



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 05:46 am GMT

PDB ID : 4HN8  
Title : Crystal structure of a putative D-glucarate dehydratase from *Pseudomonas mendocina* ymp  
Authors : Hegde, R.P.; Toro, R.; Burley, S.K.; Almo, S.C.; Ramagopal, U.A.; New York Structural Genomics Research Consortium (NYSGRG)  
Deposited on : 2012-10-19  
Resolution : 2.20 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

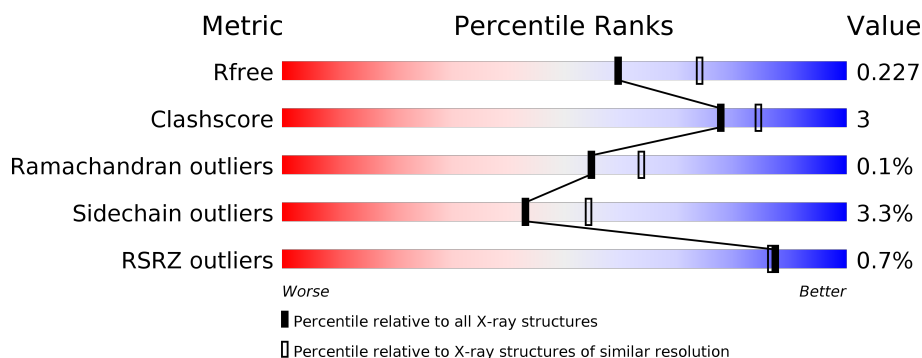
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	473	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 84%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>84%</span> <span>6%</span> <span>10%</span> </div> </div>
1	B	473	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 83%, yellow 6%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>83%</span> <span>6%</span> <span>10%</span> </div> </div>
1	C	473	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 83%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>83%</span> <span>7%</span> <span>10%</span> </div> </div>
1	D	473	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 83%, yellow 7%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>83%</span> <span>7%</span> <span>10%</span> </div> </div>
1	E	473	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 81%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>81%</span> <span>8%</span> <span>10%</span> </div> </div>
1	F	473	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 84%, yellow 6%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>84%</span> <span>6%</span> <span>10%</span> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	473	 % 79% 10% 10%
1	H	473	 % 82% 7% 10%

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
2	GOL	A	502	-	-	-	X
2	GOL	B	501	-	-	-	X
2	GOL	C	501	-	-	-	X
2	GOL	D	501	-	-	-	X
2	GOL	E	501	-	-	-	X
2	GOL	F	501	-	-	-	X
2	GOL	G	501	-	-	X	X
2	GOL	H	501	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 27164 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 D-glucarate dehydratase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	Se	0	4	0
			3328	2089	617	602	6	14			
1	B	426	Total	C	N	O	S	Se	0	1	0
			3301	2071	611	599	6	14			
1	C	427	Total	C	N	O	S	Se	0	7	0
			3348	2102	620	606	6	14			
1	D	427	Total	C	N	O	S	Se	0	6	0
			3333	2095	612	606	6	14			
1	E	426	Total	C	N	O	S	Se	0	3	0
			3312	2079	614	599	6	14			
1	F	426	Total	C	N	O	S	Se	0	3	0
			3319	2082	619	598	6	14			
1	G	426	Total	C	N	O	S	Se	0	3	0
			3314	2079	615	600	6	14			
1	H	427	Total	C	N	O	S	Se	0	2	0
			3313	2081	612	600	6	14			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
A	0	SER	-	EXPRESSION TAG	UNP A4XRL3
A	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
A	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
A	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
A	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
A	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
A	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
A	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
A	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
A	471	HIS	-	EXPRESSION TAG	UNP A4XRL3
B	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
B	0	SER	-	EXPRESSION TAG	UNP A4XRL3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
B	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
B	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
B	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
B	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
B	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
B	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
B	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
B	471	HIS	-	EXPRESSION TAG	UNP A4XRL3
C	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
C	0	SER	-	EXPRESSION TAG	UNP A4XRL3
C	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
C	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
C	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
C	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
C	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
C	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
C	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
C	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
C	471	HIS	-	EXPRESSION TAG	UNP A4XRL3
D	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
D	0	SER	-	EXPRESSION TAG	UNP A4XRL3
D	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
D	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
D	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
D	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
D	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
D	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
D	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
D	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
D	471	HIS	-	EXPRESSION TAG	UNP A4XRL3
E	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
E	0	SER	-	EXPRESSION TAG	UNP A4XRL3
E	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
E	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
E	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
E	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
E	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
E	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
E	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
E	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
E	471	HIS	-	EXPRESSION TAG	UNP A4XRL3

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
F	0	SER	-	EXPRESSION TAG	UNP A4XRL3
F	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
F	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
F	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
F	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
F	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
F	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
F	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
F	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
F	471	HIS	-	EXPRESSION TAG	UNP A4XRL3
G	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
G	0	SER	-	EXPRESSION TAG	UNP A4XRL3
G	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
G	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
G	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
G	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
G	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
G	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
G	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
G	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
G	471	HIS	-	EXPRESSION TAG	UNP A4XRL3
H	-1	MSE	-	EXPRESSION TAG	UNP A4XRL3
H	0	SER	-	EXPRESSION TAG	UNP A4XRL3
H	1	LEU	-	EXPRESSION TAG	UNP A4XRL3
H	464	GLU	-	EXPRESSION TAG	UNP A4XRL3
H	465	GLY	-	EXPRESSION TAG	UNP A4XRL3
H	466	HIS	-	EXPRESSION TAG	UNP A4XRL3
H	467	HIS	-	EXPRESSION TAG	UNP A4XRL3
H	468	HIS	-	EXPRESSION TAG	UNP A4XRL3
H	469	HIS	-	EXPRESSION TAG	UNP A4XRL3
H	470	HIS	-	EXPRESSION TAG	UNP A4XRL3
H	471	HIS	-	EXPRESSION TAG	UNP A4XRL3

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
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	E	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	103	Total	O	0	0
			103	103		
3	B	84	Total	O	0	0
			84	84		
3	C	71	Total	O	0	0
			71	71		

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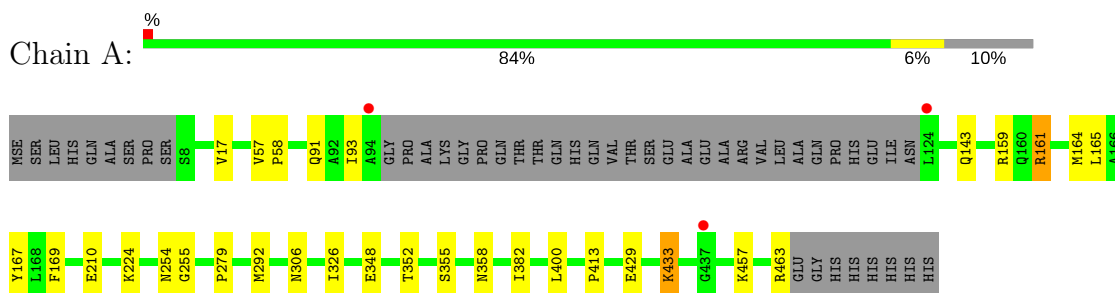
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	64	Total 64	O 64	0	0
3	E	61	Total 61	O 61	0	0
3	F	32	Total 32	O 32	0	0
3	G	53	Total 53	O 53	0	0
3	H	74	Total 74	O 74	0	0



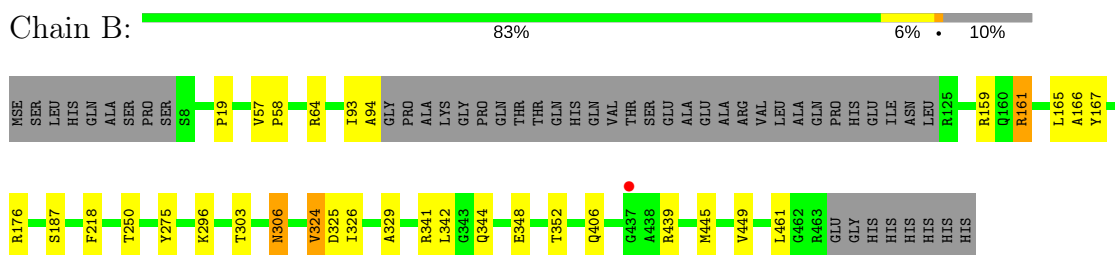
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

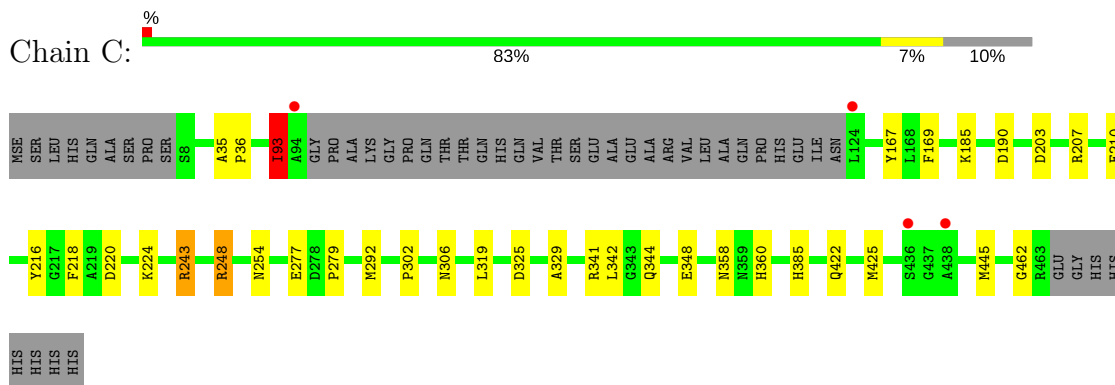
- Molecule 1: D-glucarate dehydratase



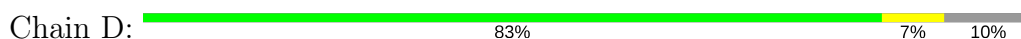
- Molecule 1: D-glucarate dehydratase

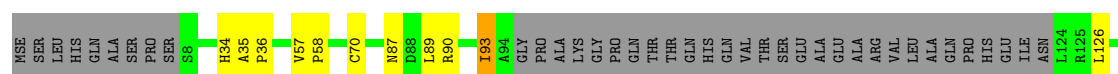


- Molecule 1: D-glucarate dehydratase

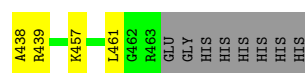
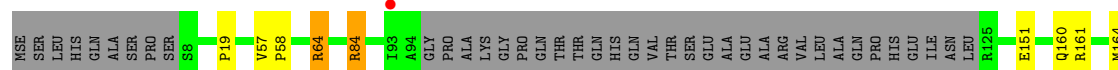
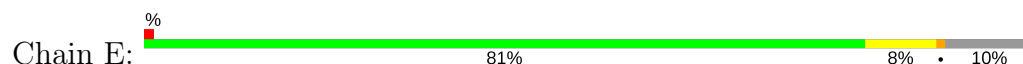


- Molecule 1: D-glucarate dehydratase

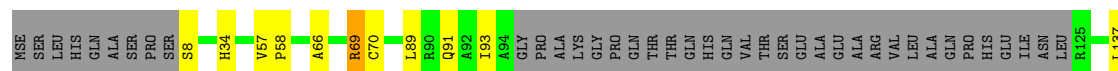
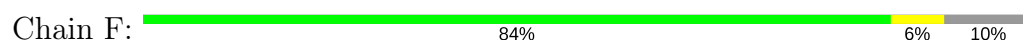




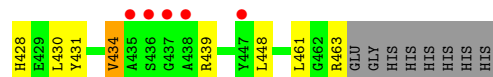
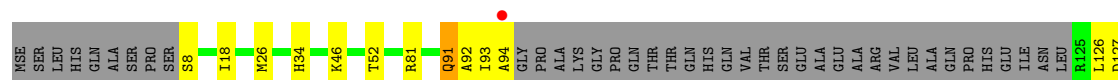
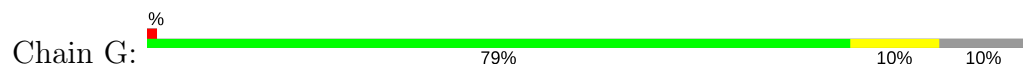
• Molecule 1: D-glucarate dehydratase



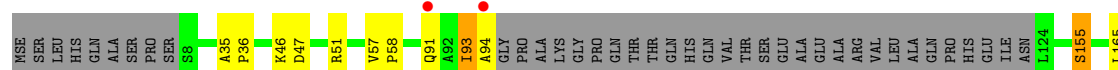
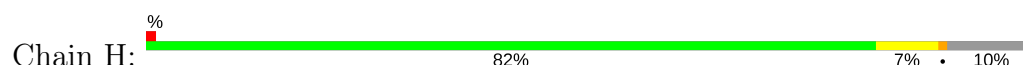
• Molecule 1: D-glucarate dehydratase

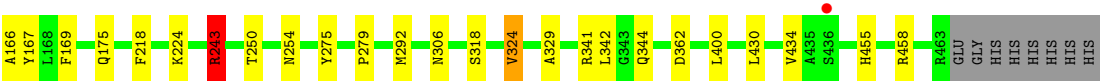


• Molecule 1: D-glucarate dehydratase



• Molecule 1: D-glucarate dehydratase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	126.36Å 148.84Å 198.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.18 – 2.20 46.18 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.18-2.20) 99.7 (46.18-2.20)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.68 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.176 , 0.222 0.185 , 0.227	Depositor DCC
$R_{free}$ test set	9478 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.4	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 30.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	27164	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL

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	0.59	0/3394	0.69	0/4574
1	B	0.59	0/3361	0.68	0/4530
1	C	0.57	0/3423	0.69	0/4613
1	D	0.57	0/3409	0.70	0/4597
1	E	0.54	0/3379	0.69	0/4555
1	F	0.53	0/3386	0.66	0/4563
1	G	0.53	0/3377	0.69	1/4551 (0.0%)
1	H	0.56	0/3377	0.69	1/4553 (0.0%)
All	All	0.56	0/27106	0.69	2/36536 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	243	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	G	139	ASP	CB-CG-OD1	5.49	123.24	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	3328	0	3295	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3301	0	3262	32	0
1	C	3348	0	3320	15	0
1	D	3333	0	3299	21	0
1	E	3312	0	3279	32	0
1	F	3319	0	3293	16	0
1	G	3314	0	3276	28	0
1	H	3313	0	3279	21	0
2	A	12	0	16	3	0
2	B	6	0	8	3	0
2	C	6	0	8	1	0
2	D	6	0	8	3	0
2	E	6	0	8	1	0
2	F	6	0	8	2	0
2	G	6	0	8	4	0
2	H	6	0	8	0	0
3	A	103	0	0	2	0
3	B	84	0	0	1	0
3	C	71	0	0	0	0
3	D	64	0	0	0	0
3	E	61	0	0	3	0
3	F	32	0	0	0	0
3	G	53	0	0	1	0
3	H	74	0	0	1	0
All	All	27164	0	26375	181	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 3.

All (181) 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:E:213[A]:ARG:HH11	1:E:213[A]:ARG:HG2	1.08	1.15
1:B:161:ARG:HG3	1:B:161:ARG:HH11	1.27	1.00
1:D:445:MSE:HE1	1:D:461:LEU:CD1	1.94	0.97
1:E:84:ARG:HG3	1:E:84:ARG:HH11	1.26	0.96
1:B:445:MSE:HE1	1:B:461:LEU:HD11	1.46	0.96
1:F:445:MSE:HE1	1:F:461:LEU:HD21	1.52	0.92
1:C:93:ILE:HG22	1:C:93:ILE:O	1.73	0.86
1:D:445:MSE:HE1	1:D:461:LEU:HD11	1.58	0.86
1:E:213[A]:ARG:NH1	1:E:213[A]:ARG:HG2	1.83	0.85
1:F:445:MSE:HE1	1:F:461:LEU:CD2	2.07	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:243:ARG:HG2	1:H:243:ARG:HH11	1.46	0.80
1:B:445:MSE:HE1	1:B:461:LEU:CD1	2.12	0.80
1:B:161:ARG:CG	1:B:161:ARG:HH11	2.01	0.74
1:D:358:ASN:HD22	2:D:501:GOL:C1	2.03	0.71
1:C:93:ILE:CG2	1:C:93:ILE:O	2.39	0.70
1:B:445:MSE:HE3	1:B:449:VAL:HG23	1.75	0.69
1:B:176:ARG:HH11	1:B:176:ARG:HG2	1.59	0.68
1:G:34:HIS:HE1	2:G:501:GOL:H31	1.60	0.67
1:G:424:ILE:HG22	1:G:425:MSE:HE2	1.76	0.66
1:H:155:SER:OG	3:H:661:HOH:O	2.13	0.65
1:G:309:ALA:HB1	1:G:315:MSE:CE	2.26	0.65
1:G:309:ALA:HB1	1:G:315:MSE:HE2	1.77	0.65
1:E:439:ARG:HH22	2:E:501:GOL:H12	1.61	0.65
1:G:224:LYS:HD2	3:G:635:HOH:O	1.98	0.64
1:E:57:VAL:HB	1:E:58:PRO:HD2	1.80	0.64
1:D:358:ASN:HD22	2:D:501:GOL:H12	1.61	0.63
1:E:84:ARG:HG3	1:E:84:ARG:NH1	2.04	0.63
1:G:34:HIS:CE1	2:G:501:GOL:H31	2.34	0.62
1:G:424:ILE:HG22	1:G:425:MSE:CE	2.30	0.62
1:H:455:HIS:CG	1:H:458:ARG:HD2	2.35	0.62
1:B:19:PRO:HG3	1:B:64:ARG:HD3	1.80	0.62
1:H:243:ARG:CG	1:H:243:ARG:HH11	2.12	0.61
1:F:34:HIS:CE1	2:F:501:GOL:H11	2.36	0.60
1:E:255:GLY:O	1:E:457:LYS:HE3	2.01	0.60
1:E:84:ARG:HH11	1:E:84:ARG:CG	2.06	0.60
1:C:210:GLU:OE1	1:C:243:ARG:NH1	2.35	0.60
1:G:296:LYS:HE3	1:G:325:ASP:OD2	2.03	0.59
1:G:91:GLN:HG3	1:G:92:ALA:N	2.18	0.58
1:F:93:ILE:N	1:F:93:ILE:HD12	2.18	0.58
1:E:250:THR:HG21	1:E:275:TYR:CZ	2.39	0.57
1:A:429:GLU:O	1:A:433:LYS:HD2	2.04	0.57
1:D:93:ILE:HG21	1:D:126:LEU:HD13	1.85	0.57
1:D:445:MSE:HE1	1:D:461:LEU:HD13	1.82	0.57
1:B:445:MSE:HE3	1:B:449:VAL:CG2	2.33	0.57
1:D:164:MSE:HE2	1:D:382:ILE:HD11	1.87	0.57
1:G:93:ILE:CG2	1:G:93:ILE:O	2.53	0.57
1:A:164:MSE:HE2	1:A:382:ILE:HD11	1.85	0.56
1:H:169:PHE:CE2	1:H:224:LYS:HG2	2.41	0.56
1:A:358:ASN:OD1	2:A:502:GOL:C1	2.54	0.55
1:E:57:VAL:HB	1:E:58:PRO:CD	2.35	0.54
1:G:385:HIS:NE2	2:G:501:GOL:H32	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:318:SER:OG	1:H:324:VAL:HG13	2.07	0.54
1:E:277:GLU:OE2	1:E:306:ASN:ND2	2.42	0.53
1:E:250:THR:HG21	1:E:275:TYR:CE2	2.43	0.53
1:A:169:PHE:CE2	1:A:224:LYS:HG2	2.43	0.53
1:B:445:MSE:CE	1:B:461:LEU:CD1	2.85	0.53
1:F:358:ASN:O	1:F:360:HIS:HD2	1.92	0.53
1:F:89:LEU:O	1:F:93:ILE:HD13	2.08	0.53
1:A:17:VAL:O	2:A:501:GOL:H31	2.08	0.52
1:B:166:ALA:HB2	1:B:218:PHE:CD2	2.45	0.52
1:F:445:MSE:HE1	1:F:461:LEU:HD23	1.91	0.52
1:C:341:ARG:HA	1:C:344:GLN:HE21	1.73	0.52
1:D:35:ALA:HB1	1:D:36:PRO:HD2	1.91	0.52
1:B:161:ARG:NH1	1:B:161:ARG:HG3	2.08	0.52
1:F:57:VAL:HB	1:F:58:PRO:HD2	1.92	0.52
1:B:176:ARG:NH1	1:B:176:ARG:HG2	2.25	0.51
1:B:296:LYS:HE3	1:B:325:ASP:OD1	2.11	0.51
1:E:213[A]:ARG:CG	1:E:213[A]:ARG:HH11	1.98	0.51
1:B:348:GLU:HG2	1:D:312:TRP:HB2	1.93	0.51
1:D:93:ILE:HG21	1:D:126:LEU:CD1	2.41	0.51
1:E:341:ARG:HA	1:E:344:GLN:HE21	1.77	0.50
1:A:355:SER:HB3	1:A:382:ILE:HB	1.93	0.50
1:F:70:CYS:SG	1:F:93:ILE:HD11	2.52	0.50
1:G:18:ILE:HG21	1:G:425:MSE:HE1	1.94	0.49
1:H:243:ARG:HG2	1:H:243:ARG:NH1	2.17	0.49
1:G:93:ILE:HG22	1:G:126:LEU:HD13	1.93	0.49
1:E:161:ARG:NH1	3:E:616:HOH:O	2.37	0.49
1:F:34:HIS:HE1	2:F:501:GOL:H11	1.76	0.49
1:G:93:ILE:O	1:G:94:ALA:HB2	2.12	0.49
1:A:93:ILE:N	1:A:93:ILE:HD12	2.27	0.49
1:E:296:LYS:HE2	1:E:325:ASP:OD1	2.13	0.49
1:E:279:PRO:HD2	1:E:292:MSE:SE	2.63	0.49
1:E:213[A]:ARG:CG	1:E:213[A]:ARG:NH1	2.64	0.49
1:H:279:PRO:HD2	1:H:292:MSE:SE	2.63	0.48
1:D:296:LYS:HE3	1:D:325:ASP:OD1	2.12	0.48
1:D:70:CYS:SG	1:D:89:LEU:HD22	2.53	0.48
1:D:169:PHE:CE2	1:D:224:LYS:HG2	2.49	0.48
1:D:279:PRO:HD2	1:D:292:MSE:SE	2.64	0.48
1:B:306[A]:ASN:ND2	3:B:612:HOH:O	2.47	0.48
1:D:93:ILE:CG2	1:D:126:LEU:HD13	2.43	0.48
1:B:57:VAL:HB	1:B:58:PRO:HD2	1.95	0.47
1:F:66:ALA:O	1:F:69[B]:ARG:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:277:GLU:OE2	1:C:306[A]:ASN:ND2	2.48	0.47
1:G:439:ARG:HH22	2:G:501:GOL:H2	1.78	0.47
1:B:439:ARG:HH22	2:B:501:GOL:H32	1.80	0.47
1:B:161:ARG:NH1	1:B:161:ARG:CG	2.68	0.47
1:C:302:PRO:HA	1:C:325:ASP:OD2	2.15	0.47
1:A:93:ILE:N	1:A:93:ILE:CD1	2.78	0.46
1:B:445:MSE:CE	1:B:449:VAL:HG23	2.43	0.46
1:C:35:ALA:HB1	1:C:36:PRO:HD2	1.97	0.46
1:E:19:PRO:HG3	1:E:64:ARG:HG3	1.97	0.46
1:E:169:PHE:CE2	1:E:224:LYS:HG2	2.50	0.46
1:D:34:HIS:CE1	2:D:501:GOL:H2	2.50	0.46
1:H:166:ALA:HB2	1:H:218:PHE:CD2	2.50	0.46
1:A:224:LYS:HD2	3:A:643:HOH:O	2.15	0.46
1:B:445:MSE:CE	1:B:449:VAL:CG2	2.93	0.46
1:H:93:ILE:O	1:H:94:ALA:C	2.54	0.46
1:B:93:ILE:O	1:B:93:ILE:HG22	2.15	0.46
1:C:279:PRO:HD2	1:C:292:MSE:SE	2.66	0.46
1:H:250:THR:HG21	1:H:275:TYR:CZ	2.50	0.46
1:H:329:ALA:HB2	1:H:342:LEU:HD23	1.97	0.46
1:H:341:ARG:HA	1:H:344:GLN:HE21	1.81	0.46
1:A:358:ASN:OD1	2:A:502:GOL:H12	2.16	0.46
1:G:93:ILE:O	1:G:93:ILE:HG23	2.16	0.45
1:B:93:ILE:O	1:B:94:ALA:HB3	2.17	0.45
1:B:19:PRO:HG3	1:B:64:ARG:CD	2.46	0.45
1:E:84:ARG:CG	1:E:84:ARG:NH1	2.69	0.45
1:B:341:ARG:HA	1:B:344:GLN:HE21	1.81	0.45
1:A:159[A]:ARG:HD3	1:A:161:ARG:O	2.16	0.45
1:E:224:LYS:HD2	3:E:647:HOH:O	2.16	0.45
1:E:344:GLN:NE2	3:E:661:HOH:O	2.45	0.45
1:G:279:PRO:HD2	1:G:292:MSE:SE	2.67	0.45
1:B:165:LEU:HD12	1:B:165:LEU:C	2.37	0.45
1:E:430:LEU:O	1:E:434:VAL:HG22	2.17	0.44
1:A:326:ILE:HA	1:A:352:THR:O	2.18	0.44
1:E:173:GLU:HB3	1:E:176:ARG:HG3	1.98	0.44
1:H:35:ALA:HB1	1:H:36:PRO:HD2	1.99	0.44
1:C:220:ASP:OD1	1:C:248:ARG:HD2	2.18	0.44
1:G:326:ILE:HA	1:G:352:THR:O	2.18	0.43
1:C:358:ASN:O	1:C:360:HIS:HD2	2.01	0.43
1:C:93:ILE:CD1	1:C:93:ILE:N	2.81	0.43
1:F:93:ILE:CD1	1:F:93:ILE:N	2.81	0.43
1:E:358:ASN:O	1:E:360:HIS:HD2	2.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:87:ASN:OD1	1:D:90:ARG:NH1	2.52	0.43
1:F:253:PRO:HD2	1:F:277:GLU:O	2.18	0.43
1:D:362:ASP:N	1:D:362:ASP:OD1	2.48	0.43
1:G:159[A]:ARG:HD3	1:G:161:ARG:O	2.18	0.43
1:H:155:SER:OG	1:H:155:SER:O	2.36	0.43
1:A:306[A]:ASN:ND2	3:A:635:HOH:O	2.51	0.43
1:C:216:TYR:HB2	1:C:218:PHE:CE2	2.54	0.43
1:D:250:THR:HG21	1:D:275:TYR:CE2	2.53	0.43
1:A:143:GLN:HA	1:A:413:PRO:HB2	2.00	0.42
1:G:26:MSE:HE2	1:G:448:LEU:HD11	2.01	0.42
1:E:164:MSE:HE2	1:E:382:ILE:HD11	2.00	0.42
1:B:329:ALA:HB2	1:B:342:LEU:HD23	2.01	0.42
1:H:47:ASP:OD1	1:H:47:ASP:C	2.57	0.42
1:D:255:GLY:O	1:D:457:LYS:HE3	2.19	0.42
1:F:169:PHE:CE2	1:F:224:LYS:HG2	2.53	0.42
1:H:46:LYS:HA	1:H:51:ARG:O	2.19	0.42
1:H:250:THR:HG21	1:H:275:TYR:CE2	2.54	0.42
1:H:165:LEU:C	1:H:165:LEU:HD12	2.40	0.42
1:H:57:VAL:HB	1:H:58:PRO:HD2	2.01	0.42
1:G:430:LEU:O	1:G:434:VAL:HB	2.20	0.42
1:A:279:PRO:HD2	1:A:292:MSE:SE	2.69	0.42
1:B:326:ILE:HA	1:B:352:THR:O	2.20	0.42
1:C:329:ALA:HB2	1:C:342:LEU:HD23	2.02	0.42
1:B:439:ARG:HH12	2:B:501:GOL:C3	2.33	0.41
1:E:315:MSE:HE2	1:E:319:LEU:HG	2.03	0.41
1:A:255:GLY:O	1:A:457:LYS:HE3	2.21	0.41
1:G:277:GLU:OE2	1:G:306[A]:ASN:ND2	2.52	0.41
1:F:137:LEU:HD23	1:F:137:LEU:HA	1.92	0.41
1:E:151:GLU:HG3	1:G:81:ARG:NH1	2.35	0.41
1:A:165:LEU:HD12	1:A:165:LEU:C	2.41	0.41
1:A:57:VAL:HB	1:A:58:PRO:HD2	2.01	0.41
1:E:169:PHE:CD2	1:E:224:LYS:HG2	2.55	0.41
1:E:411:ASP:C	1:E:411:ASP:OD1	2.57	0.41
1:B:439:ARG:NH2	2:B:501:GOL:H32	2.36	0.41
1:B:159:ARG:HD3	1:B:159:ARG:HH11	1.73	0.41
1:F:164:MSE:HE3	1:F:407:VAL:CG2	2.50	0.41
1:G:309:ALA:HB1	1:G:315:MSE:HE1	2.02	0.41
1:H:430:LEU:O	1:H:434:VAL:HG22	2.21	0.41
1:G:341:ARG:HA	1:G:344:GLN:HE21	1.86	0.41
1:G:428:HIS:O	1:G:431:TYR:HB3	2.21	0.41
1:B:250:THR:HG21	1:B:275:TYR:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:169:PHE:CE2	1:C:224:LYS:HG2	2.55	0.41
1:B:303:THR:OG1	1:B:324:VAL:HA	2.21	0.40
1:D:57:VAL:HB	1:D:58:PRO:HD2	2.04	0.40
1:G:304:ALA:HA	1:G:326:ILE:O	2.21	0.40
1:C:385:HIS:NE2	2:C:501:GOL:O2	2.43	0.40
1:G:46:LYS:HD2	1:G:52:THR:OG1	2.22	0.40
1:E:355:SER:HB3	1:E:382:ILE:HB	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	427/473 (90%)	417 (98%)	10 (2%)	0	100	100
1	B	423/473 (89%)	414 (98%)	9 (2%)	0	100	100
1	C	430/473 (91%)	420 (98%)	8 (2%)	2 (0%)	32	34
1	D	429/473 (91%)	419 (98%)	10 (2%)	0	100	100
1	E	425/473 (90%)	414 (97%)	10 (2%)	1 (0%)	51	58
1	F	425/473 (90%)	408 (96%)	14 (3%)	3 (1%)	25	24
1	G	425/473 (90%)	415 (98%)	10 (2%)	0	100	100
1	H	425/473 (90%)	419 (99%)	6 (1%)	0	100	100
All	All	3409/3784 (90%)	3326 (98%)	77 (2%)	6 (0%)	55	58

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	462	GLY
1	E	438	ALA
1	F	187	SER

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Mol	Chain	Res	Type
1	C	93	ILE
1	F	69[A]	ARG
1	F	69[B]	ARG

### 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	335/356 (94%)	326 (97%)	9 (3%)	50	62
1	B	332/356 (93%)	325 (98%)	7 (2%)	59	72
1	C	338/356 (95%)	323 (96%)	15 (4%)	33	40
1	D	336/356 (94%)	327 (97%)	9 (3%)	50	62
1	E	333/356 (94%)	321 (96%)	12 (4%)	40	50
1	F	334/356 (94%)	323 (97%)	11 (3%)	43	54
1	G	333/356 (94%)	317 (95%)	16 (5%)	30	36
1	H	333/356 (94%)	322 (97%)	11 (3%)	43	54
All	All	2674/2848 (94%)	2584 (97%)	90 (3%)	43	53

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	161	ARG
1	A	167	TYR
1	A	210	GLU
1	A	254	ASN
1	A	348	GLU
1	A	400	LEU
1	A	433	LYS
1	A	463	ARG
1	B	161	ARG
1	B	167	TYR
1	B	187	SER

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Mol	Chain	Res	Type
1	B	306[A]	ASN
1	B	306[B]	ASN
1	B	324	VAL
1	B	406	GLN
1	C	93	ILE
1	C	167	TYR
1	C	185	LYS
1	C	190	ASP
1	C	203	ASP
1	C	207	ARG
1	C	243	ARG
1	C	248	ARG
1	C	254	ASN
1	C	319	LEU
1	C	348[A]	GLU
1	C	348[B]	GLU
1	C	422	GLN
1	C	425	MSE
1	C	445	MSE
1	D	93	ILE
1	D	167	TYR
1	D	243	ARG
1	D	306	ASN
1	D	348[A]	GLU
1	D	348[B]	GLU
1	D	400	LEU
1	D	422	GLN
1	D	461	LEU
1	E	64	ARG
1	E	84	ARG
1	E	160	GLN
1	E	167	TYR
1	E	176	ARG
1	E	213[A]	ARG
1	E	213[B]	ARG
1	E	241	LYS
1	E	324	VAL
1	E	327	PRO
1	E	436	SER
1	E	461	LEU
1	F	8	SER
1	F	91	GLN

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Mol	Chain	Res	Type
1	F	161	ARG
1	F	167	TYR
1	F	215	ARG
1	F	241	LYS
1	F	254	ASN
1	F	267	LYS
1	F	306	ASN
1	F	319	LEU
1	F	425	MSE
1	G	8	SER
1	G	91	GLN
1	G	127	ASP
1	G	167	TYR
1	G	176	ARG
1	G	187	SER
1	G	189	ASP
1	G	190	ASP
1	G	243	ARG
1	G	248	ARG
1	G	254	ASN
1	G	348	GLU
1	G	400	LEU
1	G	434	VAL
1	G	461	LEU
1	G	463	ARG
1	H	91	GLN
1	H	93	ILE
1	H	155	SER
1	H	167	TYR
1	H	175	GLN
1	H	243	ARG
1	H	254	ASN
1	H	306	ASN
1	H	324	VAL
1	H	362	ASP
1	H	400	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	B	77	GLN

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Mol	Chain	Res	Type
1	B	344	GLN
1	B	408	GLN
1	C	269	GLN
1	C	284	ASN
1	C	344	GLN
1	D	128	ASN
1	D	254	ASN
1	D	306	ASN
1	D	358	ASN
1	D	453	GLN
1	E	128	ASN
1	E	337	GLN
1	E	344	GLN
1	F	306	ASN
1	F	408	GLN
1	G	344	GLN
1	H	65	GLN
1	H	344	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

9 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	A	501	-	5,5,5	0.26	0	5,5,5	0.26	0
2	GOL	A	502	-	5,5,5	0.70	0	5,5,5	1.19	0
2	GOL	B	501	-	5,5,5	0.79	0	5,5,5	0.79	0
2	GOL	C	501	-	5,5,5	0.66	0	5,5,5	0.64	0
2	GOL	D	501	-	5,5,5	0.65	0	5,5,5	1.08	0
2	GOL	E	501	-	5,5,5	0.61	0	5,5,5	1.30	1 (20%)
2	GOL	F	501	-	5,5,5	0.62	0	5,5,5	0.77	0
2	GOL	G	501	-	5,5,5	0.67	0	5,5,5	0.72	0
2	GOL	H	501	-	5,5,5	0.95	0	5,5,5	1.22	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	A	501	-	-	0/4/4/4	0/0/0/0
2	GOL	A	502	-	-	0/4/4/4	0/0/0/0
2	GOL	B	501	-	-	0/4/4/4	0/0/0/0
2	GOL	C	501	-	-	0/4/4/4	0/0/0/0
2	GOL	D	501	-	-	0/4/4/4	0/0/0/0
2	GOL	E	501	-	-	0/4/4/4	0/0/0/0
2	GOL	F	501	-	-	0/4/4/4	0/0/0/0
2	GOL	G	501	-	-	0/4/4/4	0/0/0/0
2	GOL	H	501	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	501	GOL	O2-C2-C1	-2.13	98.80	108.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 17 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	GOL	1	0
2	A	502	GOL	2	0
2	B	501	GOL	3	0
2	C	501	GOL	1	0
2	D	501	GOL	3	0
2	E	501	GOL	1	0
2	F	501	GOL	2	0
2	G	501	GOL	4	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	413/473 (87%)	-0.48	3 (0%) 87 86	11, 18, 39, 70	0
1	B	412/473 (87%)	-0.58	1 (0%) 94 94	10, 20, 41, 66	0
1	C	413/473 (87%)	-0.49	4 (0%) 82 81	11, 21, 45, 88	0
1	D	413/473 (87%)	-0.41	1 (0%) 94 94	10, 23, 47, 82	0
1	E	412/473 (87%)	-0.44	3 (0%) 87 86	14, 25, 46, 74	0
1	F	412/473 (87%)	-0.40	1 (0%) 94 94	14, 28, 50, 77	0
1	G	412/473 (87%)	-0.44	7 (1%) 70 68	14, 24, 47, 73	0
1	H	413/473 (87%)	-0.47	3 (0%) 87 86	11, 22, 48, 91	0
All	All	3300/3784 (87%)	-0.47	23 (0%) 87 86	10, 23, 46, 91	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	94	ALA	4.5
1	C	94	ALA	4.2
1	H	436	SER	4.1
1	A	124	LEU	3.8
1	D	436	SER	3.8
1	C	124	LEU	3.7
1	E	437	GLY	3.4
1	G	437	GLY	3.3
1	G	178	ASP	3.3
1	C	436	SER	3.2
1	E	93	ILE	3.1
1	A	94	ALA	2.7
1	G	94	ALA	2.7
1	G	438	ALA	2.7
1	E	436	SER	2.4
1	A	437	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	G	447	TYR	2.4
1	H	91	GLN	2.2
1	G	435	ALA	2.2
1	B	437	GLY	2.1
1	F	186	GLY	2.1
1	G	436	SER	2.1
1	C	438	ALA	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	GOL	F	501	6/6	0.80	0.27	12.48	29,37,45,46	0
2	GOL	B	501	6/6	0.94	0.14	7.32	21,27,28,36	0
2	GOL	C	501	6/6	0.90	0.19	4.77	24,32,34,43	0
2	GOL	D	501	6/6	0.89	0.23	4.76	29,32,41,42	0
2	GOL	H	501	6/6	0.91	0.20	3.59	21,26,32,37	0
2	GOL	E	501	6/6	0.94	0.17	3.42	30,32,38,44	0
2	GOL	G	501	6/6	0.86	0.19	2.36	23,32,37,52	0
2	GOL	A	502	6/6	0.93	0.16	2.29	20,26,28,37	0
2	GOL	A	501	6/6	0.92	0.12	0.87	35,36,44,45	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.