



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 09:23 am GMT

PDB ID : 4LXV  
Title : Crystal Structure of the Hemagglutinin from a H1N1pdm  
A/WASHINGTON/5/2011 virus  
Authors : Yang, H.; Chang, J.C.; Guo, Z.; Carney, P.J.; Shore, D.A.; Donis, R.O.; Cox,  
N.J.; Villanueva, J.M.; Klimov, A.I.; Stevens, J.  
Deposited on : 2013-07-30  
Resolution : 3.00 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

<https://www.wwpdb.org/validation/2017/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 : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : 1.13  
EDS : trunk30967  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30967

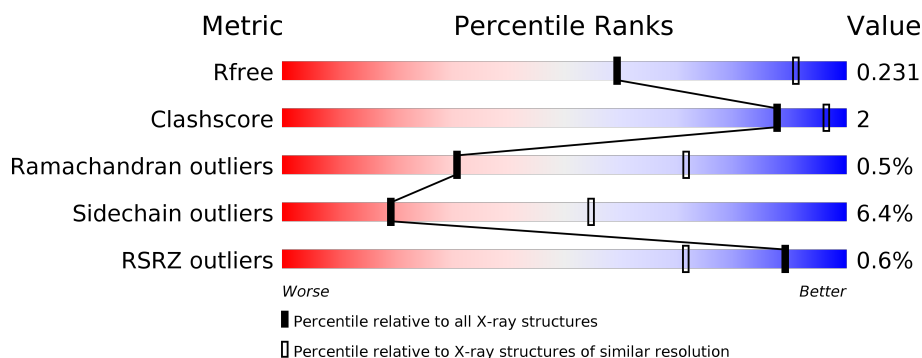
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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))
$R_{free}$	111664	1851 (3.00-3.00)
Clashscore	122126	2167 (3.00-3.00)
Ramachandran outliers	120053	2101 (3.00-3.00)
Sidechain outliers	120020	2104 (3.00-3.00)
RSRZ outliers	108989	1751 (3.00-3.00)

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	332	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div>• •</div> </div> </div>
1	C	332	<div> <div></div> <div>86%</div> <div>10%</div> <div>• •</div> </div>
1	E	332	<div> <div></div> <div>86%</div> <div>9%</div> <div>• •</div> </div>
1	G	332	<div> <div></div> <div>86%</div> <div>11%</div> <div>• •</div> </div>
1	I	332	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div>• •</div> </div> </div>
1	K	332	<div> <div></div> <div>84%</div> <div>11%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
2	B	182	 % 85% 9% 6%
2	D	182	 2% 84% 10% • 6%
2	F	182	 2% 85% 8% • 6%
2	H	182	 85% 9% 6%
2	J	182	 84% 10% • 6%
2	L	182	 % 84% 9% • 6%

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
3	NAG	A	403	X	-	-	-
3	NAG	E	404	X	-	-	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 23740 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2523	1595	435	482	11			
1	C	323	Total	C	N	O	S	0	0	0
			2523	1595	435	482	11			
1	E	323	Total	C	N	O	S	0	0	0
			2523	1595	435	482	11			
1	G	323	Total	C	N	O	S	0	0	0
			2523	1595	435	482	11			
1	I	323	Total	C	N	O	S	0	0	0
			2523	1595	435	482	11			
1	K	323	Total	C	N	O	S	0	0	0
			2523	1595	435	482	11			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	ALA	-	EXPRESSION TAG	UNP J7MFR5
A	-3	ASP	-	EXPRESSION TAG	UNP J7MFR5
A	-2	LEU	-	EXPRESSION TAG	UNP J7MFR5
A	-1	GLY	-	EXPRESSION TAG	UNP J7MFR5
A	0	SER	-	EXPRESSION TAG	UNP J7MFR5
C	-4	ALA	-	EXPRESSION TAG	UNP J7MFR5
C	-3	ASP	-	EXPRESSION TAG	UNP J7MFR5
C	-2	LEU	-	EXPRESSION TAG	UNP J7MFR5
C	-1	GLY	-	EXPRESSION TAG	UNP J7MFR5
C	0	SER	-	EXPRESSION TAG	UNP J7MFR5
E	-4	ALA	-	EXPRESSION TAG	UNP J7MFR5
E	-3	ASP	-	EXPRESSION TAG	UNP J7MFR5
E	-2	LEU	-	EXPRESSION TAG	UNP J7MFR5
E	-1	GLY	-	EXPRESSION TAG	UNP J7MFR5
E	0	SER	-	EXPRESSION TAG	UNP J7MFR5
G	-4	ALA	-	EXPRESSION TAG	UNP J7MFR5
G	-3	ASP	-	EXPRESSION TAG	UNP J7MFR5

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-2	LEU	-	EXPRESSION TAG	UNP J7MFR5
G	-1	GLY	-	EXPRESSION TAG	UNP J7MFR5
G	0	SER	-	EXPRESSION TAG	UNP J7MFR5
I	-4	ALA	-	EXPRESSION TAG	UNP J7MFR5
I	-3	ASP	-	EXPRESSION TAG	UNP J7MFR5
I	-2	LEU	-	EXPRESSION TAG	UNP J7MFR5
I	-1	GLY	-	EXPRESSION TAG	UNP J7MFR5
I	0	SER	-	EXPRESSION TAG	UNP J7MFR5
K	-4	ALA	-	EXPRESSION TAG	UNP J7MFR5
K	-3	ASP	-	EXPRESSION TAG	UNP J7MFR5
K	-2	LEU	-	EXPRESSION TAG	UNP J7MFR5
K	-1	GLY	-	EXPRESSION TAG	UNP J7MFR5
K	0	SER	-	EXPRESSION TAG	UNP J7MFR5

- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	171	Total	C	N	O	S	0	0	0
			1387	871	236	274	6			
2	D	171	Total	C	N	O	S	0	0	0
			1387	871	236	274	6			
2	F	171	Total	C	N	O	S	0	0	0
			1387	871	236	274	6			
2	H	171	Total	C	N	O	S	0	0	0
			1387	871	236	274	6			
2	J	171	Total	C	N	O	S	0	0	0
			1387	871	236	274	6			
2	L	171	Total	C	N	O	S	0	0	0
			1387	871	236	274	6			

There are 48 discrepancies between the modelled and reference sequences:

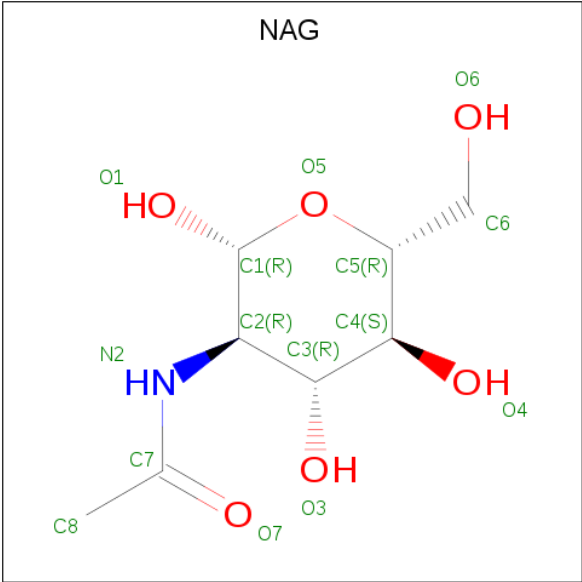
Chain	Residue	Modelled	Actual	Comment	Reference
B	175	SER	-	EXPRESSION TAG	UNP J7MFR5
B	176	GLY	-	EXPRESSION TAG	UNP J7MFR5
B	177	ARG	-	EXPRESSION TAG	UNP J7MFR5
B	178	LEU	-	EXPRESSION TAG	UNP J7MFR5
B	179	VAL	-	EXPRESSION TAG	UNP J7MFR5
B	180	PRO	-	EXPRESSION TAG	UNP J7MFR5
B	181	ARG	-	EXPRESSION TAG	UNP J7MFR5
B	182	GLY	-	EXPRESSION TAG	UNP J7MFR5
D	175	SER	-	EXPRESSION TAG	UNP J7MFR5

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Chain	Residue	Modelled	Actual	Comment	Reference
D	176	GLY	-	EXPRESSION TAG	UNP J7MFR5
D	177	ARG	-	EXPRESSION TAG	UNP J7MFR5
D	178	LEU	-	EXPRESSION TAG	UNP J7MFR5
D	179	VAL	-	EXPRESSION TAG	UNP J7MFR5
D	180	PRO	-	EXPRESSION TAG	UNP J7MFR5
D	181	ARG	-	EXPRESSION TAG	UNP J7MFR5
D	182	GLY	-	EXPRESSION TAG	UNP J7MFR5
F	175	SER	-	EXPRESSION TAG	UNP J7MFR5
F	176	GLY	-	EXPRESSION TAG	UNP J7MFR5
F	177	ARG	-	EXPRESSION TAG	UNP J7MFR5
F	178	LEU	-	EXPRESSION TAG	UNP J7MFR5
F	179	VAL	-	EXPRESSION TAG	UNP J7MFR5
F	180	PRO	-	EXPRESSION TAG	UNP J7MFR5
F	181	ARG	-	EXPRESSION TAG	UNP J7MFR5
F	182	GLY	-	EXPRESSION TAG	UNP J7MFR5
H	175	SER	-	EXPRESSION TAG	UNP J7MFR5
H	176	GLY	-	EXPRESSION TAG	UNP J7MFR5
H	177	ARG	-	EXPRESSION TAG	UNP J7MFR5
H	178	LEU	-	EXPRESSION TAG	UNP J7MFR5
H	179	VAL	-	EXPRESSION TAG	UNP J7MFR5
H	180	PRO	-	EXPRESSION TAG	UNP J7MFR5
H	181	ARG	-	EXPRESSION TAG	UNP J7MFR5
H	182	GLY	-	EXPRESSION TAG	UNP J7MFR5
J	175	SER	-	EXPRESSION TAG	UNP J7MFR5
J	176	GLY	-	EXPRESSION TAG	UNP J7MFR5
J	177	ARG	-	EXPRESSION TAG	UNP J7MFR5
J	178	LEU	-	EXPRESSION TAG	UNP J7MFR5
J	179	VAL	-	EXPRESSION TAG	UNP J7MFR5
J	180	PRO	-	EXPRESSION TAG	UNP J7MFR5
J	181	ARG	-	EXPRESSION TAG	UNP J7MFR5
J	182	GLY	-	EXPRESSION TAG	UNP J7MFR5
L	175	SER	-	EXPRESSION TAG	UNP J7MFR5
L	176	GLY	-	EXPRESSION TAG	UNP J7MFR5
L	177	ARG	-	EXPRESSION TAG	UNP J7MFR5
L	178	LEU	-	EXPRESSION TAG	UNP J7MFR5
L	179	VAL	-	EXPRESSION TAG	UNP J7MFR5
L	180	PRO	-	EXPRESSION TAG	UNP J7MFR5
L	181	ARG	-	EXPRESSION TAG	UNP J7MFR5
L	182	GLY	-	EXPRESSION TAG	UNP J7MFR5

- Molecule 3 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		

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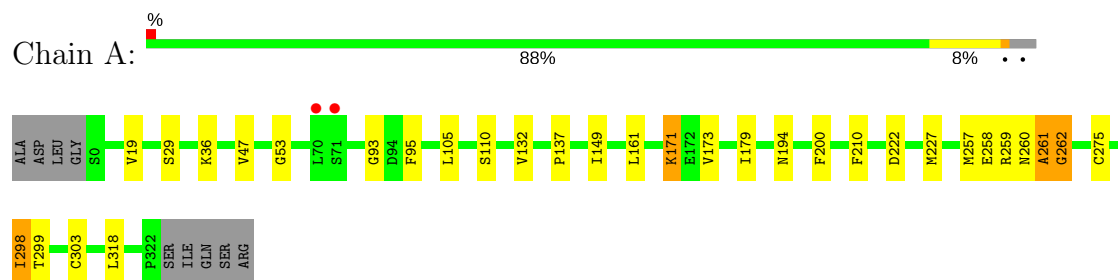
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	I	1	Total	C	N	O	0	0
			14	8	1	5		
3	I	1	Total	C	N	O	0	0
			14	8	1	5		
3	I	1	Total	C	N	O	0	0
			14	8	1	5		
3	K	1	Total	C	N	O	0	0
			14	8	1	5		
3	K	1	Total	C	N	O	0	0
			14	8	1	5		



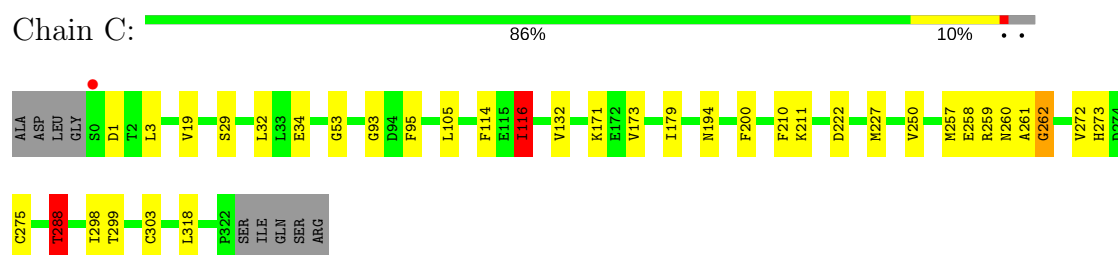
### 3 Residue-property plots [i](#)

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.

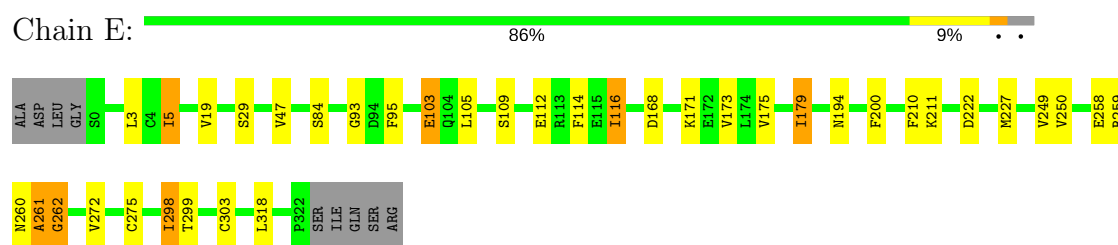
#### • Molecule 1: Hemagglutinin



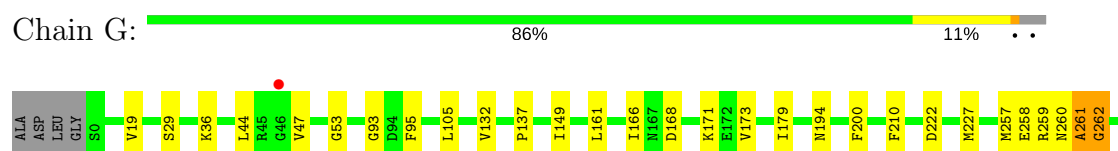
#### • Molecule 1: Hemagglutinin

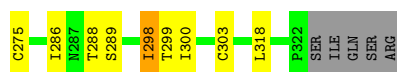


#### • Molecule 1: Hemagglutinin

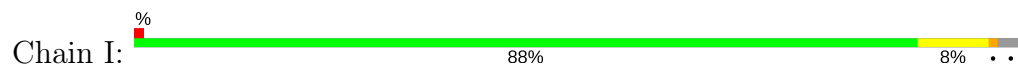


#### • Molecule 1: Hemagglutinin

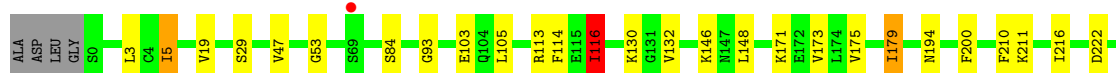
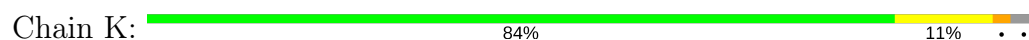




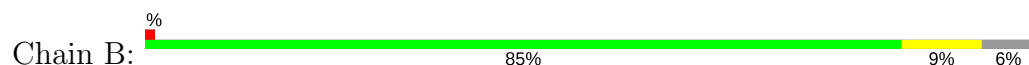
• Molecule 1: Hemagglutinin



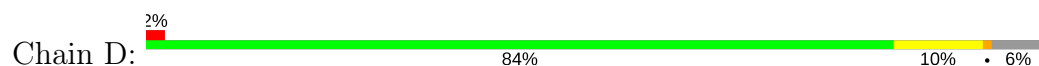
• Molecule 1: Hemagglutinin



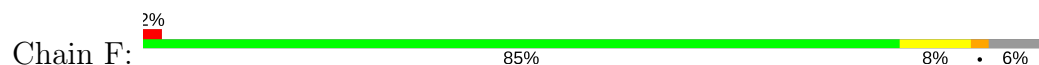
• Molecule 2: Hemagglutinin



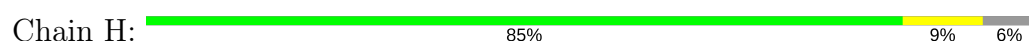
• Molecule 2: Hemagglutinin



• Molecule 2: Hemagglutinin

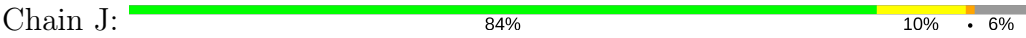


• Molecule 2: Hemagglutinin

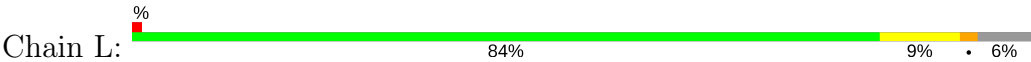




● Molecule 2: Hemagglutinin



● Molecule 2: Hemagglutinin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.24Å 226.00Å 271.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.78 – 3.00 49.73 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.2 (49.78-3.00) 98.2 (49.73-3.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.99 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.205 , 0.230 0.207 , 0.231	Depositor DCC
$R_{free}$ test set	4492 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	59.6	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 26.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	23740	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	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.*

<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

### 5.1 Standard geometry

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

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	0/2586	0.76	1/3515 (0.0%)
1	C	0.61	0/2586	0.79	3/3515 (0.1%)
1	E	0.61	0/2586	0.79	3/3515 (0.1%)
1	G	0.58	0/2586	0.76	2/3515 (0.1%)
1	I	0.60	0/2586	0.77	2/3515 (0.1%)
1	K	0.58	0/2586	0.78	3/3515 (0.1%)
2	B	0.53	0/1415	0.67	0/1906
2	D	0.54	0/1415	0.70	0/1906
2	F	0.54	0/1415	0.74	4/1906 (0.2%)
2	H	0.56	0/1415	0.68	0/1906
2	J	0.59	0/1415	0.71	0/1906
2	L	0.57	0/1415	0.75	4/1906 (0.2%)
All	All	0.58	0/24006	0.75	22/32526 (0.1%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	145	ASP	CB-CG-OD2	-7.78	111.30	118.30
2	L	145	ASP	CB-CG-OD2	-7.66	111.41	118.30
2	L	145	ASP	N-CA-CB	-7.57	96.97	110.60
2	F	145	ASP	N-CA-CB	-7.31	97.44	110.60
1	E	275	CYS	N-CA-CB	-7.21	97.61	110.60
1	G	149	ILE	CG1-CB-CG2	-7.11	95.77	111.40
1	A	149	ILE	CG1-CB-CG2	-6.94	96.13	111.40
1	C	288	THR	CB-CA-C	-6.76	93.36	111.60
1	K	275	CYS	CA-CB-SG	6.65	125.97	114.00
1	I	288	THR	CB-CA-C	-6.62	93.73	111.60
1	C	1	ASP	CB-CG-OD2	6.46	124.11	118.30
1	I	1	ASP	CB-CG-OD2	6.42	124.08	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	275	CYS	N-CA-CB	-6.38	99.11	110.60
2	L	145	ASP	CB-CG-OD1	6.12	123.81	118.30
2	F	168	LEU	CA-CB-CG	5.99	129.07	115.30
2	L	168	LEU	CA-CB-CG	5.94	128.96	115.30
2	F	145	ASP	CB-CG-OD1	5.90	123.61	118.30
1	E	275	CYS	CA-CB-SG	5.86	124.55	114.00
1	E	168	ASP	CB-CG-OD2	5.16	122.94	118.30
1	C	116	ILE	CB-CA-C	-5.14	101.32	111.60
1	G	168	ASP	CB-CG-OD2	5.10	122.89	118.30
1	K	116	ILE	CB-CA-C	-5.08	101.44	111.60

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	2523	0	2474	11	0
1	C	2523	0	2475	11	0
1	E	2523	0	2474	13	0
1	G	2523	0	2473	15	0
1	I	2523	0	2475	7	0
1	K	2523	0	2475	15	0
2	B	1387	0	1311	5	0
2	D	1387	0	1311	7	0
2	F	1387	0	1311	6	0
2	H	1387	0	1311	4	0
2	J	1387	0	1311	7	0
2	L	1387	0	1311	7	0
3	A	56	0	51	1	0
3	C	42	0	38	0	0
3	E	56	0	51	1	0
3	G	56	0	51	1	0
3	I	42	0	38	0	0
3	K	28	0	26	0	0
All	All	23740	0	22967	103	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 (103) 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:109:SER:HA	1:G:166:ILE:HD11	1.55	0.88
1:E:109:SER:HA	1:G:166:ILE:CD1	2.29	0.62
2:D:164:GLU:HA	2:D:167:LYS:HG2	1.83	0.61
1:C:200:PHE:CD2	1:C:200:PHE:C	2.76	0.57
1:A:298:ILE:HG13	1:A:298:ILE:O	2.04	0.56
1:K:114:PHE:CE1	1:K:116:ILE:HG23	2.39	0.56
1:E:298:ILE:HG13	1:E:298:ILE:O	2.03	0.56
2:J:164:GLU:HA	2:J:167:LYS:HG2	1.86	0.56
1:G:298:ILE:HG13	1:G:298:ILE:O	2.04	0.56
1:E:200:PHE:C	1:E:200:PHE:CD1	2.78	0.56
1:C:298:ILE:O	1:C:298:ILE:HG22	2.05	0.56
1:E:114:PHE:CE1	1:E:116:ILE:HG23	2.40	0.56
1:E:109:SER:HB2	1:G:166:ILE:HG13	1.88	0.55
1:I:298:ILE:HG22	1:I:298:ILE:O	2.07	0.55
1:K:298:ILE:O	1:K:298:ILE:HG13	2.07	0.55
1:K:200:PHE:CD1	1:K:200:PHE:C	2.81	0.53
1:C:261:ALA:O	1:C:262:GLY:C	2.46	0.53
1:C:114:PHE:CE1	1:C:116:ILE:HG23	2.43	0.53
1:I:200:PHE:CD2	1:I:200:PHE:C	2.81	0.53
1:G:137:PRO:HD2	3:G:401:NAG:H83	1.91	0.53
1:I:114:PHE:CE1	1:I:116:ILE:HG23	2.44	0.52
2:J:71:ASN:OD1	2:J:74:GLU:HG3	2.10	0.52
2:D:30:GLN:NE2	2:D:145:ASP:HB2	2.26	0.51
2:F:145:ASP:OD1	2:F:146:ASN:N	2.44	0.50
1:G:261:ALA:O	1:G:262:GLY:C	2.49	0.50
1:K:261:ALA:O	1:K:262:GLY:C	2.50	0.50
1:A:200:PHE:CD2	1:A:200:PHE:C	2.85	0.50
2:B:71:ASN:OD1	2:B:74:GLU:HG3	2.10	0.50
1:A:261:ALA:O	1:A:262:GLY:C	2.49	0.50
2:L:71:ASN:OD1	2:L:74:GLU:HG3	2.13	0.49
2:J:30:GLN:NE2	2:J:145:ASP:HB2	2.27	0.49
1:I:261:ALA:O	1:I:262:GLY:C	2.51	0.49
2:D:71:ASN:OD1	2:D:74:GLU:HG3	2.13	0.49
3:E:404:NAG:C1	3:E:404:NAG:C8	2.90	0.49
2:H:71:ASN:OD1	2:H:74:GLU:HG3	2.13	0.49
1:I:105:LEU:HD23	1:I:257:MET:HE1	1.95	0.48
2:F:159:TYR:HB3	2:F:160:PRO:HD3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:159:TYR:HB3	2:B:160:PRO:HD3	1.96	0.47
2:F:71:ASN:OD1	2:F:74:GLU:HG3	2.14	0.47
1:C:32:LEU:HD21	2:D:55:VAL:HG11	1.97	0.47
1:G:200:PHE:CD2	1:G:200:PHE:C	2.89	0.47
2:H:159:TYR:HB3	2:H:160:PRO:HD3	1.97	0.47
2:D:76:ARG:HG2	1:E:103:GLU:HG3	1.97	0.46
1:G:105:LEU:HD23	1:G:257:MET:HE1	1.97	0.46
1:C:105:LEU:HD23	1:C:257:MET:HE1	1.97	0.46
2:D:159:TYR:HB3	2:D:160:PRO:HD3	1.96	0.46
1:E:261:ALA:O	1:E:262:GLY:C	2.54	0.46
2:L:145:ASP:OD1	2:L:146:ASN:N	2.49	0.46
2:J:159:TYR:HB3	2:J:160:PRO:HD3	1.98	0.46
1:G:161:LEU:C	1:G:161:LEU:HD23	2.37	0.45
2:J:9:PHE:O	2:J:135:ASN:HA	2.16	0.45
1:K:179:ILE:HD12	1:K:210:PHE:HB3	1.99	0.45
2:L:159:TYR:HB3	2:L:160:PRO:HD3	1.97	0.45
1:E:179:ILE:HD12	1:E:210:PHE:HB3	1.98	0.45
1:C:179:ILE:CD1	1:C:210:PHE:CG	3.00	0.44
2:H:9:PHE:O	2:H:135:ASN:HA	2.17	0.44
1:C:179:ILE:HD12	1:C:210:PHE:CG	2.53	0.44
1:A:179:ILE:CD1	1:A:210:PHE:CG	3.01	0.44
1:A:105:LEU:HD23	1:A:257:MET:HE1	1.99	0.44
1:G:179:ILE:HD12	1:G:210:PHE:CG	2.53	0.44
2:L:2:LEU:HG	2:L:3:PHE:N	2.33	0.43
1:C:299:THR:HB	1:C:303:CYS:SG	2.58	0.43
1:G:179:ILE:CD1	1:G:210:PHE:CG	3.01	0.43
2:J:160:PRO:O	2:J:161:LYS:C	2.56	0.43
1:K:93:GLY:HA3	1:K:227:MET:O	2.18	0.43
1:A:161:LEU:C	1:A:161:LEU:HD23	2.39	0.43
1:A:179:ILE:HD12	1:A:210:PHE:CG	2.54	0.43
1:G:93:GLY:HA3	1:G:227:MET:O	2.19	0.43
2:B:9:PHE:O	2:B:135:ASN:HA	2.19	0.43
1:E:93:GLY:HA3	1:E:227:MET:O	2.19	0.43
1:I:34:GLU:HB2	1:I:288:THR:HG21	2.00	0.42
1:K:130:LYS:HD2	1:K:130:LYS:HA	1.84	0.42
2:F:94:TYR:CZ	2:F:98:LEU:HD11	2.55	0.42
2:F:2:LEU:HG	2:F:3:PHE:N	2.35	0.42
2:D:9:PHE:O	2:D:135:ASN:HA	2.20	0.42
2:F:126:LEU:HD21	2:F:152:VAL:HG21	2.02	0.42
1:C:34:GLU:HB2	1:C:288:THR:HG21	2.01	0.42
1:K:113:ARG:NH2	1:K:146:LYS:HE3	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:93:GLY:HA3	1:I:227:MET:O	2.19	0.42
1:A:299:THR:HB	1:A:303:CYS:SG	2.59	0.41
1:G:44:LEU:HD21	1:G:300:ILE:HG22	2.02	0.41
1:K:5:ILE:C	1:K:5:ILE:HD12	2.41	0.41
1:C:93:GLY:HA3	1:C:227:MET:O	2.20	0.41
1:E:112:GLU:HG3	1:K:216:ILE:HD13	2.00	0.41
2:H:2:LEU:HG	2:H:3:PHE:N	2.34	0.41
2:L:9:PHE:O	2:L:135:ASN:HA	2.19	0.41
1:K:257:MET:HE2	1:K:257:MET:HB2	1.92	0.41
1:G:299:THR:HB	1:G:303:CYS:SG	2.60	0.41
1:E:299:THR:HB	1:E:303:CYS:SG	2.61	0.41
1:G:286:ILE:CG2	1:G:288:THR:HG22	2.51	0.41
1:K:114:PHE:CZ	1:K:116:ILE:HG23	2.56	0.41
1:K:299:THR:HB	1:K:303:CYS:SG	2.61	0.41
1:A:110:SER:OG	1:A:171:LYS:HE3	2.21	0.40
1:A:93:GLY:HA3	1:A:227:MET:O	2.20	0.40
1:E:5:ILE:HD12	1:E:5:ILE:C	2.41	0.40
2:B:94:TYR:CZ	2:B:98:LEU:HD11	2.56	0.40
2:J:124:ASN:O	2:J:127:LYS:HE3	2.21	0.40
2:L:126:LEU:HD21	2:L:152:VAL:HG21	2.03	0.40
1:A:137:PRO:HD2	3:A:401:NAG:H83	2.03	0.40
2:B:160:PRO:O	2:B:161:LYS:C	2.60	0.40
1:K:148:LEU:HB3	1:K:249:VAL:HG23	2.03	0.40
1:K:286:ILE:CG2	1:K:288:THR:HG22	2.52	0.40
2:L:124:ASN:O	2:L:127:LYS:HE3	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	321/332 (97%)	303 (94%)	15 (5%)	3 (1%)	19 59

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	321/332 (97%)	301 (94%)	18 (6%)	2 (1%)	27	67
1	E	321/332 (97%)	300 (94%)	19 (6%)	2 (1%)	27	67
1	G	321/332 (97%)	302 (94%)	16 (5%)	3 (1%)	19	59
1	I	321/332 (97%)	302 (94%)	17 (5%)	2 (1%)	27	67
1	K	321/332 (97%)	302 (94%)	16 (5%)	3 (1%)	19	59
2	B	169/182 (93%)	164 (97%)	5 (3%)	0	100	100
2	D	169/182 (93%)	163 (96%)	6 (4%)	0	100	100
2	F	169/182 (93%)	164 (97%)	5 (3%)	0	100	100
2	H	169/182 (93%)	163 (96%)	6 (4%)	0	100	100
2	J	169/182 (93%)	164 (97%)	5 (3%)	0	100	100
2	L	169/182 (93%)	164 (97%)	5 (3%)	0	100	100
All	All	2940/3084 (95%)	2792 (95%)	133 (4%)	15 (0%)	31	71

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	262	GLY
1	C	262	GLY
1	E	262	GLY
1	G	262	GLY
1	I	262	GLY
1	K	262	GLY
1	A	261	ALA
1	E	261	ALA
1	G	261	ALA
1	I	261	ALA
1	K	261	ALA
1	A	53	GLY
1	C	53	GLY
1	K	53	GLY
1	G	53	GLY

### 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	283/290 (98%)	267 (94%)	16 (6%)	23	60
1	C	283/290 (98%)	263 (93%)	20 (7%)	16	50
1	E	283/290 (98%)	258 (91%)	25 (9%)	11	39
1	G	283/290 (98%)	266 (94%)	17 (6%)	21	57
1	I	283/290 (98%)	262 (93%)	21 (7%)	15	47
1	K	283/290 (98%)	258 (91%)	25 (9%)	11	39
2	B	148/156 (95%)	141 (95%)	7 (5%)	29	67
2	D	148/156 (95%)	140 (95%)	8 (5%)	24	62
2	F	148/156 (95%)	142 (96%)	6 (4%)	33	71
2	H	148/156 (95%)	140 (95%)	8 (5%)	24	62
2	J	148/156 (95%)	141 (95%)	7 (5%)	29	67
2	L	148/156 (95%)	142 (96%)	6 (4%)	33	71
All	All	2586/2676 (97%)	2420 (94%)	166 (6%)	19	55

All (166) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	VAL
1	A	29	SER
1	A	36	LYS
1	A	47	VAL
1	A	95	PHE
1	A	132	VAL
1	A	171	LYS
1	A	173	VAL
1	A	194	ASN
1	A	222	ASP
1	A	258	GLU
1	A	259	ARG
1	A	260	ASN
1	A	275	CYS
1	A	298	ILE
1	A	318	LEU
2	B	39	LYS
2	B	61	THR
2	B	64	THR

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Mol	Chain	Res	Type
2	B	102	LEU
2	B	105	GLU
2	B	108	LEU
2	B	164	GLU
1	C	3	LEU
1	C	19	VAL
1	C	29	SER
1	C	95	PHE
1	C	116	ILE
1	C	132	VAL
1	C	171	LYS
1	C	173	VAL
1	C	194	ASN
1	C	211	LYS
1	C	222	ASP
1	C	250	VAL
1	C	258	GLU
1	C	259	ARG
1	C	260	ASN
1	C	272	VAL
1	C	273	HIS
1	C	275	CYS
1	C	288	THR
1	C	318	LEU
2	D	38	LEU
2	D	39	LYS
2	D	64	THR
2	D	102	LEU
2	D	105	GLU
2	D	108	LEU
2	D	161	LYS
2	D	167	LYS
1	E	3	LEU
1	E	5	ILE
1	E	19	VAL
1	E	29	SER
1	E	47	VAL
1	E	84	SER
1	E	95	PHE
1	E	103	GLU
1	E	105	LEU
1	E	116	ILE

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Mol	Chain	Res	Type
1	E	171	LYS
1	E	173	VAL
1	E	175	VAL
1	E	179	ILE
1	E	194	ASN
1	E	211	LYS
1	E	222	ASP
1	E	249	VAL
1	E	250	VAL
1	E	258	GLU
1	E	259	ARG
1	E	260	ASN
1	E	272	VAL
1	E	298	ILE
1	E	318	LEU
2	F	64	THR
2	F	105	GLU
2	F	108	LEU
2	F	121	LYS
2	F	152	VAL
2	F	168	LEU
1	G	19	VAL
1	G	29	SER
1	G	36	LYS
1	G	47	VAL
1	G	95	PHE
1	G	132	VAL
1	G	171	LYS
1	G	173	VAL
1	G	194	ASN
1	G	222	ASP
1	G	258	GLU
1	G	259	ARG
1	G	260	ASN
1	G	275	CYS
1	G	289	SER
1	G	298	ILE
1	G	318	LEU
2	H	39	LYS
2	H	61	THR
2	H	64	THR
2	H	102	LEU

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Mol	Chain	Res	Type
2	H	105	GLU
2	H	108	LEU
2	H	161	LYS
2	H	164	GLU
1	I	3	LEU
1	I	19	VAL
1	I	29	SER
1	I	95	PHE
1	I	116	ILE
1	I	132	VAL
1	I	171	LYS
1	I	173	VAL
1	I	194	ASN
1	I	211	LYS
1	I	222	ASP
1	I	250	VAL
1	I	258	GLU
1	I	259	ARG
1	I	260	ASN
1	I	272	VAL
1	I	273	HIS
1	I	275	CYS
1	I	288	THR
1	I	318	LEU
1	I	322	PRO
2	J	38	LEU
2	J	39	LYS
2	J	64	THR
2	J	102	LEU
2	J	105	GLU
2	J	108	LEU
2	J	167	LYS
1	K	3	LEU
1	K	5	ILE
1	K	19	VAL
1	K	29	SER
1	K	47	VAL
1	K	84	SER
1	K	103	GLU
1	K	105	LEU
1	K	116	ILE
1	K	132	VAL

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Mol	Chain	Res	Type
1	K	171	LYS
1	K	173	VAL
1	K	175	VAL
1	K	179	ILE
1	K	194	ASN
1	K	211	LYS
1	K	222	ASP
1	K	249	VAL
1	K	250	VAL
1	K	258	GLU
1	K	259	ARG
1	K	260	ASN
1	K	272	VAL
1	K	298	ILE
1	K	318	LEU
2	L	64	THR
2	L	105	GLU
2	L	108	LEU
2	L	121	LYS
2	L	152	VAL
2	L	168	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

20 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	$\# Z  > 2$
3	NAG	A	401	1,3	14,14,15	0.72	0	17,19,21	2.36	6 (35%)
3	NAG	A	402	3	14,14,15	0.77	0	17,19,21	1.74	3 (17%)
3	NAG	A	403	1	14,14,15	0.28	0	17,19,21	0.66	0
3	NAG	A	404	1	14,14,15	0.77	1 (7%)	17,19,21	1.47	2 (11%)
3	NAG	C	401	1,3	14,14,15	1.03	1 (7%)	17,19,21	1.59	3 (17%)
3	NAG	C	402	3	14,14,15	0.76	0	17,19,21	1.29	4 (23%)
3	NAG	C	403	1	14,14,15	0.91	0	17,19,21	2.25	6 (35%)
3	NAG	E	401	1,3	14,14,15	0.44	0	17,19,21	1.86	3 (17%)
3	NAG	E	402	3	14,14,15	0.71	0	17,19,21	1.53	3 (17%)
3	NAG	E	403	-	14,14,15	0.96	1 (7%)	17,19,21	1.83	3 (17%)
3	NAG	E	404	1	14,14,15	0.63	0	17,19,21	1.21	2 (11%)
3	NAG	G	401	1,3	14,14,15	0.58	0	17,19,21	1.70	5 (29%)
3	NAG	G	402	3	14,14,15	0.71	0	17,19,21	2.02	7 (41%)
3	NAG	G	403	1	14,14,15	0.30	0	17,19,21	0.65	0
3	NAG	G	404	1	14,14,15	0.91	0	17,19,21	2.13	3 (17%)
3	NAG	I	401	1,3	14,14,15	0.57	0	17,19,21	1.76	2 (11%)
3	NAG	I	402	3	14,14,15	0.70	0	17,19,21	1.55	2 (11%)
3	NAG	I	403	1	14,14,15	0.75	0	17,19,21	1.81	3 (17%)
3	NAG	K	401	1	14,14,15	0.51	0	17,19,21	2.46	8 (47%)
3	NAG	K	402	1	14,14,15	0.69	0	17,19,21	2.77	8 (47%)

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
3	NAG	A	401	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	402	3	-	0/6/23/26	0/1/1/1
3	NAG	A	403	1	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	404	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	401	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	402	3	-	0/6/23/26	0/1/1/1
3	NAG	C	403	1	-	0/6/23/26	0/1/1/1
3	NAG	E	401	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	402	3	-	0/6/23/26	0/1/1/1
3	NAG	E	403	-	-	0/6/23/26	0/1/1/1
3	NAG	E	404	1	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	G	401	1,3	-	0/6/23/26	0/1/1/1
3	NAG	G	402	3	-	0/6/23/26	0/1/1/1
3	NAG	G	403	1	-	0/6/23/26	0/1/1/1
3	NAG	G	404	1	-	0/6/23/26	0/1/1/1
3	NAG	I	401	1,3	-	0/6/23/26	0/1/1/1
3	NAG	I	402	3	-	0/6/23/26	0/1/1/1
3	NAG	I	403	1	-	0/6/23/26	0/1/1/1
3	NAG	K	401	1	-	0/6/23/26	0/1/1/1
3	NAG	K	402	1	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	401	NAG	O5-C1	-3.13	1.38	1.43
3	A	404	NAG	C1-C2	2.29	1.55	1.52
3	E	403	NAG	C1-C2	2.46	1.55	1.52

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	NAG	O5-C1-C2	-6.49	102.56	111.52
3	I	401	NAG	O5-C1-C2	-5.04	104.56	111.52
3	E	401	NAG	O5-C1-C2	-4.63	105.13	111.52
3	C	401	NAG	O5-C1-C2	-4.62	105.14	111.52
3	K	401	NAG	C4-C3-C2	-3.66	105.65	111.02
3	C	403	NAG	C1-C2-N2	-3.62	104.31	110.49
3	G	402	NAG	C3-C4-C5	-3.54	103.91	110.24
3	E	404	NAG	O5-C1-C2	-3.22	107.07	111.52
3	A	402	NAG	C1-C2-N2	-3.00	105.36	110.49
3	I	402	NAG	C1-C2-N2	-2.96	105.43	110.49
3	G	402	NAG	C4-C3-C2	-2.95	106.70	111.02
3	A	401	NAG	O7-C7-C8	-2.86	116.89	122.07
3	A	401	NAG	C3-C4-C5	-2.73	105.36	110.24
3	E	402	NAG	O3-C3-C4	-2.49	104.52	110.34
3	K	402	NAG	O7-C7-C8	-2.41	117.71	122.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	401	NAG	C3-C4-C5	-2.40	105.94	110.24
3	K	401	NAG	C3-C4-C5	-2.36	106.02	110.24
3	C	401	NAG	O5-C5-C6	-2.34	103.45	107.15
3	G	401	NAG	C4-C3-C2	-2.32	107.62	111.02
3	C	402	NAG	O7-C7-C8	-2.32	117.88	122.07
3	C	402	NAG	O5-C5-C4	-2.25	105.34	110.83
3	K	401	NAG	O7-C7-C8	-2.22	118.05	122.07
3	E	404	NAG	C2-N2-C7	-2.21	119.72	122.94
3	G	401	NAG	C2-N2-C7	-2.21	119.72	122.94
3	C	402	NAG	C3-C4-C5	-2.09	106.50	110.24
3	G	401	NAG	O7-C7-C8	-2.05	118.37	122.07
3	G	404	NAG	C2-N2-C7	2.05	125.94	122.94
3	G	404	NAG	O4-C4-C5	2.08	114.51	109.31
3	C	402	NAG	O4-C4-C5	2.11	114.59	109.31
3	A	401	NAG	C1-O5-C5	2.14	115.13	112.19
3	I	403	NAG	O5-C1-C2	2.17	114.52	111.52
3	G	402	NAG	O3-C3-C2	2.22	114.14	109.39
3	C	403	NAG	O6-C6-C5	2.24	119.11	111.29
3	K	401	NAG	O7-C7-N2	2.27	126.24	121.94
3	C	403	NAG	O4-C4-C5	2.28	115.01	109.31
3	K	402	NAG	C8-C7-N2	2.38	120.27	116.10
3	G	402	NAG	C1-O5-C5	2.39	115.47	112.19
3	A	402	NAG	O5-C5-C6	2.61	111.28	107.15
3	E	403	NAG	C1-O5-C5	2.61	115.78	112.19
3	G	401	NAG	C1-C2-N2	2.65	115.02	110.49
3	K	401	NAG	C1-C2-N2	2.74	115.17	110.49
3	K	402	NAG	C1-C2-N2	2.77	115.23	110.49
3	K	401	NAG	O4-C4-C3	2.79	116.85	110.34
3	C	403	NAG	O5-C5-C6	2.80	111.58	107.15
3	A	401	NAG	C1-C2-N2	2.83	115.32	110.49
3	K	402	NAG	C3-C4-C5	2.92	115.47	110.24
3	E	401	NAG	O5-C5-C6	2.96	111.84	107.15
3	E	402	NAG	O5-C5-C6	2.99	111.88	107.15
3	G	402	NAG	C1-C2-N2	3.01	115.62	110.49
3	A	404	NAG	C1-C2-N2	3.05	115.70	110.49
3	E	401	NAG	C1-O5-C5	3.06	116.40	112.19
3	G	402	NAG	O5-C5-C6	3.21	112.23	107.15
3	I	401	NAG	C1-O5-C5	3.23	116.63	112.19
3	G	401	NAG	O5-C5-C6	3.24	112.28	107.15
3	K	402	NAG	O5-C1-C2	3.24	116.00	111.52
3	E	403	NAG	O5-C5-C6	3.26	112.31	107.15
3	K	402	NAG	C4-C3-C2	3.48	116.12	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	402	NAG	C4-C3-C2	3.61	116.31	111.02
3	A	401	NAG	O7-C7-N2	3.63	128.80	121.94
3	K	401	NAG	O3-C3-C4	3.65	118.86	110.34
3	G	402	NAG	O4-C4-C5	3.66	118.47	109.31
3	C	403	NAG	C4-C3-C2	3.68	116.41	111.02
3	I	402	NAG	C1-O5-C5	3.85	117.48	112.19
3	A	404	NAG	C1-O5-C5	3.87	117.52	112.19
3	K	402	NAG	O5-C5-C6	3.89	113.31	107.15
3	I	403	NAG	C4-C3-C2	4.23	117.22	111.02
3	E	403	NAG	O5-C1-C2	4.43	117.64	111.52
3	A	402	NAG	C4-C3-C2	4.82	118.09	111.02
3	I	403	NAG	O5-C5-C6	5.11	115.24	107.15
3	C	403	NAG	C1-O5-C5	5.77	120.13	112.19
3	K	401	NAG	C1-O5-C5	6.18	120.69	112.19
3	G	404	NAG	O5-C5-C6	6.85	118.00	107.15
3	K	402	NAG	C1-O5-C5	7.57	122.60	112.19

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	E	404	NAG	C1
3	A	403	NAG	C1

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	NAG	1	0
3	E	404	NAG	1	0
3	G	401	NAG	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, 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	323/332 (97%)	-0.28	2 (0%) 89 71	34, 59, 90, 153	0
1	C	323/332 (97%)	-0.31	1 (0%) 93 83	38, 57, 89, 138	0
1	E	323/332 (97%)	-0.34	0 100 100	37, 61, 89, 107	0
1	G	323/332 (97%)	-0.34	1 (0%) 93 83	41, 63, 96, 138	0
1	I	323/332 (97%)	-0.29	3 (0%) 84 62	38, 62, 92, 133	0
1	K	323/332 (97%)	-0.30	1 (0%) 93 83	43, 65, 95, 140	0
2	B	171/182 (93%)	-0.27	1 (0%) 89 71	39, 71, 105, 136	0
2	D	171/182 (93%)	0.13	4 (2%) 60 31	44, 78, 124, 159	0
2	F	171/182 (93%)	-0.17	3 (1%) 68 39	31, 77, 128, 152	0
2	H	171/182 (93%)	-0.19	0 100 100	44, 71, 109, 133	0
2	J	171/182 (93%)	-0.36	0 100 100	42, 62, 95, 136	0
2	L	171/182 (93%)	-0.31	1 (0%) 89 71	43, 67, 112, 134	0
All	All	2964/3084 (96%)	-0.27	17 (0%) 89 71	31, 64, 107, 159	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	172	GLU	5.4
1	A	71	SER	5.2
2	D	169	ASN	3.5
2	D	168	LEU	3.1
2	B	172	GLU	3.0
2	L	165	GLU	2.9
1	G	46	GLY	2.4
2	D	164	GLU	2.4
1	K	69	SER	2.3
1	I	156	ASN	2.2
2	F	140	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
2	F	29	GLU	2.1
1	I	46	GLY	2.1
2	F	19	ASP	2.1
1	A	70	LEU	2.0
1	C	0	SER	2.0
1	I	275	CYS	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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	E	403	14/15	0.61	0.27	90,116,128,134	0
3	NAG	A	404	14/15	0.62	0.39	92,115,124,127	0
3	NAG	G	403	14/15	0.72	0.34	58,98,103,104	0
3	NAG	E	404	14/15	0.78	0.24	106,126,139,140	0
3	NAG	A	403	14/15	0.78	0.33	48,92,100,101	0
3	NAG	C	402	14/15	0.79	0.33	82,92,105,109	0
3	NAG	A	402	14/15	0.81	0.37	80,93,104,106	0
3	NAG	G	404	14/15	0.81	0.36	84,93,98,102	0
3	NAG	K	402	14/15	0.81	0.36	83,100,110,116	0
3	NAG	I	403	14/15	0.82	0.36	92,111,118,125	0
3	NAG	C	403	14/15	0.84	0.29	67,74,76,78	0
3	NAG	G	402	14/15	0.84	0.29	68,93,106,107	0
3	NAG	E	402	14/15	0.87	0.15	57,67,80,82	0
3	NAG	I	402	14/15	0.88	0.28	64,75,81,83	0
3	NAG	K	401	14/15	0.89	0.26	63,79,82,85	0
3	NAG	A	401	14/15	0.93	0.25	56,65,72,75	0
3	NAG	G	401	14/15	0.94	0.23	62,69,84,90	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	I	401	14/15	0.95	0.30	56,61,63,68	0
3	NAG	C	401	14/15	0.96	0.20	54,65,70,80	0
3	NAG	E	401	14/15	0.97	0.13	35,44,50,53	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.