



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:23 AM UTC

PDB ID : 3N2L / pdb_00003n2l
Title : 2.1 Angstrom resolution crystal structure of an Orotate Phosphoribosyltransferase (pyrE) from *Vibrio cholerae* O1 biovar eltor str. N16961
Authors : Halavaty, A.S.; Minasov, G.; Shuvalova, L.; Dubrovskaya, I.; Winsor, J.; Kwon, K.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2010-05-18
Resolution : 2.10 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

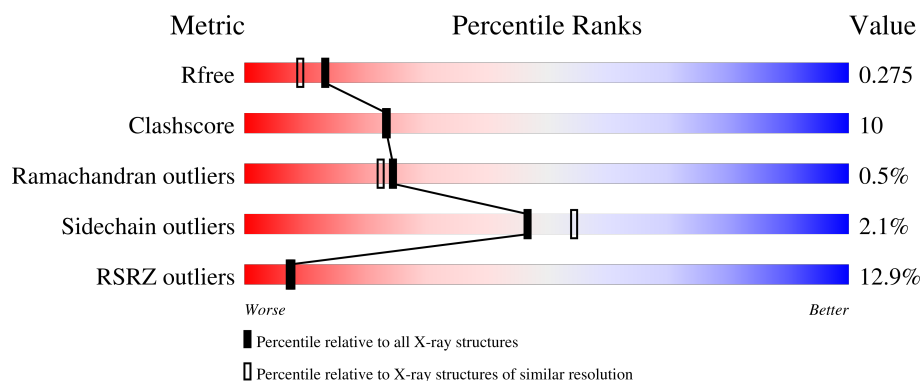
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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.



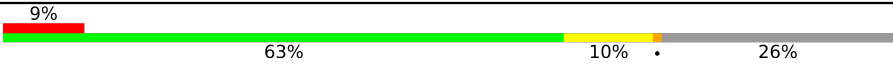
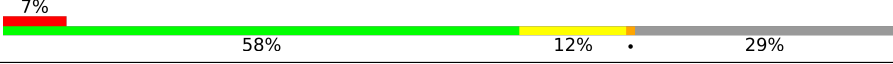
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	238	13% 66% 16% • 14%
1	B	238	7% 57% 15% • 26%
1	C	238	12% 58% 22% • 19%
1	D	238	8% 56% 14% • 29%
1	E	238	12% 66% 18% • 14%

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Mol	Chain	Length	Quality of chain
1	F	238	
1	G	238	
1	H	238	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12436 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 Orotate phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	0	3	0
			1626	1034	278	309	5			
1	B	175	Total	C	N	O	S	0	1	0
			1377	884	229	258	6			
1	C	193	Total	C	N	O	S	0	3	0
			1546	988	261	291	6			
1	D	170	Total	C	N	O	S	0	0	0
			1330	855	222	248	5			
1	E	205	Total	C	N	O	S	0	3	0
			1626	1037	275	309	5			
1	F	177	Total	C	N	O	S	0	1	0
			1391	893	232	261	5			
1	G	194	Total	C	N	O	S	0	3	0
			1554	991	268	290	5			
1	H	169	Total	C	N	O	S	0	1	0
			1335	860	220	250	5			

There are 200 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP Q9KVD5
A	-22	HIS	-	expression tag	UNP Q9KVD5
A	-21	HIS	-	expression tag	UNP Q9KVD5
A	-20	HIS	-	expression tag	UNP Q9KVD5
A	-19	HIS	-	expression tag	UNP Q9KVD5
A	-18	HIS	-	expression tag	UNP Q9KVD5
A	-17	HIS	-	expression tag	UNP Q9KVD5
A	-16	SER	-	expression tag	UNP Q9KVD5
A	-15	SER	-	expression tag	UNP Q9KVD5
A	-14	GLY	-	expression tag	UNP Q9KVD5
A	-13	VAL	-	expression tag	UNP Q9KVD5
A	-12	ASP	-	expression tag	UNP Q9KVD5
A	-11	LEU	-	expression tag	UNP Q9KVD5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLY	-	expression tag	UNP Q9KVD5
A	-9	THR	-	expression tag	UNP Q9KVD5
A	-8	GLU	-	expression tag	UNP Q9KVD5
A	-7	ASN	-	expression tag	UNP Q9KVD5
A	-6	LEU	-	expression tag	UNP Q9KVD5
A	-5	TYR	-	expression tag	UNP Q9KVD5
A	-4	PHE	-	expression tag	UNP Q9KVD5
A	-3	GLN	-	expression tag	UNP Q9KVD5
A	-2	SER	-	expression tag	UNP Q9KVD5
A	-1	ASN	-	expression tag	UNP Q9KVD5
A	0	ALA	-	expression tag	UNP Q9KVD5
A	1	MET	-	expression tag	UNP Q9KVD5
B	-23	MET	-	expression tag	UNP Q9KVD5
B	-22	HIS	-	expression tag	UNP Q9KVD5
B	-21	HIS	-	expression tag	UNP Q9KVD5
B	-20	HIS	-	expression tag	UNP Q9KVD5
B	-19	HIS	-	expression tag	UNP Q9KVD5
B	-18	HIS	-	expression tag	UNP Q9KVD5
B	-17	HIS	-	expression tag	UNP Q9KVD5
B	-16	SER	-	expression tag	UNP Q9KVD5
B	-15	SER	-	expression tag	UNP Q9KVD5
B	-14	GLY	-	expression tag	UNP Q9KVD5
B	-13	VAL	-	expression tag	UNP Q9KVD5
B	-12	ASP	-	expression tag	UNP Q9KVD5
B	-11	LEU	-	expression tag	UNP Q9KVD5
B	-10	GLY	-	expression tag	UNP Q9KVD5
B	-9	THR	-	expression tag	UNP Q9KVD5
B	-8	GLU	-	expression tag	UNP Q9KVD5
B	-7	ASN	-	expression tag	UNP Q9KVD5
B	-6	LEU	-	expression tag	UNP Q9KVD5
B	-5	TYR	-	expression tag	UNP Q9KVD5
B	-4	PHE	-	expression tag	UNP Q9KVD5
B	-3	GLN	-	expression tag	UNP Q9KVD5
B	-2	SER	-	expression tag	UNP Q9KVD5
B	-1	ASN	-	expression tag	UNP Q9KVD5
B	0	ALA	-	expression tag	UNP Q9KVD5
B	1	MET	-	expression tag	UNP Q9KVD5
C	-23	MET	-	expression tag	UNP Q9KVD5
C	-22	HIS	-	expression tag	UNP Q9KVD5
C	-21	HIS	-	expression tag	UNP Q9KVD5
C	-20	HIS	-	expression tag	UNP Q9KVD5
C	-19	HIS	-	expression tag	UNP Q9KVD5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-18	HIS	-	expression tag	UNP Q9KVD5
C	-17	HIS	-	expression tag	UNP Q9KVD5
C	-16	SER	-	expression tag	UNP Q9KVD5
C	-15	SER	-	expression tag	UNP Q9KVD5
C	-14	GLY	-	expression tag	UNP Q9KVD5
C	-13	VAL	-	expression tag	UNP Q9KVD5
C	-12	ASP	-	expression tag	UNP Q9KVD5
C	-11	LEU	-	expression tag	UNP Q9KVD5
C	-10	GLY	-	expression tag	UNP Q9KVD5
C	-9	THR	-	expression tag	UNP Q9KVD5
C	-8	GLU	-	expression tag	UNP Q9KVD5
C	-7	ASN	-	expression tag	UNP Q9KVD5
C	-6	LEU	-	expression tag	UNP Q9KVD5
C	-5	TYR	-	expression tag	UNP Q9KVD5
C	-4	PHE	-	expression tag	UNP Q9KVD5
C	-3	GLN	-	expression tag	UNP Q9KVD5
C	-2	SER	-	expression tag	UNP Q9KVD5
C	-1	ASN	-	expression tag	UNP Q9KVD5
C	0	ALA	-	expression tag	UNP Q9KVD5
C	1	MET	-	expression tag	UNP Q9KVD5
D	-23	MET	-	expression tag	UNP Q9KVD5
D	-22	HIS	-	expression tag	UNP Q9KVD5
D	-21	HIS	-	expression tag	UNP Q9KVD5
D	-20	HIS	-	expression tag	UNP Q9KVD5
D	-19	HIS	-	expression tag	UNP Q9KVD5
D	-18	HIS	-	expression tag	UNP Q9KVD5
D	-17	HIS	-	expression tag	UNP Q9KVD5
D	-16	SER	-	expression tag	UNP Q9KVD5
D	-15	SER	-	expression tag	UNP Q9KVD5
D	-14	GLY	-	expression tag	UNP Q9KVD5
D	-13	VAL	-	expression tag	UNP Q9KVD5
D	-12	ASP	-	expression tag	UNP Q9KVD5
D	-11	LEU	-	expression tag	UNP Q9KVD5
D	-10	GLY	-	expression tag	UNP Q9KVD5
D	-9	THR	-	expression tag	UNP Q9KVD5
D	-8	GLU	-	expression tag	UNP Q9KVD5
D	-7	ASN	-	expression tag	UNP Q9KVD5
D	-6	LEU	-	expression tag	UNP Q9KVD5
D	-5	TYR	-	expression tag	UNP Q9KVD5
D	-4	PHE	-	expression tag	UNP Q9KVD5
D	-3	GLN	-	expression tag	UNP Q9KVD5
D	-2	SER	-	expression tag	UNP Q9KVD5

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	ASN	-	expression tag	UNP Q9KVD5
D	0	ALA	-	expression tag	UNP Q9KVD5
D	1	MET	-	expression tag	UNP Q9KVD5
E	-23	MET	-	expression tag	UNP Q9KVD5
E	-22	HIS	-	expression tag	UNP Q9KVD5
E	-21	HIS	-	expression tag	UNP Q9KVD5
E	-20	HIS	-	expression tag	UNP Q9KVD5
E	-19	HIS	-	expression tag	UNP Q9KVD5
E	-18	HIS	-	expression tag	UNP Q9KVD5
E	-17	HIS	-	expression tag	UNP Q9KVD5
E	-16	SER	-	expression tag	UNP Q9KVD5
E	-15	SER	-	expression tag	UNP Q9KVD5
E	-14	GLY	-	expression tag	UNP Q9KVD5
E	-13	VAL	-	expression tag	UNP Q9KVD5
E	-12	ASP	-	expression tag	UNP Q9KVD5
E	-11	LEU	-	expression tag	UNP Q9KVD5
E	-10	GLY	-	expression tag	UNP Q9KVD5
E	-9	THR	-	expression tag	UNP Q9KVD5
E	-8	GLU	-	expression tag	UNP Q9KVD5
E	-7	ASN	-	expression tag	UNP Q9KVD5
E	-6	LEU	-	expression tag	UNP Q9KVD5
E	-5	TYR	-	expression tag	UNP Q9KVD5
E	-4	PHE	-	expression tag	UNP Q9KVD5
E	-3	GLN	-	expression tag	UNP Q9KVD5
E	-2	SER	-	expression tag	UNP Q9KVD5
E	-1	ASN	-	expression tag	UNP Q9KVD5
E	0	ALA	-	expression tag	UNP Q9KVD5
E	1	MET	-	expression tag	UNP Q9KVD5
F	-23	MET	-	expression tag	UNP Q9KVD5
F	-22	HIS	-	expression tag	UNP Q9KVD5
F	-21	HIS	-	expression tag	UNP Q9KVD5
F	-20	HIS	-	expression tag	UNP Q9KVD5
F	-19	HIS	-	expression tag	UNP Q9KVD5
F	-18	HIS	-	expression tag	UNP Q9KVD5
F	-17	HIS	-	expression tag	UNP Q9KVD5
F	-16	SER	-	expression tag	UNP Q9KVD5
F	-15	SER	-	expression tag	UNP Q9KVD5
F	-14	GLY	-	expression tag	UNP Q9KVD5
F	-13	VAL	-	expression tag	UNP Q9KVD5
F	-12	ASP	-	expression tag	UNP Q9KVD5
F	-11	LEU	-	expression tag	UNP Q9KVD5
F	-10	GLY	-	expression tag	UNP Q9KVD5

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-9	THR	-	expression tag	UNP Q9KVD5
F	-8	GLU	-	expression tag	UNP Q9KVD5
F	-7	ASN	-	expression tag	UNP Q9KVD5
F	-6	LEU	-	expression tag	UNP Q9KVD5
F	-5	TYR	-	expression tag	UNP Q9KVD5
F	-4	PHE	-	expression tag	UNP Q9KVD5
F	-3	GLN	-	expression tag	UNP Q9KVD5
F	-2	SER	-	expression tag	UNP Q9KVD5
F	-1	ASN	-	expression tag	UNP Q9KVD5
F	0	ALA	-	expression tag	UNP Q9KVD5
F	1	MET	-	expression tag	UNP Q9KVD5
G	-23	MET	-	expression tag	UNP Q9KVD5
G	-22	HIS	-	expression tag	UNP Q9KVD5
G	-21	HIS	-	expression tag	UNP Q9KVD5
G	-20	HIS	-	expression tag	UNP Q9KVD5
G	-19	HIS	-	expression tag	UNP Q9KVD5
G	-18	HIS	-	expression tag	UNP Q9KVD5
G	-17	HIS	-	expression tag	UNP Q9KVD5
G	-16	SER	-	expression tag	UNP Q9KVD5
G	-15	SER	-	expression tag	UNP Q9KVD5
G	-14	GLY	-	expression tag	UNP Q9KVD5
G	-13	VAL	-	expression tag	UNP Q9KVD5
G	-12	ASP	-	expression tag	UNP Q9KVD5
G	-11	LEU	-	expression tag	UNP Q9KVD5
G	-10	GLY	-	expression tag	UNP Q9KVD5
G	-9	THR	-	expression tag	UNP Q9KVD5
G	-8	GLU	-	expression tag	UNP Q9KVD5
G	-7	ASN	-	expression tag	UNP Q9KVD5
G	-6	LEU	-	expression tag	UNP Q9KVD5
G	-5	TYR	-	expression tag	UNP Q9KVD5
G	-4	PHE	-	expression tag	UNP Q9KVD5
G	-3	GLN	-	expression tag	UNP Q9KVD5
G	-2	SER	-	expression tag	UNP Q9KVD5
G	-1	ASN	-	expression tag	UNP Q9KVD5
G	0	ALA	-	expression tag	UNP Q9KVD5
G	1	MET	-	expression tag	UNP Q9KVD5
H	-23	MET	-	expression tag	UNP Q9KVD5
H	-22	HIS	-	expression tag	UNP Q9KVD5
H	-21	HIS	-	expression tag	UNP Q9KVD5
H	-20	HIS	-	expression tag	UNP Q9KVD5
H	-19	HIS	-	expression tag	UNP Q9KVD5
H	-18	HIS	-	expression tag	UNP Q9KVD5

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-17	HIS	-	expression tag	UNP Q9KVD5
H	-16	SER	-	expression tag	UNP Q9KVD5
H	-15	SER	-	expression tag	UNP Q9KVD5
H	-14	GLY	-	expression tag	UNP Q9KVD5
H	-13	VAL	-	expression tag	UNP Q9KVD5
H	-12	ASP	-	expression tag	UNP Q9KVD5
H	-11	LEU	-	expression tag	UNP Q9KVD5
H	-10	GLY	-	expression tag	UNP Q9KVD5
H	-9	THR	-	expression tag	UNP Q9KVD5
H	-8	GLU	-	expression tag	UNP Q9KVD5
H	-7	ASN	-	expression tag	UNP Q9KVD5
H	-6	LEU	-	expression tag	UNP Q9KVD5
H	-5	TYR	-	expression tag	UNP Q9KVD5
H	-4	PHE	-	expression tag	UNP Q9KVD5
H	-3	GLN	-	expression tag	UNP Q9KVD5
H	-2	SER	-	expression tag	UNP Q9KVD5
H	-1	ASN	-	expression tag	UNP Q9KVD5
H	0	ALA	-	expression tag	UNP Q9KVD5
H	1	MET	-	expression tag	UNP Q9KVD5

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	H	2	Total Cl 2 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	95	Total O 96 96	0	1
3	B	82	Total O 83 83	0	2
3	C	78	Total O 79 79	0	1
3	D	71	Total O 74 74	0	3
3	E	100	Total O 101 101	0	3
3	F	68	Total O 69 69	0	1

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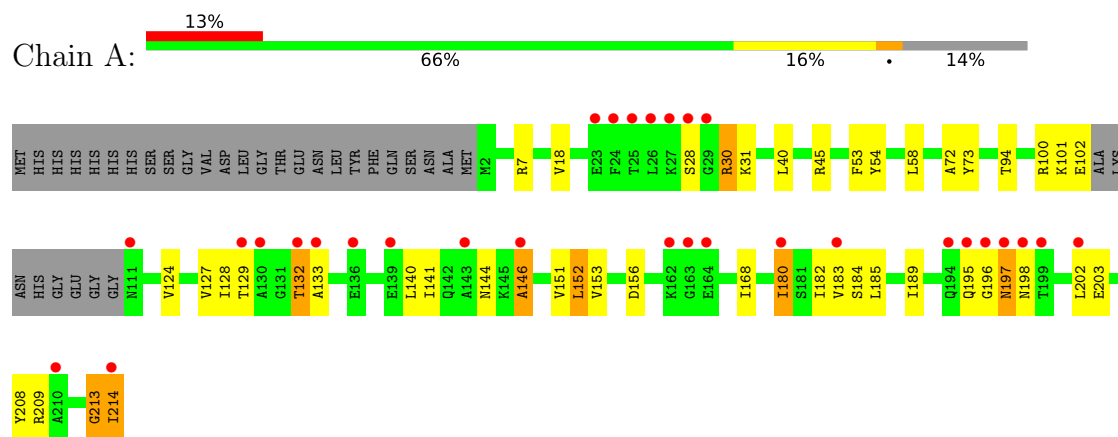
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	74	Total 74	O 74	0	1
3	H	71	Total 72	O 72	0	1

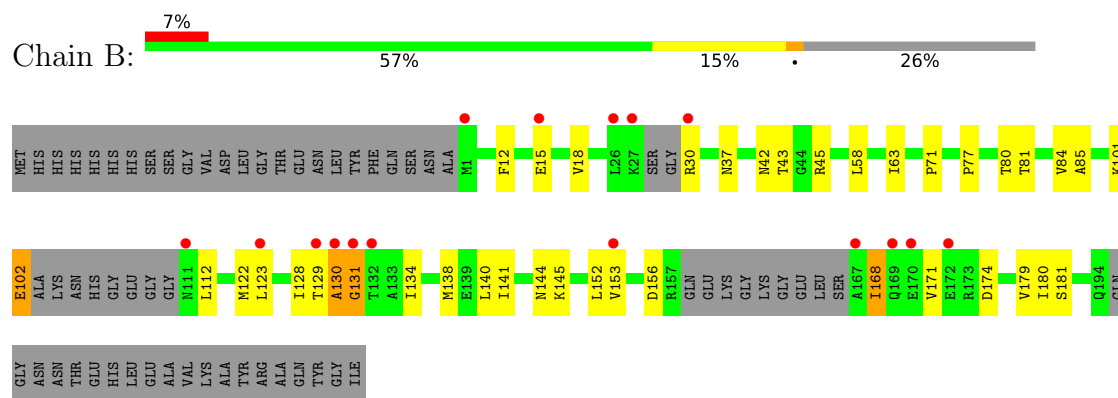
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

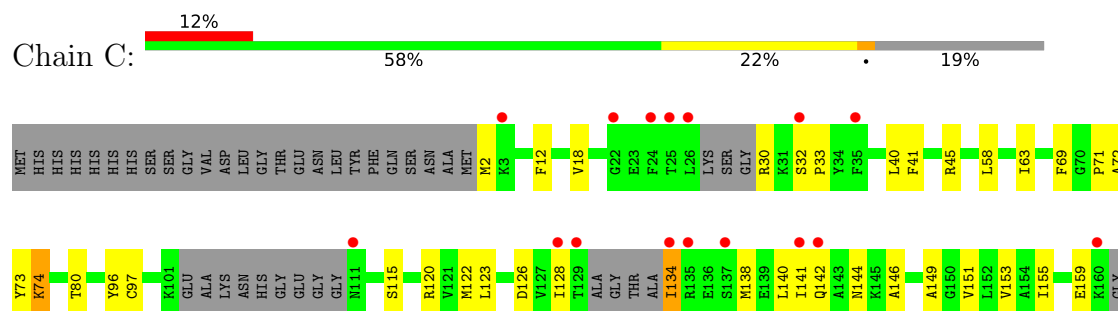
• Molecule 1: Orotate phosphoribosyltransferase



• Molecule 1: Orotate phosphoribosyltransferase

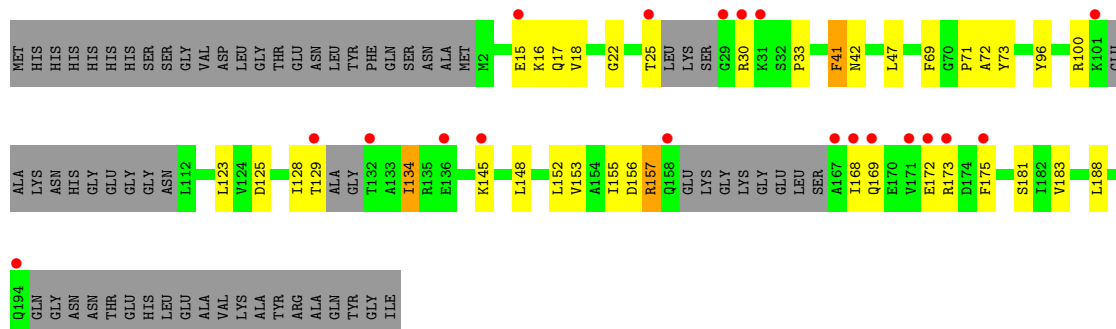


• Molecule 1: Orotate phosphoribosyltransferase

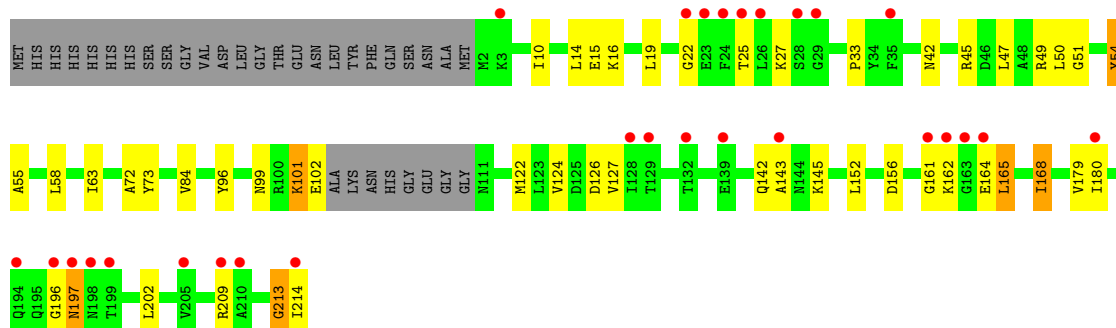




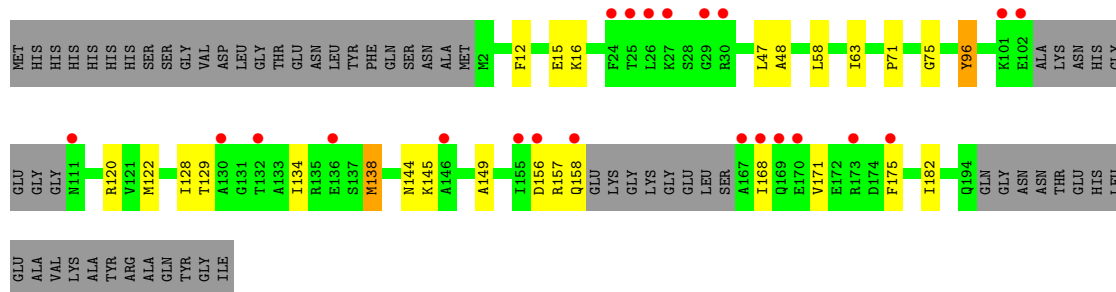
• Molecule 1: Orotate phosphoribosyltransferase



• Molecule 1: Orotate phosphoribosyltransferase

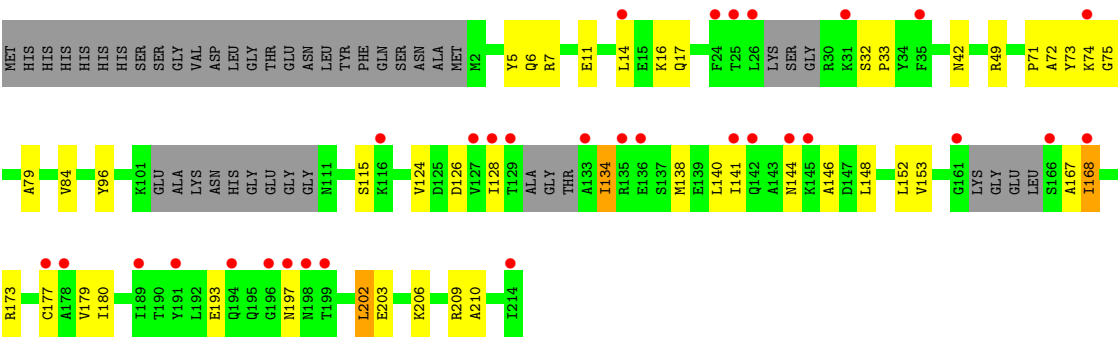


• Molecule 1: Orotate phosphoribosyltransferase

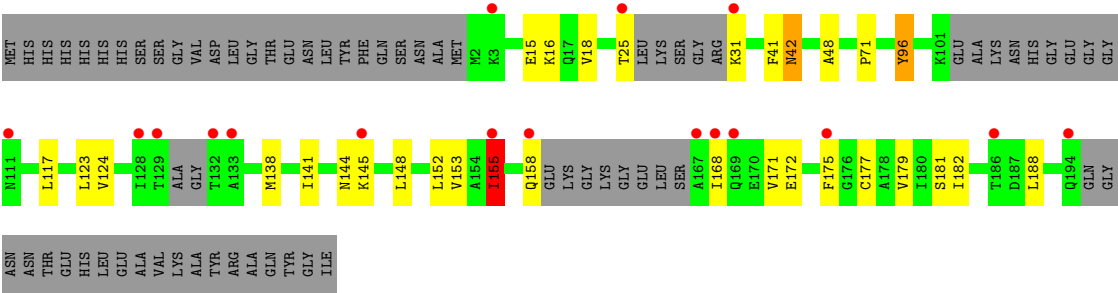


• Molecule 1: Orotate phosphoribosyltransferase





● Molecule 1: Orotate phosphoribosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	96.59Å 76.71Å 133.94Å 90.00° 92.63° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 30.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.00-2.10) 99.6 (30.00-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.222 , 0.269 0.257 , 0.275	Depositor DCC
R_{free} test set	5712 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.573	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.052 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12436	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 89.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.8806e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: CL

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	0.96	1/1651 (0.1%)	1.08	4/2222 (0.2%)
1	B	0.91	2/1397 (0.1%)	0.99	3/1882 (0.2%)
1	C	0.84	0/1568	0.96	1/2109 (0.0%)
1	D	0.85	0/1349	0.99	2/1817 (0.1%)
1	E	0.95	3/1652 (0.2%)	1.07	6/2224 (0.3%)
1	F	0.90	0/1413	1.00	4/1906 (0.2%)
1	G	0.85	1/1576 (0.1%)	0.95	1/2118 (0.0%)
1	H	0.84	1/1355 (0.1%)	1.01	4/1827 (0.2%)
All	All	0.89	8/11961 (0.1%)	1.01	25/16105 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	45	ARG	C-O	-5.85	1.17	1.24
1	E	51	GLY	N-CA	-5.67	1.38	1.45
1	B	85	ALA	CA-CB	5.59	1.62	1.53
1	E	54	TYR	C-O	-5.53	1.17	1.24
1	A	53	PHE	CA-C	5.25	1.59	1.52
1	H	155	ILE	CA-CB	-5.25	1.50	1.55
1	G	49	ARG	C-O	-5.11	1.18	1.24
1	E	55	ALA	CA-CB	5.01	1.61	1.53

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	164	GLU	CB-CA-C	-9.74	91.04	110.42
1	A	213	GLY	N-CA-C	9.59	122.88	112.33
1	E	164	GLU	N-CA-C	8.12	128.10	110.80
1	A	30	ARG	N-CA-C	7.53	120.29	110.43
1	E	213	GLY	N-CA-C	7.41	121.62	111.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	71	PRO	CB-CA-C	-7.11	102.86	111.46
1	H	42	ASN	N-CA-C	7.06	121.61	112.92
1	B	130	ALA	N-CA-C	-6.58	104.25	113.20
1	A	146	ALA	N-CA-C	-6.56	100.50	110.14
1	E	165	LEU	N-CA-CB	6.38	122.55	111.13
1	E	165	LEU	N-CA-C	-6.30	98.35	109.06
1	D	41	PHE	N-CA-C	-6.25	101.68	110.50
1	G	168	ILE	CB-CA-C	-6.21	104.03	111.97
1	F	175	PHE	N-CA-C	-6.09	100.69	110.14
1	F	129	THR	N-CA-C	-5.96	105.50	112.89
1	B	168	ILE	N-CA-C	-5.71	104.95	110.72
1	E	101	LYS	N-CA-C	-5.66	105.01	111.07
1	F	182	ILE	N-CA-C	-5.52	105.34	110.53
1	A	198	ASN	N-CA-C	5.38	116.83	110.97
1	H	71	PRO	N-CA-C	5.33	118.84	110.80
1	F	156	ASP	CB-CA-C	-5.32	98.80	110.67
1	C	32	SER	N-CA-C	5.31	117.36	110.40
1	D	15	GLU	CB-CA-C	-5.24	102.04	110.74
1	B	174	ASP	N-CA-C	5.23	119.66	113.28
1	H	182	ILE	N-CA-C	-5.04	105.48	110.62

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	1626	0	1623	44	0
1	B	1377	0	1383	28	0
1	C	1546	0	1540	46	0
1	D	1330	0	1332	23	0
1	E	1626	0	1619	37	0
1	F	1391	0	1391	20	0
1	G	1554	0	1552	36	0
1	H	1335	0	1330	25	0
2	A	1	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	2	0	0	0	0
3	A	96	0	0	4	0
3	B	83	0	0	0	0
3	C	79	0	0	1	0
3	D	74	0	0	0	0
3	E	101	0	0	1	0
3	F	69	0	0	0	0
3	G	74	0	0	1	0
3	H	72	0	0	1	0
All	All	12436	0	11770	241	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (241) 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:69:PHE:HD1	1:C:97[B]:CYS:HG	0.96	0.90
1:C:69:PHE:HD1	1:C:97[B]:CYS:SG	1.99	0.85
1:E:15:GLU:OE1	1:E:16:LYS:HE2	1.80	0.81
1:G:140:LEU:O	1:G:144:ASN:ND2	2.21	0.73
1:B:101:LYS:O	1:B:102:GLU:C	2.32	0.71
1:D:25:THR:C	1:D:30:ARG:O	2.38	0.67
1:E:84:VAL:HG13	1:F:48:ALA:HB2	1.76	0.67
1:H:145:LYS:O	1:H:145:LYS:HG2	1.96	0.66
1:A:196:GLY:O	1:A:197:ASN:C	2.40	0.64
1:A:18:VAL:HG22	1:A:40:LEU:HB2	1.80	0.64
1:A:214:ILE:N	1:A:214:ILE:HD12	2.12	0.63
1:G:202:LEU:HD23	1:G:202:LEU:C	2.23	0.63
1:E:15:GLU:OE1	1:E:49:ARG:NH2	2.32	0.63
1:E:58:LEU:HG	1:E:152:LEU:HD22	1.80	0.63
1:A:152:LEU:HD23	1:A:153:VAL:O	1.99	0.63
1:D:168:ILE:HG13	1:D:169:GLN:N	2.12	0.63
1:D:30:ARG:HD3	1:D:129:THR:HG21	1.81	0.62
1:E:10:ILE:O	1:E:14:LEU:HG	1.98	0.62
1:C:189:ILE:C	1:C:189:ILE:HD12	2.24	0.62
1:A:152:LEU:HD23	1:A:153:VAL:N	2.15	0.61
1:F:138:MET:HE1	1:F:171:VAL:HG13	1.82	0.61
1:E:84:VAL:HG13	1:F:48:ALA:CB	2.30	0.61
1:F:157:ARG:O	1:F:158:GLN:HB2	2.01	0.60
1:E:101:LYS:O	1:E:102:GLU:C	2.44	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:202:LEU:C	1:C:202:LEU:HD23	2.26	0.60
1:C:195:GLN:O	1:C:196:GLY:O	2.19	0.60
1:G:202:LEU:HD21	1:G:206:LYS:HE3	1.83	0.60
1:D:148:LEU:HD13	1:D:175:PHE:HB3	1.84	0.59
1:A:182:ILE:HG22	1:A:183:VAL:HG13	1.84	0.59
1:E:45:ARG:NH1	3:E:596:HOH:O	2.28	0.59
1:G:128:ILE:HG22	1:G:167:ALA:HB1	1.84	0.59
1:F:63:ILE:HG21	1:F:122:MET:HE1	1.85	0.58
1:B:63:ILE:HD12	1:B:180:ILE:HD13	1.84	0.58
1:B:112:LEU:HD21	1:B:140:LEU:HD21	1.85	0.58
1:C:189:ILE:HD12	1:C:190:THR:N	2.18	0.58
1:B:168:ILE:HG23	1:B:179:VAL:HB	1.85	0.58
1:A:124:VAL:HG12	1:A:152:LEU:HB3	1.85	0.58
1:D:71:PRO:HG3	1:D:123:LEU:HD22	1.86	0.58
1:E:209:ARG:O	1:E:213:GLY:N	2.37	0.58
1:G:134:ILE:CG2	1:G:138:MET:HE2	2.33	0.58
1:B:128:ILE:HD12	1:B:134:ILE:HD13	1.86	0.57
1:A:152:LEU:HA	1:A:180:ILE:O	2.03	0.57
1:H:155:ILE:CD1	1:H:188:LEU:HD11	2.33	0.57
1:C:128:ILE:HD12	1:C:168:ILE:HG12	1.86	0.57
1:C:173:ARG:NH2	1:C:174:ASP:OD1	2.37	0.57
1:G:16:LYS:O	1:G:17:GLN:HB2	2.04	0.57
1:H:158:GLN:NE2	3:H:239:HOH:O	2.37	0.57
1:G:202:LEU:C	1:G:202:LEU:CD2	2.78	0.57
1:C:140:LEU:O	1:C:144:ASN:ND2	2.25	0.56
1:D:128:ILE:HG12	1:D:134:ILE:HD12	1.88	0.55
1:E:168:ILE:HG12	1:E:179:VAL:HG11	1.87	0.55
1:C:30:ARG:NH2	1:C:159:GLU:OE1	2.39	0.55
1:H:144:ASN:O	1:H:145:LYS:HB3	2.06	0.55
1:G:202:LEU:HD23	1:G:202:LEU:O	2.06	0.55
1:E:156:ASP:CG	1:E:168:ILE:HD12	2.32	0.55
1:B:123:LEU:HD11	1:B:141:ILE:HD12	1.88	0.55
1:E:161:GLY:O	1:E:162:LYS:C	2.48	0.54
1:G:16:LYS:HE3	3:G:227:HOH:O	2.07	0.54
1:G:74[A]:LYS:HG2	1:G:126:ASP:CG	2.33	0.54
1:A:202:LEU:HD23	1:A:203:GLU:OE1	2.08	0.54
1:G:42:ASN:HA	1:H:96[B]:TYR:CE2	2.43	0.53
1:F:138:MET:HE1	1:F:171:VAL:CG1	2.38	0.53
1:C:128:ILE:HD12	1:C:179:VAL:HG11	1.90	0.53
1:B:128:ILE:C	1:B:130:ALA:H	2.16	0.53
1:C:134:ILE:HG23	1:C:138:MET:HE2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LEU:HG	1:A:152:LEU:HD12	1.91	0.53
1:E:202:LEU:HD23	1:E:202:LEU:O	2.09	0.53
1:E:162:LYS:HA	1:E:162:LYS:HE2	1.91	0.53
1:C:209:ARG:O	1:C:213:GLY:N	2.40	0.52
1:B:138:MET:HE1	1:B:171:VAL:CG1	2.40	0.52
1:G:84:VAL:HG13	1:H:48:ALA:HB2	1.91	0.52
1:A:140:LEU:O	1:A:144:ASN:ND2	2.32	0.52
1:E:14:LEU:CD2	1:E:19:LEU:HD23	2.39	0.52
1:D:153:VAL:O	1:D:181:SER:HA	2.09	0.52
1:D:69:PHE:CD2	1:D:123:LEU:HD23	2.45	0.52
1:B:153:VAL:O	1:B:181:SER:HA	2.09	0.52
1:D:168:ILE:HG13	1:D:169:GLN:H	1.73	0.51
1:E:25:THR:HG23	1:E:25:THR:O	2.10	0.51
1:B:129:THR:C	1:B:131:GLY:H	2.19	0.51
1:C:71:PRO:CG	1:C:123:LEU:HD21	2.41	0.51
1:E:124:VAL:HG12	1:E:152:LEU:HB3	1.92	0.51
1:G:128:ILE:HD12	1:G:179:VAL:HG11	1.92	0.51
1:B:134:ILE:HG22	1:B:138:MET:SD	2.51	0.51
1:A:7:ARG:NH2	3:A:498:HOH:O	2.38	0.51
1:C:186:THR:O	1:C:189:ILE:HG13	2.10	0.50
1:H:25:THR:HA	1:H:31:LYS:HA	1.93	0.50
1:C:138:MET:HG3	1:C:175:PHE:CE1	2.46	0.50
1:B:138:MET:HE1	1:B:171:VAL:HG13	1.94	0.50
1:C:45:ARG:NH2	3:C:595:HOH:O	2.44	0.50
1:C:151:VAL:HG23	1:C:177:CYS:SG	2.51	0.50
1:E:202:LEU:HD11	1:G:193:GLU:HG2	1.94	0.50
1:E:22:GLY:O	1:E:33:PRO:HA	2.11	0.50
1:B:128:ILE:C	1:B:130:ALA:N	2.69	0.50
1:G:42:ASN:C	1:H:96[B]:TYR:CE2	2.90	0.50
1:E:63:ILE:HD12	1:E:180:ILE:HD11	1.94	0.49
1:H:123:LEU:HD11	1:H:138:MET:HG3	1.94	0.49
1:B:123:LEU:HD11	1:B:141:ILE:CD1	2.42	0.49
1:B:71:PRO:HG3	1:B:123:LEU:HD23	1.94	0.49
1:A:182:ILE:C	1:A:183:VAL:HG13	2.37	0.49
1:C:128:ILE:HG13	1:C:153:VAL:HG11	1.94	0.49
1:E:42:ASN:HA	1:F:96[B]:TYR:CE2	2.48	0.49
1:A:45[B]:ARG:NH1	3:A:246:HOH:O	2.46	0.49
1:E:14:LEU:HD21	1:E:19:LEU:HD23	1.94	0.49
1:G:11:GLU:HA	1:G:14:LEU:HD12	1.93	0.48
1:A:101:LYS:O	1:A:102:GLU:HG3	2.13	0.48
1:C:165:LEU:HD12	1:C:165:LEU:C	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:128:ILE:O	1:F:128:ILE:HG13	2.14	0.48
1:H:148:LEU:HD13	1:H:175:PHE:HB3	1.96	0.48
1:A:30:ARG:HD3	1:C:212:TYR:CE1	2.49	0.48
1:C:128:ILE:CD1	1:C:179:VAL:HG11	2.43	0.48
1:C:202:LEU:C	1:C:202:LEU:CD2	2.86	0.48
1:A:152:LEU:CD2	1:A:153:VAL:O	2.61	0.48
1:A:213:GLY:C	1:A:214:ILE:HD12	2.39	0.48
1:C:138:MET:HG3	1:C:175:PHE:CZ	2.49	0.47
1:B:128:ILE:HD12	1:B:134:ILE:CD1	2.44	0.47
1:E:214:ILE:HD12	1:E:214:ILE:N	2.29	0.47
1:A:127:VAL:HG23	1:A:129:THR:HG23	1.96	0.47
1:A:156:ASP:CG	1:A:168:ILE:HD12	2.40	0.47
1:C:155:ILE:HA	1:C:183:VAL:O	2.15	0.47
1:C:2:MET:HE1	1:C:191:TYR:HB2	1.97	0.47
1:A:101:LYS:O	1:A:102:GLU:CG	2.63	0.47
1:C:165:LEU:HD12	1:C:165:LEU:O	2.15	0.47
1:E:27:LYS:H	1:E:27:LYS:HD2	1.80	0.47
1:E:96:TYR:CZ	1:F:47:LEU:HD11	2.49	0.47
1:B:30:ARG:NH1	1:B:30:ARG:HB3	2.30	0.47
1:B:80:THR:O	1:B:84:VAL:HG23	2.15	0.47
1:H:155:ILE:HD11	1:H:188:LEU:HD11	1.97	0.47
1:H:138:MET:HE1	1:H:171:VAL:CG1	2.45	0.46
1:A:54:TYR:CD2	1:A:152:LEU:HD22	2.49	0.46
1:H:153:VAL:O	1:H:181:SER:HA	2.15	0.46
1:A:45[B]:ARG:NE	3:A:218:HOH:O	2.36	0.46
1:A:168:ILE:HG13	3:A:261:HOH:O	2.16	0.46
1:A:152:LEU:HG	1:A:180:ILE:HG22	1.97	0.46
1:G:202:LEU:HD22	1:G:203:GLU:CD	2.41	0.46
1:C:58:LEU:HD21	1:C:122:MET:HG2	1.98	0.46
1:D:145:LYS:HB2	1:D:145:LYS:HE3	1.69	0.46
1:G:84:VAL:HG13	1:H:48:ALA:CB	2.46	0.46
1:C:141:ILE:HG23	1:C:146:ALA:HB3	1.98	0.46
1:A:54:TYR:CE2	1:A:152:LEU:HD22	2.52	0.45
1:G:148:LEU:HD21	1:G:177:CYS:SG	2.57	0.45
1:E:202:LEU:HD11	1:G:193:GLU:CG	2.46	0.45
1:B:58:LEU:HD11	1:B:122:MET:HG2	1.98	0.45
1:A:72:ALA:HA	1:A:73:TYR:HA	1.82	0.45
1:A:128:ILE:HG13	1:A:153:VAL:HG11	1.98	0.45
1:A:132:THR:HG23	1:A:133:ALA:H	1.82	0.45
1:B:12:PHE:O	1:B:15[B]:GLU:HG2	2.15	0.45
1:F:157:ARG:O	1:F:158:GLN:CB	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:196:GLY:O	1:E:197:ASN:C	2.59	0.45
1:H:18:VAL:HG21	1:H:41:PHE:CE1	2.51	0.45
1:A:209:ARG:O	1:A:213:GLY:N	2.47	0.45
1:B:152:LEU:HD13	1:B:180:ILE:HG13	1.99	0.45
1:C:202:LEU:HD23	1:C:202:LEU:O	2.16	0.45
1:D:18:VAL:HG21	1:D:41:PHE:CD1	2.52	0.45
1:H:123:LEU:HD21	1:H:141:ILE:CD1	2.47	0.45
1:E:145:LYS:HD3	1:E:145:LYS:N	2.32	0.44
1:F:128:ILE:HG21	1:F:168:ILE:HG12	1.98	0.44
1:G:128:ILE:HB	1:G:168:ILE:HD12	2.00	0.44
1:H:15:GLU:HG3	1:H:16:LYS:HE3	1.99	0.44
1:H:155:ILE:HD12	1:H:188:LEU:HD11	1.99	0.44
1:B:156:ASP:C	1:B:156:ASP:OD1	2.60	0.44
1:H:117:LEU:HD12	1:H:117:LEU:HA	1.83	0.44
1:H:172:GLU:HG2	1:H:177:CYS:O	2.18	0.44
1:C:74[A]:LYS:HG2	1:C:126:ASP:CG	2.42	0.44
1:E:202:LEU:HD23	1:E:202:LEU:C	2.42	0.44
1:C:18:VAL:HG22	1:C:40:LEU:HB2	1.99	0.44
1:C:69:PHE:CD1	1:C:97[B]:CYS:SG	2.90	0.44
1:H:168:ILE:CG2	1:H:179:VAL:HG11	2.47	0.44
1:A:156:ASP:OD2	1:A:156:ASP:C	2.60	0.44
1:G:16:LYS:O	1:G:17:GLN:CB	2.66	0.43
1:A:152:LEU:CD2	1:A:152:LEU:C	2.91	0.43
1:A:180:ILE:HD13	1:A:180:ILE:HA	1.79	0.43
1:G:72:ALA:HA	1:G:73:TYR:HA	1.82	0.43
1:H:18:VAL:HG21	1:H:41:PHE:CD1	2.52	0.43
1:F:12:PHE:O	1:F:16:LYS:HG3	2.19	0.43
1:F:128:ILE:HD12	1:F:134:ILE:HD13	2.01	0.43
1:C:74[A]:LYS:HZ2	1:C:126:ASP:HB3	1.84	0.43
1:G:209:ARG:O	1:G:210:ALA:C	2.60	0.43
1:E:58:LEU:HD21	1:E:122:MET:HG2	2.00	0.43
1:E:72:ALA:HA	1:E:73:TYR:HA	1.75	0.43
1:F:144:ASN:O	1:F:145:LYS:HB2	2.18	0.43
1:B:77:PRO:O	1:B:81:THR:HG22	2.18	0.43
1:D:156:ASP:HB2	1:D:157:ARG:NH2	2.33	0.43
1:G:71:PRO:HD2	1:G:75:GLY:HA3	2.00	0.43
1:G:134:ILE:HG21	1:G:138:MET:HE2	1.99	0.43
1:A:94:THR:HG23	1:A:94:THR:O	2.18	0.43
1:A:141:ILE:HG23	1:A:146:ALA:HB3	2.00	0.43
1:G:152:LEU:HD13	1:G:180:ILE:HG13	2.01	0.43
1:D:152:LEU:HD12	1:D:181:SER:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:VAL:HG13	1:B:37:ASN:O	2.19	0.42
1:G:5:TYR:CE1	1:G:6:GLN:HG3	2.54	0.42
1:G:128:ILE:CD1	1:G:179:VAL:HG11	2.49	0.42
1:C:63:ILE:HG21	1:C:122:MET:HE1	2.01	0.42
1:E:47:LEU:HD11	1:F:96[B]:TYR:CE2	2.54	0.42
1:F:58:LEU:HD11	1:F:122:MET:HG2	2.01	0.42
1:A:100:ARG:NH1	2:A:215:CL:CL	2.79	0.42
1:A:128:ILE:HD11	1:A:151:VAL:HG11	2.00	0.42
1:A:184:SER:O	1:A:185:LEU:C	2.62	0.42
1:C:80:THR:HG23	1:D:47:LEU:HD13	2.01	0.42
1:E:165:LEU:HD23	1:E:165:LEU:HA	1.94	0.42
1:G:32:SER:OG	1:G:33:PRO:HD2	2.18	0.42
1:G:79:ALA:HA	1:G:124:VAL:HG21	2.01	0.42
1:G:115:SER:HB2	1:H:42:ASN:HB2	2.01	0.42
1:C:12:PHE:CE2	1:C:41:PHE:CE1	3.08	0.42
1:A:31:LYS:N	1:A:31:LYS:HD2	2.34	0.42
1:A:141:ILE:HG22	1:A:146:ALA:O	2.19	0.42
1:C:138:MET:HG3	1:C:175:PHE:CD1	2.55	0.42
1:D:16:LYS:O	1:D:17:GLN:HB2	2.20	0.42
1:A:189:ILE:CD1	1:C:209:ARG:HD3	2.50	0.41
1:B:42:ASN:OD1	1:B:43:THR:HG23	2.20	0.41
1:F:120:ARG:HB3	1:F:149:ALA:HB2	2.02	0.41
1:H:148:LEU:CD1	1:H:175:PHE:HB3	2.49	0.41
1:C:134:ILE:CG2	1:C:138:MET:HE2	2.49	0.41
1:D:72:ALA:HA	1:D:73:TYR:HA	1.83	0.41
1:E:50:LEU:HD11	1:E:54:TYR:HE1	1.85	0.41
1:H:124:VAL:HG12	1:H:152:LEU:HB3	2.01	0.41
1:D:69:PHE:HD2	1:D:123:LEU:HD23	1.83	0.41
1:B:128:ILE:HG21	1:B:168:ILE:HG13	2.02	0.41
1:C:72:ALA:HA	1:C:73:TYR:HA	1.82	0.41
1:B:144:ASN:O	1:B:145:LYS:HB2	2.20	0.41
1:F:15:GLU:OE2	1:F:16:LYS:NZ	2.53	0.41
1:F:63:ILE:CG2	1:F:122:MET:HE1	2.51	0.41
1:G:128:ILE:HG13	1:G:153:VAL:HG11	2.02	0.41
1:A:203:GLU:OE1	1:A:203:GLU:HA	2.21	0.41
1:A:208:TYR:CD1	1:C:33:PRO:HD3	2.56	0.41
1:C:120:ARG:HB3	1:C:149:ALA:HB2	2.02	0.41
1:D:125:ASP:O	1:D:153:VAL:HA	2.20	0.41
1:E:126:ASP:OD1	1:E:127:VAL:N	2.49	0.41
1:G:141:ILE:HG23	1:G:146:ALA:HB3	2.02	0.41
1:G:5:TYR:CD1	1:G:5:TYR:C	2.99	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:72:ALA:HB2	1:E:99:ASN:O	2.20	0.40
1:E:142:GLN:HG3	1:E:143:ALA:N	2.35	0.40
1:C:74[B]:LYS:HE3	1:D:100:ARG:HH22	1.86	0.40
1:D:22:GLY:O	1:D:33:PRO:HA	2.22	0.40
1:G:202:LEU:HD21	1:G:206:LYS:CE	2.50	0.40
1:C:71:PRO:HG3	1:C:123:LEU:HD21	2.02	0.40
1:C:115:SER:HB2	1:D:42:ASN:HB2	2.03	0.40
1:F:71:PRO:HD2	1:F:75:GLY:HA3	2.04	0.40
1:B:168:ILE:HG23	1:B:179:VAL:CB	2.49	0.40
1:D:155:ILE:HD12	1:D:188:LEU:HD11	2.03	0.40
1:A:101:LYS:C	1:A:102:GLU:HG3	2.46	0.40
1:D:155:ILE:HA	1:D:183:VAL:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	204/238 (86%)	197 (97%)	5 (2%)	2 (1%)	12	9
1	B	168/238 (71%)	164 (98%)	3 (2%)	1 (1%)	21	18
1	C	186/238 (78%)	182 (98%)	2 (1%)	2 (1%)	11	8
1	D	160/238 (67%)	157 (98%)	3 (2%)	0	100	100
1	E	204/238 (86%)	196 (96%)	7 (3%)	1 (0%)	24	22
1	F	172/238 (72%)	171 (99%)	1 (1%)	0	100	100
1	G	187/238 (79%)	185 (99%)	1 (0%)	1 (0%)	24	22
1	H	160/238 (67%)	159 (99%)	1 (1%)	0	100	100
All	All	1441/1904 (76%)	1411 (98%)	23 (2%)	7 (0%)	24	22

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	C	196	GLY
1	A	197	ASN
1	E	197	ASN
1	G	197	ASN
1	C	166	SER
1	B	131	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	168/191 (88%)	163 (97%)	5 (3%)	36	41
1	B	144/191 (75%)	143 (99%)	1 (1%)	76	83
1	C	162/191 (85%)	156 (96%)	6 (4%)	30	33
1	D	139/191 (73%)	134 (96%)	5 (4%)	31	34
1	E	168/191 (88%)	167 (99%)	1 (1%)	78	86
1	F	145/191 (76%)	142 (98%)	3 (2%)	47	54
1	G	161/191 (84%)	155 (96%)	6 (4%)	30	33
1	H	140/191 (73%)	137 (98%)	3 (2%)	47	54
All	All	1227/1528 (80%)	1197 (98%)	30 (2%)	47	49

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	132	THR
1	A	152	LEU
1	A	180	ILE
1	A	195	GLN
1	A	214	ILE
1	B	102	GLU
1	C	74[A]	LYS
1	C	74[B]	LYS
1	C	96	TYR

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Mol	Chain	Res	Type
1	C	134	ILE
1	C	142	GLN
1	C	189	ILE
1	D	96	TYR
1	D	134	ILE
1	D	157	ARG
1	D	172	GLU
1	D	173	ARG
1	E	168	ILE
1	F	96[A]	TYR
1	F	96[B]	TYR
1	F	138	MET
1	G	7[A]	ARG
1	G	7[B]	ARG
1	G	96	TYR
1	G	134	ILE
1	G	173	ARG
1	G	202	LEU
1	H	96[A]	TYR
1	H	96[B]	TYR
1	H	155	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	GLN
1	C	17	GLN
1	C	158	GLN
1	C	211	GLN
1	E	6	GLN
1	E	142	GLN
1	E	195	GLN
1	G	158	GLN
1	G	169	GLN
1	G	195	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	205/238 (86%)	0.88	30 (14%) 6 6	14, 40, 86, 96	3 (1%)
1	B	175/238 (73%)	0.66	16 (9%) 15 15	15, 38, 70, 86	1 (0%)
1	C	193/238 (81%)	1.03	29 (15%) 5 5	18, 48, 78, 92	3 (1%)
1	D	170/238 (71%)	0.77	19 (11%) 10 10	16, 41, 74, 87	0
1	E	205/238 (86%)	1.00	28 (13%) 6 7	18, 41, 84, 95	3 (1%)
1	F	177/238 (74%)	0.62	22 (12%) 8 8	14, 39, 76, 96	1 (0%)
1	G	194/238 (81%)	1.03	31 (15%) 5 5	15, 50, 79, 93	3 (1%)
1	H	169/238 (71%)	0.76	17 (10%) 12 13	15, 41, 66, 84	1 (0%)
All	All	1488/1904 (78%)	0.85	192 (12%) 7 8	14, 42, 80, 96	15 (1%)

All (192) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	132	THR	6.7
1	C	214	ILE	6.6
1	B	130	ALA	6.6
1	D	168	ILE	6.0
1	C	129	THR	5.5
1	D	30	ARG	5.3
1	G	31	LYS	4.8
1	A	214	ILE	4.7
1	H	168	ILE	4.7
1	E	132	THR	4.6
1	G	133	ALA	4.5
1	A	162	LYS	4.4
1	E	29	GLY	4.3
1	C	142	GLN	4.3
1	A	129	THR	4.2
1	D	25	THR	4.2

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Mol	Chain	Res	Type	RSRZ
1	E	214	ILE	4.2
1	F	167	ALA	4.2
1	G	189	ILE	4.1
1	D	158	GLN	4.1
1	H	25	THR	4.1
1	F	30	ARG	4.1
1	A	143	ALA	4.0
1	B	167	ALA	4.0
1	F	158	GLN	4.0
1	A	27	LYS	3.9
1	E	164	GLU	3.9
1	C	165	LEU	3.9
1	A	163	GLY	3.8
1	F	26	LEU	3.8
1	F	130	ALA	3.8
1	F	156	ASP	3.8
1	E	129	THR	3.8
1	B	26	LEU	3.7
1	D	171	VAL	3.7
1	E	196	GLY	3.7
1	G	24	PHE	3.7
1	H	31	LYS	3.7
1	D	175	PHE	3.7
1	B	27	LYS	3.7
1	A	24	PHE	3.6
1	G	166	SER	3.6
1	F	27	LYS	3.6
1	D	129	THR	3.6
1	G	197	ASN	3.6
1	A	196	GLY	3.6
1	C	26	LEU	3.5
1	H	145	LYS	3.5
1	E	25	THR	3.5
1	H	158	GLN	3.4
1	G	127	VAL	3.4
1	C	24	PHE	3.4
1	C	25	THR	3.4
1	B	170	GLU	3.3
1	B	172	GLU	3.3
1	A	29	GLY	3.3
1	A	130	ALA	3.2
1	B	1	MET	3.2

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Mol	Chain	Res	Type	RSRZ
1	E	197	ASN	3.2
1	C	35	PHE	3.2
1	E	210	ALA	3.1
1	E	24	PHE	3.1
1	A	132	THR	3.1
1	A	183	VAL	3.1
1	C	134	ILE	3.1
1	G	141	ILE	3.1
1	G	161	GLY	3.0
1	F	173	ARG	3.0
1	G	26	LEU	3.0
1	G	74[A]	LYS	3.0
1	A	25	THR	3.0
1	G	35	PHE	3.0
1	C	167	ALA	3.0
1	D	167	ALA	3.0
1	D	173	ARG	3.0
1	A	164	GLU	3.0
1	E	163	GLY	3.0
1	A	133	ALA	2.9
1	E	35[A]	PHE	2.9
1	E	28	SER	2.9
1	G	142	GLN	2.9
1	G	129	THR	2.9
1	E	180	ILE	2.8
1	F	155	ILE	2.8
1	A	28	SER	2.8
1	A	194	GLN	2.8
1	G	168	ILE	2.8
1	H	175	PHE	2.8
1	A	23	GLU	2.8
1	E	205	VAL	2.8
1	A	180	ILE	2.8
1	C	111	ASN	2.8
1	C	197	ASN	2.7
1	A	146	ALA	2.7
1	H	155	ILE	2.7
1	C	160	LYS	2.7
1	G	145	LYS	2.7
1	G	25	THR	2.7
1	C	177	CYS	2.7
1	E	162	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	22	GLY	2.7
1	G	196	GLY	2.6
1	A	210	ALA	2.6
1	A	198	ASN	2.6
1	H	167	ALA	2.6
1	A	202	LEU	2.6
1	D	31	LYS	2.6
1	B	30	ARG	2.6
1	B	153	VAL	2.6
1	F	168	ILE	2.6
1	D	136	GLU	2.6
1	C	3	LYS	2.6
1	B	123	LEU	2.6
1	G	14	LEU	2.6
1	A	195	GLN	2.5
1	G	214	ILE	2.5
1	E	161	GLY	2.5
1	A	197	ASN	2.5
1	C	178	ALA	2.5
1	G	178	ALA	2.5
1	G	191	TYR	2.5
1	G	116	LYS	2.5
1	G	136	GLU	2.5
1	B	131	GLY	2.5
1	G	198	ASN	2.5
1	C	128	ILE	2.4
1	H	169	GLN	2.4
1	C	171	VAL	2.4
1	C	32	SER	2.4
1	F	25	THR	2.4
1	F	29	GLY	2.4
1	A	111	ASN	2.4
1	A	199	THR	2.4
1	C	173	ARG	2.4
1	A	26	LEU	2.4
1	F	111	ASN	2.4
1	E	128	ILE	2.4
1	D	101	LYS	2.4
1	E	199	THR	2.4
1	A	139	GLU	2.3
1	F	146	ALA	2.3
1	E	198	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	186	THR	2.3
1	G	144	ASN	2.3
1	C	135	ARG	2.3
1	C	141	ILE	2.3
1	H	186	THR	2.3
1	E	139	GLU	2.2
1	D	169	GLN	2.2
1	E	194	GLN	2.2
1	B	132	THR	2.2
1	D	15	GLU	2.2
1	B	169	GLN	2.2
1	D	194	GLN	2.2
1	G	135	ARG	2.2
1	H	128	ILE	2.2
1	E	26	LEU	2.2
1	F	132	THR	2.2
1	H	129	THR	2.2
1	C	196	GLY	2.2
1	D	29	GLY	2.2
1	D	145	LYS	2.2
1	E	3[A]	LYS	2.2
1	H	3	LYS	2.2
1	B	129	THR	2.2
1	B	15[A]	GLU	2.1
1	F	24	PHE	2.1
1	F	136	GLU	2.1
1	C	194	GLN	2.1
1	H	133	ALA	2.1
1	G	177	CYS	2.1
1	A	136	GLU	2.1
1	F	102	GLU	2.1
1	G	194	GLN	2.1
1	F	175	PHE	2.1
1	C	22	GLY	2.1
1	E	143	ALA	2.1
1	F	101	LYS	2.1
1	C	166	SER	2.1
1	E	23	GLU	2.1
1	D	172	GLU	2.1
1	F	170	GLU	2.0
1	H	111	ASN	2.0
1	F	169	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	132	THR	2.0
1	G	199	THR	2.0
1	G	128	ILE	2.0
1	C	137	SER	2.0
1	B	111	ASN	2.0
1	H	194	GLN	2.0
1	E	209	ARG	2.0
1	C	212	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CL	H	215	1/1	0.91	0.17	62,62,62,62	0
2	CL	H	216	1/1	0.93	0.15	61,61,61,61	0
2	CL	A	215	1/1	0.98	0.06	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.