



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:44 PM UTC

PDB ID : 4DDF / pdb_00004ddf
Title : Computationally Designed Self-assembling Octahedral Cage protein, O333,
Crystallized in space group P4
Authors : Sawaya, M.R.; King, N.P.; Sheffler, W.; Baker, D.; Yeates, T.O.
Deposited on : 2012-01-18
Resolution : 3.15 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

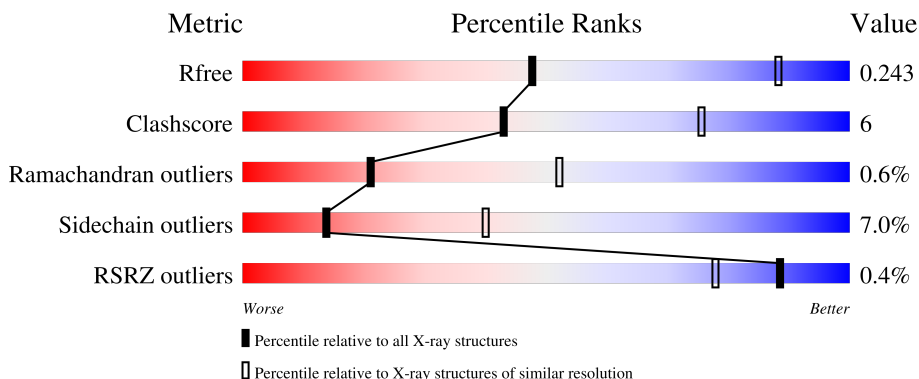
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	192	 72% 21% • 5%
1	B	192	 77% 18% • 5%
1	C	192	 71% 23% • 5%
1	D	192	 71% 21% • 5%
1	E	192	 73% 18% • 5%

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Mol	Chain	Length	Quality of chain
1	F	192	<p>% 79% 15% • 5%</p>
1	G	192	<p>76% 18% •• 5%</p>
1	H	192	<p>% 79% 15% • 5%</p>
1	I	192	<p>75% 18% • 5%</p>
1	J	192	<p>70% 22% •• 5%</p>
1	K	192	<p>70% 19% 5% • 5%</p>
1	L	192	<p>% 77% 16% •• 5%</p>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 15798 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 Propanediol utilization polyhedral body protein PduT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	183	1310	828	222	251	9	0	0	0
1	B	183	1310	828	222	251	9	0	0	0
1	C	183	1310	828	222	251	9	0	0	0
1	D	183	1310	828	222	251	9	0	0	0
1	E	183	1310	828	222	251	9	0	0	0
1	F	183	1310	828	222	251	9	0	0	0
1	G	183	1310	828	222	251	9	0	0	0
1	H	183	1310	828	222	251	9	0	0	0
1	I	183	1310	828	222	251	9	0	0	0
1	J	183	1310	828	222	251	9	0	0	0
1	K	183	1310	828	222	251	9	0	0	0
1	L	183	1310	828	222	251	9	0	0	0

There are 216 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	ALA	LYS	engineered mutation	UNP E7V033
A	38	SER	CYS	engineered mutation	UNP E7V033
A	67	LEU	MET	engineered mutation	UNP E7V033
A	148	ALA	ASN	engineered mutation	UNP E7V033
A	149	LEU	ASN	engineered mutation	UNP E7V033

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Chain	Residue	Modelled	Actual	Comment	Reference
A	156	SER	GLU	engineered mutation	UNP E7V033
A	160	ALA	GLU	engineered mutation	UNP E7V033
A	161	TYR	LYS	engineered mutation	UNP E7V033
A	167	ALA	ARG	engineered mutation	UNP E7V033
A	169	LEU	VAL	engineered mutation	UNP E7V033
A	185	LEU	-	expression tag	UNP E7V033
A	186	GLU	-	expression tag	UNP E7V033
A	187	HIS	-	expression tag	UNP E7V033
A	188	HIS	-	expression tag	UNP E7V033
A	189	HIS	-	expression tag	UNP E7V033
A	190	HIS	-	expression tag	UNP E7V033
A	191	HIS	-	expression tag	UNP E7V033
A	192	HIS	-	expression tag	UNP E7V033
B	15	ALA	LYS	engineered mutation	UNP E7V033
B	38	SER	CYS	engineered mutation	UNP E7V033
B	67	LEU	MET	engineered mutation	UNP E7V033
B	148	ALA	ASN	engineered mutation	UNP E7V033
B	149	LEU	ASN	engineered mutation	UNP E7V033
B	156	SER	GLU	engineered mutation	UNP E7V033
B	160	ALA	GLU	engineered mutation	UNP E7V033
B	161	TYR	LYS	engineered mutation	UNP E7V033
B	167	ALA	ARG	engineered mutation	UNP E7V033
B	169	LEU	VAL	engineered mutation	UNP E7V033
B	185	LEU	-	expression tag	UNP E7V033
B	186	GLU	-	expression tag	UNP E7V033
B	187	HIS	-	expression tag	UNP E7V033
B	188	HIS	-	expression tag	UNP E7V033
B	189	HIS	-	expression tag	UNP E7V033
B	190	HIS	-	expression tag	UNP E7V033
B	191	HIS	-	expression tag	UNP E7V033
B	192	HIS	-	expression tag	UNP E7V033
C	15	ALA	LYS	engineered mutation	UNP E7V033
C	38	SER	CYS	engineered mutation	UNP E7V033
C	67	LEU	MET	engineered mutation	UNP E7V033
C	148	ALA	ASN	engineered mutation	UNP E7V033
C	149	LEU	ASN	engineered mutation	UNP E7V033
C	156	SER	GLU	engineered mutation	UNP E7V033
C	160	ALA	GLU	engineered mutation	UNP E7V033
C	161	TYR	LYS	engineered mutation	UNP E7V033
C	167	ALA	ARG	engineered mutation	UNP E7V033
C	169	LEU	VAL	engineered mutation	UNP E7V033
C	185	LEU	-	expression tag	UNP E7V033

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Chain	Residue	Modelled	Actual	Comment	Reference
C	186	GLU	-	expression tag	UNP E7V033
C	187	HIS	-	expression tag	UNP E7V033
C	188	HIS	-	expression tag	UNP E7V033
C	189	HIS	-	expression tag	UNP E7V033
C	190	HIS	-	expression tag	UNP E7V033
C	191	HIS	-	expression tag	UNP E7V033
C	192	HIS	-	expression tag	UNP E7V033
D	15	ALA	LYS	engineered mutation	UNP E7V033
D	38	SER	CYS	engineered mutation	UNP E7V033
D	67	LEU	MET	engineered mutation	UNP E7V033
D	148	ALA	ASN	engineered mutation	UNP E7V033
D	149	LEU	ASN	engineered mutation	UNP E7V033
D	156	SER	GLU	engineered mutation	UNP E7V033
D	160	ALA	GLU	engineered mutation	UNP E7V033
D	161	TYR	LYS	engineered mutation	UNP E7V033
D	167	ALA	ARG	engineered mutation	UNP E7V033
D	169	LEU	VAL	engineered mutation	UNP E7V033
D	185	LEU	-	expression tag	UNP E7V033
D	186	GLU	-	expression tag	UNP E7V033
D	187	HIS	-	expression tag	UNP E7V033
D	188	HIS	-	expression tag	UNP E7V033
D	189	HIS	-	expression tag	UNP E7V033
D	190	HIS	-	expression tag	UNP E7V033
D	191	HIS	-	expression tag	UNP E7V033
D	192	HIS	-	expression tag	UNP E7V033
E	15	ALA	LYS	engineered mutation	UNP E7V033
E	38	SER	CYS	engineered mutation	UNP E7V033
E	67	LEU	MET	engineered mutation	UNP E7V033
E	148	ALA	ASN	engineered mutation	UNP E7V033
E	149	LEU	ASN	engineered mutation	UNP E7V033
E	156	SER	GLU	engineered mutation	UNP E7V033
E	160	ALA	GLU	engineered mutation	UNP E7V033
E	161	TYR	LYS	engineered mutation	UNP E7V033
E	167	ALA	ARG	engineered mutation	UNP E7V033
E	169	LEU	VAL	engineered mutation	UNP E7V033
E	185	LEU	-	expression tag	UNP E7V033
E	186	GLU	-	expression tag	UNP E7V033
E	187	HIS	-	expression tag	UNP E7V033
E	188	HIS	-	expression tag	UNP E7V033
E	189	HIS	-	expression tag	UNP E7V033
E	190	HIS	-	expression tag	UNP E7V033
E	191	HIS	-	expression tag	UNP E7V033

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Chain	Residue	Modelled	Actual	Comment	Reference
E	192	HIS	-	expression tag	UNP E7V033
F	15	ALA	LYS	engineered mutation	UNP E7V033
F	38	SER	CYS	engineered mutation	UNP E7V033
F	67	LEU	MET	engineered mutation	UNP E7V033
F	148	ALA	ASN	engineered mutation	UNP E7V033
F	149	LEU	ASN	engineered mutation	UNP E7V033
F	156	SER	GLU	engineered mutation	UNP E7V033
F	160	ALA	GLU	engineered mutation	UNP E7V033
F	161	TYR	LYS	engineered mutation	UNP E7V033
F	167	ALA	ARG	engineered mutation	UNP E7V033
F	169	LEU	VAL	engineered mutation	UNP E7V033
F	185	LEU	-	expression tag	UNP E7V033
F	186	GLU	-	expression tag	UNP E7V033
F	187	HIS	-	expression tag	UNP E7V033
F	188	HIS	-	expression tag	UNP E7V033
F	189	HIS	-	expression tag	UNP E7V033
F	190	HIS	-	expression tag	UNP E7V033
F	191	HIS	-	expression tag	UNP E7V033
F	192	HIS	-	expression tag	UNP E7V033
G	15	ALA	LYS	engineered mutation	UNP E7V033
G	38	SER	CYS	engineered mutation	UNP E7V033
G	67	LEU	MET	engineered mutation	UNP E7V033
G	148	ALA	ASN	engineered mutation	UNP E7V033
G	149	LEU	ASN	engineered mutation	UNP E7V033
G	156	SER	GLU	engineered mutation	UNP E7V033
G	160	ALA	GLU	engineered mutation	UNP E7V033
G	161	TYR	LYS	engineered mutation	UNP E7V033
G	167	ALA	ARG	engineered mutation	UNP E7V033
G	169	LEU	VAL	engineered mutation	UNP E7V033
G	185	LEU	-	expression tag	UNP E7V033
G	186	GLU	-	expression tag	UNP E7V033
G	187	HIS	-	expression tag	UNP E7V033
G	188	HIS	-	expression tag	UNP E7V033
G	189	HIS	-	expression tag	UNP E7V033
G	190	HIS	-	expression tag	UNP E7V033
G	191	HIS	-	expression tag	UNP E7V033
G	192	HIS	-	expression tag	UNP E7V033
H	15	ALA	LYS	engineered mutation	UNP E7V033
H	38	SER	CYS	engineered mutation	UNP E7V033
H	67	LEU	MET	engineered mutation	UNP E7V033
H	148	ALA	ASN	engineered mutation	UNP E7V033
H	149	LEU	ASN	engineered mutation	UNP E7V033

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Chain	Residue	Modelled	Actual	Comment	Reference
H	156	SER	GLU	engineered mutation	UNP E7V033
H	160	ALA	GLU	engineered mutation	UNP E7V033
H	161	TYR	LYS	engineered mutation	UNP E7V033
H	167	ALA	ARG	engineered mutation	UNP E7V033
H	169	LEU	VAL	engineered mutation	UNP E7V033
H	185	LEU	-	expression tag	UNP E7V033
H	186	GLU	-	expression tag	UNP E7V033
H	187	HIS	-	expression tag	UNP E7V033
H	188	HIS	-	expression tag	UNP E7V033
H	189	HIS	-	expression tag	UNP E7V033
H	190	HIS	-	expression tag	UNP E7V033
H	191	HIS	-	expression tag	UNP E7V033
H	192	HIS	-	expression tag	UNP E7V033
I	15	ALA	LYS	engineered mutation	UNP E7V033
I	38	SER	CYS	engineered mutation	UNP E7V033
I	67	LEU	MET	engineered mutation	UNP E7V033
I	148	ALA	ASN	engineered mutation	UNP E7V033
I	149	LEU	ASN	engineered mutation	UNP E7V033
I	156	SER	GLU	engineered mutation	UNP E7V033
I	160	ALA	GLU	engineered mutation	UNP E7V033
I	161	TYR	LYS	engineered mutation	UNP E7V033
I	167	ALA	ARG	engineered mutation	UNP E7V033
I	169	LEU	VAL	engineered mutation	UNP E7V033
I	185	LEU	-	expression tag	UNP E7V033
I	186	GLU	-	expression tag	UNP E7V033
I	187	HIS	-	expression tag	UNP E7V033
I	188	HIS	-	expression tag	UNP E7V033
I	189	HIS	-	expression tag	UNP E7V033
I	190	HIS	-	expression tag	UNP E7V033
I	191	HIS	-	expression tag	UNP E7V033
I	192	HIS	-	expression tag	UNP E7V033
J	15	ALA	LYS	engineered mutation	UNP E7V033
J	38	SER	CYS	engineered mutation	UNP E7V033
J	67	LEU	MET	engineered mutation	UNP E7V033
J	148	ALA	ASN	engineered mutation	UNP E7V033
J	149	LEU	ASN	engineered mutation	UNP E7V033
J	156	SER	GLU	engineered mutation	UNP E7V033
J	160	ALA	GLU	engineered mutation	UNP E7V033
J	161	TYR	LYS	engineered mutation	UNP E7V033
J	167	ALA	ARG	engineered mutation	UNP E7V033
J	169	LEU	VAL	engineered mutation	UNP E7V033
J	185	LEU	-	expression tag	UNP E7V033

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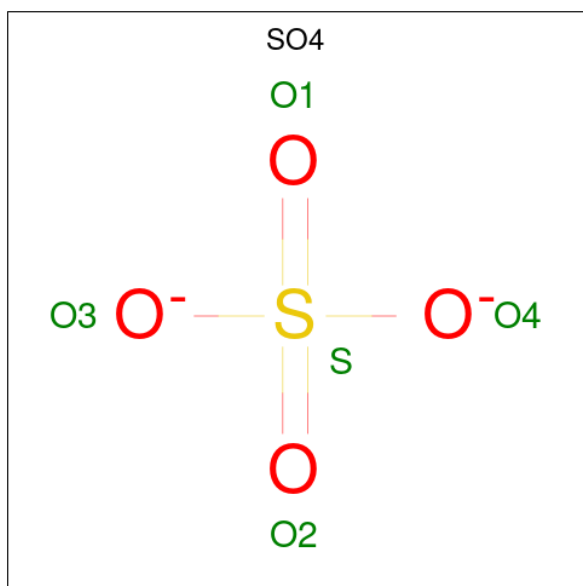
Chain	Residue	Modelled	Actual	Comment	Reference
J	186	GLU	-	expression tag	UNP E7V033
J	187	HIS	-	expression tag	UNP E7V033
J	188	HIS	-	expression tag	UNP E7V033
J	189	HIS	-	expression tag	UNP E7V033
J	190	HIS	-	expression tag	UNP E7V033
J	191	HIS	-	expression tag	UNP E7V033
J	192	HIS	-	expression tag	UNP E7V033
K	15	ALA	LYS	engineered mutation	UNP E7V033
K	38	SER	CYS	engineered mutation	UNP E7V033
K	67	LEU	MET	engineered mutation	UNP E7V033
K	148	ALA	ASN	engineered mutation	UNP E7V033
K	149	LEU	ASN	engineered mutation	UNP E7V033
K	156	SER	GLU	engineered mutation	UNP E7V033
K	160	ALA	GLU	engineered mutation	UNP E7V033
K	161	TYR	LYS	engineered mutation	UNP E7V033
K	167	ALA	ARG	engineered mutation	UNP E7V033
K	169	LEU	VAL	engineered mutation	UNP E7V033
K	185	LEU	-	expression tag	UNP E7V033
K	186	GLU	-	expression tag	UNP E7V033
K	187	HIS	-	expression tag	UNP E7V033
K	188	HIS	-	expression tag	UNP E7V033
K	189	HIS	-	expression tag	UNP E7V033
K	190	HIS	-	expression tag	UNP E7V033
K	191	HIS	-	expression tag	UNP E7V033
K	192	HIS	-	expression tag	UNP E7V033
L	15	ALA	LYS	engineered mutation	UNP E7V033
L	38	SER	CYS	engineered mutation	UNP E7V033
L	67	LEU	MET	engineered mutation	UNP E7V033
L	148	ALA	ASN	engineered mutation	UNP E7V033
L	149	LEU	ASN	engineered mutation	UNP E7V033
L	156	SER	GLU	engineered mutation	UNP E7V033
L	160	ALA	GLU	engineered mutation	UNP E7V033
L	161	TYR	LYS	engineered mutation	UNP E7V033
L	167	ALA	ARG	engineered mutation	UNP E7V033
L	169	LEU	VAL	engineered mutation	UNP E7V033
L	185	LEU	-	expression tag	UNP E7V033
L	186	GLU	-	expression tag	UNP E7V033
L	187	HIS	-	expression tag	UNP E7V033
L	188	HIS	-	expression tag	UNP E7V033
L	189	HIS	-	expression tag	UNP E7V033
L	190	HIS	-	expression tag	UNP E7V033
L	191	HIS	-	expression tag	UNP E7V033

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Chain	Residue	Modelled	Actual	Comment	Reference
L	192	HIS	-	expression tag	UNP E7V033

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	D	1	5	4	1	0	0
2	E	1	5	4	1	0	0
2	G	1	5	4	1	0	0
2	H	1	5	4	1	0	0
2	I	1	5	4	1	0	0
2	J	1	5	4	1	0	0
2	K	1	5	4	1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	2	Total	Cl	0	0
			2	2		
3	C	1	Total	Cl	0	0
			1	1		
3	E	1	Total	Cl	0	0
			1	1		
3	F	1	Total	Cl	0	0
			1	1		
3	G	1	Total	Cl	0	0
			1	1		
3	H	1	Total	Cl	0	0
			1	1		
3	I	1	Total	Cl	0	0
			1	1		
3	J	1	Total	Cl	0	0
			1	1		
3	K	1	Total	Cl	0	0
			1	1		
3	L	1	Total	Cl	0	0
			1	1		

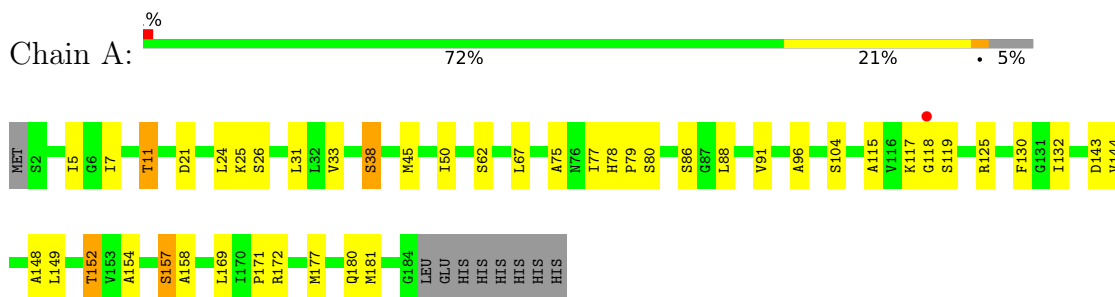
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		
4	D	1	Total	O	0	0
			1	1		
4	E	1	Total	O	0	0
			1	1		
4	K	1	Total	O	0	0
			1	1		
4	L	1	Total	O	0	0
			1	1		

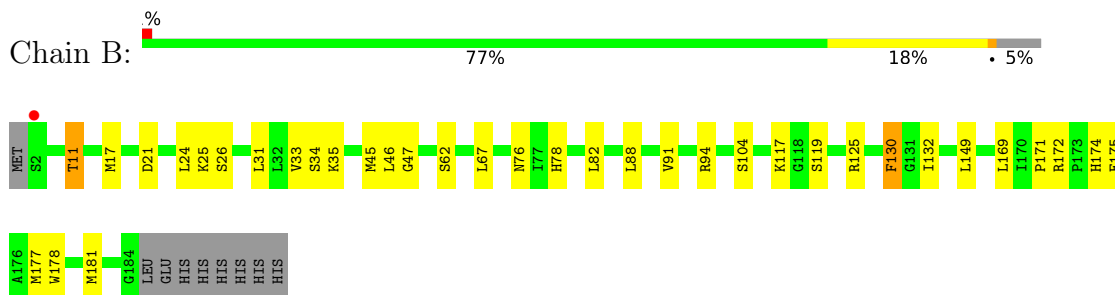
3 Residue-property plots [i](#)

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.

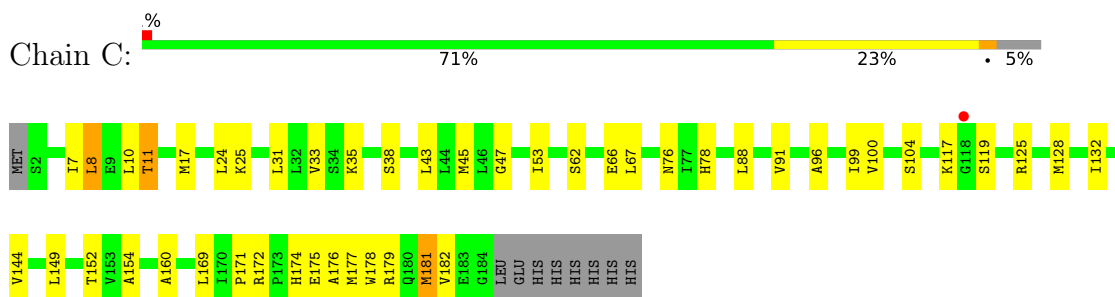
- Molecule 1: Propanediol utilization polyhedral body protein PduT



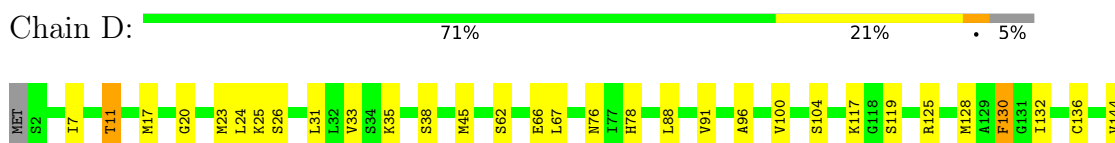
- Molecule 1: Propanediol utilization polyhedral body protein PduT



- Molecule 1: Propanediol utilization polyhedral body protein PduT



- Molecule 1: Propanediol utilization polyhedral body protein PduT

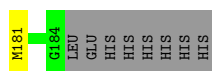
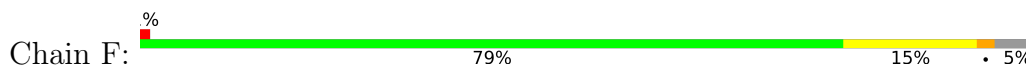




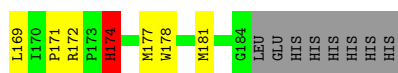
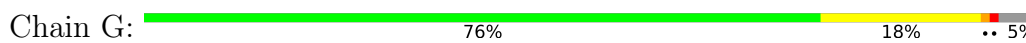
• Molecule 1: Propanediol utilization polyhedral body protein PduT



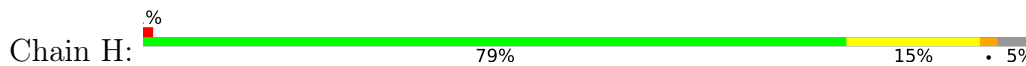
• Molecule 1: Propanediol utilization polyhedral body protein PduT



• Molecule 1: Propanediol utilization polyhedral body protein PduT



• Molecule 1: Propanediol utilization polyhedral body protein PduT



• Molecule 1: Propanediol utilization polyhedral body protein PduT



4 Data and refinement statistics i

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	169.46Å 169.46Å 119.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.84 – 3.15 19.84 – 3.15	Depositor EDS
% Data completeness (in resolution range)	90.3 (19.84-3.15) 89.9 (19.84-3.15)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.93 (at 3.15Å)	Xtrriage
Refinement program	BUSTER-TNT, BUSTER 2.10.0	Depositor
R, R_{free}	0.183 , 0.221 0.201 , 0.243	Depositor DCC
R_{free} test set	2704 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	63.7	Xtrriage
Anisotropy	0.026	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 75.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.043 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15798	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.02 % 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 1.4041e-04. 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: CL, SO4

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	0.97	2/1326 (0.2%)	1.44	12/1801 (0.7%)
1	B	0.97	0/1326	1.41	8/1801 (0.4%)
1	C	0.88	1/1326 (0.1%)	1.41	7/1801 (0.4%)
1	D	0.81	1/1326 (0.1%)	1.39	14/1801 (0.8%)
1	E	0.84	0/1326	1.44	14/1801 (0.8%)
1	F	0.77	0/1326	1.39	9/1801 (0.5%)
1	G	0.76	0/1326	1.38	10/1801 (0.6%)
1	H	0.74	0/1326	1.33	7/1801 (0.4%)
1	I	0.78	0/1326	1.37	6/1801 (0.3%)
1	J	1.00	1/1326 (0.1%)	1.48	14/1801 (0.8%)
1	K	0.91	2/1326 (0.2%)	1.41	10/1801 (0.6%)
1	L	0.82	1/1326 (0.1%)	1.37	8/1801 (0.4%)
All	All	0.86	8/15912 (0.1%)	1.40	119/21612 (0.6%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	172	ARG	CA-CB	7.83	1.58	1.52
1	K	172	ARG	CA-C	6.58	1.58	1.53
1	J	24	LEU	CA-C	6.00	1.60	1.52
1	L	172	ARG	CA-C	5.58	1.59	1.52
1	C	172	ARG	CA-C	5.12	1.59	1.52
1	D	172	ARG	CA-C	5.10	1.59	1.52
1	A	158	ALA	CA-C	5.09	1.59	1.52
1	A	118	GLY	C-N	5.03	1.39	1.33

All (119) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	ALA	CA-C-N	8.62	131.72	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	ALA	C-N-CA	8.62	131.72	120.60
1	E	91	VAL	CA-C-N	8.12	131.51	120.38
1	E	91	VAL	C-N-CA	8.12	131.51	120.38
1	C	132	ILE	N-CA-C	7.66	120.63	111.05
1	H	132	ILE	N-CA-C	7.62	120.57	111.05
1	I	132	ILE	N-CA-C	7.51	120.44	111.05
1	B	132	ILE	N-CA-C	7.51	120.44	111.05
1	B	119	SER	N-CA-C	7.49	119.08	108.74
1	A	132	ILE	N-CA-C	7.40	120.30	111.05
1	F	132	ILE	N-CA-C	7.38	120.28	111.05
1	K	132	ILE	N-CA-C	7.36	120.25	111.05
1	E	132	ILE	N-CA-C	7.35	120.24	111.05
1	G	132	ILE	N-CA-C	7.34	120.23	111.05
1	H	119	SER	N-CA-C	7.15	118.61	108.74
1	L	132	ILE	N-CA-C	7.15	119.99	111.05
1	J	132	ILE	N-CA-C	7.10	119.93	111.05
1	J	119	SER	N-CA-C	7.09	118.52	108.74
1	K	91	VAL	CA-C-N	7.07	130.08	120.54
1	K	91	VAL	C-N-CA	7.07	130.08	120.54
1	K	119	SER	N-CA-C	7.00	118.40	108.74
1	D	91	VAL	CA-C-N	6.79	129.68	120.38
1	D	91	VAL	C-N-CA	6.79	129.68	120.38
1	D	119	SER	N-CA-C	6.75	118.06	108.74
1	C	119	SER	N-CA-C	6.66	117.93	108.74
1	G	91	VAL	CA-C-N	6.60	129.42	120.38
1	G	91	VAL	C-N-CA	6.60	129.42	120.38
1	A	91	VAL	CA-C-N	6.57	129.38	120.38
1	A	91	VAL	C-N-CA	6.57	129.38	120.38
1	E	25	LYS	CA-C-N	6.53	129.33	120.38
1	E	25	LYS	C-N-CA	6.53	129.33	120.38
1	F	91	VAL	CA-C-N	6.48	129.26	120.38
1	F	91	VAL	C-N-CA	6.48	129.26	120.38
1	K	88	LEU	N-CA-C	6.44	120.08	108.48
1	B	91	VAL	CA-C-N	6.44	129.20	120.38
1	B	91	VAL	C-N-CA	6.44	129.20	120.38
1	F	119	SER	N-CA-C	6.43	117.61	108.74
1	K	25	LYS	CA-C-N	6.39	128.84	120.28
1	K	25	LYS	C-N-CA	6.39	128.84	120.28
1	J	91	VAL	CA-C-N	6.38	129.13	120.38
1	J	91	VAL	C-N-CA	6.38	129.13	120.38
1	L	119	SER	N-CA-C	6.37	117.53	108.74
1	I	91	VAL	CA-C-N	6.33	129.09	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	91	VAL	C-N-CA	6.33	129.09	120.54
1	C	91	VAL	CA-C-N	6.25	128.94	120.38
1	C	91	VAL	C-N-CA	6.25	128.94	120.38
1	L	91	VAL	CA-C-N	6.19	128.86	120.38
1	L	91	VAL	C-N-CA	6.19	128.86	120.38
1	A	119	SER	N-CA-C	6.13	117.46	108.14
1	E	178	TRP	CA-C-N	6.03	128.36	120.28
1	E	178	TRP	C-N-CA	6.03	128.36	120.28
1	L	25	LYS	CA-C-N	6.00	128.31	120.28
1	L	25	LYS	C-N-CA	6.00	128.31	120.28
1	D	25	LYS	CA-C-N	5.92	128.22	120.28
1	D	25	LYS	C-N-CA	5.92	128.22	120.28
1	D	132	ILE	N-CA-C	5.87	119.27	111.17
1	G	25	LYS	CA-C-N	5.85	128.12	120.28
1	G	25	LYS	C-N-CA	5.85	128.12	120.28
1	J	32	LEU	N-CA-C	5.82	117.71	111.36
1	F	25	LYS	CA-C-N	5.81	128.06	120.28
1	F	25	LYS	C-N-CA	5.81	128.06	120.28
1	A	157	SER	CA-C-N	5.79	128.03	120.28
1	A	157	SER	C-N-CA	5.79	128.03	120.28
1	H	25	LYS	CA-C-N	5.72	127.94	120.28
1	H	25	LYS	C-N-CA	5.72	127.94	120.28
1	J	25	LYS	CA-C-N	5.71	128.20	120.38
1	J	25	LYS	C-N-CA	5.71	128.20	120.38
1	A	25	LYS	CA-C-N	5.70	127.92	120.28
1	A	25	LYS	C-N-CA	5.70	127.92	120.28
1	F	174	HIS	CA-C-N	5.62	128.08	120.38
1	F	174	HIS	C-N-CA	5.62	128.08	120.38
1	C	25	LYS	CA-C-N	5.61	128.07	120.38
1	C	25	LYS	C-N-CA	5.61	128.07	120.38
1	D	182	VAL	N-CA-C	5.56	116.62	111.45
1	I	25	LYS	CA-C-N	5.55	127.72	120.28
1	I	25	LYS	C-N-CA	5.55	127.72	120.28
1	E	26	SER	N-CA-C	5.50	117.99	111.33
1	K	79	PRO	CA-C-N	5.42	128.29	120.38
1	K	79	PRO	C-N-CA	5.42	128.29	120.38
1	C	181	MET	N-CA-C	5.42	117.27	111.36
1	F	143	ASP	CA-CB-CG	5.42	118.02	112.60
1	D	158	ALA	CA-C-N	5.38	125.95	119.98
1	D	158	ALA	C-N-CA	5.38	125.95	119.98
1	J	130	PHE	CA-CB-CG	5.35	119.15	113.80
1	G	43	LEU	N-CA-C	5.34	117.61	108.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	26	SER	N-CA-C	5.33	117.09	111.28
1	A	26	SER	N-CA-C	5.32	117.08	111.28
1	G	26	SER	N-CA-C	5.31	117.07	111.28
1	E	97	VAL	CA-C-N	5.29	126.19	122.33
1	E	97	VAL	C-N-CA	5.29	126.19	122.33
1	B	25	LYS	CA-C-N	5.28	127.36	120.28
1	B	25	LYS	C-N-CA	5.28	127.36	120.28
1	A	143	ASP	CA-CB-CG	5.25	117.86	112.60
1	H	143	ASP	CA-CB-CG	5.24	117.84	112.60
1	D	130	PHE	CA-CB-CG	5.22	119.02	113.80
1	I	143	ASP	CA-CB-CG	5.21	117.81	112.60
1	J	30	ASP	CA-CB-CG	5.19	117.79	112.60
1	B	26	SER	N-CA-C	5.18	116.92	111.28
1	L	174	HIS	N-CA-C	5.17	117.20	110.53
1	J	157	SER	CA-C-N	5.14	127.12	120.44
1	J	157	SER	C-N-CA	5.14	127.12	120.44
1	E	143	ASP	CA-CB-CG	5.12	117.72	112.60
1	D	26	SER	N-CA-C	5.10	116.84	111.28
1	E	130	PHE	CA-CB-CG	5.09	118.89	113.80
1	L	143	ASP	CA-CB-CG	5.09	117.69	112.60
1	H	158	ALA	CA-C-N	5.09	125.63	119.98
1	H	158	ALA	C-N-CA	5.09	125.63	119.98
1	G	143	ASP	CA-CB-CG	5.08	117.68	112.60
1	E	120	ASN	N-CA-C	-5.08	104.71	111.87
1	E	78	HIS	N-CA-C	5.06	116.21	109.93
1	D	157	SER	CA-C-N	5.05	127.01	120.44
1	D	157	SER	C-N-CA	5.05	127.01	120.44
1	J	143	ASP	CA-CB-CG	5.05	117.65	112.60
1	J	60	GLY	CA-C-N	5.04	127.04	120.28
1	J	60	GLY	C-N-CA	5.04	127.04	120.28
1	D	174	HIS	N-CA-C	5.04	117.93	110.48
1	B	130	PHE	CA-CB-CG	5.03	118.83	113.80
1	G	174	HIS	CA-C-N	5.01	128.34	120.82
1	G	174	HIS	C-N-CA	5.01	128.34	120.82

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	1310	0	1355	21	0
1	B	1310	0	1355	16	0
1	C	1310	0	1355	27	0
1	D	1310	0	1355	24	0
1	E	1310	0	1355	17	0
1	F	1310	0	1355	13	0
1	G	1310	0	1355	18	0
1	H	1310	0	1355	16	0
1	I	1310	0	1355	18	0
1	J	1310	0	1355	22	0
1	K	1310	0	1355	29	0
1	L	1310	0	1355	16	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	10	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	G	5	0	0	0	0
2	H	5	0	0	0	0
2	I	5	0	0	0	0
2	J	5	0	0	0	0
2	K	5	0	0	0	0
2	L	5	0	0	0	0
3	A	1	0	0	1	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	1	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
4	A	2	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	K	1	0	0	0	0
4	L	1	0	0	0	0
All	All	15798	0	16260	199	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 6.

All (199) 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:99:ILE:HD11	1:C:181:MET:SD	2.09	0.93
1:C:152:THR:HG23	1:D:152:THR:HG23	1.52	0.91
1:B:24:LEU:HD21	1:B:31:LEU:HB2	1.57	0.87
1:B:177:MET:HB3	1:B:181:MET:HE3	1.61	0.83
1:I:152:THR:HG23	1:J:152:THR:HG23	1.62	0.82
1:A:148:ALA:O	1:A:152:THR:HG23	1.88	0.73
1:J:78:HIS:CE1	1:J:117:LYS:HB2	2.28	0.68
1:E:88:LEU:HD23	1:E:125:ARG:HB3	1.75	0.67
1:A:177:MET:HB3	1:A:181:MET:HE3	1.75	0.67
1:C:178:TRP:CE2	1:C:182:VAL:HG21	2.32	0.65
1:J:4:ALA:HB2	1:J:50:ILE:HD13	1.79	0.65
1:D:117:LYS:HG3	1:K:93:LYS:HZ2	1.62	0.65
1:C:169:LEU:HD23	1:C:171:PRO:HD3	1.79	0.64
1:F:177:MET:HB3	1:F:181:MET:HE3	1.78	0.64
1:J:177:MET:HB3	1:J:181:MET:HE3	1.79	0.64
1:A:79:PRO:HG2	3:A:202:CL:CL	2.35	0.64
1:C:174:HIS:HD2	1:C:176:ALA:H	1.44	0.64
1:I:177:MET:HB3	1:I:181:MET:HE3	1.81	0.62
1:B:33:VAL:HB	1:B:45:MET:HB2	1.82	0.62
1:B:117:LYS:HE3	1:J:92:ASP:HB3	1.82	0.62
1:G:177:MET:HB3	1:G:181:MET:HE3	1.80	0.62
1:C:177:MET:O	1:C:181:MET:HG2	2.00	0.61
1:J:2:SER:HB2	1:J:76:ASN:HA	1.83	0.61
1:B:21:ASP:HB2	1:C:177:MET:HG3	1.83	0.60
1:E:17:MET:HB3	1:F:181:MET:HE2	1.84	0.59
1:H:177:MET:HB3	1:H:181:MET:HE3	1.84	0.58
1:K:17:MET:HB3	1:L:181:MET:HE2	1.84	0.58
1:A:5:ILE:HD11	1:A:77:ILE:HG21	1.85	0.58
1:G:78:HIS:CE1	1:G:117:LYS:HB2	2.39	0.57
1:C:177:MET:HB3	1:C:181:MET:HE2	1.86	0.57
1:E:24:LEU:HD21	1:E:31:LEU:HB2	1.86	0.57
1:I:115:ALA:O	1:I:119:SER:HB3	2.03	0.57
1:K:78:HIS:CE1	1:K:117:LYS:HB2	2.40	0.57
1:D:78:HIS:CE1	1:D:117:LYS:HB2	2.40	0.57
1:E:173:PRO:HG2	1:E:178:TRP:HE3	1.70	0.57
1:H:24:LEU:HD21	1:H:31:LEU:HB2	1.88	0.56
1:H:78:HIS:CE1	1:H:117:LYS:HB2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:78:HIS:CE1	1:L:117:LYS:HB2	2.41	0.56
1:L:24:LEU:HD21	1:L:31:LEU:HB2	1.88	0.55
1:I:24:LEU:HD21	1:I:31:LEU:HB2	1.90	0.54
1:H:25:LYS:HG2	1:I:174:HIS:CD2	2.43	0.54
1:K:24:LEU:HD21	1:K:31:LEU:HB2	1.90	0.54
1:C:96:ALA:HB2	1:C:144:VAL:HG22	1.89	0.54
1:D:38:SER:HA	1:E:130:PHE:CZ	2.43	0.54
1:E:174:HIS:C	1:E:176:ALA:H	2.15	0.54
1:G:24:LEU:HD21	1:G:31:LEU:HB2	1.90	0.54
1:J:24:LEU:HB3	1:K:174:HIS:CD2	2.43	0.54
1:D:169:LEU:HD23	1:D:171:PRO:HD3	1.91	0.53
1:B:169:LEU:HD23	1:B:171:PRO:HD3	1.89	0.53
1:K:170:ILE:HG21	1:K:173:PRO:HB3	1.90	0.53
1:E:43:LEU:HD21	1:E:128:MET:HE1	1.89	0.53
1:I:169:LEU:HD23	1:I:171:PRO:HD3	1.91	0.52
1:L:169:LEU:HD23	1:L:171:PRO:HD3	1.91	0.52
1:J:169:LEU:HD23	1:J:171:PRO:HD3	1.90	0.52
1:A:24:LEU:HD21	1:A:31:LEU:HB2	1.92	0.52
1:K:25:LYS:NZ	1:L:172:ARG:HH21	2.08	0.52
1:F:33:VAL:HB	1:F:45:MET:HB2	1.92	0.52
1:I:33:VAL:HB	1:I:45:MET:HB2	1.92	0.52
1:A:78:HIS:CE1	1:A:117:LYS:HB2	2.45	0.51
1:C:24:LEU:HD21	1:C:31:LEU:HB2	1.92	0.51
1:C:33:VAL:HB	1:C:45:MET:HB2	1.91	0.51
1:L:33:VAL:HB	1:L:45:MET:HB2	1.93	0.51
1:G:169:LEU:HD23	1:G:171:PRO:HD3	1.91	0.51
1:K:94:ARG:HD3	1:K:178:TRP:CE2	2.46	0.51
1:D:17:MET:HB3	1:E:181:MET:HE2	1.92	0.51
1:D:20:GLY:HA2	1:D:23:MET:HE2	1.93	0.51
1:F:78:HIS:CE1	1:F:117:LYS:HB2	2.45	0.50
1:D:33:VAL:HB	1:D:45:MET:HB2	1.92	0.50
1:L:88:LEU:HD23	1:L:125:ARG:HB3	1.92	0.50
1:H:17:MET:HB3	1:I:181:MET:HE2	1.93	0.50
1:G:17:MET:HB3	1:H:181:MET:HE2	1.93	0.50
1:E:33:VAL:HB	1:E:45:MET:HB2	1.93	0.50
1:D:24:LEU:HD21	1:D:31:LEU:HB2	1.94	0.49
1:F:24:LEU:HD21	1:F:31:LEU:HB2	1.93	0.49
1:G:99:ILE:HG12	1:G:139:VAL:HG22	1.93	0.49
1:A:33:VAL:HB	1:A:45:MET:HB2	1.92	0.49
1:J:17:MET:HG2	1:K:177:MET:HE1	1.94	0.49
1:K:96:ALA:HB2	1:K:144:VAL:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:33:VAL:HB	1:G:45:MET:HB2	1.94	0.49
1:C:88:LEU:HD23	1:C:125:ARG:HB3	1.94	0.49
1:E:176:ALA:HA	1:E:179:ARG:HD3	1.93	0.49
1:G:25:LYS:NZ	1:H:172:ARG:HE	2.10	0.49
1:K:97:VAL:HG23	1:K:170:ILE:HB	1.93	0.49
1:A:169:LEU:HD23	1:A:171:PRO:HD3	1.95	0.49
1:D:88:LEU:HD23	1:D:125:ARG:HB3	1.94	0.49
1:K:38:SER:HA	1:L:130:PHE:CZ	2.48	0.48
1:L:177:MET:HB3	1:L:181:MET:HE3	1.95	0.48
1:F:169:LEU:HD23	1:F:171:PRO:HD3	1.95	0.48
1:J:88:LEU:HD23	1:J:125:ARG:HB3	1.95	0.48
1:H:88:LEU:HD23	1:H:125:ARG:HB3	1.95	0.48
1:J:33:VAL:HB	1:J:45:MET:HB2	1.96	0.48
1:A:50:ILE:HD13	1:A:75:ALA:HB2	1.96	0.48
1:C:8:LEU:HD23	1:C:10:LEU:HD21	1.95	0.48
1:G:177:MET:HG3	1:I:21:ASP:HB2	1.95	0.47
1:G:25:LYS:HZ3	1:H:172:ARG:HE	1.62	0.47
1:L:174:HIS:ND1	1:L:175:GLU:OE1	2.48	0.47
1:B:17:MET:HB3	1:C:177:MET:HE2	1.96	0.47
1:D:128:MET:HG3	1:D:136:CYS:SG	2.55	0.47
1:D:174:HIS:CD2	1:F:25:LYS:HG2	2.50	0.47
1:J:32:LEU:HD22	1:J:86:SER:HB3	1.97	0.47
1:A:180:GLN:HE21	1:C:17:MET:HE2	1.79	0.47
1:H:33:VAL:HB	1:H:45:MET:HB2	1.96	0.47
1:J:10:LEU:HD23	1:J:16:GLY:HA2	1.97	0.47
1:K:170:ILE:HG12	1:K:181:MET:HE1	1.97	0.47
1:F:96:ALA:HB2	1:F:144:VAL:HG22	1.97	0.46
1:D:125:ARG:CZ	1:D:184:GLY:HA2	2.46	0.46
1:K:33:VAL:HB	1:K:45:MET:HB2	1.97	0.46
1:J:47:GLY:HA3	1:J:82:LEU:HD23	1.97	0.46
1:A:154:ALA:O	1:A:157:SER:HB3	2.16	0.46
1:H:78:HIS:HB3	3:H:202:CL:CL	2.53	0.46
1:D:117:LYS:HG3	1:K:93:LYS:NZ	2.28	0.46
1:G:88:LEU:HD23	1:G:125:ARG:HB3	1.97	0.46
1:I:88:LEU:HD23	1:I:125:ARG:HB3	1.97	0.46
1:A:96:ALA:HB2	1:A:144:VAL:HG22	1.98	0.45
1:B:31:LEU:HD21	1:B:34:SER:HB2	1.98	0.45
1:K:7:ILE:HG12	1:K:45:MET:HG2	1.98	0.45
1:C:171:PRO:HB2	1:K:172:ARG:NH2	2.31	0.45
1:I:96:ALA:HB2	1:I:144:VAL:HG22	1.98	0.45
1:G:38:SER:HA	1:H:130:PHE:CZ	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:100:VAL:HG21	1:J:154:ALA:HB3	1.99	0.45
1:I:10:LEU:HD23	1:I:16:GLY:HA2	1.98	0.45
1:K:170:ILE:CG2	1:K:173:PRO:HB3	2.47	0.45
1:B:88:LEU:HD23	1:B:125:ARG:HB3	1.99	0.44
1:K:11:THR:HG22	1:K:67:LEU:HA	2.00	0.44
1:J:7:ILE:HB	1:J:72:LEU:HB3	1.99	0.44
1:J:181:MET:HE2	1:L:17:MET:HB3	1.99	0.44
1:C:78:HIS:CE1	1:C:117:LYS:HB2	2.53	0.44
1:K:97:VAL:CG2	1:K:173:PRO:HG3	2.48	0.44
1:K:182:VAL:HG23	1:K:183:GLU:HG3	2.00	0.44
1:G:174:HIS:HD2	1:I:25:LYS:HE2	1.83	0.44
1:H:7:ILE:HG12	1:H:45:MET:HG2	2.00	0.44
1:K:171:PRO:C	1:K:172:ARG:HG3	2.42	0.44
1:L:11:THR:HG22	1:L:67:LEU:HA	2.00	0.44
1:A:7:ILE:HG12	1:A:45:MET:HG2	2.00	0.44
1:A:181:MET:HE2	1:C:17:MET:HB3	1.99	0.44
1:E:7:ILE:HG12	1:E:45:MET:HG2	2.00	0.44
1:H:38:SER:HA	1:I:130:PHE:CZ	2.53	0.44
1:K:47:GLY:C	1:K:53:ILE:HD11	2.44	0.43
1:B:94:ARG:HG2	1:B:178:TRP:CD2	2.53	0.43
1:I:11:THR:HG22	1:I:67:LEU:HA	2.00	0.43
1:K:178:TRP:NE1	1:K:182:VAL:HG21	2.33	0.43
1:B:67:LEU:HD11	1:D:160:ALA:HA	2.00	0.43
1:D:100:VAL:HG21	1:D:154:ALA:HB3	1.99	0.43
1:F:7:ILE:HG12	1:F:45:MET:HG2	2.01	0.43
1:K:100:VAL:HG21	1:K:154:ALA:HB3	2.01	0.43
1:B:78:HIS:CE1	1:B:117:LYS:HB2	2.53	0.43
1:C:160:ALA:HA	1:F:67:LEU:HD11	1.99	0.43
1:B:47:GLY:HA3	1:B:82:LEU:HD23	2.00	0.43
1:D:17:MET:HB3	1:E:181:MET:CE	2.49	0.43
1:K:99:ILE:O	1:K:167:ALA:HA	2.19	0.43
1:H:67:LEU:HD11	1:J:160:ALA:HA	2.01	0.42
1:C:11:THR:HG22	1:C:67:LEU:HA	2.02	0.42
1:G:11:THR:HG22	1:G:67:LEU:HA	2.02	0.42
1:D:7:ILE:HG12	1:D:45:MET:HG2	2.01	0.42
1:E:91:VAL:HG11	1:E:178:TRP:HH2	1.85	0.42
1:K:144:VAL:HG13	1:K:169:LEU:HD21	2.02	0.42
1:A:38:SER:HA	1:B:130:PHE:CZ	2.55	0.42
1:C:88:LEU:HD21	1:C:125:ARG:HD3	2.01	0.42
1:G:96:ALA:HB2	1:G:144:VAL:HG22	2.01	0.42
1:J:96:ALA:HB2	1:J:144:VAL:HG22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:100:VAL:HG21	1:G:154:ALA:HB3	2.02	0.42
1:C:88:LEU:CD2	1:C:125:ARG:HD3	2.50	0.42
1:D:96:ALA:HB2	1:D:144:VAL:HG22	2.02	0.42
1:I:100:VAL:HG21	1:I:154:ALA:HB3	2.02	0.42
1:G:7:ILE:HG12	1:G:45:MET:HG2	2.02	0.41
1:G:94:ARG:HG2	1:G:178:TRP:CD2	2.55	0.41
1:A:88:LEU:HD23	1:A:125:ARG:HB3	2.02	0.41
1:C:43:LEU:HD21	1:C:128:MET:HE1	2.01	0.41
1:D:178:TRP:O	1:D:182:VAL:HG22	2.21	0.41
1:L:7:ILE:HG12	1:L:45:MET:HG2	2.02	0.41
1:L:94:ARG:HG2	1:L:178:TRP:CD2	2.55	0.41
1:D:144:VAL:HG13	1:D:169:LEU:HD21	2.03	0.41
1:D:177:MET:HE2	1:D:181:MET:HE3	2.03	0.41
1:E:11:THR:HG22	1:E:67:LEU:HA	2.02	0.41
1:D:11:THR:HG22	1:D:67:LEU:HA	2.02	0.41
1:H:114:ARG:HD2	1:H:157:SER:HB2	2.02	0.41
1:C:100:VAL:HG21	1:C:154:ALA:HB3	2.01	0.41
1:L:100:VAL:HG21	1:L:154:ALA:HB3	2.03	0.41
1:C:7:ILE:HG12	1:C:45:MET:HG2	2.01	0.41
1:E:174:HIS:C	1:E:176:ALA:N	2.79	0.41
1:F:11:THR:HG22	1:F:67:LEU:HA	2.02	0.41
1:I:94:ARG:HG2	1:I:178:TRP:CD2	2.56	0.41
1:A:11:THR:HG22	1:A:67:LEU:HA	2.03	0.41
1:A:130:PHE:CZ	1:C:38:SER:HA	2.56	0.41
1:B:11:THR:HG22	1:B:67:LEU:HA	2.02	0.41
1:E:178:TRP:CZ2	1:E:182:VAL:HG21	2.56	0.41
1:A:144:VAL:HG13	1:A:169:LEU:HD21	2.03	0.41
1:K:79:PRO:O	1:K:83:PRO:HD2	2.21	0.41
1:J:17:MET:HB3	1:K:177:MET:HE2	2.03	0.41
1:E:114:ARG:HD2	1:E:157:SER:HB2	2.03	0.40
1:C:47:GLY:C	1:C:53:ILE:HD11	2.46	0.40
1:F:114:ARG:HD2	1:F:157:SER:HB2	2.04	0.40
1:I:114:ARG:HD2	1:I:157:SER:HB2	2.03	0.40
1:J:114:ARG:HD2	1:J:157:SER:HB2	2.03	0.40
1:L:43:LEU:CD2	1:L:128:MET:HE1	2.51	0.40
1:A:5:ILE:HG12	1:A:77:ILE:HG13	2.02	0.40
1:J:144:VAL:HG13	1:J:169:LEU:HD21	2.04	0.40
1:A:21:ASP:HB2	1:B:177:MET:HG3	2.02	0.40
1:D:130:PHE:CZ	1:F:38:SER:HA	2.57	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	181/192 (94%)	167 (92%)	13 (7%)	1 (1%)	21	52
1	B	181/192 (94%)	174 (96%)	6 (3%)	1 (1%)	21	52
1	C	181/192 (94%)	172 (95%)	8 (4%)	1 (1%)	21	52
1	D	181/192 (94%)	172 (95%)	8 (4%)	1 (1%)	21	52
1	E	181/192 (94%)	167 (92%)	12 (7%)	2 (1%)	11	39
1	F	181/192 (94%)	171 (94%)	10 (6%)	0	100	100
1	G	181/192 (94%)	169 (93%)	12 (7%)	0	100	100
1	H	181/192 (94%)	169 (93%)	10 (6%)	2 (1%)	11	39
1	I	181/192 (94%)	171 (94%)	9 (5%)	1 (1%)	21	52
1	J	181/192 (94%)	172 (95%)	7 (4%)	2 (1%)	11	39
1	K	181/192 (94%)	170 (94%)	10 (6%)	1 (1%)	21	52
1	L	181/192 (94%)	172 (95%)	8 (4%)	1 (1%)	21	52
All	All	2172/2304 (94%)	2046 (94%)	113 (5%)	13 (1%)	21	52

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	175	GLU
1	K	96	ALA
1	E	94	ARG
1	L	175	GLU
1	J	24	LEU
1	C	104	SER
1	A	104	SER
1	B	104	SER
1	D	104	SER
1	I	104	SER
1	J	104	SER
1	E	124	VAL

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Mol	Chain	Res	Type
1	H	171	PRO

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	139/148 (94%)	131 (94%)	8 (6%)	18	46
1	B	139/148 (94%)	130 (94%)	9 (6%)	15	42
1	C	139/148 (94%)	130 (94%)	9 (6%)	15	42
1	D	139/148 (94%)	132 (95%)	7 (5%)	22	50
1	E	139/148 (94%)	125 (90%)	14 (10%)	7	26
1	F	139/148 (94%)	130 (94%)	9 (6%)	15	42
1	G	139/148 (94%)	131 (94%)	8 (6%)	18	46
1	H	139/148 (94%)	130 (94%)	9 (6%)	15	42
1	I	139/148 (94%)	130 (94%)	9 (6%)	15	42
1	J	139/148 (94%)	129 (93%)	10 (7%)	13	39
1	K	139/148 (94%)	124 (89%)	15 (11%)	6	23
1	L	139/148 (94%)	130 (94%)	9 (6%)	15	42
All	All	1668/1776 (94%)	1552 (93%)	116 (7%)	14	40

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	THR
1	A	38	SER
1	A	62	SER
1	A	80	SER
1	A	86	SER
1	A	149	LEU
1	A	152	THR
1	A	172	ARG
1	B	11	THR

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Mol	Chain	Res	Type
1	B	35	LYS
1	B	46	LEU
1	B	62	SER
1	B	76	ASN
1	B	149	LEU
1	B	172	ARG
1	B	174	HIS
1	B	175	GLU
1	C	8	LEU
1	C	11	THR
1	C	35	LYS
1	C	62	SER
1	C	66	GLU
1	C	76	ASN
1	C	149	LEU
1	C	175	GLU
1	C	179	ARG
1	D	11	THR
1	D	35	LYS
1	D	62	SER
1	D	66	GLU
1	D	76	ASN
1	D	149	LEU
1	D	172	ARG
1	E	2	SER
1	E	11	THR
1	E	26	SER
1	E	41	LYS
1	E	62	SER
1	E	76	ASN
1	E	93	LYS
1	E	121	VAL
1	E	149	LEU
1	E	169	LEU
1	E	174	HIS
1	E	175	GLU
1	E	177	MET
1	E	182	VAL
1	F	11	THR
1	F	35	LYS
1	F	62	SER
1	F	76	ASN

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Mol	Chain	Res	Type
1	F	132	ILE
1	F	149	LEU
1	F	172	ARG
1	F	175	GLU
1	F	179	ARG
1	G	11	THR
1	G	35	LYS
1	G	62	SER
1	G	76	ASN
1	G	93	LYS
1	G	149	LEU
1	G	172	ARG
1	G	174	HIS
1	H	11	THR
1	H	35	LYS
1	H	62	SER
1	H	76	ASN
1	H	93	LYS
1	H	132	ILE
1	H	149	LEU
1	H	172	ARG
1	H	174	HIS
1	I	11	THR
1	I	35	LYS
1	I	62	SER
1	I	66	GLU
1	I	76	ASN
1	I	132	ILE
1	I	149	LEU
1	I	172	ARG
1	I	174	HIS
1	J	8	LEU
1	J	10	LEU
1	J	11	THR
1	J	29	VAL
1	J	50	ILE
1	J	62	SER
1	J	66	GLU
1	J	117	LYS
1	J	149	LEU
1	J	172	ARG
1	K	11	THR

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Mol	Chain	Res	Type
1	K	17	MET
1	K	62	SER
1	K	66	GLU
1	K	76	ASN
1	K	80	SER
1	K	90	SER
1	K	94	ARG
1	K	97	VAL
1	K	99	ILE
1	K	122	THR
1	K	149	LEU
1	K	170	ILE
1	K	172	ARG
1	K	182	VAL
1	L	11	THR
1	L	35	LYS
1	L	38	SER
1	L	62	SER
1	L	76	ASN
1	L	132	ILE
1	L	149	LEU
1	L	172	ARG
1	L	175	GLU

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	3	GLN
1	A	78	HIS
1	A	120	ASN
1	C	120	ASN
1	C	174	HIS
1	D	3	GLN
1	D	120	ASN
1	E	180	GLN
1	F	120	ASN
1	G	3	GLN
1	G	120	ASN
1	G	174	HIS
1	H	120	ASN
1	I	120	ASN
1	J	3	GLN

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Mol	Chain	Res	Type
1	J	55	GLN
1	J	76	ASN
1	K	3	GLN
1	K	95	GLN
1	K	120	ASN
1	L	120	ASN

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](#)

Of 24 ligands modelled in this entry, 12 are monoatomic - leaving 12 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	SO4	H	201	-	4,4,4	0.37	0	6,6,6	0.17	0
2	SO4	C	202	-	4,4,4	0.43	0	6,6,6	0.19	0
2	SO4	J	201	-	4,4,4	0.45	0	6,6,6	0.37	0
2	SO4	L	201	-	4,4,4	0.26	0	6,6,6	0.20	0
2	SO4	I	201	-	4,4,4	0.27	0	6,6,6	0.18	0
2	SO4	C	201	-	4,4,4	0.21	0	6,6,6	0.20	0
2	SO4	B	201	-	4,4,4	0.24	0	6,6,6	0.42	0
2	SO4	E	201	-	4,4,4	0.24	0	6,6,6	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	201	-	4,4,4	0.33	0	6,6,6	0.22	0
2	SO4	G	201	-	4,4,4	0.28	0	6,6,6	0.19	0
2	SO4	D	201	-	4,4,4	0.21	0	6,6,6	0.29	0
2	SO4	K	201	-	4,4,4	0.31	0	6,6,6	0.12	0

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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	183/192 (95%)	-0.53	1 (0%) 87 75	24, 50, 93, 116	0
1	B	183/192 (95%)	-0.59	1 (0%) 87 75	21, 45, 85, 103	0
1	C	183/192 (95%)	-0.51	1 (0%) 87 75	32, 58, 95, 117	0
1	D	183/192 (95%)	-0.23	0 100 100	45, 81, 122, 131	0
1	E	183/192 (95%)	-0.32	1 (0%) 87 75	51, 77, 110, 122	0
1	F	183/192 (95%)	-0.18	1 (0%) 87 75	55, 83, 123, 141	0
1	G	183/192 (95%)	-0.08	0 100 100	65, 89, 124, 143	0
1	H	183/192 (95%)	-0.01	1 (0%) 87 75	63, 94, 128, 149	0
1	I	183/192 (95%)	-0.13	0 100 100	53, 89, 123, 136	0
1	J	183/192 (95%)	-0.47	0 100 100	28, 58, 95, 109	0
1	K	183/192 (95%)	-0.52	0 100 100	35, 58, 91, 112	0
1	L	183/192 (95%)	-0.27	2 (1%) 78 60	38, 74, 107, 121	0
All	All	2196/2304 (95%)	-0.32	8 (0%) 88 78	21, 74, 116, 149	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	118	GLY	3.2
1	C	118	GLY	3.2
1	F	26	SER	2.5
1	L	26	SER	2.4
1	H	60	GLY	2.2
1	E	118	GLY	2.2
1	B	2	SER	2.1
1	L	2	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	J	201	5/5	0.77	0.10	127,127,127,128	0
2	SO4	C	202	5/5	0.80	0.17	144,144,145,145	0
2	SO4	B	201	5/5	0.80	0.10	112,112,113,113	0
2	SO4	I	201	5/5	0.83	0.09	147,147,148,148	0
2	SO4	G	201	5/5	0.83	0.09	134,134,134,135	0
2	SO4	D	201	5/5	0.84	0.13	124,124,125,125	0
2	SO4	E	201	5/5	0.87	0.08	116,116,117,118	0
3	CL	I	202	1/1	0.87	0.16	80,80,80,80	0
3	CL	G	202	1/1	0.88	0.16	90,90,90,90	0
3	CL	C	203	1/1	0.88	0.16	79,79,79,79	0
3	CL	J	202	1/1	0.89	0.20	48,48,48,48	0
3	CL	H	202	1/1	0.90	0.17	114,114,114,114	0
3	CL	F	201	1/1	0.90	0.08	79,79,79,79	0
2	SO4	K	201	5/5	0.90	0.07	113,113,114,115	0
3	CL	A	202	1/1	0.91	0.09	53,53,53,53	0
2	SO4	C	201	5/5	0.91	0.07	120,122,123,123	0
3	CL	E	202	1/1	0.91	0.09	65,65,65,65	0
2	SO4	A	201	5/5	0.91	0.09	109,109,109,110	0
2	SO4	L	201	5/5	0.92	0.09	96,97,98,98	0
3	CL	L	202	1/1	0.92	0.14	79,79,79,79	0
2	SO4	H	201	5/5	0.93	0.09	116,116,116,116	0
3	CL	B	203	1/1	0.94	0.11	35,35,35,35	0
3	CL	K	202	1/1	0.95	0.08	58,58,58,58	0
3	CL	B	202	1/1	0.96	0.07	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.