3763	0	3703	67	0
1	C	3806	0	3747	119	1
1	D	3700	0	3641	74	0
1	E	3735	0	3669	85	0
1	F	3660	0	3607	105	0
2	A	166	0	0	7	1
2	B	214	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	156	0	0	6	0
2	D	152	0	0	3	0
2	E	136	0	0	1	0
2	F	112	0	0	6	0
All	All	23360	0	22064	486	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 11.

All (486) 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:293:GLU:CG	1:C:480:ARG:NH2	1.97	1.26
1:B:385:GLY:O	1:B:389:GLN:HG3	1.38	1.24
1:C:293:GLU:HG3	1:C:480:ARG:NH2	1.48	1.23
1:F:500:SER:N	1:F:503:ARG:HH21	1.47	1.10
1:F:482:ARG:H	1:F:482:ARG:HD3	1.20	1.06
1:C:472:LYS:HZ3	1:C:475:LEU:HD23	1.16	1.05
1:F:500:SER:H	1:F:503:ARG:NH2	1.54	1.04
1:C:293:GLU:HG2	1:C:480:ARG:NH2	1.71	1.03
1:C:293:GLU:HG2	1:C:480:ARG:CZ	1.88	1.03
1:F:221:THR:HG22	1:F:224:GLN:HG2	1.40	1.02
1:B:385:GLY:O	1:B:389:GLN:CG	2.07	1.02
1:C:114:MSE:HE3	1:C:156:TYR:HB3	1.38	1.01
1:D:221:THR:HG22	1:D:224:GLN:HG3	1.39	0.99
1:E:298:ASP:HB2	1:E:343:ARG:CZ	1.92	0.98
1:C:293:GLU:HG3	1:C:480:ARG:HH22	1.26	0.97
1:B:445:VAL:HA	1:F:154:ILE:HD11	1.46	0.94
1:C:472:LYS:NZ	1:C:475:LEU:HD23	1.83	0.93
1:C:293:GLU:CG	1:C:480:ARG:HH22	1.74	0.90
1:A:445:VAL:HA	1:D:154:ILE:HD11	1.52	0.90
1:C:199:LYS:NZ	2:C:601:HOH:O	2.04	0.90
1:E:213:SER:HB2	1:E:219:GLN:HA	1.54	0.90
1:F:500:SER:H	1:F:503:ARG:HH21	0.94	0.89
1:B:434:ARG:HD2	1:B:497:MSE:HE3	1.51	0.89
1:F:221:THR:N	2:F:601:HOH:O	2.08	0.86
1:F:434:ARG:HD2	1:F:497:MSE:HE3	1.57	0.86
1:D:438:TRP:CZ2	1:D:504:THR:HG21	2.11	0.85
1:D:438:TRP:HZ2	1:D:504:THR:HG21	1.41	0.85
1:A:299:MSE:HE3	1:A:374:PHE:HB3	1.58	0.84
1:E:375:MSE:HE1	1:E:422:TYR:HD1	1.42	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:MSE:HE1	1:B:414:THR:HG21	1.61	0.82
1:E:88:VAL:O	1:E:121:LYS:NZ	2.12	0.82
1:C:451:ALA:O	1:C:452:VAL:HG13	1.81	0.81
1:F:127:ASP:HA	1:F:130:MSE:HE2	1.61	0.81
1:C:375:MSE:HE1	1:C:422:TYR:HA	1.62	0.80
1:C:293:GLU:N	1:C:293:GLU:OE2	2.16	0.79
1:D:479:TYR:N	2:D:601:HOH:O	2.16	0.78
1:E:298:ASP:OD1	1:E:341:GLN:NE2	2.15	0.78
1:A:139:ILE:HG23	1:A:178:MSE:HE3	1.65	0.77
1:C:293:GLU:O	1:C:415:ARG:NH2	2.17	0.77
1:E:33:VAL:HG13	1:E:34:LEU:HD12	1.67	0.77
1:A:375:MSE:HE1	1:A:422:TYR:HD1	1.49	0.77
1:C:471:VAL:HG13	1:C:472:LYS:H	1.50	0.77
1:C:375:MSE:HE2	1:C:417:ALA:HA	1.66	0.77
1:B:114:MSE:HE1	1:B:119:GLY:HA3	1.66	0.77
1:C:116:GLU:OE1	1:D:74:ARG:NH2	2.18	0.76
1:C:221:THR:HG23	1:C:224:GLN:H	1.51	0.75
1:B:20:ARG:NH2	1:C:484:GLY:O	2.20	0.75
1:C:473:GLU:N	1:C:473:GLU:OE2	2.19	0.74
1:C:199:LYS:HZ2	1:C:200:ASP:H	1.35	0.74
1:A:475:LEU:N	2:A:603:HOH:O	2.21	0.74
1:C:477:ASP:O	1:C:479:TYR:N	2.21	0.73
1:F:500:SER:H	1:F:503:ARG:CZ	2.02	0.73
1:B:147:ILE:HD11	1:F:451:ALA:HB1	1.68	0.73
1:A:434:ARG:HD2	1:A:497:MSE:HE3	1.71	0.72
1:B:114:MSE:HE2	1:B:156:TYR:HB3	1.71	0.72
1:B:114:MSE:HE1	1:B:119:GLY:CA	2.20	0.72
1:A:311:ASP:OD1	2:A:601:HOH:O	2.08	0.71
1:C:74:ARG:NH2	1:D:116:GLU:OE2	2.20	0.70
1:D:220:VAL:HG23	1:D:224:GLN:HB2	1.73	0.70
1:F:500:SER:N	1:F:503:ARG:NH2	2.25	0.69
1:C:478:ASP:HB3	1:C:481:ARG:HD2	1.74	0.69
1:F:482:ARG:HD3	1:F:482:ARG:N	2.03	0.69
1:C:204:MSE:HE2	1:E:392:ILE:HD13	1.75	0.69
1:F:417:ALA:HB3	1:F:443:ILE:HA	1.74	0.68
1:C:449:ASN:C	1:C:449:ASN:OD1	2.36	0.68
1:D:481:ARG:H	1:D:481:ARG:HD2	1.58	0.68
1:E:80:ILE:HD12	1:E:81:PRO:HD2	1.75	0.68
1:B:52:ARG:NH2	1:B:66:GLU:OE2	2.27	0.68
1:C:116:GLU:HB3	1:C:152:MSE:CE	2.23	0.67
1:E:378:PRO:HB3	1:E:416:LYS:HD2	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:471:VAL:HG13	1:C:472:LYS:N	2.09	0.67
1:C:445:VAL:HA	1:E:154:ILE:HD11	1.77	0.67
1:C:476:VAL:C	1:C:478:ASP:H	2.03	0.67
1:E:208:GLY:O	1:E:212:VAL:HG23	1.95	0.67
1:C:246:GLN:OE1	2:C:602:HOH:O	2.13	0.66
1:F:499:ILE:HA	1:F:503:ARG:NH2	2.09	0.66
1:D:215:VAL:HG12	1:D:216:MSE:HE3	1.78	0.66
1:D:272:HIS:NE2	1:D:274:CYS:HB2	2.10	0.66
1:C:293:GLU:HG2	1:C:480:ARG:NH1	2.10	0.66
1:B:114:MSE:HE2	1:B:156:TYR:CB	2.26	0.65
1:B:503:ARG:NH2	1:B:507:GLU:OE1	2.28	0.65
1:F:480:ARG:N	2:F:605:HOH:O	2.29	0.65
1:A:382:PRO:O	2:A:602:HOH:O	2.13	0.65
1:E:316:HIS:CB	1:E:356:LYS:HE3	2.26	0.65
1:C:476:VAL:O	1:C:478:ASP:N	2.29	0.65
1:A:375:MSE:HE1	1:A:422:TYR:CD1	2.31	0.64
1:B:130:MSE:HE1	1:B:163:ARG:HG2	1.79	0.64
1:C:199:LYS:NZ	1:C:200:ASP:H	1.94	0.64
1:A:323:ILE:O	1:A:356:LYS:NZ	2.30	0.64
1:C:436:MSE:HE3	1:C:507:GLU:HB2	1.77	0.64
1:C:375:MSE:CE	1:C:422:TYR:HA	2.27	0.64
1:E:298:ASP:OD2	1:E:300:ALA:HB3	1.97	0.64
1:B:405:SER:OG	1:B:409:LYS:HE3	1.98	0.64
1:E:24:LEU:HD21	1:F:503:ARG:NH2	2.13	0.63
1:F:114:MSE:HE3	1:F:118:PHE:CD1	2.34	0.63
1:E:375:MSE:HE1	1:E:422:TYR:CD1	2.30	0.62
1:C:114:MSE:HE3	1:C:156:TYR:CB	2.25	0.61
1:D:285:LEU:HA	1:D:288:VAL:HG13	1.81	0.61
1:D:373:THR:HB	1:D:411:THR:HG23	1.83	0.61
1:F:406:THR:O	1:F:518:LYS:HE2	1.99	0.61
1:F:126:TYR:O	1:F:130:MSE:HG3	2.01	0.61
1:D:272:HIS:CE1	1:D:331:GLU:HA	2.36	0.61
1:C:429:GLN:OE1	2:C:603:HOH:O	2.16	0.61
1:A:15:THR:HG22	1:B:481:ARG:HH22	1.65	0.61
1:A:197:MSE:HE3	1:A:238:ALA:HB2	1.82	0.60
1:F:480:ARG:HG2	1:F:482:ARG:NH1	2.16	0.60
1:F:114:MSE:HE3	1:F:118:PHE:HD1	1.65	0.60
1:C:148:GLN:OE1	1:C:148:GLN:N	2.29	0.60
1:C:472:LYS:HZ3	1:C:475:LEU:CD2	2.04	0.60
1:E:126:TYR:CE2	1:E:136:VAL:HG11	2.37	0.60
1:F:72:ARG:HB2	1:F:82:ARG:HH21	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:36:GLU:OE2	1:F:50:THR:HG21	2.01	0.60
1:F:500:SER:H	1:F:503:ARG:NE	1.99	0.60
1:C:183:GLY:O	1:C:186:VAL:HG22	2.02	0.59
1:F:316:HIS:HB2	1:F:356:LYS:HE3	1.83	0.59
1:C:293:GLU:CG	1:C:480:ARG:CZ	2.61	0.59
1:E:350:ASP:HB2	1:E:386:GLN:HE22	1.67	0.59
1:D:387:GLU:HA	1:D:391:ILE:HG22	1.85	0.59
1:F:316:HIS:CB	1:F:356:LYS:HE3	2.32	0.59
1:C:76:VAL:HA	1:C:82:ARG:HD2	1.84	0.58
1:F:500:SER:H	1:F:503:ARG:HE	1.50	0.58
1:A:375:MSE:HE2	1:A:413:ILE:HG12	1.85	0.58
1:E:316:HIS:HB2	1:E:356:LYS:HE3	1.84	0.58
1:B:16:SER:O	1:B:20:ARG:HG3	2.03	0.58
1:B:148:GLN:H	1:B:148:GLN:CD	2.07	0.58
1:B:159:GLU:O	1:B:162:VAL:HG12	2.02	0.58
1:A:385:GLY:O	1:A:389:GLN:HG3	2.03	0.58
1:C:475:LEU:HG	1:C:475:LEU:O	2.03	0.58
1:D:438:TRP:HE1	1:D:504:THR:CG2	2.16	0.58
1:E:298:ASP:CG	1:E:300:ALA:H	2.11	0.58
1:F:178:MSE:HE1	1:F:249:ILE:HG13	1.86	0.58
1:C:436:MSE:HG2	1:C:497:MSE:HE2	1.85	0.57
1:C:470:ALA:HA	1:C:473:GLU:OE1	2.03	0.57
1:C:213:SER:OG	1:C:218:GLU:O	2.15	0.57
1:D:52:ARG:HH22	1:D:66:GLU:CD	2.13	0.57
1:E:299:MSE:HE3	1:E:302:VAL:HB	1.86	0.56
1:F:246:GLN:CD	1:F:246:GLN:H	2.14	0.56
1:C:470:ALA:HA	1:C:473:GLU:CD	2.30	0.56
1:E:212:VAL:HG21	1:E:225:LEU:HD21	1.87	0.56
1:C:473:GLU:HG2	1:C:474:ASN:H	1.71	0.56
1:A:176:LEU:HB3	1:A:178:MSE:HE2	1.88	0.55
1:A:99:GLN:OE1	1:B:514:SER:HB2	2.05	0.55
1:B:197:MSE:SE	1:B:204:MSE:HE2	2.56	0.55
1:E:350:ASP:HB2	1:E:386:GLN:NE2	2.21	0.55
1:A:285:LEU:O	1:A:288:VAL:HG12	2.07	0.55
1:D:376:ASP:OD1	1:D:416:LYS:HE2	2.06	0.55
1:D:221:THR:HG23	1:D:223:GLU:H	1.70	0.55
1:B:30:GLU:HA	1:B:34:LEU:HD12	1.89	0.55
1:B:448:ALA:HA	1:B:479:TYR:CD2	2.42	0.55
1:B:506:TYR:CZ	1:B:510:ARG:HD2	2.42	0.55
1:E:139:ILE:HG12	1:E:176:LEU:HD12	1.87	0.55
1:E:316:HIS:HB3	1:E:356:LYS:HE3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:381:LEU:HD12	1:E:382:PRO:HD2	1.88	0.54
1:A:146:ARG:HD2	1:A:149:GLU:OE1	2.07	0.54
1:C:116:GLU:HB3	1:C:152:MSE:HE3	1.89	0.54
1:F:151:VAL:HA	1:F:154:ILE:HD12	1.89	0.54
1:B:114:MSE:HE3	1:B:115:GLY:O	2.07	0.54
1:E:99:GLN:OE1	1:F:514:SER:HB2	2.07	0.54
1:F:481:ARG:H	1:F:482:ARG:NH2	2.06	0.54
1:F:482:ARG:H	1:F:482:ARG:CD	2.06	0.54
1:A:23:ASP:OD2	1:A:27:ARG:NH1	2.40	0.54
1:A:197:MSE:HE2	1:A:204:MSE:SE	2.57	0.54
1:E:351:ILE:H	1:E:386:GLN:HE22	1.54	0.53
1:F:500:SER:N	1:F:503:ARG:HE	2.06	0.53
1:B:345:LEU:O	1:B:348:VAL:HG22	2.08	0.53
1:A:38:LYS:HD2	1:A:38:LYS:C	2.33	0.53
1:D:297:TYR:OH	1:D:376:ASP:OD2	2.21	0.53
1:D:510:ARG:NH2	1:F:62:GLY:O	2.42	0.53
1:E:298:ASP:CB	1:E:343:ARG:CZ	2.78	0.53
1:C:73:HIS:CE1	1:C:82:ARG:HG2	2.43	0.53
1:C:88:VAL:HB	1:C:118:PHE:CE2	2.44	0.53
1:D:428:ARG:HH21	1:D:496:ASP:CG	2.16	0.53
1:B:127:ASP:HA	1:B:130:MSE:HE3	1.91	0.53
1:E:293:GLU:OE1	1:E:480:ARG:NH1	2.42	0.53
1:F:197:MSE:HE3	1:F:204:MSE:HG3	1.91	0.53
1:F:340:ASN:HD21	1:F:425:MSE:HE2	1.74	0.53
1:F:16:SER:O	1:F:20:ARG:HG3	2.07	0.53
1:A:299:MSE:HE3	1:A:374:PHE:CB	2.37	0.53
1:F:340:ASN:ND2	1:F:425:MSE:HE2	2.24	0.53
1:C:41:ASP:O	1:C:45:LEU:HD13	2.09	0.52
1:C:473:GLU:HG2	1:C:474:ASN:N	2.25	0.52
1:C:285:LEU:O	1:C:288:VAL:HG12	2.09	0.52
1:B:43:GLN:HG2	1:B:49:LEU:O	2.09	0.52
1:C:479:TYR:CE2	1:C:483:PHE:CD2	2.97	0.52
1:F:319:PHE:O	1:F:356:LYS:NZ	2.30	0.52
1:A:23:ASP:O	1:A:27:ARG:HG3	2.09	0.52
1:D:197:MSE:SE	1:D:204:MSE:HE2	2.60	0.52
1:F:197:MSE:SE	1:F:204:MSE:HE2	2.59	0.52
1:F:375:MSE:HE1	1:F:422:TYR:HD1	1.74	0.52
1:C:345:LEU:O	1:C:348:VAL:HG22	2.10	0.52
1:C:375:MSE:HE3	1:C:377:VAL:HG12	1.91	0.52
1:A:43:GLN:NE2	1:A:49:LEU:O	2.39	0.52
1:E:299:MSE:CE	1:E:302:VAL:HB	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:293:GLU:HG3	1:C:480:ARG:HH21	1.62	0.52
1:D:481:ARG:HD2	1:D:481:ARG:N	2.24	0.51
1:E:375:MSE:SE	1:E:422:TYR:HA	2.61	0.51
1:E:507:GLU:OE1	1:E:507:GLU:HA	2.10	0.51
1:F:435:VAL:HG12	1:F:495:VAL:HG12	1.92	0.51
1:B:406:THR:O	1:B:518:LYS:HE2	2.10	0.51
1:C:316:HIS:HB2	1:C:356:LYS:HE2	1.92	0.51
1:F:350:ASP:HB2	1:F:386:GLN:HE22	1.75	0.51
1:C:29:GLU:CD	1:C:33:VAL:HG21	2.36	0.51
1:F:220:VAL:CG1	1:F:224:GLN:HG3	2.41	0.51
1:F:220:VAL:HG13	1:F:224:GLN:HG3	1.93	0.51
1:B:52:ARG:HG3	2:B:705:HOH:O	2.11	0.51
1:E:88:VAL:HB	1:E:118:PHE:CE2	2.46	0.51
1:E:139:ILE:HG23	1:E:178:MSE:SE	2.61	0.51
1:A:88:VAL:HB	1:A:118:PHE:CE2	2.46	0.51
1:C:320:ALA:HB2	1:C:352:ASP:HB3	1.92	0.51
1:E:298:ASP:HB2	1:E:343:ARG:NH2	2.24	0.51
1:F:101:CYS:HB3	1:F:125:ILE:HG23	1.93	0.51
1:B:385:GLY:O	1:B:389:GLN:HG2	2.04	0.50
1:C:473:GLU:C	1:C:475:LEU:H	2.20	0.50
1:A:152:MSE:HE3	1:D:492:HIS:CE1	2.46	0.50
1:E:24:LEU:HD21	1:F:503:ARG:HH22	1.76	0.50
1:E:33:VAL:CG1	1:E:34:LEU:HD12	2.41	0.50
1:A:117:ALA:O	1:A:121:LYS:HG3	2.12	0.50
1:B:433:ASP:OD1	1:B:518:LYS:NZ	2.36	0.50
1:D:316:HIS:HB3	1:D:356:LYS:HE3	1.93	0.50
1:D:68:ASP:CB	1:D:121:LYS:HD2	2.42	0.50
1:D:176:LEU:HD23	1:D:196:VAL:HB	1.93	0.50
1:D:323:ILE:O	1:D:356:LYS:NZ	2.44	0.50
1:C:375:MSE:HE1	1:C:422:TYR:CA	2.36	0.50
1:D:506:TYR:CE2	1:D:510:ARG:HD2	2.47	0.50
1:C:116:GLU:HB3	1:C:152:MSE:HE1	1.91	0.50
1:C:231:HIS:CD2	1:C:237:ASN:HD22	2.30	0.50
1:D:499:ILE:HG22	2:D:603:HOH:O	2.12	0.50
1:C:86:ASP:OD2	1:C:117:ALA:HB3	2.12	0.49
1:F:500:SER:HB3	1:F:503:ARG:NE	2.27	0.49
1:F:38:LYS:O	1:F:42:ARG:HG3	2.12	0.49
1:B:497:MSE:HE1	2:B:611:HOH:O	2.12	0.49
1:D:427:SER:O	1:D:430:ILE:HG22	2.12	0.49
1:A:116:GLU:OE2	1:F:74:ARG:NH2	2.43	0.49
1:D:481:ARG:NH2	2:D:601:HOH:O	2.21	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:136:VAL:HG23	1:E:171:ILE:HD12	1.95	0.49
1:E:183:GLY:O	1:E:186:VAL:HG22	2.13	0.49
1:C:477:ASP:OD1	1:C:477:ASP:C	2.56	0.49
1:E:220:VAL:HG21	1:E:225:LEU:HD13	1.95	0.49
1:B:43:GLN:HG3	1:B:48:LYS:HB2	1.93	0.49
1:F:427:SER:O	1:F:430:ILE:HG22	2.12	0.49
1:D:99:GLN:OE1	1:E:514:SER:HB2	2.13	0.49
1:D:16:SER:O	1:D:20:ARG:HG3	2.13	0.48
1:F:480:ARG:HD2	1:F:483:PHE:HD2	1.78	0.48
1:E:23:ASP:O	1:E:27:ARG:HG3	2.12	0.48
1:F:500:SER:HB3	1:F:503:ARG:HE	1.78	0.48
1:A:206:VAL:N	1:D:387:GLU:OE2	2.42	0.48
1:D:71:VAL:HB	1:D:121:LYS:HD3	1.95	0.48
1:D:439:PRO:HD3	1:D:499:ILE:O	2.13	0.48
1:D:342:PRO:HA	1:D:347:GLY:N	2.28	0.48
1:C:204:MSE:HB2	1:C:231:HIS:CE1	2.48	0.48
1:E:497:MSE:HG2	1:E:499:ILE:HG23	1.95	0.48
1:A:28:HIS:CE1	1:A:32:VAL:HG21	2.49	0.48
1:D:29:GLU:OE2	1:D:34:LEU:HD21	2.14	0.48
1:A:116:GLU:CD	1:F:74:ARG:HH22	2.21	0.48
1:D:125:ILE:HD12	1:D:125:ILE:HA	1.80	0.48
1:B:177:ILE:HB	1:B:197:MSE:HG2	1.96	0.47
1:C:375:MSE:CE	1:C:417:ALA:HA	2.41	0.47
1:B:178:MSE:HE1	1:B:249:ILE:HG13	1.96	0.47
1:E:43:GLN:NE2	1:E:49:LEU:O	2.44	0.47
1:E:231:HIS:HA	1:E:235:SER:OG	2.14	0.47
1:E:338:VAL:O	1:E:373:THR:HA	2.14	0.47
1:F:387:GLU:HA	1:F:391:ILE:HG22	1.96	0.47
1:B:43:GLN:HG3	1:B:48:LYS:CB	2.44	0.47
1:B:438:TRP:CZ2	1:B:501:PRO:HB3	2.49	0.47
1:D:181:CYS:O	1:D:204:MSE:HA	2.14	0.47
1:F:273:ASP:OD1	1:F:273:ASP:N	2.47	0.47
1:A:445:VAL:HG13	1:A:446:MSE:HG2	1.96	0.47
1:B:126:TYR:O	1:B:130:MSE:HG3	2.14	0.47
1:C:43:GLN:OE1	1:C:48:LYS:HE2	2.15	0.47
1:A:159:GLU:OE1	1:A:163:ARG:NH1	2.47	0.47
1:C:436:MSE:CE	1:C:504:THR:HA	2.44	0.47
1:D:68:ASP:HB2	1:D:121:LYS:HD2	1.97	0.47
1:A:126:TYR:CE2	1:A:136:VAL:HG11	2.50	0.47
1:B:506:TYR:CE2	1:B:510:ARG:HD2	2.49	0.47
1:D:495:VAL:HG12	1:D:497:MSE:H	1.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:86:ASP:OD2	1:E:117:ALA:HB3	2.14	0.47
1:F:320:ALA:HB2	1:F:352:ASP:HB3	1.96	0.47
1:A:320:ALA:HB2	1:A:352:ASP:HB3	1.96	0.47
1:B:316:HIS:HB3	1:B:356:LYS:HE3	1.97	0.47
1:C:497:MSE:HE3	1:C:499:ILE:HG21	1.97	0.47
1:F:375:MSE:HE1	1:F:422:TYR:CD1	2.50	0.47
1:A:375:MSE:HE3	1:A:417:ALA:HB2	1.97	0.47
1:B:435:VAL:HG12	1:B:495:VAL:HG12	1.96	0.47
1:B:320:ALA:HB2	1:B:352:ASP:HB3	1.96	0.46
1:F:375:MSE:HE2	1:F:413:ILE:HG23	1.97	0.46
1:A:95:ILE:HG13	1:A:100:VAL:HG21	1.97	0.46
1:A:293:GLU:HA	1:A:415:ARG:NH2	2.30	0.46
1:D:316:HIS:CB	1:D:356:LYS:HE3	2.45	0.46
1:F:73:HIS:CE1	1:F:82:ARG:HG2	2.50	0.46
1:A:220:VAL:HG22	1:A:224:GLN:HB2	1.97	0.46
1:F:386:GLN:NE2	2:F:603:HOH:O	2.26	0.46
1:C:74:ARG:HG3	1:C:116:GLU:OE1	2.15	0.46
1:F:183:GLY:O	1:F:186:VAL:HG22	2.16	0.46
1:F:279:THR:HG22	1:F:281:ALA:H	1.80	0.46
1:C:220:VAL:HG22	1:C:224:GLN:HB2	1.97	0.46
1:C:473:GLU:O	1:C:476:VAL:HG22	2.15	0.46
1:D:104:SER:HA	1:D:139:ILE:HB	1.98	0.46
1:F:481:ARG:H	1:F:482:ARG:HH21	1.62	0.46
1:B:23:ASP:O	1:B:27:ARG:HG3	2.16	0.46
1:A:249:ILE:O	1:A:253:GLN:HG3	2.16	0.46
1:C:148:GLN:H	1:C:148:GLN:CD	2.19	0.46
1:D:62:GLY:O	1:E:510:ARG:NH2	2.49	0.46
1:B:99:GLN:OE1	1:C:514:SER:HB2	2.16	0.46
1:C:439:PRO:HD3	1:C:499:ILE:O	2.15	0.46
1:D:272:HIS:NE2	1:D:331:GLU:HA	2.30	0.46
1:F:148:GLN:H	1:F:148:GLN:CD	2.13	0.46
1:C:242:GLY:HA2	1:C:247:ASP:OD2	2.16	0.46
1:C:472:LYS:HZ1	1:C:476:VAL:HG12	1.79	0.46
1:D:220:VAL:CG2	1:D:224:GLN:HB2	2.42	0.46
1:E:244:ASP:OD1	1:E:247:ASP:N	2.39	0.46
1:E:387:GLU:HA	1:E:391:ILE:HG22	1.97	0.46
1:C:392:ILE:HD13	1:E:206:VAL:HG23	1.96	0.45
1:F:286:ASP:OD1	1:F:502:SER:HB3	2.15	0.45
1:A:68:ASP:HB2	1:A:121:LYS:HE3	1.99	0.45
1:C:117:ALA:O	1:C:121:LYS:HG3	2.17	0.45
1:F:204:MSE:HB2	1:F:231:HIS:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:231:HIS:HA	1:C:235:SER:OG	2.16	0.45
1:C:376:ASP:OD1	1:C:416:LYS:HB2	2.17	0.45
1:F:177:ILE:HB	1:F:197:MSE:HG2	1.97	0.45
1:A:448:ALA:O	1:A:476:VAL:HG22	2.17	0.45
1:E:82:ARG:O	1:E:82:ARG:CG	2.65	0.45
1:B:448:ALA:HA	1:B:479:TYR:CE2	2.51	0.45
1:D:221:THR:HG23	1:D:223:GLU:N	2.31	0.45
1:E:249:ILE:O	1:E:253:GLN:HG3	2.17	0.45
1:F:86:ASP:OD2	1:F:117:ALA:HB3	2.17	0.45
1:F:439:PRO:HD3	1:F:499:ILE:O	2.17	0.45
1:B:117:ALA:O	1:B:121:LYS:HG3	2.17	0.45
1:C:44:HIS:HE1	2:C:718:HOH:O	1.99	0.45
1:F:231:HIS:HD2	1:F:235:SER:OG	2.00	0.45
1:F:431:GLY:HA3	2:F:686:HOH:O	2.16	0.45
1:C:221:THR:HG22	1:C:224:GLN:OE1	2.17	0.45
1:D:437:ALA:O	1:D:498:VAL:HA	2.16	0.45
1:E:202:SER:O	1:E:228:PRO:HD3	2.17	0.45
1:C:471:VAL:CG1	1:C:472:LYS:N	2.79	0.45
1:F:210:GLU:CD	1:F:210:GLU:H	2.25	0.45
1:F:350:ASP:HB2	1:F:386:GLN:NE2	2.32	0.45
1:A:73:HIS:CE1	1:A:82:ARG:HG2	2.52	0.45
1:B:213:SER:HB2	1:B:219:GLN:OE1	2.17	0.45
1:C:344:HIS:ND1	1:C:345:LEU:HG	2.32	0.45
1:C:471:VAL:CG1	1:C:472:LYS:H	2.25	0.45
1:C:436:MSE:HE1	1:C:504:THR:HA	1.99	0.44
1:A:183:GLY:O	1:A:186:VAL:HG22	2.17	0.44
1:A:446:MSE:SE	1:D:147:ILE:HD11	2.66	0.44
1:C:29:GLU:HA	1:C:33:VAL:HG22	1.99	0.44
1:E:285:LEU:O	1:E:288:VAL:HG12	2.17	0.44
1:C:48:LYS:NZ	2:C:610:HOH:O	2.44	0.44
1:C:48:LYS:NZ	1:C:141:ASP:OD1	2.44	0.44
1:F:344:HIS:ND1	1:F:345:LEU:HG	2.32	0.44
1:F:50:THR:HG22	1:F:51:ALA:N	2.32	0.44
1:F:424:VAL:O	1:F:426:GLY:N	2.49	0.44
1:A:199:LYS:O	1:A:228:PRO:HG2	2.18	0.44
1:C:438:TRP:N	1:C:438:TRP:CD1	2.83	0.44
1:D:338:VAL:O	1:D:373:THR:HA	2.18	0.44
1:D:45:LEU:HD23	1:D:45:LEU:HA	1.62	0.44
1:E:104:SER:HA	1:E:139:ILE:HB	2.00	0.44
1:E:251:TRP:CD1	1:E:313:LEU:HD11	2.53	0.44
1:C:476:VAL:C	1:C:478:ASP:N	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:249:ILE:O	1:D:253:GLN:HG3	2.18	0.44
1:E:329:ARG:HA	1:E:333:HIS:O	2.17	0.44
1:F:176:LEU:HD22	1:F:178:MSE:HE3	2.00	0.44
1:B:439:PRO:HD3	1:B:499:ILE:O	2.17	0.44
1:E:205:PHE:CZ	1:E:209:PRO:HD3	2.53	0.44
1:E:298:ASP:OD1	1:E:343:ARG:HD2	2.18	0.44
1:E:427:SER:O	1:E:430:ILE:HG22	2.18	0.44
1:A:329:ARG:HA	1:A:333:HIS:O	2.18	0.44
1:D:88:VAL:HB	1:D:118:PHE:CE2	2.52	0.44
1:E:208:GLY:C	1:E:212:VAL:HG23	2.42	0.44
1:E:361:ILE:HD12	1:E:398:LEU:HD11	2.00	0.44
1:F:416:LYS:HD2	1:F:418:TYR:CZ	2.53	0.44
1:B:316:HIS:HB2	1:B:356:LYS:HE2	2.00	0.43
1:C:29:GLU:OE1	1:C:33:VAL:HG21	2.18	0.43
1:C:375:MSE:HE3	1:C:377:VAL:CG1	2.48	0.43
1:D:272:HIS:CE1	1:D:274:CYS:SG	3.11	0.43
1:E:16:SER:HB2	1:F:291:ASP:OD1	2.18	0.43
1:E:181:CYS:O	1:E:204:MSE:HA	2.18	0.43
1:F:114:MSE:CE	1:F:119:GLY:HA2	2.48	0.43
1:A:16:SER:O	1:A:20:ARG:HG3	2.18	0.43
1:B:86:ASP:OD2	1:B:117:ALA:HB3	2.19	0.43
1:D:438:TRP:CE2	1:D:504:THR:HG21	2.53	0.43
1:D:272:HIS:CD2	1:D:331:GLU:HA	2.53	0.43
1:C:436:MSE:HE2	1:C:499:ILE:HD11	2.00	0.43
1:F:202:SER:O	1:F:228:PRO:HD3	2.18	0.43
1:B:162:VAL:HG23	1:F:403:ALA:HB1	1.99	0.43
1:F:279:THR:HG22	1:F:280:GLU:N	2.33	0.43
1:F:349:LEU:HD11	1:F:425:MSE:SE	2.69	0.43
1:B:285:LEU:HD21	1:B:302:VAL:HA	2.00	0.43
1:E:381:LEU:HD12	1:E:382:PRO:CD	2.49	0.43
1:A:71:VAL:HA	1:B:491:ALA:HA	2.01	0.43
1:C:38:LYS:HE3	1:C:38:LYS:HB3	1.74	0.43
1:C:378:PRO:HB3	1:C:416:LYS:HD2	1.99	0.43
1:E:147:ILE:HD12	1:E:147:ILE:H	1.83	0.43
1:E:299:MSE:HG2	1:E:339:ALA:HB1	2.01	0.43
1:F:204:MSE:O	1:F:231:HIS:HE1	2.02	0.43
1:A:72:ARG:HB2	1:A:82:ARG:HH21	1.83	0.43
1:B:74:ARG:HG3	1:B:116:GLU:OE1	2.18	0.43
1:E:72:ARG:CZ	1:E:82:ARG:HE	2.32	0.43
1:E:293:GLU:OE2	1:E:480:ARG:CZ	2.67	0.43
1:B:338:VAL:O	1:B:373:THR:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:473:GLU:C	1:C:475:LEU:N	2.76	0.43
1:D:231:HIS:HA	1:D:235:SER:OG	2.19	0.43
1:F:375:MSE:SE	1:F:422:TYR:HA	2.68	0.43
1:B:350:ASP:HB2	1:B:386:GLN:OE1	2.18	0.42
1:C:338:VAL:HG12	1:C:425:MSE:HE1	2.01	0.42
1:D:197:MSE:HE3	1:D:204:MSE:HG3	2.01	0.42
1:D:375:MSE:HE2	1:D:425:MSE:O	2.19	0.42
1:A:71:VAL:CG1	1:A:117:ALA:HB1	2.49	0.42
1:A:321:ARG:NH1	2:A:613:HOH:O	2.50	0.42
1:D:179:GLY:HA2	1:D:201:ILE:O	2.19	0.42
1:D:439:PRO:HG3	1:F:24:LEU:HD22	2.01	0.42
1:F:348:VAL:HG12	1:F:378:PRO:HG2	2.01	0.42
1:A:359:ARG:HD3	2:A:636:HOH:O	2.20	0.42
1:A:505:ARG:NH2	2:A:618:HOH:O	2.52	0.42
1:D:501:PRO:O	1:D:504:THR:HG23	2.19	0.42
1:F:499:ILE:CA	1:F:503:ARG:NH2	2.80	0.42
1:C:348:VAL:CG1	1:C:378:PRO:HG2	2.49	0.42
1:E:213:SER:HA	1:E:218:GLU:C	2.44	0.42
1:E:439:PRO:HD3	1:E:499:ILE:O	2.19	0.42
1:F:221:THR:HB	2:F:601:HOH:O	2.18	0.42
1:A:416:LYS:HD2	1:A:418:TYR:CZ	2.54	0.42
1:B:38:LYS:HE3	1:B:38:LYS:HB3	1.82	0.42
1:C:316:HIS:HB3	1:C:356:LYS:HE3	2.01	0.42
1:C:417:ALA:HB3	1:C:443:ILE:HA	2.01	0.42
1:C:497:MSE:HE1	1:C:507:GLU:CG	2.50	0.42
1:A:417:ALA:HB3	1:A:443:ILE:HA	2.02	0.42
1:B:215:VAL:HG12	1:B:216:MSE:CE	2.50	0.42
1:C:338:VAL:O	1:C:373:THR:HA	2.20	0.42
1:C:244:ASP:HB2	2:C:602:HOH:O	2.20	0.42
1:C:472:LYS:O	1:C:476:VAL:HG13	2.20	0.42
1:E:299:MSE:HE3	1:E:299:MSE:O	2.20	0.42
1:E:358:ALA:O	1:E:362:ARG:HG3	2.20	0.42
1:C:68:ASP:OD2	1:C:124:LYS:HE3	2.19	0.41
1:D:88:VAL:HG13	1:D:121:LYS:HE3	2.02	0.41
1:D:198:VAL:HG12	1:D:201:ILE:HG12	2.02	0.41
1:E:299:MSE:HG2	1:E:340:ASN:O	2.20	0.41
1:E:322:ASN:OD1	1:E:322:ASN:N	2.52	0.41
1:F:95:ILE:HG13	1:F:100:VAL:HG21	2.01	0.41
1:F:205:PHE:CZ	1:F:209:PRO:HD3	2.55	0.41
1:F:316:HIS:HB3	1:F:356:LYS:HE3	2.00	0.41
1:A:338:VAL:O	1:A:373:THR:HA	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:213:SER:HB3	1:F:219:GLN:HA	2.02	0.41
1:F:435:VAL:CG1	1:F:495:VAL:HG12	2.50	0.41
1:B:242:GLY:HA2	1:B:247:ASP:OD2	2.20	0.41
1:D:438:TRP:HE1	1:D:504:THR:HG22	1.83	0.41
1:F:231:HIS:HA	1:F:235:SER:OG	2.21	0.41
1:F:480:ARG:HA	1:F:482:ARG:NH2	2.36	0.41
1:A:376:ASP:OD1	1:A:415:ARG:HB3	2.20	0.41
1:E:436:MSE:HE2	1:E:497:MSE:SE	2.69	0.41
1:C:436:MSE:CE	1:C:507:GLU:HB2	2.49	0.41
1:D:23:ASP:O	1:D:27:ARG:HG3	2.20	0.41
1:D:77:GLU:HB2	1:D:80:ILE:HG13	2.02	0.41
1:B:289:ILE:HD11	1:B:438:TRP:HZ3	1.86	0.41
1:E:151:VAL:HA	1:E:154:ILE:HD12	2.01	0.41
1:A:86:ASP:OD2	1:A:117:ALA:HB3	2.21	0.41
1:C:376:ASP:CG	1:C:415:ARG:HB3	2.45	0.41
1:D:202:SER:O	1:D:228:PRO:HD3	2.21	0.41
1:A:298:ASP:HB2	1:A:343:ARG:HD2	2.03	0.41
1:C:231:HIS:HD2	1:C:237:ASN:HD22	1.68	0.41
1:C:427:SER:O	1:C:430:ILE:HG22	2.21	0.41
1:B:307:LEU:HD11	1:B:337:VAL:HG21	2.02	0.41
1:D:424:VAL:O	1:D:426:GLY:N	2.48	0.41
1:E:298:ASP:HB3	1:E:301:ASP:H	1.86	0.41
1:F:125:ILE:HD13	1:F:125:ILE:HA	1.81	0.41
1:F:525:ARG:NH2	2:F:616:HOH:O	2.51	0.41
1:C:141:ASP:OD1	1:C:179:GLY:HA3	2.21	0.41
1:E:74:ARG:HD3	1:E:116:GLU:OE2	2.21	0.41
1:E:299:MSE:CG	1:E:339:ALA:HB1	2.51	0.41
1:F:231:HIS:CD2	1:F:237:ASN:HD22	2.39	0.41
1:A:15:THR:HG22	1:B:481:ARG:NH2	2.32	0.40
1:C:497:MSE:HE1	1:C:507:GLU:HG3	2.03	0.40
1:A:140:ASN:ND2	2:A:605:HOH:O	2.31	0.40
1:A:148:GLN:H	1:A:148:GLN:CD	2.29	0.40
1:B:358:ALA:O	1:B:362:ARG:HG3	2.21	0.40
1:B:416:LYS:HD2	1:B:418:TYR:CZ	2.56	0.40
1:C:221:THR:HG22	1:C:224:GLN:HG3	2.03	0.40
1:E:251:TRP:NE1	1:E:313:LEU:HD11	2.35	0.40
1:F:182:ALA:HA	1:F:205:PHE:O	2.21	0.40
1:F:232:ALA:O	1:F:319:PHE:HB2	2.21	0.40
1:E:213:SER:HA	1:E:218:GLU:O	2.22	0.40
1:A:381:LEU:HA	1:A:382:PRO:HD3	1.79	0.40
1:B:448:ALA:O	1:B:449:ASN:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:226:GLY:O	1:C:231:HIS:CE1	2.75	0.40
1:D:52:ARG:NH2	1:D:66:GLU:OE2	2.55	0.40
1:B:114:MSE:HE2	1:B:156:TYR:HB2	2.01	0.40
1:D:438:TRP:NE1	1:D:504:THR:CG2	2.84	0.40
1:E:29:GLU:O	1:E:33:VAL:HG12	2.22	0.40
1:E:98:ARG:NH1	2:E:617:HOH:O	2.54	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:C:469:ALA:N	2:A:764:HOH:O[1_655]	2.02	0.18

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	490/570 (86%)	470 (96%)	18 (4%)	2 (0%)	30	37
1	B	493/570 (86%)	478 (97%)	15 (3%)	0	100	100
1	C	499/570 (88%)	480 (96%)	16 (3%)	3 (1%)	21	27
1	D	484/570 (85%)	467 (96%)	17 (4%)	0	100	100
1	E	485/570 (85%)	469 (97%)	15 (3%)	1 (0%)	43	54
1	F	475/570 (83%)	459 (97%)	16 (3%)	0	100	100
All	All	2926/3420 (86%)	2823 (96%)	97 (3%)	6 (0%)	43	54

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	477	ASP
1	C	478	ASP

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Mol	Chain	Res	Type
1	A	383	GLY
1	C	474	ASN
1	A	452	VAL
1	E	141	ASP

### 5.3.2 Protein sidechains [i](#)

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	395/442 (89%)	392 (99%)	3 (1%)	73	83
1	B	394/442 (89%)	388 (98%)	6 (2%)	57	70
1	C	398/442 (90%)	396 (100%)	2 (0%)	81	87
1	D	386/442 (87%)	386 (100%)	0	100	100
1	E	391/442 (88%)	389 (100%)	2 (0%)	81	87
1	F	383/442 (87%)	381 (100%)	2 (0%)	81	87
All	All	2347/2652 (88%)	2332 (99%)	15 (1%)	78	86

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LEU
1	A	126	TYR
1	A	148	GLN
1	B	126	TYR
1	B	389	GLN
1	B	409	LYS
1	B	445	VAL
1	B	452	VAL
1	B	497	MSE
1	C	178	MSE
1	C	450	SER
1	E	100	VAL
1	E	162	VAL
1	F	94	THR

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Mol	Chain	Res	Type
1	F	224	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	237	ASN
1	A	316	HIS
1	A	344	HIS
1	A	492	HIS
1	B	164	ASN
1	C	164	ASN
1	C	231	HIS
1	C	237	ASN
1	C	492	HIS
1	D	140	ASN
1	D	344	HIS
1	D	492	HIS
1	E	237	ASN
1	E	246	GLN
1	E	386	GLN
1	F	43	GLN
1	F	92	HIS
1	F	164	ASN
1	F	224	GLN
1	F	231	HIS
1	F	386	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	483/570 (84%)	-0.03	15 (3%) 51 52	25, 42, 77, 104	0
1	B	484/570 (84%)	-0.14	15 (3%) 51 52	23, 39, 73, 112	0
1	C	490/570 (85%)	-0.01	24 (4%) 35 33	25, 41, 89, 141	0
1	D	475/570 (83%)	0.01	8 (1%) 69 71	25, 45, 78, 130	0
1	E	481/570 (84%)	0.14	16 (3%) 49 50	27, 50, 85, 112	0
1	F	470/570 (82%)	0.15	13 (2%) 55 56	28, 50, 80, 102	0
All	All	2883/3420 (84%)	0.02	91 (3%) 50 51	23, 44, 81, 141	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	215	VAL	6.6
1	C	469	ALA	5.9
1	C	452	VAL	5.8
1	A	454	ILE	5.7
1	F	452	VAL	5.7
1	A	452	VAL	5.6
1	B	452	VAL	5.4
1	D	448	ALA	4.9
1	C	471	VAL	4.9
1	F	214	ALA	4.5
1	C	214	ALA	4.3
1	B	475	LEU	4.2
1	C	448	ALA	4.1
1	C	470	ALA	4.0
1	E	76	VAL	4.0
1	E	217	GLY	3.9
1	D	449	ASN	3.9
1	C	80	ILE	3.9
1	C	14	PRO	3.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	451	ALA	3.8
1	F	15	THR	3.8
1	C	479	TYR	3.7
1	B	451	ALA	3.6
1	E	214	ALA	3.6
1	D	294	GLN	3.6
1	F	480	ARG	3.6
1	E	15	THR	3.5
1	E	14	PRO	3.5
1	C	474	ASN	3.5
1	B	476	VAL	3.5
1	E	75	THR	3.4
1	E	298	ASP	3.3
1	C	294	GLN	3.3
1	C	79	GLY	3.3
1	C	472	LYS	3.2
1	F	211	VAL	3.2
1	C	473	GLU	3.1
1	E	476	VAL	3.1
1	F	503	ARG	3.0
1	F	482	ARG	3.0
1	B	78	ALA	2.9
1	A	475	LEU	2.9
1	C	449	ASN	2.8
1	E	80	ILE	2.7
1	B	478	ASP	2.7
1	D	76	VAL	2.7
1	C	78	ALA	2.7
1	C	215	VAL	2.6
1	E	475	LEU	2.6
1	F	81	PRO	2.5
1	B	34	LEU	2.5
1	E	474	ASN	2.5
1	B	14	PRO	2.4
1	C	475	LEU	2.4
1	B	15	THR	2.4
1	B	450	SER	2.4
1	A	34	LEU	2.4
1	A	45	LEU	2.4
1	B	479	TYR	2.4
1	B	477	ASP	2.3
1	F	45	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	215	VAL	2.3
1	A	37	LYS	2.3
1	C	76	VAL	2.3
1	C	15	THR	2.3
1	B	80	ILE	2.3
1	E	418	TYR	2.3
1	F	147	ILE	2.2
1	A	211	VAL	2.2
1	A	453	PRO	2.2
1	F	448	ALA	2.2
1	C	126	TYR	2.2
1	A	383	GLY	2.2
1	D	481	ARG	2.2
1	B	76	VAL	2.2
1	D	447	GLY	2.2
1	E	213	SER	2.2
1	C	210	GLU	2.2
1	E	211	VAL	2.1
1	B	449	ASN	2.1
1	D	78	ALA	2.1
1	A	214	ALA	2.1
1	A	217	GLY	2.1
1	E	478	ASP	2.1
1	E	212	VAL	2.1
1	A	81	PRO	2.1
1	D	272	HIS	2.1
1	F	279	THR	2.1
1	A	294	GLN	2.0
1	C	476	VAL	2.0
1	A	273	ASP	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.