



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 03:43 PM UTC

PDB ID : 3VCD / pdb_00003vcd
Title : Computationally Designed Self-assembling Octahedral Cage protein, O333,
Crystallized in space group R32
Authors : Sawaya, M.R.; King, N.P.; Sheffler, W.; Baker, D.; Yeates, T.O.
Deposited on : 2012-01-03
Resolution : 2.35 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

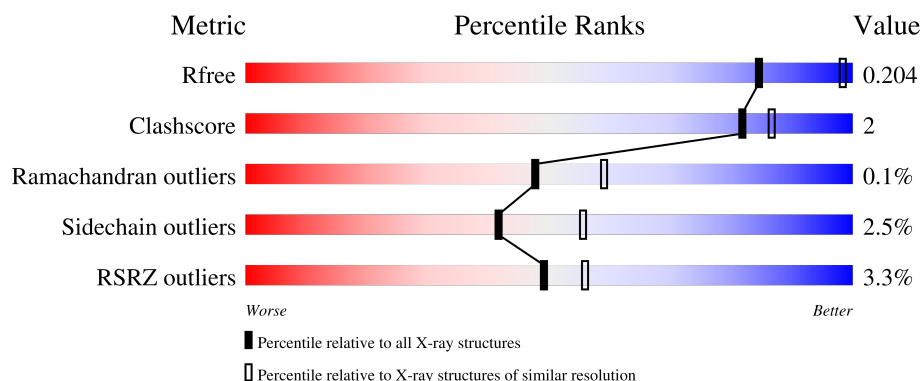
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	3% 88% 8% •
1	B	192	4% 83% 11% • 5%
1	C	192	5% 88% 6% • 5%
1	D	192	2% 81% 13% • 5%
1	E	192	4% 86% 8% • •

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Mol	Chain	Length	Quality of chain
1	F	192	5% 90% 5% • 5%
1	G	192	2% 92% • • 5%
1	H	192	% 90% 5% • 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11002 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
1	A	184	Total	C	N	O	S	0	0	0
			1318	834	223	252	9			
1	B	183	Total	C	N	O	S	0	0	0
			1310	828	222	251	9			
1	C	183	Total	C	N	O	S	0	0	0
			1310	828	222	251	9			
1	D	183	Total	C	N	O	S	0	0	0
			1310	828	222	251	9			
1	E	184	Total	C	N	O	S	0	0	0
			1318	834	223	252	9			
1	F	183	Total	C	N	O	S	0	0	0
			1310	828	222	251	9			
1	G	183	Total	C	N	O	S	0	0	0
			1310	828	222	251	9			
1	H	183	Total	C	N	O	S	0	0	0
			1310	828	222	251	9			

There are 144 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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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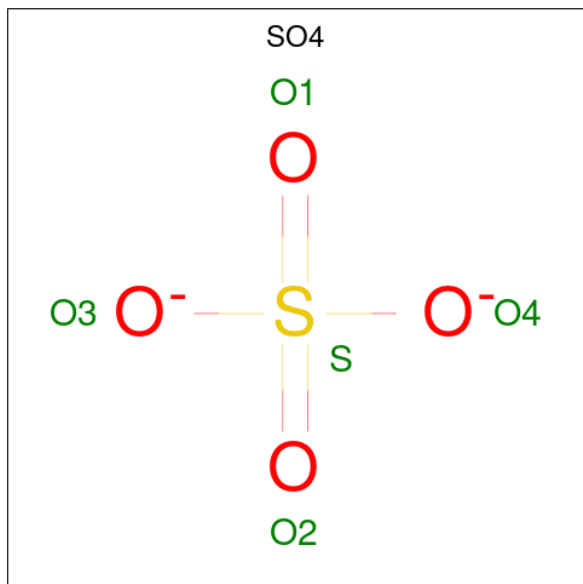
Chain	Residue	Modelled	Actual	Comment	Reference
F	161	TYR	LYS	engineered mutation	UNP E7V033
F	167	ALA	ARG	engineered mutation	UNP E7V033
F	169	LEU	VAL	engineered mutation	UNP E7V033
F	185	LEU	-	expression tag	UNP E7V033
F	186	GLU	-	expression tag	UNP E7V033
F	187	HIS	-	expression tag	UNP E7V033
F	188	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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	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	H	1	Total	Cl	0	0
			1	1		

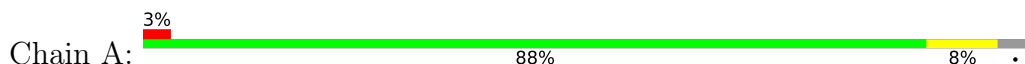
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	36	Total 36	O 36	0	0
4	B	47	Total 47	O 47	0	0
4	C	46	Total 46	O 46	0	0
4	D	56	Total 56	O 56	0	0
4	E	41	Total 41	O 41	0	0
4	F	35	Total 35	O 35	0	0
4	G	48	Total 48	O 48	0	0
4	H	66	Total 66	O 66	0	0

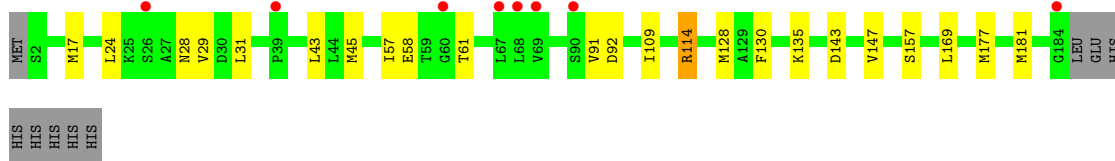
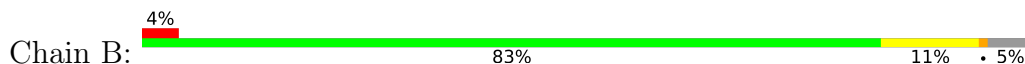
3 Residue-property plots [i](#)

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.

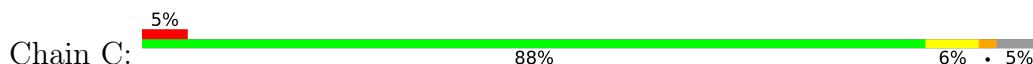
- Molecule 1: Propanediol utilization polyhedral body protein PduT



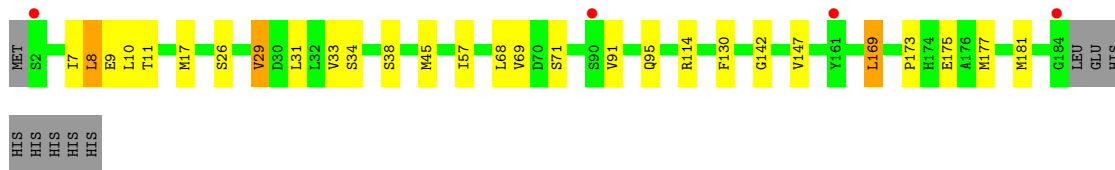
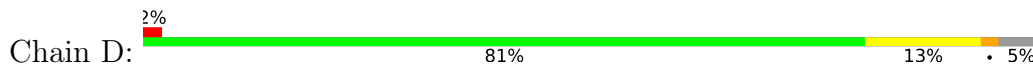
- Molecule 1: Propanediol utilization polyhedral body protein PduT



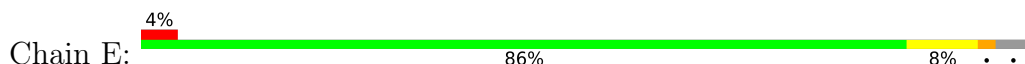
- Molecule 1: Propanediol utilization polyhedral body protein PduT



- Molecule 1: Propanediol utilization polyhedral body protein PduT

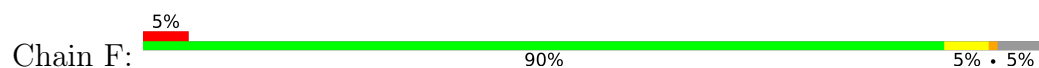


- Molecule 1: Propanediol utilization polyhedral body protein PduT

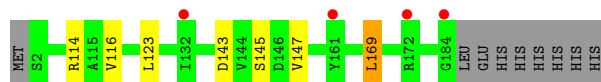
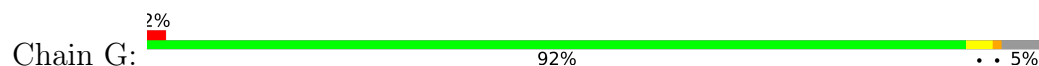




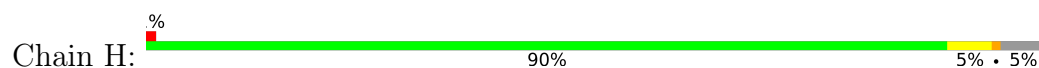
- Molecule 1: Propanediol utilization polyhedral body protein PduT



- Molecule 1: Propanediol utilization polyhedral body protein PduT



- Molecule 1: Propanediol utilization polyhedral body protein PduT



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	137.92Å 137.92Å 560.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	81.74 – 2.35 81.74 – 2.35	Depositor EDS
% Data completeness (in resolution range)	95.0 (81.74-2.35) 95.0 (81.74-2.35)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 2.34Å)	Xtriage
Refinement program	BUSTER-TNT, BUSTER 2.10.0	Depositor
R, R_{free}	0.180 , 0.213 0.178 , 0.204	Depositor DCC
R_{free} test set	4055 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.136	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 61.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11002	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, 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	0.78	0/1334	1.23	6/1812 (0.3%)
1	B	0.88	0/1326	1.19	3/1801 (0.2%)
1	C	0.82	0/1326	1.21	2/1801 (0.1%)
1	D	0.91	0/1326	1.24	3/1801 (0.2%)
1	E	0.79	0/1334	1.24	5/1812 (0.3%)
1	F	0.80	0/1326	1.22	5/1801 (0.3%)
1	G	0.79	0/1326	1.18	1/1801 (0.1%)
1	H	0.86	0/1326	1.19	1/1801 (0.1%)
All	All	0.83	0/10624	1.21	26/14430 (0.2%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	91	VAL	CA-C-N	7.07	130.08	120.54
1	E	91	VAL	C-N-CA	7.07	130.08	120.54
1	A	91	VAL	CA-C-N	6.84	129.77	120.54
1	A	91	VAL	C-N-CA	6.84	129.77	120.54
1	A	69	VAL	N-CA-C	-6.77	105.16	111.45
1	F	69	VAL	N-CA-C	-6.74	105.19	111.45
1	B	29	VAL	N-CA-CB	6.39	120.23	111.41
1	E	178	TRP	CA-C-N	6.09	128.76	120.54
1	E	178	TRP	C-N-CA	6.09	128.76	120.54
1	F	91	VAL	CA-C-N	5.95	128.57	120.54
1	F	91	VAL	C-N-CA	5.95	128.57	120.54
1	C	174	HIS	CA-C-N	5.79	128.36	120.54
1	C	174	HIS	C-N-CA	5.79	128.36	120.54
1	F	174	HIS	CA-C-N	5.48	127.94	120.54
1	F	174	HIS	C-N-CA	5.48	127.94	120.54
1	A	174	HIS	CA-C-N	5.48	127.93	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	HIS	C-N-CA	5.48	127.93	120.54
1	H	35	LYS	N-CA-C	5.39	117.03	108.79
1	A	143	ASP	CA-CB-CG	5.31	117.91	112.60
1	D	173	PRO	CA-C-N	5.12	128.14	120.87
1	D	173	PRO	C-N-CA	5.12	128.14	120.87
1	E	182	VAL	N-CA-CB	5.10	118.47	110.31
1	D	175	GLU	CB-CG-CD	5.09	121.25	112.60
1	B	143	ASP	CA-CB-CG	5.08	117.68	112.60
1	B	28	ASN	N-CA-C	5.06	118.42	111.54
1	G	143	ASP	CA-CB-CG	5.00	117.61	112.60

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	1318	0	1366	7	0
1	B	1310	0	1355	11	0
1	C	1310	0	1355	9	0
1	D	1310	0	1355	13	0
1	E	1318	0	1366	10	0
1	F	1310	0	1355	5	0
1	G	1310	0	1355	3	0
1	H	1310	0	1355	4	0
2	A	10	0	0	1	0
2	B	15	0	0	0	0
2	C	15	0	0	0	0
2	D	25	0	0	0	0
2	E	20	0	0	0	0
2	F	15	0	0	0	0
2	G	10	0	0	0	0
2	H	20	0	0	0	0
3	H	1	0	0	0	0
4	A	36	0	0	0	0
4	B	47	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	46	0	0	0	0
4	D	56	0	0	0	0
4	E	41	0	0	0	0
4	F	35	0	0	1	0
4	G	48	0	0	0	0
4	H	66	0	0	1	0
All	All	11002	0	10862	53	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 (53) 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:175:GLU:H	1:C:175:GLU:CD	1.96	0.70
1:B:43:LEU:HD22	1:B:128:MET:HE1	1.74	0.69
1:D:8:LEU:HD13	1:D:57:ILE:HD12	1.81	0.63
1:B:57:ILE:O	1:B:61:THR:HG23	2.00	0.62
1:E:43:LEU:HD22	1:E:128:MET:HE1	1.83	0.61
1:H:147:VAL:HG12	1:H:169:LEU:HD12	1.83	0.60
1:E:145:SER:HB2	1:G:145:SER:HB2	1.84	0.60
1:C:38:SER:HB3	1:C:130:PHE:CZ	2.36	0.60
1:D:147:VAL:HG12	1:D:169:LEU:HD12	1.83	0.59
1:B:43:LEU:CD2	1:B:128:MET:HE1	2.33	0.58
1:A:147:VAL:HG12	1:A:169:LEU:HD12	1.87	0.56
1:G:147:VAL:HG12	1:G:169:LEU:HD12	1.88	0.56
1:D:38:SER:HA	1:E:130:PHE:CZ	2.41	0.56
1:B:147:VAL:HG12	1:B:169:LEU:HD12	1.91	0.52
1:A:130:PHE:CZ	1:C:38:SER:HA	2.43	0.52
1:F:147:VAL:HG12	1:F:169:LEU:HD12	1.90	0.52
1:E:99:ILE:CG2	1:E:168:SER:OG	2.57	0.52
1:E:43:LEU:CD2	1:E:128:MET:HE1	2.40	0.51
1:E:147:VAL:HG12	1:E:169:LEU:HD12	1.93	0.50
1:C:147:VAL:HG12	1:C:169:LEU:HD12	1.94	0.49
1:A:116:VAL:CG1	1:A:123:LEU:HB2	2.43	0.48
1:D:17:MET:HE2	1:E:180:GLN:HG2	1.96	0.48
1:C:177:MET:HB3	1:C:181:MET:HE3	1.95	0.48
1:E:175:GLU:O	1:E:178:TRP:HB3	2.13	0.48
1:D:7:ILE:O	1:D:71:SER:HA	2.13	0.48
1:C:38:SER:HB3	1:C:130:PHE:HZ	1.79	0.47
1:F:172:ARG:NH1	4:F:326:HOH:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:MET:HE2	1:C:180:GLN:HG2	1.95	0.47
1:B:24:LEU:HD21	1:B:31:LEU:HB2	1.95	0.47
1:B:177:MET:HB3	1:B:181:MET:HE3	1.97	0.46
1:A:177:MET:HB3	1:A:181:MET:HE3	1.97	0.46
1:G:116:VAL:CG1	1:G:123:LEU:HB2	2.46	0.46
1:D:9:GLU:HB3	1:D:69:VAL:HB	1.99	0.45
1:D:130:PHE:CZ	1:F:38:SER:HA	2.52	0.44
1:F:177:MET:HB3	1:F:181:MET:HE3	1.97	0.44
1:D:26:SER:HB2	1:D:29:VAL:HG22	2.00	0.44
1:D:177:MET:HB3	1:D:181:MET:HE3	2.00	0.44
1:D:11:THR:O	1:E:135:LYS:NZ	2.47	0.43
1:H:177:MET:HB3	1:H:181:MET:HE3	2.00	0.43
1:D:95:GLN:NE2	1:D:142:GLY:HA2	2.33	0.43
1:C:94:ARG:HG2	1:C:178:TRP:CD2	2.54	0.43
1:H:33:VAL:HB	1:H:45:MET:HB2	2.00	0.43
1:E:114:ARG:HD2	1:E:157:SER:HB2	2.00	0.43
1:H:35:LYS:HE2	4:H:303:HOH:O	2.19	0.42
1:F:94:ARG:HG2	1:F:178:TRP:CD2	2.55	0.42
1:D:33:VAL:HB	1:D:45:MET:HB2	2.02	0.41
1:A:132:ILE:N	2:A:202:SO4:O1	2.53	0.41
1:A:11:THR:O	1:B:135:LYS:NZ	2.53	0.41
1:C:116:VAL:CG1	1:C:123:LEU:HB2	2.50	0.41
1:B:114:ARG:HD2	1:B:157:SER:HB2	2.02	0.40
1:B:45:MET:HE1	1:B:109:ILE:HD11	2.04	0.40
1:A:38:SER:HA	1:B:130:PHE:CZ	2.56	0.40
1:D:31:LEU:HD21	1:D:34:SER:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	182/192 (95%)	174 (96%)	8 (4%)	0	100	100
1	B	181/192 (94%)	173 (96%)	8 (4%)	0	100	100
1	C	181/192 (94%)	174 (96%)	7 (4%)	0	100	100
1	D	181/192 (94%)	176 (97%)	5 (3%)	0	100	100
1	E	182/192 (95%)	175 (96%)	7 (4%)	0	100	100
1	F	181/192 (94%)	174 (96%)	7 (4%)	0	100	100
1	G	181/192 (94%)	176 (97%)	5 (3%)	0	100	100
1	H	181/192 (94%)	175 (97%)	5 (3%)	1 (1%)	21	24
All	All	1450/1536 (94%)	1397 (96%)	52 (4%)	1 (0%)	48	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	92	ASP

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	140/148 (95%)	138 (99%)	2 (1%)	59	73
1	B	139/148 (94%)	135 (97%)	4 (3%)	37	49
1	C	139/148 (94%)	135 (97%)	4 (3%)	37	49
1	D	139/148 (94%)	132 (95%)	7 (5%)	22	27
1	E	140/148 (95%)	135 (96%)	5 (4%)	31	41
1	F	139/148 (94%)	138 (99%)	1 (1%)	76	86
1	G	139/148 (94%)	137 (99%)	2 (1%)	59	73
1	H	139/148 (94%)	136 (98%)	3 (2%)	45	59
All	All	1114/1184 (94%)	1086 (98%)	28 (2%)	42	55

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	ASP
1	A	183	GLU
1	B	58	GLU
1	B	91	VAL
1	B	92	ASP
1	B	114	ARG
1	C	94	ARG
1	C	143	ASP
1	C	169	LEU
1	C	175	GLU
1	D	8	LEU
1	D	10	LEU
1	D	29	VAL
1	D	68	LEU
1	D	91	VAL
1	D	114	ARG
1	D	169	LEU
1	E	92	ASP
1	E	93	LYS
1	E	94	ARG
1	E	114	ARG
1	E	182	VAL
1	F	169	LEU
1	G	114	ARG
1	G	169	LEU
1	H	91	VAL
1	H	93	LYS
1	H	114	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	55	GLN
1	D	3	GLN
1	D	55	GLN
1	D	95	GLN
1	E	3	GLN
1	E	55	GLN
1	F	3	GLN
1	F	55	GLN
1	F	95	GLN
1	G	3	GLN

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Mol	Chain	Res	Type
1	G	55	GLN
1	H	3	GLN
1	H	55	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 27 ligands modelled in this entry, 1 is monoatomic - leaving 26 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	C	203	-	4,4,4	0.34	0	6,6,6	0.24	0
2	SO4	G	201	-	4,4,4	0.29	0	6,6,6	0.21	0
2	SO4	C	202	-	4,4,4	0.21	0	6,6,6	0.27	0
2	SO4	A	201	-	4,4,4	0.30	0	6,6,6	0.34	0
2	SO4	D	204	-	4,4,4	0.29	0	6,6,6	0.12	0
2	SO4	F	202	-	4,4,4	0.45	0	6,6,6	0.14	0
2	SO4	D	201	-	4,4,4	0.33	0	6,6,6	0.24	0
2	SO4	H	204	-	4,4,4	0.20	0	6,6,6	0.32	0
2	SO4	E	203	-	4,4,4	0.25	0	6,6,6	0.18	0
2	SO4	E	204	-	4,4,4	0.38	0	6,6,6	0.21	0
2	SO4	C	201	-	4,4,4	0.20	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	E	201	-	4,4,4	0.33	0	6,6,6	0.31	0
2	SO4	E	202	-	4,4,4	0.19	0	6,6,6	0.27	0
2	SO4	F	203	-	4,4,4	0.25	0	6,6,6	0.37	0
2	SO4	H	201	-	4,4,4	0.38	0	6,6,6	0.39	0
2	SO4	B	202	-	4,4,4	0.30	0	6,6,6	0.24	0
2	SO4	D	205	-	4,4,4	0.28	0	6,6,6	0.20	0
2	SO4	G	202	-	4,4,4	0.29	0	6,6,6	0.25	0
2	SO4	F	201	-	4,4,4	0.38	0	6,6,6	0.35	0
2	SO4	D	203	-	4,4,4	0.39	0	6,6,6	0.34	0
2	SO4	A	202	-	4,4,4	0.25	0	6,6,6	0.50	0
2	SO4	H	203	-	4,4,4	0.27	0	6,6,6	0.45	0
2	SO4	D	202	-	4,4,4	0.18	0	6,6,6	0.22	0
2	SO4	B	201	-	4,4,4	0.35	0	6,6,6	0.18	0
2	SO4	B	203	-	4,4,4	0.32	0	6,6,6	0.05	0
2	SO4	H	202	-	4,4,4	0.36	0	6,6,6	0.10	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	202	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	184/192 (95%)	0.20	5 (2%)	56	63	47, 71, 107, 134	0
1	B	183/192 (95%)	0.35	8 (4%)	39	45	43, 61, 98, 127	0
1	C	183/192 (95%)	0.29	9 (4%)	35	41	45, 68, 101, 113	0
1	D	183/192 (95%)	0.05	4 (2%)	62	67	42, 57, 90, 103	0
1	E	184/192 (95%)	0.31	7 (3%)	44	50	49, 70, 106, 118	0
1	F	183/192 (95%)	0.37	10 (5%)	30	36	48, 70, 108, 137	0
1	G	183/192 (95%)	0.31	4 (2%)	62	67	49, 68, 97, 111	0
1	H	183/192 (95%)	0.02	2 (1%)	78	82	42, 57, 83, 98	0
All	All	1466/1536 (95%)	0.24	49 (3%)	49	56	42, 66, 101, 137	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	185	LEU	6.3
1	C	184	GLY	5.8
1	G	184	GLY	5.3
1	D	184	GLY	5.1
1	F	129	ALA	4.8
1	F	184	GLY	4.8
1	F	2	SER	4.4
1	E	185	LEU	4.2
1	C	130	PHE	4.1
1	A	129	ALA	4.1
1	C	129	ALA	4.0
1	B	184	GLY	4.0
1	F	130	PHE	3.9
1	D	2	SER	3.7
1	F	132	ILE	3.5
1	D	90	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	F	160	ALA	3.3
1	B	68	LEU	3.0
1	C	26	SER	3.0
1	C	131	GLY	3.0
1	B	69	VAL	2.9
1	F	93	LYS	2.8
1	A	130	PHE	2.8
1	F	131	GLY	2.7
1	H	184	GLY	2.7
1	G	172	ARG	2.7
1	E	39	PRO	2.6
1	B	39	PRO	2.5
1	B	26	SER	2.5
1	C	176	ALA	2.5
1	B	60	GLY	2.4
1	E	132	ILE	2.4
1	C	2	SER	2.4
1	D	161	TYR	2.3
1	H	26	SER	2.3
1	F	174	HIS	2.3
1	E	171	PRO	2.3
1	E	2	SER	2.3
1	G	132	ILE	2.3
1	E	40	GLY	2.2
1	F	128	MET	2.2
1	A	131	GLY	2.2
1	C	132	ILE	2.1
1	C	128	MET	2.1
1	B	67	LEU	2.1
1	E	28	ASN	2.1
1	B	90	SER	2.0
1	G	161	TYR	2.0
1	A	2	SER	2.0

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

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

6.3 Carbohydrates ⓘ

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	H	201	5/5	0.67	0.32	69,71,72,72	5
2	SO4	E	204	5/5	0.75	0.18	75,76,78,78	5
2	SO4	A	202	5/5	0.77	0.23	67,68,70,71	5
2	SO4	D	203	5/5	0.77	0.24	78,79,80,81	5
2	SO4	E	203	5/5	0.78	0.22	96,97,98,98	5
2	SO4	D	205	5/5	0.78	0.25	64,66,67,67	5
2	SO4	G	202	5/5	0.78	0.23	83,85,86,86	5
2	SO4	E	201	5/5	0.78	0.30	68,70,71,75	5
2	SO4	H	204	5/5	0.80	0.22	70,72,73,74	5
2	SO4	H	202	5/5	0.81	0.26	75,77,79,80	5
2	SO4	H	203	5/5	0.81	0.29	75,78,82,83	5
2	SO4	A	201	5/5	0.81	0.27	68,71,73,77	5
2	SO4	D	202	5/5	0.82	0.23	69,72,73,73	5
2	SO4	B	203	5/5	0.85	0.11	89,89,91,91	5
2	SO4	C	202	5/5	0.85	0.17	82,82,86,86	5
2	SO4	F	203	5/5	0.85	0.16	67,68,72,72	5
2	SO4	E	202	5/5	0.85	0.18	82,84,84,86	5
2	SO4	C	203	5/5	0.86	0.11	66,69,70,74	5
3	CL	H	205	1/1	0.86	0.24	98,98,98,98	0
2	SO4	B	202	5/5	0.87	0.24	67,69,70,73	5
2	SO4	D	201	5/5	0.87	0.18	71,74,76,79	5
2	SO4	F	202	5/5	0.88	0.20	74,75,76,78	5
2	SO4	D	204	5/5	0.90	0.13	100,101,101,103	5
2	SO4	G	201	5/5	0.90	0.19	81,83,85,85	5
2	SO4	C	201	5/5	0.91	0.16	86,86,87,89	5
2	SO4	B	201	5/5	0.91	0.12	73,75,77,80	5
2	SO4	F	201	5/5	0.91	0.14	54,56,57,59	5

6.5 Other polymers [i](#)

There are no such residues in this entry.