



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 11:24 PM BST

PDB ID : 6K40
Title : Crystal structure of alkyl hydroperoxide reductase from *D. radiodurans* R1
Authors : Kim, M.-K.; Zhang, J.; Zhao, L.
Deposited on : 2019-05-22
Resolution : 2.27 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

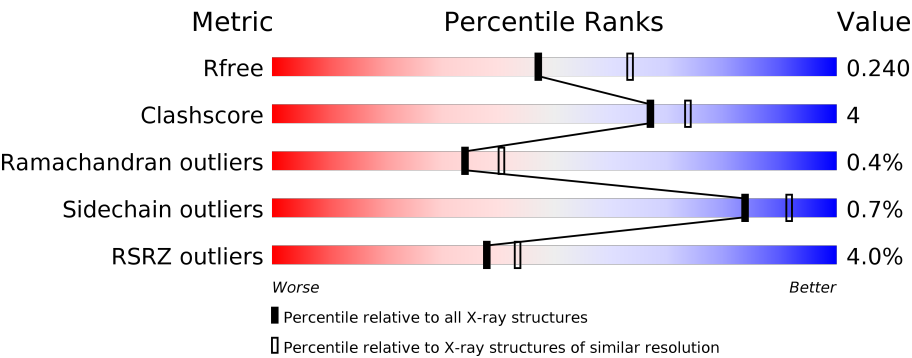
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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}	130704	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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 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	197	<div><div>3%</div><div></div><div>88%</div><div>7%</div><div>6%</div></div>
1	B	197	<div><div>3%</div><div></div><div>88%</div><div>6%</div><div>6%</div></div>
1	C	197	<div><div>4%</div><div></div><div>84%</div><div>10%</div><div>6%</div></div>
1	D	197	<div><div>2%</div><div></div><div>85%</div><div>9%</div><div>6%</div></div>
1	E	197	<div><div>5%</div><div></div><div>83%</div><div>10%</div><div>6%</div></div>
1	F	197	<div><div>4%</div><div></div><div>82%</div><div>12%</div><div>6%</div></div>

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Mol	Chain	Length	Quality of chain
1	G	197	
1	H	197	
1	I	197	
1	J	197	
1	K	197	
1	L	197	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	TRS	G	302	-	-	-	X
4	TRS	I	202	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18705 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 Alkyl hydroperoxide reductase AhpD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	186	Total	C	N	O	S	0	0	0
			1437	907	254	268	8			
1	B	185	Total	C	N	O	S	0	0	0
			1430	902	253	267	8			
1	C	186	Total	C	N	O	S	0	0	0
			1437	907	254	268	8			
1	D	185	Total	C	N	O	S	0	0	0
			1430	902	253	267	8			
1	E	185	Total	C	N	O	S	0	0	0
			1430	902	253	267	8			
1	G	186	Total	C	N	O	S	0	0	0
			1437	907	254	268	8			
1	F	185	Total	C	N	O	S	0	0	0
			1430	902	253	267	8			
1	H	185	Total	C	N	O	S	0	0	0
			1430	902	253	267	8			
1	I	185	Total	C	N	O	S	0	0	0
			1430	902	253	267	8			
1	J	185	Total	C	N	O	S	0	0	0
			1430	902	253	267	8			
1	K	186	Total	C	N	O	S	0	0	0
			1437	907	254	268	8			
1	L	186	Total	C	N	O	S	0	0	0
			1437	907	254	268	8			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	initiating methionine	UNP Q9RTJ7
A	-4	HIS	-	expression tag	UNP Q9RTJ7
A	-3	HIS	-	expression tag	UNP Q9RTJ7
A	-2	HIS	-	expression tag	UNP Q9RTJ7
A	-1	HIS	-	expression tag	UNP Q9RTJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP Q9RTJ7
A	1	HIS	-	expression tag	UNP Q9RTJ7
B	-5	MET	-	initiating methionine	UNP Q9RTJ7
B	-4	HIS	-	expression tag	UNP Q9RTJ7
B	-3	HIS	-	expression tag	UNP Q9RTJ7
B	-2	HIS	-	expression tag	UNP Q9RTJ7
B	-1	HIS	-	expression tag	UNP Q9RTJ7
B	0	HIS	-	expression tag	UNP Q9RTJ7
B	1	HIS	-	expression tag	UNP Q9RTJ7
C	-5	MET	-	initiating methionine	UNP Q9RTJ7
C	-4	HIS	-	expression tag	UNP Q9RTJ7
C	-3	HIS	-	expression tag	UNP Q9RTJ7
C	-2	HIS	-	expression tag	UNP Q9RTJ7
C	-1	HIS	-	expression tag	UNP Q9RTJ7
C	0	HIS	-	expression tag	UNP Q9RTJ7
C	1	HIS	-	expression tag	UNP Q9RTJ7
D	-5	MET	-	initiating methionine	UNP Q9RTJ7
D	-4	HIS	-	expression tag	UNP Q9RTJ7
D	-3	HIS	-	expression tag	UNP Q9RTJ7
D	-2	HIS	-	expression tag	UNP Q9RTJ7
D	-1	HIS	-	expression tag	UNP Q9RTJ7
D	0	HIS	-	expression tag	UNP Q9RTJ7
D	1	HIS	-	expression tag	UNP Q9RTJ7
E	-5	MET	-	initiating methionine	UNP Q9RTJ7
E	-4	HIS	-	expression tag	UNP Q9RTJ7
E	-3	HIS	-	expression tag	UNP Q9RTJ7
E	-2	HIS	-	expression tag	UNP Q9RTJ7
E	-1	HIS	-	expression tag	UNP Q9RTJ7
E	0	HIS	-	expression tag	UNP Q9RTJ7
E	1	HIS	-	expression tag	UNP Q9RTJ7
G	-5	MET	-	initiating methionine	UNP Q9RTJ7
G	-4	HIS	-	expression tag	UNP Q9RTJ7
G	-3	HIS	-	expression tag	UNP Q9RTJ7
G	-2	HIS	-	expression tag	UNP Q9RTJ7
G	-1	HIS	-	expression tag	UNP Q9RTJ7
G	0	HIS	-	expression tag	UNP Q9RTJ7
G	1	HIS	-	expression tag	UNP Q9RTJ7
F	-5	MET	-	initiating methionine	UNP Q9RTJ7
F	-4	HIS	-	expression tag	UNP Q9RTJ7
F	-3	HIS	-	expression tag	UNP Q9RTJ7
F	-2	HIS	-	expression tag	UNP Q9RTJ7
F	-1	HIS	-	expression tag	UNP Q9RTJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	HIS	-	expression tag	UNP Q9RTJ7
F	1	HIS	-	expression tag	UNP Q9RTJ7
H	-5	MET	-	initiating methionine	UNP Q9RTJ7
H	-4	HIS	-	expression tag	UNP Q9RTJ7
H	-3	HIS	-	expression tag	UNP Q9RTJ7
H	-2	HIS	-	expression tag	UNP Q9RTJ7
H	-1	HIS	-	expression tag	UNP Q9RTJ7
H	0	HIS	-	expression tag	UNP Q9RTJ7
H	1	HIS	-	expression tag	UNP Q9RTJ7
I	-5	MET	-	initiating methionine	UNP Q9RTJ7
I	-4	HIS	-	expression tag	UNP Q9RTJ7
I	-3	HIS	-	expression tag	UNP Q9RTJ7
I	-2	HIS	-	expression tag	UNP Q9RTJ7
I	-1	HIS	-	expression tag	UNP Q9RTJ7
I	0	HIS	-	expression tag	UNP Q9RTJ7
I	1	HIS	-	expression tag	UNP Q9RTJ7
J	-5	MET	-	initiating methionine	UNP Q9RTJ7
J	-4	HIS	-	expression tag	UNP Q9RTJ7
J	-3	HIS	-	expression tag	UNP Q9RTJ7
J	-2	HIS	-	expression tag	UNP Q9RTJ7
J	-1	HIS	-	expression tag	UNP Q9RTJ7
J	0	HIS	-	expression tag	UNP Q9RTJ7
J	1	HIS	-	expression tag	UNP Q9RTJ7
K	-5	MET	-	initiating methionine	UNP Q9RTJ7
K	-4	HIS	-	expression tag	UNP Q9RTJ7
K	-3	HIS	-	expression tag	UNP Q9RTJ7
K	-2	HIS	-	expression tag	UNP Q9RTJ7
K	-1	HIS	-	expression tag	UNP Q9RTJ7
K	0	HIS	-	expression tag	UNP Q9RTJ7
K	1	HIS	-	expression tag	UNP Q9RTJ7
L	-5	MET	-	initiating methionine	UNP Q9RTJ7
L	-4	HIS	-	expression tag	UNP Q9RTJ7
L	-3	HIS	-	expression tag	UNP Q9RTJ7
L	-2	HIS	-	expression tag	UNP Q9RTJ7
L	-1	HIS	-	expression tag	UNP Q9RTJ7
L	0	HIS	-	expression tag	UNP Q9RTJ7
L	1	HIS	-	expression tag	UNP Q9RTJ7

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



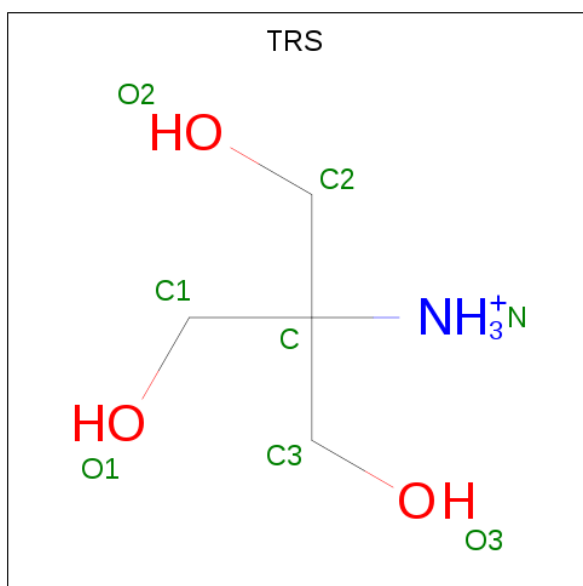
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	J	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		
2	L	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	C	O	0	0
			7	4	3		
3	J	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	G	1	Total	C	N	O	0	0
			8	4	1	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	I	1	Total	C	N	O	0	0
			8	4	1	3		

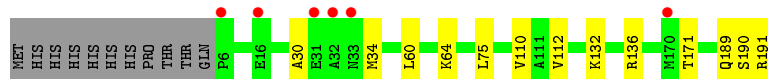
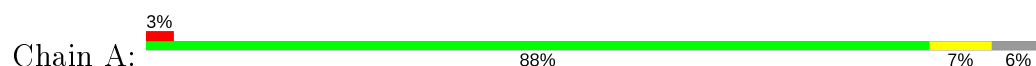
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	154	Total	O	0	0
			154	154		
5	B	148	Total	O	0	0
			148	148		
5	C	138	Total	O	0	0
			138	138		
5	D	119	Total	O	0	0
			119	119		
5	E	133	Total	O	0	0
			133	133		
5	G	147	Total	O	0	0
			147	147		
5	F	72	Total	O	0	0
			72	72		
5	H	66	Total	O	0	0
			66	66		
5	I	151	Total	O	0	0
			151	151		
5	J	86	Total	O	0	0
			86	86		
5	K	113	Total	O	0	0
			113	113		
5	L	75	Total	O	0	0
			75	75		

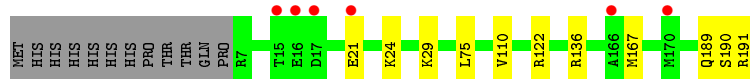
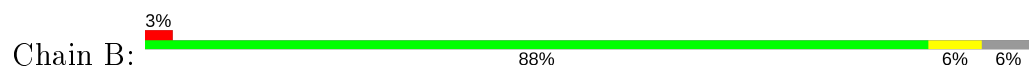
3 Residue-property plots [i](#)

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.

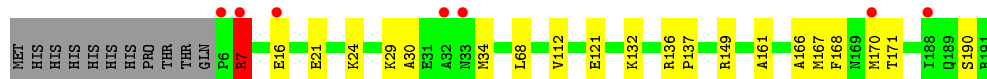
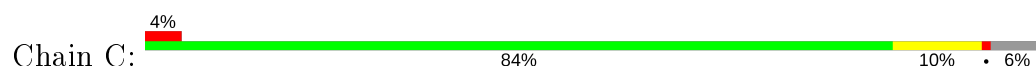
- Molecule 1: Alkyl hydroperoxide reductase AhpD



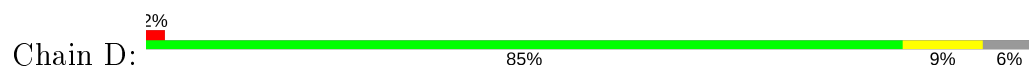
- Molecule 1: Alkyl hydroperoxide reductase AhpD



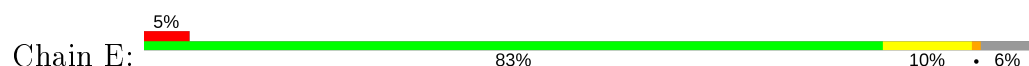
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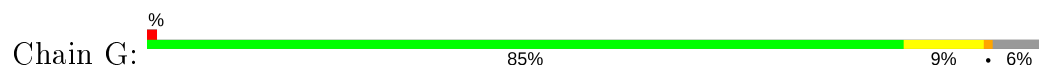
- Molecule 1: Alkyl hydroperoxide reductase AhpD



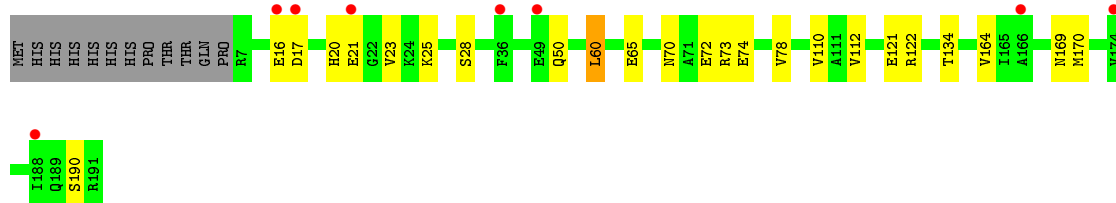
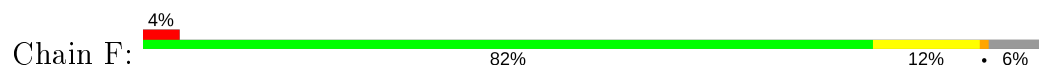
- Molecule 1: Alkyl hydroperoxide reductase AhpD



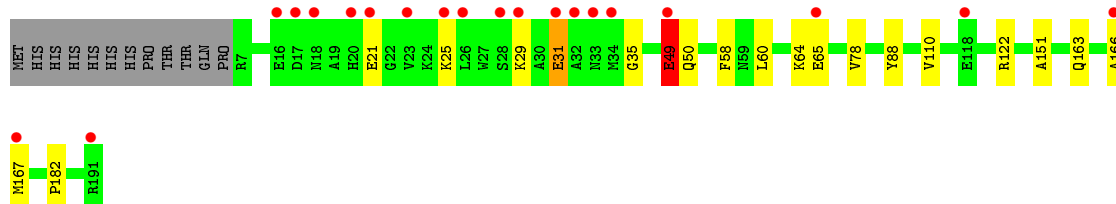
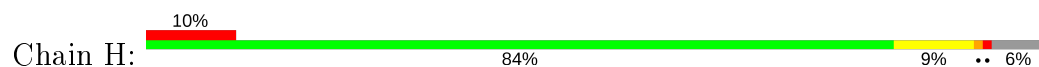
- Molecule 1: Alkyl hydroperoxide reductase AhpD



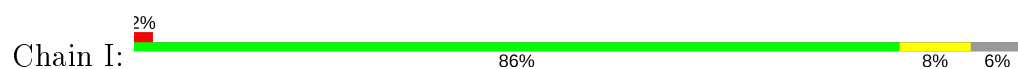
- Molecule 1: Alkyl hydroperoxide reductase AhpD



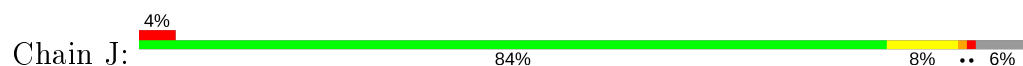
- Molecule 1: Alkyl hydroperoxide reductase AhpD



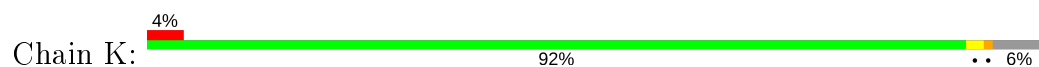
- Molecule 1: Alkyl hydroperoxide reductase AhpD



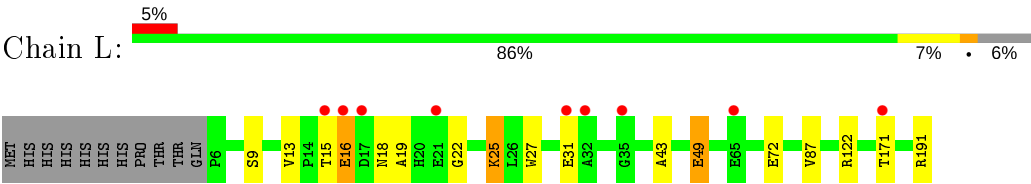
- Molecule 1: Alkyl hydroperoxide reductase AhpD



- Molecule 1: Alkyl hydroperoxide reductase AhpD



● Molecule 1: Alkyl hydroperoxide reductase AhpD



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.09Å 166.86Å 285.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.10 – 2.27 35.10 – 2.27	Depositor EDS
% Data completeness (in resolution range)	91.1 (35.10-2.27) 89.8 (35.10-2.27)	Depositor EDS
R_{merge}	0.51	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.93 (at 2.27Å)	Xtriage
Refinement program	PHENIX 1.15_3459	Depositor
R, R_{free}	0.208 , 0.239 0.209 , 0.240	Depositor DCC
R_{free} test set	1997 reflections (1.73%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	18705	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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: GOL, TRS, PEG

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	0.28	0/1466	0.47	0/1989
1	B	0.27	0/1458	0.45	1/1978 (0.1%)
1	C	0.37	0/1466	0.59	3/1989 (0.2%)
1	D	0.34	0/1458	0.59	6/1978 (0.3%)
1	E	0.35	0/1458	0.57	2/1978 (0.1%)
1	F	0.41	0/1458	0.68	8/1978 (0.4%)
1	G	0.32	0/1466	0.53	2/1989 (0.1%)
1	H	0.44	1/1458 (0.1%)	0.71	5/1978 (0.3%)
1	I	0.26	0/1458	0.46	0/1978
1	J	0.38	0/1458	0.64	4/1978 (0.2%)
1	K	0.33	0/1466	0.51	1/1989 (0.1%)
1	L	0.38	0/1466	0.67	4/1989 (0.2%)
All	All	0.35	1/17536 (0.0%)	0.58	36/23791 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	L	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	49	GLU	CB-CG	6.92	1.65	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	16	GLU	CA-CB-CG	-10.04	91.32	113.40
1	F	25	LYS	CB-CG-CD	-9.21	87.65	111.60
1	H	65	GLU	N-CA-CB	8.17	125.30	110.60
1	E	31	GLU	CA-CB-CG	-8.11	95.57	113.40
1	F	65	GLU	CA-CB-CG	7.24	129.32	113.40
1	J	24	LYS	CD-CE-NZ	-6.81	96.04	111.70
1	D	49	GLU	CA-CB-CG	6.73	128.21	113.40
1	F	121	GLU	CA-CB-CG	6.59	127.90	113.40
1	L	49	GLU	CB-CA-C	-6.57	97.26	110.40
1	D	21	GLU	CB-CA-C	6.51	123.43	110.40
1	F	25	LYS	CA-CB-CG	6.49	127.67	113.40
1	J	17	ASP	CB-CG-OD1	-6.12	112.79	118.30
1	F	60	LEU	CB-CG-CD2	-6.10	100.64	111.00
1	C	29	LYS	CB-CG-CD	-6.07	95.83	111.60
1	C	121	GLU	CA-CB-CG	6.05	126.71	113.40
1	F	16	GLU	CA-CB-CG	5.86	126.29	113.40
1	L	25	LYS	CB-CG-CD	5.82	126.72	111.60
1	H	49	GLU	CB-CA-C	-5.71	98.97	110.40
1	J	121	GLU	CA-CB-CG	5.67	125.87	113.40
1	B	21	GLU	CA-CB-CG	5.64	125.82	113.40
1	C	7	ARG	CB-CG-CD	5.61	126.20	111.60
1	H	29	LYS	CA-CB-CG	5.56	125.63	113.40
1	H	31	GLU	CB-CA-C	-5.45	99.49	110.40
1	E	31	GLU	CG-CD-OE1	-5.45	107.40	118.30
1	D	49	GLU	CB-CG-CD	-5.42	99.56	114.20
1	L	22	GLY	N-CA-C	-5.39	99.64	113.10
1	G	65	GLU	CA-CB-CG	5.36	125.19	113.40
1	F	17	ASP	CB-CG-OD2	-5.31	113.52	118.30
1	K	21	GLU	CB-CA-C	-5.30	99.80	110.40
1	F	21	GLU	CA-CB-CG	-5.29	101.77	113.40
1	J	121	GLU	CB-CA-C	5.14	120.69	110.40
1	D	17	ASP	CB-CG-OD2	-5.11	113.70	118.30
1	D	48	GLY	C-N-CA	-5.09	108.96	121.70
1	G	21	GLU	N-CA-CB	-5.09	101.44	110.60
1	D	21	GLU	N-CA-CB	-5.08	101.46	110.60
1	H	65	GLU	CA-CB-CG	5.03	124.46	113.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	16	GLU	Peptide
1	L	31	GLU	Sidechain

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	1437	0	1411	10	0
1	B	1430	0	1403	7	0
1	C	1437	0	1411	13	0
1	D	1430	0	1403	13	0
1	E	1430	0	1403	19	0
1	F	1430	0	1403	19	0
1	G	1437	0	1411	14	0
1	H	1430	0	1403	21	0
1	I	1430	0	1403	10	0
1	J	1430	0	1403	11	0
1	K	1437	0	1411	6	0
1	L	1437	0	1411	13	0
2	A	6	0	8	1	0
2	B	6	0	8	0	0
2	C	6	0	8	0	0
2	D	12	0	16	1	0
2	E	6	0	8	1	0
2	F	6	0	8	0	0
2	G	6	0	8	0	0
2	H	6	0	8	0	0
2	I	6	0	8	0	0
2	J	6	0	8	0	0
2	K	6	0	8	0	0
2	L	6	0	8	1	0
3	E	7	0	10	1	0
3	J	7	0	10	0	0
4	G	8	0	12	0	3
4	I	8	0	12	0	3
5	A	154	0	0	1	0
5	B	148	0	0	3	0
5	C	138	0	0	3	0
5	D	119	0	0	4	0
5	E	133	0	0	4	0
5	F	72	0	0	0	0
5	G	147	0	0	2	0
5	H	66	0	0	0	0
5	I	151	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	J	86	0	0	0	0
5	K	113	0	0	0	0
5	L	75	0	0	1	0
All	All	18705	0	17024	132	3

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 4.

All (132) 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:F:60:LEU:HD21	1:H:49:GLU:HG2	1.25	1.11
1:F:60:LEU:HD21	1:H:49:GLU:CG	1.82	1.07
1:F:60:LEU:HD11	1:H:49:GLU:CD	1.84	0.97
1:D:134:THR:HG22	1:D:169:ASN:HD21	1.36	0.88
1:K:21:GLU:N	1:K:21:GLU:OE1	2.07	0.88
1:E:31:GLU:OE1	1:E:36:PHE:N	2.09	0.85
1:J:21:GLU:O	1:J:25:LYS:HB2	1.79	0.83
1:I:134:THR:HG22	1:I:169:ASN:HD21	1.44	0.81
1:C:7:ARG:HD3	5:C:303:HOH:O	1.81	0.80
1:H:88:TYR:HA	1:H:182:PRO:HG3	1.63	0.79
1:A:191:ARG:HG2	1:G:191:ARG:HG2	1.68	0.76
1:H:58:PHE:HD2	1:H:167:MET:SD	2.10	0.73
1:F:60:LEU:HD11	1:H:49:GLU:OE2	1.89	0.73
1:H:60:LEU:HA	1:H:64:LYS:HD3	1.72	0.71
1:A:189:GLN:O	1:A:191:ARG:N	2.24	0.70
1:E:121:GLU:OE2	5:E:301:HOH:O	2.10	0.69
1:E:31:GLU:OE1	1:E:35:GLY:C	2.31	0.69
1:E:48:GLY:O	1:E:52:LEU:HD22	1.94	0.68
1:A:171:THR:HB	2:A:201:GOL:H12	1.75	0.67
1:G:29:LYS:HE3	1:G:33:ASN:OD1	1.96	0.66
1:A:136:ARG:NH1	5:A:301:HOH:O	2.31	0.64
1:F:60:LEU:CD2	1:H:49:GLU:HG2	2.15	0.64
1:L:171:THR:HB	2:L:201:GOL:H11	1.79	0.64
1:B:29:LYS:HA	1:J:188:ILE:HG21	1.80	0.64
1:E:31:GLU:OE1	1:E:36:PHE:CA	2.47	0.62
1:D:136:ARG:NH1	5:D:305:HOH:O	2.30	0.62
1:I:130:ALA:O	1:I:134:THR:HG23	1.99	0.62
2:D:201:GOL:O1	5:D:301:HOH:O	2.15	0.62
1:B:75:LEU:HD12	1:B:110:VAL:HG11	1.81	0.61
1:C:7:ARG:CD	5:C:303:HOH:O	2.44	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:21:GLU:HG2	1:G:22:GLY:N	2.14	0.60
1:D:130:ALA:O	1:D:134:THR:HG23	2.01	0.60
1:I:7:ARG:N	5:I:303:HOH:O	2.34	0.60
1:C:7:ARG:NE	5:C:303:HOH:O	2.35	0.60
1:G:29:LYS:NZ	5:G:401:HOH:O	2.28	0.60
1:G:75:LEU:HD12	1:G:110:VAL:HG11	1.83	0.59
1:F:60:LEU:HD21	1:H:49:GLU:HG3	1.79	0.59
1:B:24:LYS:NZ	5:B:302:HOH:O	2.35	0.59
1:E:75:LEU:HD12	1:E:110:VAL:HG11	1.85	0.59
1:F:60:LEU:CD1	1:H:49:GLU:CD	2.66	0.58
1:C:30:ALA:O	1:C:34:MET:HG2	2.05	0.57
1:C:149:ARG:NH2	5:E:302:HOH:O	2.32	0.56
1:A:75:LEU:HD12	1:A:110:VAL:HG11	1.87	0.56
1:F:20:HIS:HB3	1:F:23:VAL:HG23	1.88	0.56
1:J:21:GLU:OE2	1:J:24:LYS:NZ	2.39	0.56
1:D:75:LEU:HD12	1:D:110:VAL:HG11	1.88	0.55
1:H:21:GLU:O	1:H:25:LYS:HG3	2.05	0.55
1:G:14:PRO:HG3	1:G:46:LEU:HA	1.89	0.55
1:E:93:HIS:NE2	2:E:201:GOL:H32	2.21	0.54
1:A:189:GLN:C	1:A:191:ARG:H	2.09	0.54
1:E:29:LYS:HD3	1:E:33:ASN:ND2	2.22	0.54
1:C:137:PRO:HD2	3:E:202:PEG:H42	1.90	0.54
1:C:21:GLU:HA	1:C:24:LYS:HD3	1.89	0.54
1:J:75:LEU:HD12	1:J:110:VAL:HG11	1.90	0.53
1:E:48:GLY:O	1:E:52:LEU:CD2	2.55	0.53
1:L:122:ARG:NH2	5:L:304:HOH:O	2.42	0.52
1:D:166:ALA:HB1	1:G:170:MET:HG3	1.92	0.51
1:F:170:MET:HG2	1:H:166:ALA:HB1	1.93	0.51
1:C:166:ALA:HB1	1:E:170:MET:HG3	1.91	0.51
1:F:60:LEU:HD11	1:H:49:GLU:OE1	2.10	0.50
1:D:30:ALA:O	1:D:34:MET:HG2	2.11	0.50
1:B:136:ARG:NH1	5:B:306:HOH:O	2.43	0.50
1:E:7:ARG:N	5:E:310:HOH:O	2.45	0.50
1:A:132:LYS:HG3	1:A:136:ARG:NH2	2.27	0.50
1:K:20:HIS:ND1	1:K:21:GLU:OE1	2.45	0.49
1:D:191:ARG:NH1	5:D:307:HOH:O	2.36	0.49
1:I:134:THR:HG22	1:I:169:ASN:ND2	2.21	0.49
1:E:21:GLU:O	1:E:25:LYS:HG3	2.13	0.48
1:I:142:GLU:OE2	1:K:7:ARG:NH2	2.37	0.48
1:E:21:GLU:CD	1:E:21:GLU:H	2.16	0.48
1:L:15:THR:N	1:L:18:ASN:OD1	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:149:ARG:NH2	5:D:311:HOH:O	2.46	0.47
1:D:134:THR:HG22	1:D:169:ASN:ND2	2.16	0.47
1:C:167:MET:HG2	1:E:170:MET:HE3	1.96	0.47
1:F:60:LEU:CD1	1:H:49:GLU:OE1	2.63	0.46
1:C:68:LEU:HD11	1:C:161:ALA:HB2	1.98	0.46
1:E:125:ALA:HB2	5:E:301:HOH:O	2.16	0.46
1:H:78:VAL:HG11	1:H:110:VAL:HG23	1.98	0.46
1:A:30:ALA:O	1:A:34:MET:HG2	2.15	0.46
1:D:166:ALA:CB	1:G:170:MET:HG3	2.45	0.46
1:G:136:ARG:NH1	5:G:407:HOH:O	2.34	0.46
1:I:20:HIS:HB3	1:I:23:VAL:HG23	1.98	0.46
1:I:166:ALA:HB1	1:K:170:MET:HG3	1.98	0.46
1:E:31:GLU:HG3	1:E:31:GLU:O	2.14	0.45
1:B:122:ARG:NH2	5:B:308:HOH:O	2.45	0.45
1:C:132:LYS:HG3	1:C:136:ARG:NH2	2.31	0.45
1:J:120:SER:N	1:J:123:GLU:OE1	2.34	0.45
1:L:13:VAL:HG12	1:L:43:ALA:HA	1.98	0.45
1:E:68:LEU:HD11	1:E:161:ALA:HB2	1.99	0.45
1:L:16:GLU:CG	1:L:27:TRP:HZ3	2.29	0.45
1:H:60:LEU:HA	1:H:64:LYS:HB2	1.98	0.45
1:E:29:LYS:HD3	1:E:33:ASN:HD21	1.81	0.45
1:F:73:ARG:HG2	1:F:164:VAL:HG21	1.99	0.45
1:I:64:LYS:O	1:I:73:ARG:NH1	2.45	0.45
1:L:72:GLU:OE2	1:L:122:ARG:NE	2.45	0.45
1:J:15:THR:C	1:J:17:ASP:H	2.19	0.44
1:B:191:ARG:HD3	1:B:191:ARG:HA	1.82	0.44
1:K:20:HIS:CE1	1:K:21:GLU:OE1	2.71	0.44
1:F:72:GLU:OE2	1:F:122:ARG:NE	2.49	0.43
1:G:14:PRO:HD2	1:G:42:ARG:O	2.19	0.43
1:F:50:GLN:HB3	1:H:163:GLN:HE22	1.83	0.43
1:I:16:GLU:O	1:I:24:LYS:HE2	2.18	0.43
1:L:15:THR:O	1:L:19:ALA:HB2	2.18	0.43
1:F:134:THR:HG23	1:F:169:ASN:OD1	2.18	0.43
1:H:31:GLU:OE2	1:H:35:GLY:HA2	2.18	0.43
1:L:9:SER:HB2	1:L:43:ALA:HB1	2.01	0.43
1:F:78:VAL:HG11	1:F:110:VAL:HG23	2.01	0.43
1:L:191:ARG:HA	1:L:191:ARG:HD3	1.73	0.42
1:F:112:VAL:HB	1:L:87:VAL:HG22	2.01	0.42
1:H:49:GLU:HB3	1:H:50:GLN:H	1.60	0.42
1:L:16:GLU:HG3	1:L:27:TRP:HZ3	1.84	0.42
1:B:189:GLN:O	1:B:191:ARG:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:60:LEU:HD21	1:L:49:GLU:CB	2.50	0.42
1:K:20:HIS:HB3	1:K:23:VAL:HG23	2.02	0.42
1:C:168:PHE:HA	1:C:171:THR:OG1	2.20	0.42
1:G:29:LYS:CE	1:G:33:ASN:OD1	2.64	0.42
1:J:30:ALA:O	1:J:34:MET:HG2	2.20	0.42
1:C:112:VAL:HB	1:D:87:VAL:HG22	2.02	0.41
1:I:68:LEU:HD11	1:I:161:ALA:HB2	2.02	0.41
1:J:98:ARG:HD2	1:J:191:ARG:HA	2.02	0.41
1:J:60:LEU:HD21	1:L:49:GLU:HB3	2.03	0.41
1:A:112:VAL:HB	1:G:87:VAL:HG22	2.01	0.41
1:E:31:GLU:OE1	1:E:36:PHE:HA	2.21	0.41
1:D:68:LEU:HD11	1:D:161:ALA:HB2	2.02	0.41
1:D:191:ARG:HD3	1:D:191:ARG:HA	1.88	0.41
1:A:60:LEU:HA	1:A:64:LYS:HD3	2.03	0.40
1:F:70:ASN:O	1:F:74:GLU:HG2	2.21	0.40
1:G:132:LYS:HG3	1:G:136:ARG:NH2	2.36	0.40
1:J:75:LEU:CD1	1:J:110:VAL:HG11	2.52	0.40
1:G:69:SER:O	1:G:73:ARG:HD2	2.21	0.40
1:H:122:ARG:HG3	1:H:151:ALA:HB1	2.04	0.40
1:F:60:LEU:CD2	1:H:49:GLU:CG	2.75	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:302:TRS:C2	4:I:202:TRS:C2[4_445]	1.49	0.71
4:G:302:TRS:C2	4:I:202:TRS:O2[4_445]	1.92	0.28
4:G:302:TRS:O2	4:I:202:TRS:O2[4_445]	2.14	0.06

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	184/197 (93%)	183 (100%)	0	1 (0%)	29	34
1	B	183/197 (93%)	181 (99%)	1 (0%)	1 (0%)	29	34
1	C	184/197 (93%)	181 (98%)	1 (0%)	2 (1%)	14	14
1	D	183/197 (93%)	181 (99%)	2 (1%)	0	100	100
1	E	183/197 (93%)	182 (100%)	1 (0%)	0	100	100
1	F	183/197 (93%)	182 (100%)	0	1 (0%)	29	34
1	G	184/197 (93%)	183 (100%)	0	1 (0%)	29	34
1	H	183/197 (93%)	180 (98%)	2 (1%)	1 (0%)	29	34
1	I	183/197 (93%)	182 (100%)	1 (0%)	0	100	100
1	J	183/197 (93%)	181 (99%)	1 (0%)	1 (0%)	29	34
1	K	184/197 (93%)	183 (100%)	1 (0%)	0	100	100
1	L	184/197 (93%)	182 (99%)	2 (1%)	0	100	100
All	All	2201/2364 (93%)	2181 (99%)	12 (0%)	8 (0%)	34	40

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	SER
1	B	190	SER
1	C	7	ARG
1	C	190	SER
1	G	190	SER
1	H	49	GLU
1	F	190	SER
1	J	190	SER

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	147/158 (93%)	147 (100%)	0	100	100
1	B	146/158 (92%)	145 (99%)	1 (1%)	84	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	147/158 (93%)	145 (99%)	2 (1%)	67	79
1	D	146/158 (92%)	146 (100%)	0	100	100
1	E	146/158 (92%)	143 (98%)	3 (2%)	53	68
1	F	146/158 (92%)	145 (99%)	1 (1%)	84	91
1	G	147/158 (93%)	147 (100%)	0	100	100
1	H	146/158 (92%)	146 (100%)	0	100	100
1	I	146/158 (92%)	145 (99%)	1 (1%)	84	91
1	J	146/158 (92%)	142 (97%)	4 (3%)	44	59
1	K	147/158 (93%)	147 (100%)	0	100	100
1	L	147/158 (93%)	146 (99%)	1 (1%)	84	91
All	All	1757/1896 (93%)	1744 (99%)	13 (1%)	84	91

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	167	MET
1	C	7	ARG
1	C	170	MET
1	E	16	GLU
1	E	28	SER
1	E	153	LEU
1	F	28	SER
1	I	17	ASP
1	J	17	ASP
1	J	25	LYS
1	J	69	SER
1	J	167	MET
1	L	25	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	84	ASN
1	D	169	ASN
1	E	33	ASN
1	H	18	ASN
1	H	163	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

17 ligands are modelled in this entry.

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	GOL	D	202	-	5,5,5	0.99	0	5,5,5	0.98	0
2	GOL	L	201	-	5,5,5	0.87	0	5,5,5	1.02	0
2	GOL	I	201	-	5,5,5	0.84	0	5,5,5	1.00	0
3	PEG	J	202	-	6,6,6	0.48	0	5,5,5	0.22	0
2	GOL	K	201	-	5,5,5	0.88	0	5,5,5	0.89	0
2	GOL	H	201	-	5,5,5	0.92	0	5,5,5	1.02	0
2	GOL	J	201	-	5,5,5	0.88	0	5,5,5	0.92	0
3	PEG	E	202	-	6,6,6	0.48	0	5,5,5	0.30	0
2	GOL	F	201	-	5,5,5	0.98	0	5,5,5	0.88	0
4	TRS	I	202	-	7,7,7	0.36	0	9,9,9	0.52	0
4	TRS	G	302	-	7,7,7	0.29	0	9,9,9	0.39	0
2	GOL	E	201	-	5,5,5	0.90	0	5,5,5	1.02	0
2	GOL	D	201	-	5,5,5	0.92	0	5,5,5	1.02	0
2	GOL	A	201	-	5,5,5	1.01	0	5,5,5	0.89	0
2	GOL	G	301	-	5,5,5	0.84	0	5,5,5	0.99	0
2	GOL	C	201	-	5,5,5	0.90	0	5,5,5	0.99	0
2	GOL	B	201	-	5,5,5	0.98	0	5,5,5	0.92	0

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	GOL	D	202	-	-	2/4/4/4	-
2	GOL	L	201	-	-	2/4/4/4	-
2	GOL	I	201	-	-	2/4/4/4	-
3	PEG	J	202	-	-	1/4/4/4	-
2	GOL	K	201	-	-	2/4/4/4	-
2	GOL	H	201	-	-	2/4/4/4	-
2	GOL	J	201	-	-	2/4/4/4	-
3	PEG	E	202	-	-	2/4/4/4	-
2	GOL	F	201	-	-	2/4/4/4	-
4	TRS	I	202	-	-	9/9/9/9	-
4	TRS	G	302	-	-	0/9/9/9	-
2	GOL	E	201	-	-	2/4/4/4	-
2	GOL	D	201	-	-	1/4/4/4	-
2	GOL	A	201	-	-	2/4/4/4	-
2	GOL	G	301	-	-	2/4/4/4	-
2	GOL	C	201	-	-	0/4/4/4	-
2	GOL	B	201	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	202	GOL	C1-C2-C3-O3
2	L	201	GOL	O1-C1-C2-C3
2	F	201	GOL	O1-C1-C2-O2
2	F	201	GOL	O1-C1-C2-C3
2	K	201	GOL	O1-C1-C2-O2
2	K	201	GOL	O1-C1-C2-C3
2	H	201	GOL	O1-C1-C2-O2
2	H	201	GOL	O1-C1-C2-C3
4	I	202	TRS	N-C-C1-O1
2	E	201	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
2	A	201	GOL	O1-C1-C2-O2
2	A	201	GOL	O1-C1-C2-C3
2	G	301	GOL	O1-C1-C2-O2
2	G	301	GOL	O1-C1-C2-C3
2	B	201	GOL	O1-C1-C2-O2
2	B	201	GOL	O1-C1-C2-C3
2	E	201	GOL	O1-C1-C2-O2
3	E	202	PEG	O2-C3-C4-O4
3	E	202	PEG	O1-C1-C2-O2
2	I	201	GOL	O1-C1-C2-C3
2	J	201	GOL	O1-C1-C2-C3
2	I	201	GOL	O1-C1-C2-O2
4	I	202	TRS	C3-C-C1-O1
2	D	202	GOL	O2-C2-C3-O3
4	I	202	TRS	N-C-C2-O2
4	I	202	TRS	C1-C-C3-O3
4	I	202	TRS	N-C-C3-O3
3	J	202	PEG	O2-C3-C4-O4
4	I	202	TRS	C2-C-C1-O1
4	I	202	TRS	C1-C-C2-O2
4	I	202	TRS	C2-C-C3-O3
2	L	201	GOL	O1-C1-C2-O2
2	D	201	GOL	C1-C2-C3-O3
2	J	201	GOL	O1-C1-C2-O2
4	I	202	TRS	C3-C-C2-O2

There are no ring outliers.

7 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	201	GOL	1	0
3	E	202	PEG	1	0
4	I	202	TRS	0	3
4	G	302	TRS	0	3
2	E	201	GOL	1	0
2	D	201	GOL	1	0
2	A	201	GOL	1	0

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	186/197 (94%)	-0.12	6 (3%) 47 53	16, 22, 36, 53	0
1	B	185/197 (93%)	-0.13	6 (3%) 47 53	15, 23, 36, 49	0
1	C	186/197 (94%)	0.07	7 (3%) 40 45	19, 28, 48, 62	0
1	D	185/197 (93%)	-0.10	3 (1%) 72 77	17, 27, 40, 63	0
1	E	185/197 (93%)	0.11	9 (4%) 29 35	17, 28, 50, 66	0
1	F	185/197 (93%)	0.30	8 (4%) 35 40	25, 41, 58, 74	0
1	G	186/197 (94%)	-0.12	1 (0%) 91 93	17, 26, 39, 54	0
1	H	185/197 (93%)	0.59	20 (10%) 5 7	26, 42, 66, 84	0
1	I	185/197 (93%)	-0.17	3 (1%) 72 77	20, 27, 36, 47	0
1	J	185/197 (93%)	0.22	8 (4%) 35 40	21, 34, 53, 76	0
1	K	186/197 (94%)	0.13	8 (4%) 35 40	20, 29, 48, 61	0
1	L	186/197 (94%)	0.36	9 (4%) 30 36	24, 38, 55, 76	0
All	All	2225/2364 (94%)	0.09	88 (3%) 38 43	15, 29, 53, 84	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	6	PRO	5.3
1	D	17	ASP	4.5
1	H	17	ASP	4.4
1	L	21	GLU	4.2
1	H	33	ASN	3.9
1	H	16	GLU	3.9
1	C	16	GLU	3.8
1	L	17	ASP	3.8
1	K	118	GLU	3.7
1	J	17	ASP	3.7
1	E	35	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	H	49	GLU	3.6
1	H	32	ALA	3.6
1	J	190	SER	3.6
1	H	25	LYS	3.5
1	A	6	PRO	3.5
1	J	15	THR	3.5
1	E	33	ASN	3.5
1	L	16	GLU	3.5
1	F	21	GLU	3.4
1	J	16	GLU	3.3
1	H	167	MET	3.3
1	C	7	ARG	3.3
1	C	32	ALA	3.2
1	H	21	GLU	3.2
1	L	15	THR	3.1
1	A	16	GLU	3.0
1	F	49	GLU	3.0
1	H	118	GLU	2.9
1	L	171	THR	2.9
1	H	34	MET	2.9
1	B	17	ASP	2.8
1	L	35	GLY	2.8
1	E	15	THR	2.8
1	G	21	GLU	2.8
1	F	16	GLU	2.8
1	A	32	ALA	2.7
1	K	17	ASP	2.7
1	C	33	ASN	2.6
1	J	21	GLU	2.6
1	K	21	GLU	2.6
1	K	33	ASN	2.6
1	H	18	ASN	2.6
1	H	23	VAL	2.5
1	H	65	GLU	2.5
1	B	15	THR	2.5
1	H	28	SER	2.5
1	C	6	PRO	2.5
1	L	31	GLU	2.5
1	E	170	MET	2.4
1	K	32	ALA	2.4
1	I	17	ASP	2.4
1	I	170	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	I	166	ALA	2.4
1	H	31	GLU	2.4
1	B	21	GLU	2.3
1	E	31	GLU	2.3
1	H	191	ARG	2.3
1	E	21	GLU	2.3
1	L	65	GLU	2.3
1	F	166	ALA	2.3
1	H	26	LEU	2.3
1	E	16	GLU	2.3
1	F	188	ILE	2.3
1	L	32	ALA	2.2
1	A	31	GLU	2.2
1	J	33	ASN	2.2
1	F	174	VAL	2.2
1	E	167	MET	2.2
1	F	17	ASP	2.2
1	F	36	PHE	2.1
1	B	166	ALA	2.1
1	H	29	LYS	2.1
1	C	170	MET	2.1
1	J	25	LYS	2.1
1	J	22	GLY	2.1
1	D	166	ALA	2.1
1	H	166	ALA	2.1
1	K	171	THR	2.1
1	A	170	MET	2.1
1	A	33	ASN	2.1
1	K	168	PHE	2.1
1	C	188	ILE	2.1
1	B	16	GLU	2.1
1	D	16	GLU	2.1
1	B	170	MET	2.1
1	E	17	ASP	2.0
1	H	20	HIS	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 monosaccharides in this entry.

6.4 Ligands ⓘ

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	GOL	E	201	6/6	0.52	0.35	34,37,38,39	0
4	TRS	G	302	8/8	0.57	0.62	54,55,57,57	0
2	GOL	J	201	6/6	0.62	0.39	42,45,46,47	0
2	GOL	D	201	6/6	0.72	0.34	39,40,44,50	0
2	GOL	L	201	6/6	0.74	0.33	48,48,48,48	0
4	TRS	I	202	8/8	0.77	0.42	42,43,46,48	0
3	PEG	J	202	7/7	0.78	0.23	36,38,38,39	0
2	GOL	H	201	6/6	0.79	0.35	48,49,49,50	0
2	GOL	F	201	6/6	0.82	0.28	44,46,46,47	0
2	GOL	B	201	6/6	0.86	0.17	28,28,28,28	0
2	GOL	G	301	6/6	0.87	0.18	33,34,34,36	0
2	GOL	I	201	6/6	0.87	0.18	34,34,35,37	0
2	GOL	K	201	6/6	0.88	0.19	37,37,37,38	0
3	PEG	E	202	7/7	0.88	0.16	31,31,32,32	0
2	GOL	C	201	6/6	0.88	0.17	35,37,37,37	0
2	GOL	D	202	6/6	0.88	0.22	28,30,30,31	0
2	GOL	A	201	6/6	0.93	0.13	30,31,31,31	0

6.5 Other polymers ⓘ

There are no such residues in this entry.