



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:46 PM GMT

PDB ID : 4JFR
Title : Crystal structure of anabolic ornithine carbamoyltransferase from *Vibrio vulnificus* in complex with carbamoyl phosphate
Authors : Shabalin, I.G.; Winsor, J.; Grimshaw, S.; Osinski, T.; Bajor, J.; Chordia, M.D.; Shuvalova, L.; Anderson, W.F.; Minor, W.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2013-02-28
Resolution : 2.17 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

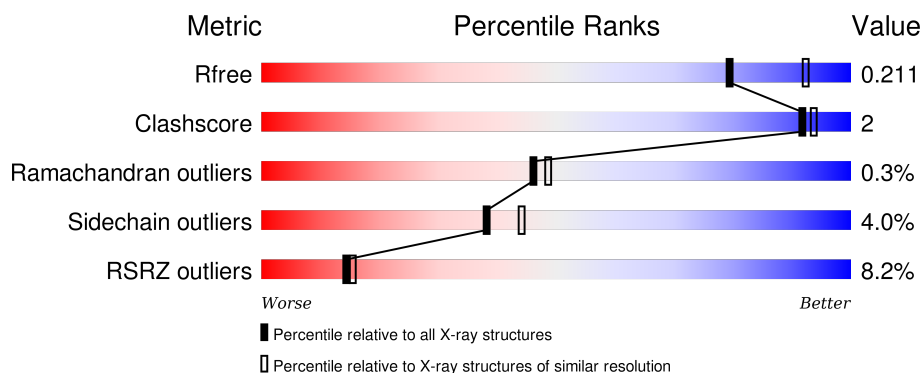
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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}	91344	5130 (2.20-2.16)
Clashscore	102246	5965 (2.20-2.16)
Ramachandran outliers	100387	5863 (2.20-2.16)
Sidechain outliers	100360	5864 (2.20-2.16)
RSRZ outliers	91569	5142 (2.20-2.16)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	358	<div> <div>2%</div> <div>86%</div> <div>7%</div> <div>6%</div> </div>
1	B	358	<div> <div>12%</div> <div>88%</div> <div>6%</div> <div>5%</div> </div>
1	C	358	<div> <div>9%</div> <div>84%</div> <div>9%</div> <div>7%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8086 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 Ornithine carbamoyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	1	0
			2600	1648	439	495	18			
1	B	340	Total	C	N	O	S	0	0	0
			2621	1661	443	499	18			
1	C	334	Total	C	N	O	S	0	0	0
			2572	1628	434	492	18			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	EXPRESSION TAG	UNP Q8DCF5
A	-22	HIS	-	EXPRESSION TAG	UNP Q8DCF5
A	-21	HIS	-	EXPRESSION TAG	UNP Q8DCF5
A	-20	HIS	-	EXPRESSION TAG	UNP Q8DCF5
A	-19	HIS	-	EXPRESSION TAG	UNP Q8DCF5
A	-18	HIS	-	EXPRESSION TAG	UNP Q8DCF5
A	-17	HIS	-	EXPRESSION TAG	UNP Q8DCF5
A	-16	SER	-	EXPRESSION TAG	UNP Q8DCF5
A	-15	SER	-	EXPRESSION TAG	UNP Q8DCF5
A	-14	GLY	-	EXPRESSION TAG	UNP Q8DCF5
A	-13	VAL	-	EXPRESSION TAG	UNP Q8DCF5
A	-12	ASP	-	EXPRESSION TAG	UNP Q8DCF5
A	-11	LEU	-	EXPRESSION TAG	UNP Q8DCF5
A	-10	GLY	-	EXPRESSION TAG	UNP Q8DCF5
A	-9	THR	-	EXPRESSION TAG	UNP Q8DCF5
A	-8	GLU	-	EXPRESSION TAG	UNP Q8DCF5
A	-7	ASN	-	EXPRESSION TAG	UNP Q8DCF5
A	-6	LEU	-	EXPRESSION TAG	UNP Q8DCF5
A	-5	TYR	-	EXPRESSION TAG	UNP Q8DCF5
A	-4	PHE	-	EXPRESSION TAG	UNP Q8DCF5
A	-3	GLN	-	EXPRESSION TAG	UNP Q8DCF5
A	-2	SER	-	EXPRESSION TAG	UNP Q8DCF5
A	-1	ASN	-	EXPRESSION TAG	UNP Q8DCF5

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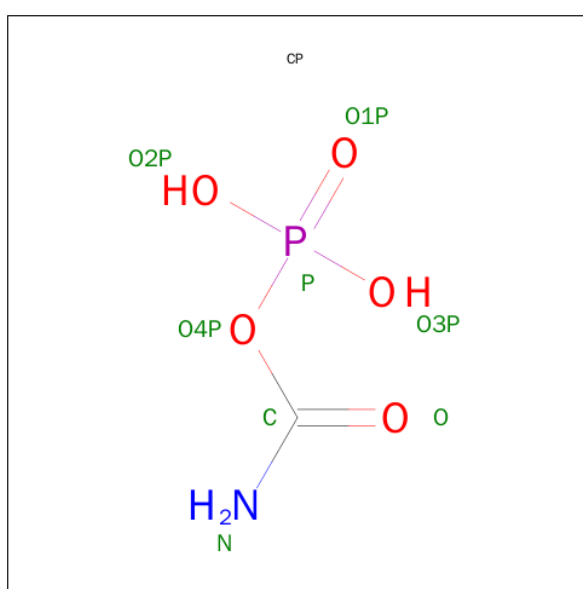
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	EXPRESSION TAG	UNP Q8DCF5
B	-23	MET	-	EXPRESSION TAG	UNP Q8DCF5
B	-22	HIS	-	EXPRESSION TAG	UNP Q8DCF5
B	-21	HIS	-	EXPRESSION TAG	UNP Q8DCF5
B	-20	HIS	-	EXPRESSION TAG	UNP Q8DCF5
B	-19	HIS	-	EXPRESSION TAG	UNP Q8DCF5
B	-18	HIS	-	EXPRESSION TAG	UNP Q8DCF5
B	-17	HIS	-	EXPRESSION TAG	UNP Q8DCF5
B	-16	SER	-	EXPRESSION TAG	UNP Q8DCF5
B	-15	SER	-	EXPRESSION TAG	UNP Q8DCF5
B	-14	GLY	-	EXPRESSION TAG	UNP Q8DCF5
B	-13	VAL	-	EXPRESSION TAG	UNP Q8DCF5
B	-12	ASP	-	EXPRESSION TAG	UNP Q8DCF5
B	-11	LEU	-	EXPRESSION TAG	UNP Q8DCF5
B	-10	GLY	-	EXPRESSION TAG	UNP Q8DCF5
B	-9	THR	-	EXPRESSION TAG	UNP Q8DCF5
B	-8	GLU	-	EXPRESSION TAG	UNP Q8DCF5
B	-7	ASN	-	EXPRESSION TAG	UNP Q8DCF5
B	-6	LEU	-	EXPRESSION TAG	UNP Q8DCF5
B	-5	TYR	-	EXPRESSION TAG	UNP Q8DCF5
B	-4	PHE	-	EXPRESSION TAG	UNP Q8DCF5
B	-3	GLN	-	EXPRESSION TAG	UNP Q8DCF5
B	-2	SER	-	EXPRESSION TAG	UNP Q8DCF5
B	-1	ASN	-	EXPRESSION TAG	UNP Q8DCF5
B	0	ALA	-	EXPRESSION TAG	UNP Q8DCF5
C	-23	MET	-	EXPRESSION TAG	UNP Q8DCF5
C	-22	HIS	-	EXPRESSION TAG	UNP Q8DCF5
C	-21	HIS	-	EXPRESSION TAG	UNP Q8DCF5
C	-20	HIS	-	EXPRESSION TAG	UNP Q8DCF5
C	-19	HIS	-	EXPRESSION TAG	UNP Q8DCF5
C	-18	HIS	-	EXPRESSION TAG	UNP Q8DCF5
C	-17	HIS	-	EXPRESSION TAG	UNP Q8DCF5
C	-16	SER	-	EXPRESSION TAG	UNP Q8DCF5
C	-15	SER	-	EXPRESSION TAG	UNP Q8DCF5
C	-14	GLY	-	EXPRESSION TAG	UNP Q8DCF5
C	-13	VAL	-	EXPRESSION TAG	UNP Q8DCF5
C	-12	ASP	-	EXPRESSION TAG	UNP Q8DCF5
C	-11	LEU	-	EXPRESSION TAG	UNP Q8DCF5
C	-10	GLY	-	EXPRESSION TAG	UNP Q8DCF5
C	-9	THR	-	EXPRESSION TAG	UNP Q8DCF5
C	-8	GLU	-	EXPRESSION TAG	UNP Q8DCF5
C	-7	ASN	-	EXPRESSION TAG	UNP Q8DCF5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	LEU	-	EXPRESSION TAG	UNP Q8DCF5
C	-5	TYR	-	EXPRESSION TAG	UNP Q8DCF5
C	-4	PHE	-	EXPRESSION TAG	UNP Q8DCF5
C	-3	GLN	-	EXPRESSION TAG	UNP Q8DCF5
C	-2	SER	-	EXPRESSION TAG	UNP Q8DCF5
C	-1	ASN	-	EXPRESSION TAG	UNP Q8DCF5
C	0	ALA	-	EXPRESSION TAG	UNP Q8DCF5

- Molecule 2 is PHOSPHORIC ACID MONO(FORMAMIDE)ESTER (three-letter code: CP) (formula: $\text{CH}_4\text{NO}_5\text{P}$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			8	1	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			8	1	1	5	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total 1	Mg 1	0	0

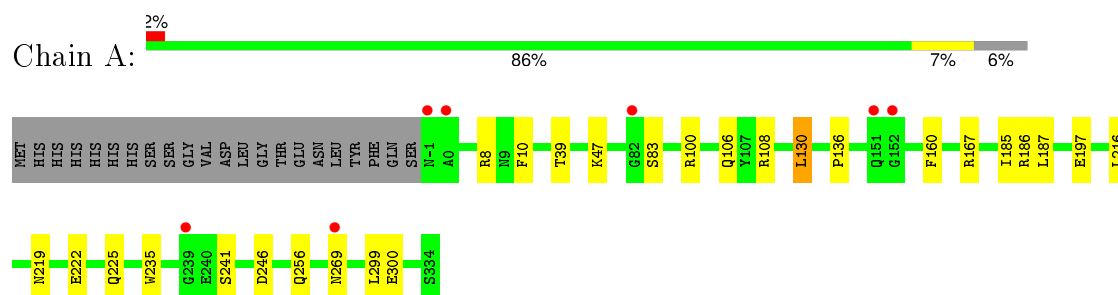
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	130	Total 130	O 130	0	0
5	B	72	Total 72	O 72	0	0
5	C	70	Total 71	O 71	0	1

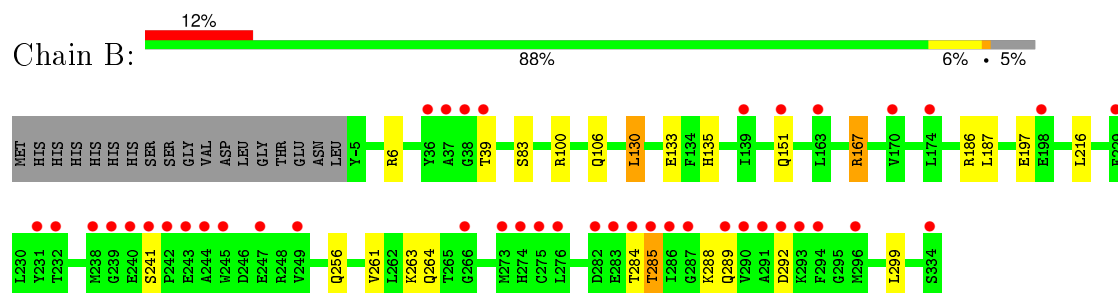
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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.

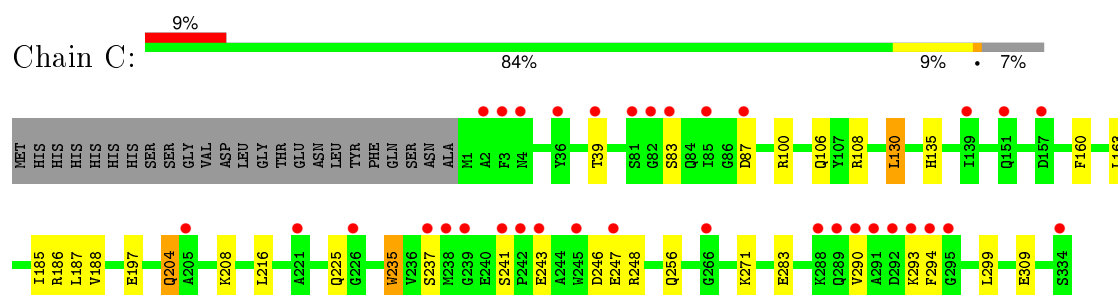
• Molecule 1: Ornithine carbamoyltransferase



• Molecule 1: Ornithine carbamoyltransferase



• Molecule 1: Ornithine carbamoyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.34Å 80.92Å 177.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.96 – 2.17 38.93 – 2.17	Depositor EDS
% Data completeness (in resolution range)	99.5 (38.96-2.17) 99.6 (38.93-2.17)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.18Å)	Xtriage
Refinement program	HKL-3000, REFMAC 5.7.0029, COOT	Depositor
R, R_{free}	0.171 , 0.205 0.178 , 0.211	Depositor DCC
R_{free} test set	3185 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	44.3	Xtriage
Anisotropy	0.445	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.3	EDS
Estimated twinning fraction	0.022 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 62540 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8086	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, CP, 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	0.82	0/2651	0.86	5/3577 (0.1%)
1	B	0.80	0/2670	0.82	5/3603 (0.1%)
1	C	0.78	0/2619	0.87	6/3536 (0.2%)
All	All	0.80	0/7940	0.85	16/10716 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	108	ARG	NE-CZ-NH1	-9.84	115.38	120.30
1	C	108	ARG	NE-CZ-NH2	8.31	124.45	120.30
1	C	100	ARG	NE-CZ-NH2	-7.35	116.63	120.30
1	C	100	ARG	NE-CZ-NH1	7.24	123.92	120.30
1	A	100	ARG	NE-CZ-NH2	-6.83	116.88	120.30
1	A	100	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	C	186	ARG	NE-CZ-NH1	6.30	123.45	120.30
1	A	108	ARG	NE-CZ-NH2	-6.29	117.15	120.30
1	B	100	ARG	NE-CZ-NH2	-5.72	117.44	120.30
1	A	8	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	C	186	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	B	133	GLU	OE1-CD-OE2	-5.32	116.92	123.30
1	A	186	ARG	NE-CZ-NH1	5.27	122.94	120.30
1	B	186	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	100	ARG	NE-CZ-NH1	5.06	122.83	120.30
1	B	186	ARG	NE-CZ-NH2	-5.02	117.79	120.30

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	2600	0	2562	9	0
1	B	2621	0	2575	10	0
1	C	2572	0	2522	14	0
2	A	8	0	2	0	0
2	B	8	0	2	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	C	1	0	0	0	0
5	A	130	0	0	2	0
5	B	72	0	0	0	0
5	C	71	0	0	1	0
All	All	8086	0	7663	32	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 2.

All (32) 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:B:285:THR:HG23	1:C:87:ASP:HA	1.55	0.86
1:B:167:ARG:C	1:B:167:ARG:HD2	2.00	0.82
1:B:288:LYS:HD2	1:B:292:ASP:OD2	1.85	0.76
1:A:197:GLU:HG3	5:A:629:HOH:O	2.01	0.61
1:C:197:GLU:N	1:C:197:GLU:OE1	2.33	0.60
1:C:243:GLU:HA	1:C:293:LYS:HE2	1.87	0.57
1:C:290:VAL:HG11	1:C:299:LEU:HD21	1.86	0.56
1:C:204:GLN:O	1:C:208:LYS:HG2	2.07	0.55
1:A:256:GLN:HG3	1:A:299:LEU:O	2.08	0.53
1:C:256:GLN:HG3	1:C:299:LEU:O	2.09	0.51
1:C:187:LEU:HD12	1:C:216:LEU:CD2	2.42	0.50
1:A:219:ASN:HB3	1:A:222:GLU:HG2	1.93	0.50
1:B:256:GLN:HG3	1:B:299:LEU:O	2.13	0.49
1:C:235:TRP:O	1:C:248:ARG:NH1	2.45	0.49
1:A:235:TRP:CZ2	1:A:300:GLU:HA	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:LEU:HD12	1:B:216:LEU:CD2	2.44	0.48
1:A:130:LEU:C	1:A:130:LEU:HD23	2.35	0.46
1:A:187:LEU:HD12	1:A:216:LEU:CD2	2.45	0.46
1:B:167:ARG:C	1:B:167:ARG:CD	2.80	0.45
1:A:160:PHE:CZ	1:A:185:ILE:HD12	2.52	0.45
1:B:261:VAL:HA	1:B:264:GLN:HE21	1.81	0.45
1:B:130:LEU:C	1:B:130:LEU:HD23	2.37	0.45
1:B:197:GLU:H	1:B:197:GLU:CD	2.21	0.43
1:C:290:VAL:CG1	1:C:299:LEU:HD21	2.49	0.43
1:A:47:LYS:HE3	5:A:501:HOH:O	2.18	0.42
1:C:160:PHE:CZ	1:C:185:ILE:HD12	2.54	0.42
1:C:225:GLN:NE2	5:C:517:HOH:O	2.44	0.42
1:C:130:LEU:HD23	1:C:130:LEU:C	2.40	0.42
1:B:167:ARG:HD2	1:B:167:ARG:O	2.16	0.42
1:C:246:ASP:HA	1:C:294:PHE:CE1	2.54	0.42
1:A:10:PHE:CE2	1:A:136:PRO:HB2	2.55	0.41
1:C:163:LEU:HA	1:C:188:VAL:O	2.21	0.41

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	335/358 (94%)	326 (97%)	8 (2%)	1 (0%)	46	48
1	B	338/358 (94%)	328 (97%)	9 (3%)	1 (0%)	46	48
1	C	332/358 (93%)	324 (98%)	7 (2%)	1 (0%)	46	48
All	All	1005/1074 (94%)	978 (97%)	24 (2%)	3 (0%)	46	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	130	LEU
1	B	130	LEU
1	C	130	LEU

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	270/292 (92%)	262 (97%)	8 (3%)	48	57
1	B	272/292 (93%)	260 (96%)	12 (4%)	35	39
1	C	266/292 (91%)	254 (96%)	12 (4%)	34	38
All	All	808/876 (92%)	776 (96%)	32 (4%)	38	44

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	THR
1	A	83	SER
1	A	106	GLN
1	A	167	ARG
1	A	225	GLN
1	A	241	SER
1	A	246	ASP
1	A	269	ASN
1	B	6	ARG
1	B	39	THR
1	B	83	SER
1	B	106	GLN
1	B	135	HIS
1	B	151	GLN
1	B	167	ARG
1	B	241	SER
1	B	263	LYS
1	B	284	THR
1	B	285	THR
1	B	289	GLN

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Mol	Chain	Res	Type
1	C	39	THR
1	C	83	SER
1	C	106	GLN
1	C	135	HIS
1	C	204	GLN
1	C	235	TRP
1	C	237	SER
1	C	241	SER
1	C	247	GLU
1	C	271	LYS
1	C	283	GLU
1	C	309	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	264	GLN
1	C	264	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length

(or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	CP	A	401	-	6,7,7	2.19	2 (33%)	6,10,10	2.36	2 (33%)
2	CP	B	401	-	6,7,7	2.45	1 (16%)	6,10,10	2.53	3 (50%)

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	CP	A	401	-	-	0/3/5/5	0/0/0/0
2	CP	B	401	-	-	0/3/5/5	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	CP	O-C	3.59	1.27	1.21
2	A	401	CP	O4P-C	3.82	1.43	1.38
2	B	401	CP	O4P-C	5.79	1.46	1.38

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	CP	O2P-P-O4P	-3.76	92.98	105.25
2	B	401	CP	O2P-P-O4P	-3.11	95.09	105.25
2	B	401	CP	O3P-P-O1P	-2.01	104.09	110.58
2	A	401	CP	O2P-P-O1P	3.83	122.92	110.58
2	B	401	CP	O3P-P-O2P	4.71	125.32	107.38

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	336/358 (93%)	0.01	7 (2%) 67 68	34, 49, 75, 94	0
1	B	340/358 (94%)	0.47	42 (12%) 5 6	36, 58, 103, 128	0
1	C	334/358 (93%)	0.45	34 (10%) 9 10	35, 60, 91, 115	0
All	All	1010/1074 (94%)	0.31	83 (8%) 14 15	34, 55, 91, 128	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	290	VAL	7.7
1	B	241	SER	7.1
1	B	245	TRP	7.0
1	B	239	GLY	6.6
1	C	239	GLY	5.3
1	C	238	MET	5.3
1	B	242	PRO	5.3
1	C	334	SER	5.2
1	B	240	GLU	5.1
1	C	294	PHE	4.8
1	B	247	GLU	4.8
1	B	287	GLY	4.6
1	C	245	TRP	4.6
1	B	286	ILE	4.3
1	C	85	ILE	4.3
1	A	239	GLY	4.2
1	C	82	GLY	4.2
1	A	152	GLY	4.1
1	B	37	ALA	4.1
1	B	289	GLN	3.9
1	B	139	ILE	3.9
1	C	292	ASP	3.9
1	B	293	LYS	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	334	SER	3.7
1	B	231	TYR	3.7
1	C	83	SER	3.7
1	C	2	ALA	3.6
1	C	139	ILE	3.5
1	B	39	THR	3.5
1	B	294	PHE	3.5
1	C	241	SER	3.4
1	C	295	GLY	3.2
1	C	237	SER	3.2
1	C	151	GLN	3.2
1	A	0	ALA	3.2
1	C	242	PRO	3.2
1	B	249	VAL	3.1
1	B	170	VAL	3.0
1	A	82	GLY	3.0
1	C	293	LYS	3.0
1	C	36	TYR	2.9
1	B	283	GLU	2.9
1	A	-1	ASN	2.9
1	B	292	ASP	2.9
1	C	247	GLU	2.9
1	B	291	ALA	2.9
1	B	274	HIS	2.8
1	C	87	ASP	2.8
1	B	284	THR	2.8
1	A	151	GLN	2.8
1	B	275	CYS	2.8
1	C	221	ALA	2.8
1	C	243	GLU	2.7
1	C	226	GLY	2.7
1	B	244	ALA	2.7
1	C	4	ASN	2.7
1	B	174	LEU	2.7
1	B	243	GLU	2.6
1	C	289	GLN	2.6
1	B	285	THR	2.6
1	B	273	MET	2.6
1	B	38	GLY	2.5
1	B	296	MET	2.5
1	C	291	ALA	2.5
1	B	238	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	282	ASP	2.4
1	C	290	VAL	2.4
1	B	151	GLN	2.3
1	C	81	SER	2.3
1	B	163	LEU	2.3
1	C	205	ALA	2.3
1	B	276	LEU	2.2
1	C	3	PHE	2.2
1	B	198	GLU	2.2
1	B	266	GLY	2.1
1	C	288	LYS	2.1
1	B	229	PHE	2.1
1	B	232	THR	2.0
1	C	39	THR	2.0
1	B	36	TYR	2.0
1	A	269	ASN	2.0
1	C	266	GLY	2.0
1	C	157	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	CL	C	401	1/1	0.98	0.15	0.26	49,49,49,49	0
3	CL	A	402	1/1	0.98	0.10	-1.16	37,37,37,37	0
3	CL	B	402	1/1	0.97	0.06	-2.03	44,44,44,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CP	A	401	8/8	0.99	0.09	-2.21	37,40,42,46	0
2	CP	B	401	8/8	0.98	0.06	-2.28	41,54,57,59	0
4	MG	C	402	1/1	0.86	0.14	-	52,52,52,52	1

6.5 Other polymers [i](#)

There are no such residues in this entry.