



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:37 PM GMT

PDB ID : 3U7U  
Title : Crystal structure of extracellular region of human epidermal growth factor receptor 4 in complex with neuregulin-1 beta  
Authors : Liu, P.; Cleveland IV, T.E.; Bouyain, S.; Longo, P.A.; Leahy, D.J.  
Deposited on : 2011-10-14  
Resolution : 3.03 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

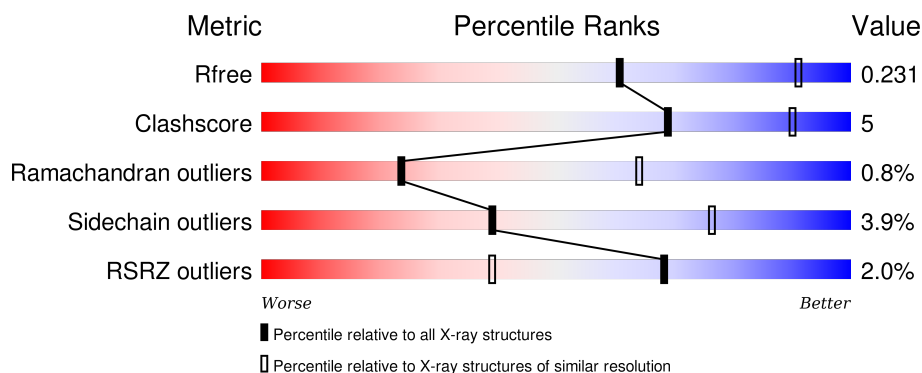
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.03 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))
$R_{free}$	91344	1995 (3.08-3.00)
Clashscore	102246	2351 (3.08-3.00)
Ramachandran outliers	100387	2272 (3.08-3.00)
Sidechain outliers	100360	2275 (3.08-3.00)
RSRZ outliers	91569	2013 (3.08-3.00)

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	615	<div> <div>4%</div> <div>77%</div> <div>11%</div> <div>•</div> <div>11%</div> </div>
1	B	615	<div> <div>83%</div> <div>14%</div> <div>••</div> </div>
1	C	615	<div> <div>3%</div> <div>84%</div> <div>12%</div> <div>••</div> </div>
1	D	615	<div> <div>84%</div> <div>13%</div> <div>••</div> </div>
1	E	615	<div> <div>2%</div> <div>80%</div> <div>15%</div> <div>••</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	615	
2	G	55	
2	H	55	
2	I	55	
2	J	55	
2	K	55	
2	L	55	

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	B	1004	-	-	-	X
3	NAG	D	1002	-	-	-	X
3	NAG	D	1004	-	-	-	X
3	NAG	D	1005	-	-	-	X
3	NAG	E	1003	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 30527 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 Receptor tyrosine-protein kinase erbB-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	550	Total	C	N	O	S	0	0	0
			4284	2658	746	829	51			
1	B	601	Total	C	N	O	S	0	0	0
			4685	2908	815	904	58			
1	C	598	Total	C	N	O	S	0	0	0
			4649	2879	812	900	58			
1	D	600	Total	C	N	O	S	0	0	0
			4683	2906	815	904	58			
1	E	593	Total	C	N	O	S	0	0	0
			4615	2856	807	894	58			
1	F	591	Total	C	N	O	S	0	0	0
			4595	2843	804	890	58			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	524	ASP	GLY	CONFLICT	UNP Q15303
B	524	ASP	GLY	CONFLICT	UNP Q15303
C	524	ASP	GLY	CONFLICT	UNP Q15303
D	524	ASP	GLY	CONFLICT	UNP Q15303
E	524	ASP	GLY	CONFLICT	UNP Q15303
F	524	ASP	GLY	CONFLICT	UNP Q15303

- Molecule 2 is a protein called Neuregulin 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	50	Total	C	N	O	S	0	0	0
			391	243	67	73	8			
2	H	51	Total	C	N	O	S	0	0	0
			398	247	68	75	8			
2	I	50	Total	C	N	O	S	0	0	0
			391	243	67	73	8			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	J	51	Total	C	N	O	S	0	0	0
			398	247	68	75	8			
2	K	49	Total	C	N	O	S	0	0	0
			385	240	66	71	8			
2	L	51	Total	C	N	O	S	0	0	0
			398	247	68	75	8			

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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	B	1	Total	C	N	O		0	0
			14	8	1	5			
3	B	1	Total	C	N	O		0	0
			14	8	1	5			
3	B	1	Total	C	N	O		0	0
			14	8	1	5			
3	B	1	Total	C	N	O		0	0
			14	8	1	5			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	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	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	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	F	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	30	Total	O	0	0
			30	30		

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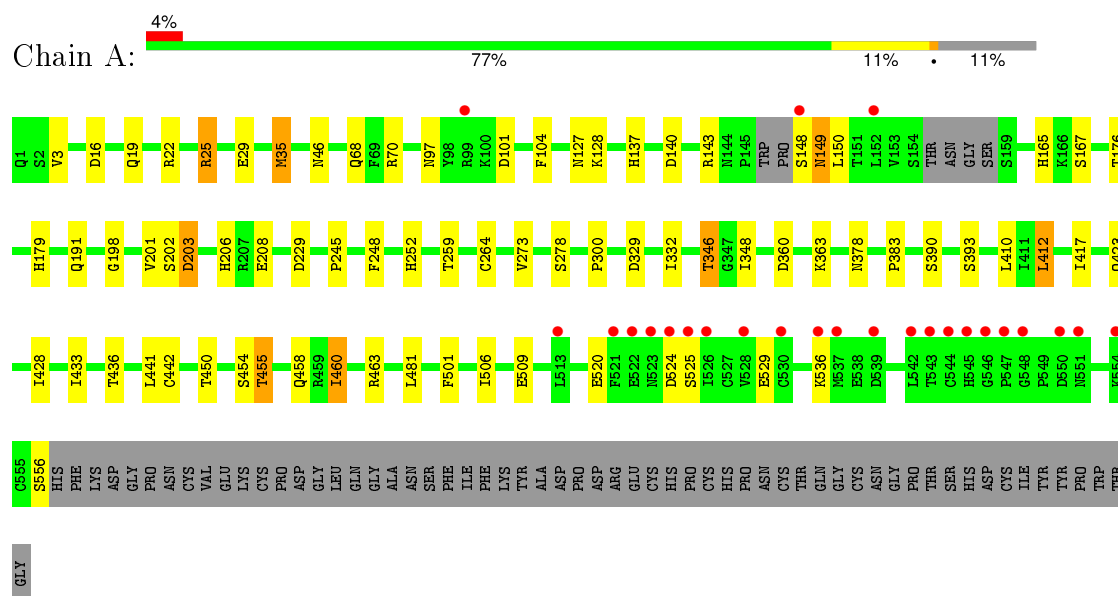
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	63	Total 63	O 63	0	0
4	C	37	Total 37	O 37	0	0
4	D	55	Total 55	O 55	0	0
4	E	33	Total 33	O 33	0	0
4	F	52	Total 52	O 52	0	0
4	G	1	Total 1	O 1	0	0
4	H	9	Total 9	O 9	0	0
4	I	2	Total 2	O 2	0	0
4	J	3	Total 3	O 3	0	0
4	K	2	Total 2	O 2	0	0
4	L	4	Total 4	O 4	0	0

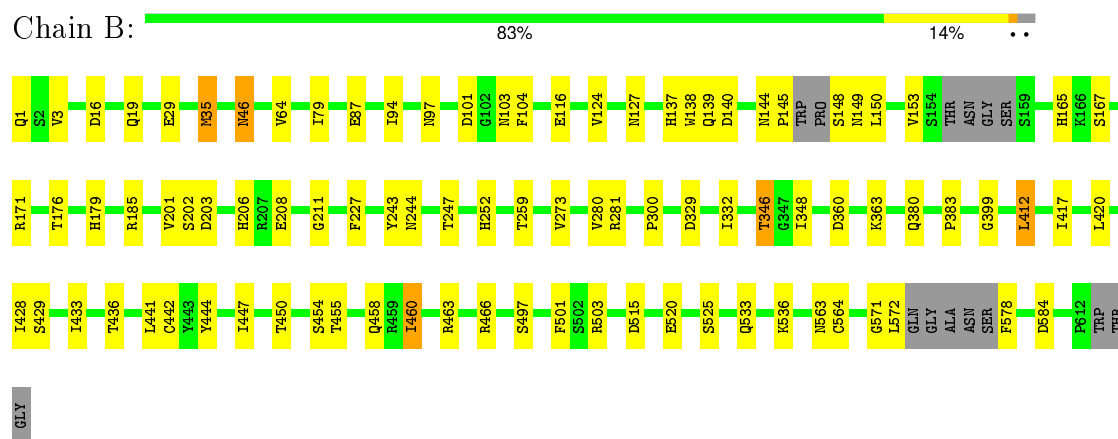
### 3 Residue-property plots

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

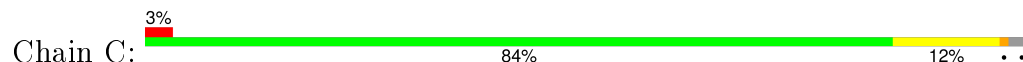
- Molecule 1: Receptor tyrosine-protein kinase erbB-4



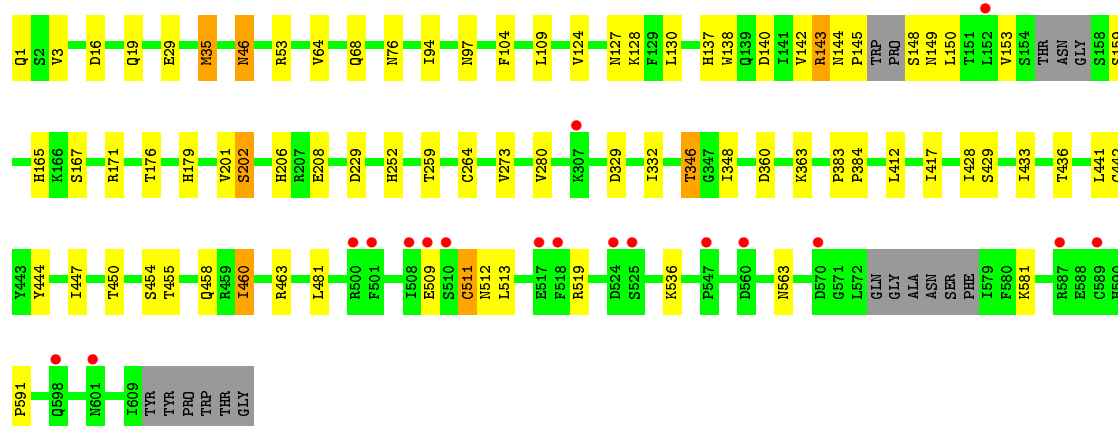
- Molecule 1: Receptor tyrosine-protein kinase erbB-4



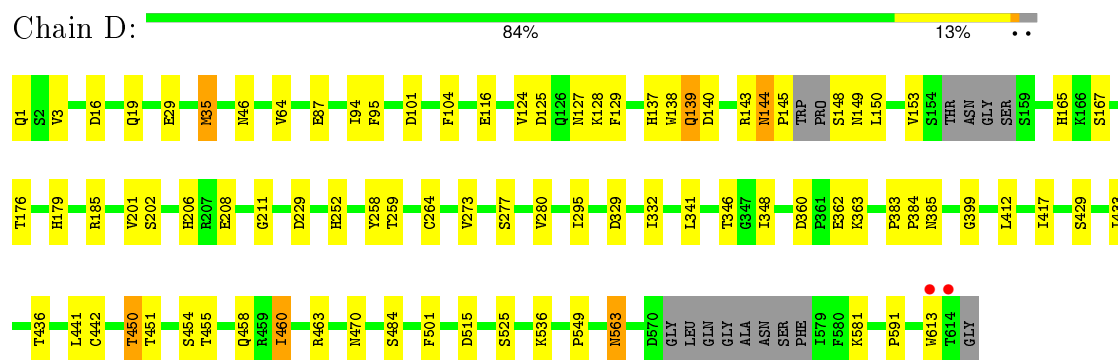
- Molecule 1: Receptor tyrosine-protein kinase erbB-4



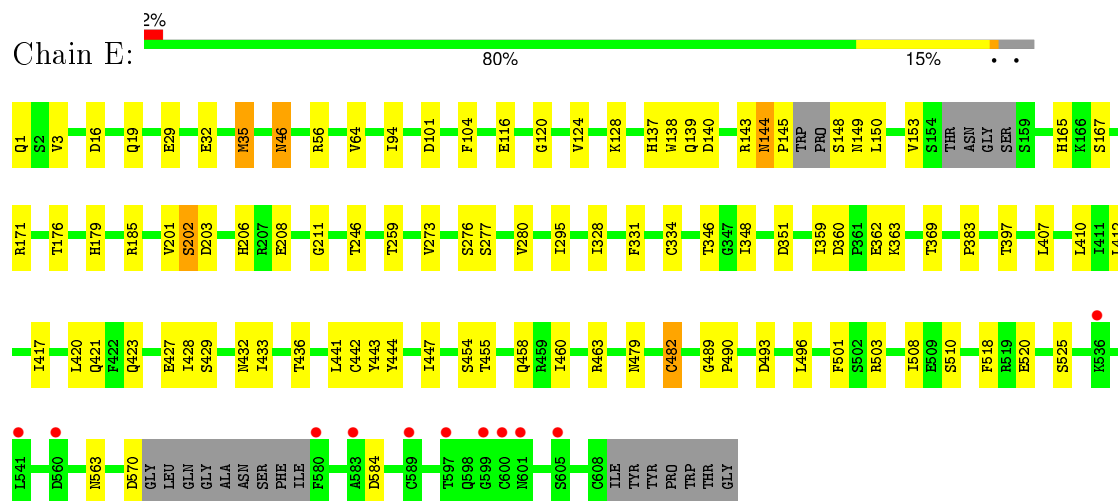




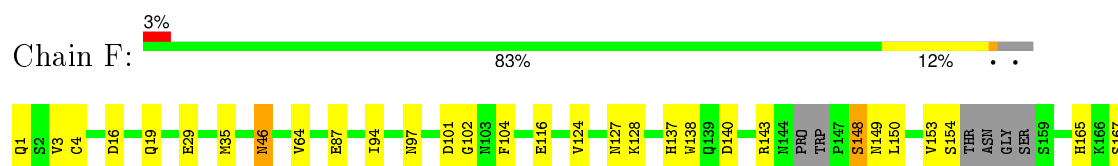
- Molecule 1: Receptor tyrosine-protein kinase erbB-4

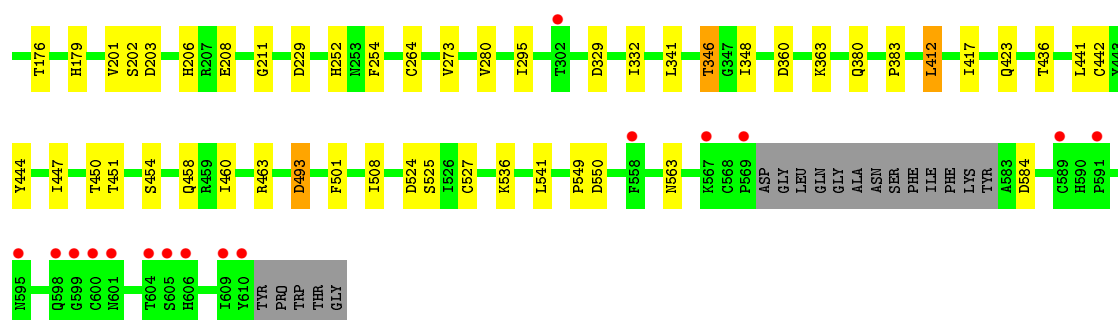


- Molecule 1: Receptor tyrosine-protein kinase erbB-4



- Molecule 1: Receptor tyrosine-protein kinase erbB-4





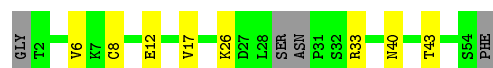
- Molecule 2: Neuregulin 1

Chain G: 85% 5% 9%



- Molecule 2: Neuregulin 1

Chain H: 78% 15% 7%



- Molecule 2: Neuregulin 1

Chain I: 73% 18% 9%



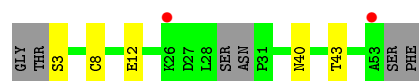
- Molecule 2: Neuregulin 1

Chain J: 76% 16% 7%



- Molecule 2: Neuregulin 1

Chain K: 4% 80% 9% 11%



- Molecule 2: Neuregulin 1

Chain L: 2% 85% 7% 7%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.73Å 223.51Å 146.92Å 90.00° 99.72° 90.00°	Depositor
Resolution (Å)	49.63 – 3.03 49.63 – 3.03	Depositor EDS
% Data completeness (in resolution range)	(Not available) (49.63-3.03) 98.7 (49.63-3.03)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.91 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, $R_{free}$	0.190 , 0.227 0.193 , 0.231	Depositor DCC
$R_{free}$ test set	5198 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	73.3	Xtriage
Anisotropy	0.361	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 69.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 103996 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	30527	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.50	0/4366	0.76	1/5908 (0.0%)
1	B	0.53	0/4783	0.75	0/6476
1	C	0.49	0/4743	0.74	0/6420
1	D	0.51	0/4782	0.75	0/6477
1	E	0.50	0/4709	0.74	0/6374
1	F	0.52	0/4688	0.75	0/6346
2	G	0.48	0/397	0.76	0/528
2	H	0.48	0/404	0.76	0/538
2	I	0.46	0/397	0.74	0/528
2	J	0.49	0/404	0.75	0/538
2	K	0.45	0/391	0.78	1/520 (0.2%)
2	L	0.49	0/404	0.78	0/538
All	All	0.51	0/30468	0.75	2/41191 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	3	SER	C-N-CA	5.80	136.19	121.70
1	A	203	ASP	CB-CG-OD2	5.18	122.96	118.30

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	4284	0	4089	38	0
1	B	4685	0	4432	49	0
1	C	4649	0	4407	45	0
1	D	4683	0	4427	51	0
1	E	4615	0	4365	45	0
1	F	4595	0	4348	47	0
2	G	391	0	373	1	0
2	H	398	0	380	4	0
2	I	391	0	373	4	0
2	J	398	0	380	5	0
2	K	385	0	368	1	0
2	L	398	0	380	2	0
3	A	56	0	51	0	0
3	B	84	0	78	2	0
3	C	28	0	26	0	0
3	D	70	0	65	6	0
3	E	42	0	39	0	0
3	F	84	0	78	2	0
4	A	30	0	0	0	0
4	B	63	0	0	2	0
4	C	37	0	0	1	0
4	D	55	0	0	2	0
4	E	33	0	0	1	0
4	F	52	0	0	2	0
4	G	1	0	0	0	0
4	H	9	0	0	0	0
4	I	2	0	0	0	0
4	J	3	0	0	0	0
4	K	2	0	0	0	0
4	L	4	0	0	0	0
All	All	30527	0	28659	270	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 5.

All (270) 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:D:143:ARG:O	1:D:145:PRO:HD3	1.69	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:145:PRO:HB2	1:F:254:PHE:HB3	1.52	0.88
1:E:3:VAL:HG13	1:E:35:MET:HG3	1.57	0.87
1:D:3:VAL:HG13	1:D:35:MET:HG3	1.56	0.87
1:D:206:HIS:HD2	1:D:208:GLU:H	1.20	0.86
1:F:46:ASN:HA	4:F:624:HOH:O	1.76	0.85
1:A:206:HIS:HD2	1:A:208:GLU:H	1.22	0.85
1:C:3:VAL:HG13	1:C:35:MET:HG3	1.58	0.85
1:A:3:VAL:HG13	1:A:35:MET:HG3	1.60	0.83
1:B:206:HIS:HD2	1:B:208:GLU:H	1.21	0.82
1:F:206:HIS:HD2	1:F:208:GLU:H	1.26	0.82
1:B:3:VAL:HG13	1:B:35:MET:HG3	1.61	0.82
1:E:206:HIS:HD2	1:E:208:GLU:H	1.27	0.79
1:E:397:THR:HG22	1:E:427:GLU:HB3	1.67	0.77
1:B:466:ARG:NH1	3:B:1005:NAG:H2	2.01	0.76
1:C:206:HIS:HD2	1:C:208:GLU:H	1.34	0.74
1:E:143:ARG:O	1:E:145:PRO:HD3	1.88	0.73
1:B:206:HIS:CD2	1:B:208:GLU:H	2.06	0.73
1:C:143:ARG:O	1:C:145:PRO:HD3	1.89	0.73
1:B:466:ARG:HH12	3:B:1005:NAG:H2	1.54	0.71
1:C:145:PRO:HB2	1:F:254:PHE:CB	2.22	0.69
1:A:206:HIS:CD2	1:A:208:GLU:H	2.10	0.68
1:D:385:ASN:HD22	3:D:1004:NAG:H83	1.58	0.68
1:F:3:VAL:HG13	1:F:35:MET:HG3	1.73	0.68
1:C:455:THR:H	1:C:458:GLN:HE21	1.40	0.68
1:D:206:HIS:CD2	1:D:208:GLU:H	2.07	0.67
1:D:143:ARG:C	1:D:145:PRO:HD3	2.14	0.67
1:F:3:VAL:HG13	1:F:35:MET:CG	2.25	0.67
1:C:97:ASN:H	1:C:127:ASN:ND2	1.95	0.65
1:B:46:ASN:HA	4:B:657:HOH:O	1.96	0.65
1:E:206:HIS:CD2	1:E:208:GLU:H	2.12	0.65
1:B:454:SER:H	1:B:458:GLN:HE22	1.45	0.65
1:C:137:HIS:HD2	1:C:140:ASP:HB2	1.62	0.64
1:A:137:HIS:HD2	1:A:140:ASP:HB2	1.62	0.64
1:A:455:THR:H	1:A:458:GLN:HE21	1.47	0.63
1:C:53:ARG:O	1:C:76:ASN:ND2	2.32	0.62
1:C:206:HIS:CD2	1:C:208:GLU:H	2.19	0.61
1:D:454:SER:H	1:D:458:GLN:NE2	1.98	0.61
1:F:423:GLN:HE21	1:F:493:ASP:HB2	1.65	0.61
1:D:167:SER:HA	4:D:632:HOH:O	1.99	0.61
1:F:417:ILE:HD12	1:F:441:LEU:HD13	1.83	0.60
1:D:148:SER:O	1:D:150:LEU:N	2.28	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:HIS:HD2	1:B:167:SER:H	1.50	0.59
1:F:206:HIS:CD2	1:F:208:GLU:H	2.14	0.59
1:E:444:TYR:HA	1:E:447:ILE:HD12	1.83	0.59
1:B:137:HIS:HD2	1:B:140:ASP:HB2	1.67	0.58
1:F:97:ASN:H	1:F:127:ASN:HD22	1.50	0.58
1:F:165:HIS:HD2	1:F:167:SER:H	1.51	0.58
1:F:3:VAL:CG1	1:F:35:MET:HG3	2.32	0.58
1:D:501:PHE:CD2	1:D:525:SER:HA	2.39	0.58
1:F:451:THR:HG22	3:F:1004:NAG:HN2	1.68	0.58
1:A:97:ASN:HD22	1:A:127:ASN:HD21	1.52	0.57
1:F:102:GLY:HA3	4:F:626:HOH:O	2.04	0.57
1:F:137:HIS:HD2	1:F:140:ASP:HB2	1.69	0.57
1:A:25:ARG:CD	3:D:1005:NAG:H62	2.34	0.57
1:B:501:PHE:CD2	1:B:525:SER:HA	2.39	0.57
1:D:139:GLN:O	1:D:185:ARG:NH2	2.37	0.57
1:A:390:SER:O	1:A:393:SER:HB2	2.04	0.57
1:A:454:SER:H	1:A:458:GLN:NE2	2.02	0.56
1:A:25:ARG:HD3	3:D:1005:NAG:H62	1.86	0.56
1:E:417:ILE:HD12	1:E:441:LEU:HD13	1.87	0.56
1:C:165:HIS:HD2	1:C:167:SER:H	1.53	0.56
1:E:165:HIS:HD2	1:E:167:SER:H	1.53	0.56
1:B:417:ILE:HD12	1:B:441:LEU:HD13	1.88	0.56
1:C:202:SER:HB3	1:D:201:VAL:HB	1.86	0.56
1:E:501:PHE:CD2	1:E:525:SER:HA	2.41	0.56
1:D:258:TYR:HA	1:D:277:SER:O	2.05	0.56
1:D:148:SER:C	1:D:150:LEU:H	2.07	0.56
1:D:360:ASP:HB3	1:D:363:LYS:HG3	1.88	0.55
1:B:571:GLY:HA2	1:B:578:PHE:HA	1.88	0.55
1:E:328:ILE:HD13	1:E:359:ILE:HD11	1.88	0.55
1:E:137:HIS:HD2	1:E:140:ASP:HB2	1.72	0.55
1:F:501:PHE:CD2	1:F:525:SER:HA	2.41	0.55
1:D:454:SER:H	1:D:458:GLN:HE22	1.54	0.55
1:A:248:PHE:HB2	1:B:281:ARG:HG3	1.87	0.55
1:F:454:SER:H	1:F:458:GLN:NE2	2.04	0.55
1:E:407:LEU:HD23	1:E:432:ASN:HB2	1.89	0.55
2:I:5:LEU:HD21	2:I:35:LEU:HD21	1.87	0.55
1:A:165:HIS:HD2	1:A:167:SER:H	1.53	0.55
1:D:124:VAL:HB	1:D:153:VAL:HG22	1.89	0.54
1:F:87:GLU:OE2	2:L:33:ARG:HD3	2.07	0.54
1:A:501:PHE:CD2	1:A:525:SER:HA	2.41	0.54
1:C:417:ILE:HD12	1:C:441:LEU:HD13	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:454:SER:H	1:B:458:GLN:NE2	2.05	0.54
1:C:124:VAL:HB	1:C:153:VAL:HG22	1.88	0.54
1:A:417:ILE:HD12	1:A:441:LEU:HD13	1.88	0.54
1:C:97:ASN:H	1:C:127:ASN:HD22	1.55	0.54
1:B:399:GLY:O	1:B:429:SER:HB2	2.06	0.54
1:D:165:HIS:HD2	1:D:167:SER:H	1.53	0.54
1:F:360:ASP:HB3	1:F:363:LYS:HG3	1.89	0.54
1:D:417:ILE:HD12	1:D:441:LEU:HD13	1.89	0.54
1:D:137:HIS:HD2	1:D:140:ASP:HB2	1.73	0.54
1:B:360:ASP:HB3	1:B:363:LYS:HG3	1.90	0.54
1:E:124:VAL:HB	1:E:153:VAL:HG22	1.89	0.54
1:A:346:THR:HG21	2:G:17:VAL:HG11	1.89	0.54
1:C:481:LEU:HD21	1:C:509:GLU:HG3	1.90	0.53
1:C:360:ASP:HB3	1:C:363:LYS:HG3	1.90	0.53
1:E:346:THR:HG23	1:E:351:ASP:HB2	1.91	0.53
1:E:454:SER:H	1:E:458:GLN:HE22	1.57	0.53
1:F:116:GLU:HB2	1:F:211:GLY:HA2	1.90	0.53
1:B:124:VAL:HB	1:B:153:VAL:HG22	1.90	0.53
1:B:97:ASN:HD22	1:B:127:ASN:HD21	1.56	0.53
1:E:116:GLU:HB2	1:E:211:GLY:HA2	1.91	0.53
1:E:479:ASN:O	1:E:482:CYS:HB2	2.09	0.53
1:A:360:ASP:HB3	1:A:363:LYS:HG3	1.90	0.52
1:C:444:TYR:HA	1:C:447:ILE:HD12	1.92	0.52
1:A:378:ASN:HA	1:A:410:LEU:HB2	1.91	0.52
1:A:97:ASN:H	1:A:127:ASN:HD22	1.58	0.51
1:E:360:ASP:HB3	1:E:363:LYS:HG3	1.92	0.51
1:E:273:VAL:HG13	1:E:280:VAL:HG23	1.92	0.51
1:F:124:VAL:HB	1:F:153:VAL:HG22	1.92	0.51
1:E:176:THR:H	1:E:179:HIS:HD2	1.58	0.51
1:D:273:VAL:HG13	1:D:280:VAL:HG23	1.93	0.51
1:C:329:ASP:O	1:C:332:ILE:HG12	2.11	0.50
1:C:346:THR:HG21	2:I:17:VAL:HG11	1.94	0.50
1:F:423:GLN:NE2	1:F:493:ASP:HB2	2.26	0.50
1:F:454:SER:H	1:F:458:GLN:HE22	1.59	0.50
1:B:346:THR:HG21	2:H:17:VAL:HG11	1.93	0.50
1:B:116:GLU:HB2	1:B:211:GLY:HA2	1.95	0.49
1:D:329:ASP:O	1:D:332:ILE:HG12	2.12	0.49
1:C:176:THR:H	1:C:179:HIS:HD2	1.57	0.49
1:A:329:ASP:O	1:A:332:ILE:HG12	2.12	0.49
1:C:176:THR:H	1:C:179:HIS:CD2	2.30	0.49
1:F:273:VAL:HG13	1:F:280:VAL:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:139:GLN:O	1:E:185:ARG:NH2	2.44	0.49
1:F:329:ASP:O	1:F:332:ILE:HG12	2.12	0.49
1:E:176:THR:H	1:E:179:HIS:CD2	2.31	0.49
1:B:444:TYR:HA	1:B:447:ILE:HD12	1.95	0.49
1:F:346:THR:HG21	2:L:17:VAL:HG11	1.95	0.48
1:B:455:THR:H	1:B:458:GLN:HE21	1.61	0.48
1:F:176:THR:H	1:F:179:HIS:HD2	1.61	0.48
2:I:6:VAL:HG13	2:I:26:LYS:HE2	1.96	0.48
1:D:176:THR:H	1:D:179:HIS:HD2	1.59	0.48
1:A:22:ARG:HD2	1:D:484:SER:HB2	1.94	0.48
1:F:4:CYS:O	1:F:35:MET:HB2	2.13	0.48
1:C:138:TRP:HE1	1:C:153:VAL:HG21	1.79	0.48
1:B:176:THR:H	1:B:179:HIS:HD2	1.62	0.48
1:F:444:TYR:HA	1:F:447:ILE:HD12	1.94	0.48
1:E:443:TYR:CG	1:E:489:GLY:HA2	2.49	0.48
1:C:171:ARG:HH21	1:F:295:ILE:CD1	2.27	0.48
1:C:97:ASN:HD22	1:C:127:ASN:HD21	1.60	0.47
1:F:138:TRP:HE1	1:F:153:VAL:HG21	1.78	0.47
1:C:273:VAL:HG13	1:C:280:VAL:HG23	1.96	0.47
1:A:97:ASN:HD22	1:A:127:ASN:ND2	2.11	0.47
1:B:329:ASP:O	1:B:332:ILE:HG12	2.15	0.47
1:C:511:CYS:HB3	1:C:513:LEU:HG	1.97	0.47
1:B:171:ARG:HH21	1:E:295:ILE:HD11	1.80	0.47
1:C:454:SER:H	1:C:458:GLN:NE2	2.11	0.47
1:A:176:THR:H	1:A:179:HIS:HD2	1.62	0.47
1:D:16:ASP:HB3	1:D:19:GLN:HB2	1.97	0.47
1:E:276:SER:OG	1:E:277:SER:N	2.48	0.47
1:D:116:GLU:HB2	1:D:211:GLY:HA2	1.96	0.47
1:C:512:ASN:HB3	1:C:519:ARG:HA	1.97	0.47
1:C:16:ASP:HB3	1:C:19:GLN:HB2	1.97	0.46
1:A:16:ASP:HB3	1:A:19:GLN:HB2	1.97	0.46
1:D:176:THR:H	1:D:179:HIS:CD2	2.32	0.46
1:D:348:ILE:HD12	1:D:383:PRO:HD3	1.97	0.46
1:E:16:ASP:HB3	1:E:19:GLN:HB2	1.98	0.46
1:A:148:SER:C	1:A:150:LEU:H	2.18	0.46
1:E:138:TRP:HE1	1:E:153:VAL:HG21	1.81	0.46
1:C:137:HIS:CD2	1:C:140:ASP:HB2	2.48	0.46
1:B:176:THR:H	1:B:179:HIS:CD2	2.34	0.45
1:A:348:ILE:HD12	1:A:383:PRO:HD3	1.98	0.45
1:D:450:THR:HG22	3:D:1005:NAG:H2	1.97	0.45
1:E:421:GLN:NE2	1:E:490:PRO:HD2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:420:LEU:HG	1:B:441:LEU:HD11	1.98	0.45
1:D:362:GLU:OE1	3:D:1004:NAG:O4	2.24	0.45
1:B:139:GLN:O	1:B:185:ARG:NH2	2.50	0.45
1:F:16:ASP:HB3	1:F:19:GLN:HB2	1.99	0.45
1:D:138:TRP:HE1	1:D:153:VAL:HG21	1.82	0.45
1:B:428:ILE:HD13	1:B:433:ILE:HD11	1.97	0.45
1:A:176:THR:H	1:A:179:HIS:CD2	2.35	0.45
1:B:16:ASP:HB3	1:B:19:GLN:HB2	1.99	0.45
1:D:455:THR:H	1:D:458:GLN:HE21	1.65	0.45
1:B:87:GLU:OE2	2:H:33:ARG:HD3	2.17	0.45
1:A:245:PRO:HG3	1:B:227:PHE:CE2	2.52	0.45
1:F:176:THR:H	1:F:179:HIS:CD2	2.34	0.44
1:B:533:GLN:HB3	1:B:564:CYS:HB2	1.99	0.44
1:C:171:ARG:NH2	1:F:295:ILE:HD12	2.32	0.44
1:B:144:ASN:HA	1:B:145:PRO:HD3	1.77	0.44
1:B:348:ILE:HD12	1:B:383:PRO:HD3	1.99	0.44
1:D:436:THR:HA	1:D:463:ARG:O	2.18	0.44
1:B:380:GLN:HG2	1:B:412:LEU:HD21	2.00	0.44
1:D:229:ASP:HB2	1:D:264:CYS:SG	2.57	0.44
1:C:348:ILE:HD12	1:C:383:PRO:HD3	1.98	0.44
1:E:202:SER:HB3	1:F:201:VAL:HB	1.98	0.44
1:F:64:VAL:O	1:F:94:ILE:HA	2.18	0.44
1:E:508:ILE:HD12	1:E:510:SER:O	2.18	0.44
1:E:46:ASN:HA	4:E:626:HOH:O	2.16	0.44
1:D:144:ASN:HD22	1:D:148:SER:HB3	1.83	0.44
1:B:148:SER:C	1:B:150:LEU:H	2.21	0.44
1:F:451:THR:HG22	3:F:1004:NAG:N2	2.32	0.44
1:A:300:PRO:HG2	1:B:300:PRO:HG2	2.00	0.44
2:J:6:VAL:HG13	2:J:26:LYS:HE2	1.99	0.44
1:F:436:THR:HA	1:F:463:ARG:O	2.18	0.43
1:E:348:ILE:HD12	1:E:383:PRO:HD3	1.99	0.43
1:E:331:PHE:HA	1:E:334:CYS:SG	2.58	0.43
1:A:137:HIS:CD2	1:A:140:ASP:HB2	2.48	0.43
1:D:549:PRO:HG2	1:D:563:ASN:HB3	1.99	0.43
1:E:421:GLN:HE21	1:E:490:PRO:HD2	1.83	0.43
1:C:148:SER:C	1:C:150:LEU:H	2.21	0.43
1:F:341:LEU:HA	1:F:341:LEU:HD23	1.94	0.43
1:B:138:TRP:HE1	1:B:153:VAL:HG21	1.82	0.43
1:A:436:THR:HA	1:A:463:ARG:O	2.18	0.43
1:E:420:LEU:HG	1:E:441:LEU:HD11	2.01	0.43
1:D:95:PHE:HA	1:D:125:ASP:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:646:HOH:O	2:J:52:MET:HE3	2.19	0.43
1:D:383:PRO:HA	1:D:384:PRO:HD3	1.89	0.43
2:H:6:VAL:HG13	2:H:26:LYS:HE2	1.99	0.43
2:I:8:CYS:HB3	2:I:12:GLU:HB2	2.01	0.43
1:C:428:ILE:HD13	1:C:433:ILE:HD11	2.01	0.43
1:E:454:SER:H	1:E:458:GLN:NE2	2.16	0.43
1:E:428:ILE:HD13	1:E:433:ILE:HD11	2.00	0.43
1:F:148:SER:C	1:F:150:LEU:H	2.23	0.43
1:D:346:THR:HG21	2:J:17:VAL:HG11	2.01	0.42
1:C:436:THR:HA	1:C:463:ARG:O	2.18	0.42
1:F:3:VAL:HG13	1:F:35:MET:HG2	1.98	0.42
1:E:120:GLY:H	1:E:144:ASN:ND2	2.17	0.42
1:A:229:ASP:HB2	1:A:264:CYS:SG	2.59	0.42
1:F:508:ILE:HD11	1:F:527:CYS:SG	2.59	0.42
1:F:549:PRO:HB2	1:F:563:ASN:HB2	2.01	0.42
1:D:206:HIS:CD2	1:D:208:GLU:HB2	2.54	0.42
1:B:436:THR:HA	1:B:463:ARG:O	2.19	0.42
1:E:148:SER:C	1:E:150:LEU:H	2.22	0.42
1:D:433:ILE:HB	1:D:460:ILE:HD12	2.01	0.42
1:A:198:GLY:HA3	1:A:203:ASP:OD2	2.20	0.42
1:B:103:ASN:HA	4:B:626:HOH:O	2.19	0.42
1:C:109:LEU:HG	1:C:130:LEU:HD11	2.01	0.42
1:D:127:ASN:O	1:D:129:PHE:N	2.49	0.42
1:B:273:VAL:HG13	1:B:280:VAL:HG23	2.00	0.42
1:E:436:THR:HA	1:E:463:ARG:O	2.20	0.42
1:C:433:ILE:HB	1:C:460:ILE:HD12	2.00	0.42
1:B:64:VAL:O	1:B:94:ILE:HA	2.20	0.42
1:C:581:LYS:HG2	1:C:591:PRO:HA	2.00	0.42
1:D:451:THR:HG22	3:D:1005:NAG:C7	2.49	0.42
1:A:481:LEU:HD21	1:A:509:GLU:HG2	2.02	0.42
1:C:138:TRP:O	1:C:142:VAL:HG23	2.19	0.42
1:C:229:ASP:HB2	1:C:264:CYS:SG	2.60	0.42
1:D:143:ARG:O	1:D:145:PRO:CD	2.55	0.41
1:C:171:ARG:HH21	1:F:295:ILE:HD12	1.85	0.41
1:C:383:PRO:HA	1:C:384:PRO:HD3	1.91	0.41
1:B:244:ASN:OD1	1:B:247:THR:HG23	2.20	0.41
1:E:64:VAL:O	1:E:94:ILE:HA	2.20	0.41
1:D:295:ILE:HD11	1:E:171:ARG:HH21	1.85	0.41
1:E:503:ARG:HD3	1:E:520:GLU:OE1	2.21	0.41
2:J:8:CYS:HB3	2:J:12:GLU:HB2	2.01	0.41
1:D:399:GLY:O	1:D:429:SER:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:32:GLU:HG2	1:E:56:ARG:HE	1.85	0.41
1:C:64:VAL:O	1:C:94:ILE:HA	2.21	0.41
1:C:46:ASN:HA	4:C:619:HOH:O	2.20	0.41
1:A:433:ILE:HB	1:A:460:ILE:HD12	2.03	0.41
2:K:8:CYS:HB3	2:K:12:GLU:HB2	2.03	0.41
1:D:501:PHE:HD2	1:D:525:SER:HA	1.84	0.41
2:H:8:CYS:HB3	2:H:12:GLU:HB2	2.02	0.41
1:A:428:ILE:HD13	1:A:433:ILE:HD11	2.03	0.41
1:A:410:LEU:HD12	1:A:412:LEU:HD22	2.02	0.40
1:D:87:GLU:OE2	2:J:33:ARG:HD3	2.21	0.40
1:E:423:GLN:HE21	1:E:493:ASP:HB2	1.86	0.40
1:D:581:LYS:HG2	1:D:591:PRO:HA	2.03	0.40
1:D:341:LEU:HD23	1:D:341:LEU:HA	1.91	0.40
1:D:64:VAL:O	1:D:94:ILE:HA	2.21	0.40
1:B:165:HIS:CD2	1:B:167:SER:H	2.35	0.40
1:B:433:ILE:HB	1:B:460:ILE:HD12	2.02	0.40
1:F:380:GLN:HG2	1:F:412:LEU:HD21	2.03	0.40
1:B:79:ILE:HA	1:B:116:GLU:O	2.21	0.40
1:A:278:SER:HB3	1:B:243:TYR:CD1	2.56	0.40
1:F:229:ASP:HB2	1:F:264:CYS:SG	2.61	0.40
1:F:348:ILE:HD12	1:F:383:PRO:HD3	2.02	0.40
1:B:503:ARG:NH1	1:B:520:GLU:OE2	2.55	0.40
1:A:520:GLU:HG2	1:A:529:GLU:HA	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	544/615 (88%)	508 (93%)	31 (6%)	5 (1%)	21 61
1	B	593/615 (96%)	558 (94%)	33 (6%)	2 (0%)	46 82

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	590/615 (96%)	548 (93%)	37 (6%)	5 (1%)	24	64
1	D	592/615 (96%)	553 (93%)	35 (6%)	4 (1%)	26	67
1	E	585/615 (95%)	544 (93%)	38 (6%)	3 (0%)	34	74
1	F	583/615 (95%)	546 (94%)	32 (6%)	5 (1%)	21	61
2	G	46/55 (84%)	43 (94%)	2 (4%)	1 (2%)	8	35
2	H	47/55 (86%)	43 (92%)	3 (6%)	1 (2%)	9	37
2	I	46/55 (84%)	43 (94%)	2 (4%)	1 (2%)	8	35
2	J	47/55 (86%)	44 (94%)	2 (4%)	1 (2%)	9	37
2	K	45/55 (82%)	41 (91%)	3 (7%)	1 (2%)	8	35
2	L	47/55 (86%)	44 (94%)	2 (4%)	1 (2%)	9	37
All	All	3765/4020 (94%)	3515 (93%)	220 (6%)	30 (1%)	24	64

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	ASN
1	B	149	ASN
1	C	149	ASN
1	C	159	SER
1	D	128	LYS
1	D	144	ASN
1	D	149	ASN
1	E	128	LYS
1	E	144	ASN
1	E	149	ASN
1	F	148	SER
1	F	149	ASN
1	A	128	LYS
1	F	128	LYS
1	A	423	GLN
1	A	536	LYS
1	B	536	LYS
1	C	128	LYS
1	D	536	LYS
2	G	40	ASN
2	H	40	ASN
2	I	40	ASN
2	J	40	ASN

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Mol	Chain	Res	Type
2	K	40	ASN
2	L	40	ASN
1	C	536	LYS
1	F	536	LYS
1	A	524	ASP
1	F	524	ASP
1	C	144	ASN

### 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	488/544 (90%)	464 (95%)	24 (5%)	31	69
1	B	534/544 (98%)	513 (96%)	21 (4%)	39	76
1	C	531/544 (98%)	512 (96%)	19 (4%)	42	78
1	D	534/544 (98%)	516 (97%)	18 (3%)	44	80
1	E	527/544 (97%)	502 (95%)	25 (5%)	32	70
1	F	525/544 (96%)	506 (96%)	19 (4%)	42	78
2	G	45/49 (92%)	44 (98%)	1 (2%)	60	87
2	H	46/49 (94%)	45 (98%)	1 (2%)	60	87
2	I	45/49 (92%)	43 (96%)	2 (4%)	35	73
2	J	46/49 (94%)	45 (98%)	1 (2%)	60	87
2	K	44/49 (90%)	43 (98%)	1 (2%)	58	86
2	L	46/49 (94%)	45 (98%)	1 (2%)	60	87
All	All	3411/3558 (96%)	3278 (96%)	133 (4%)	39	76

All (133) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	29	GLU
1	A	35	MET

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Mol	Chain	Res	Type
1	A	46	ASN
1	A	68	GLN
1	A	70	ARG
1	A	101	ASP
1	A	104	PHE
1	A	143	ARG
1	A	149	ASN
1	A	191	GLN
1	A	201	VAL
1	A	202	SER
1	A	252	HIS
1	A	259	THR
1	A	273	VAL
1	A	346	THR
1	A	412	LEU
1	A	442	CYS
1	A	450	THR
1	A	455	THR
1	A	460	ILE
1	A	506	ILE
1	A	556	SER
1	B	1	GLN
1	B	29	GLU
1	B	35	MET
1	B	46	ASN
1	B	101	ASP
1	B	104	PHE
1	B	201	VAL
1	B	202	SER
1	B	203	ASP
1	B	252	HIS
1	B	259	THR
1	B	346	THR
1	B	412	LEU
1	B	442	CYS
1	B	450	THR
1	B	460	ILE
1	B	497	SER
1	B	515	ASP
1	B	563	ASN
1	B	572	LEU
1	B	584	ASP

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Mol	Chain	Res	Type
1	C	1	GLN
1	C	29	GLU
1	C	35	MET
1	C	46	ASN
1	C	68	GLN
1	C	104	PHE
1	C	143	ARG
1	C	201	VAL
1	C	202	SER
1	C	252	HIS
1	C	259	THR
1	C	346	THR
1	C	412	LEU
1	C	429	SER
1	C	442	CYS
1	C	450	THR
1	C	460	ILE
1	C	511	CYS
1	C	563	ASN
1	D	1	GLN
1	D	29	GLU
1	D	35	MET
1	D	46	ASN
1	D	101	ASP
1	D	104	PHE
1	D	139	GLN
1	D	202	SER
1	D	252	HIS
1	D	259	THR
1	D	412	LEU
1	D	442	CYS
1	D	450	THR
1	D	460	ILE
1	D	470	ASN
1	D	515	ASP
1	D	563	ASN
1	D	613	TRP
1	E	1	GLN
1	E	29	GLU
1	E	35	MET
1	E	46	ASN
1	E	101	ASP

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Mol	Chain	Res	Type
1	E	104	PHE
1	E	201	VAL
1	E	202	SER
1	E	203	ASP
1	E	246	THR
1	E	259	THR
1	E	362	GLU
1	E	369	THR
1	E	410	LEU
1	E	412	LEU
1	E	429	SER
1	E	442	CYS
1	E	455	THR
1	E	460	ILE
1	E	482	CYS
1	E	496	LEU
1	E	518	PHE
1	E	563	ASN
1	E	570	ASP
1	E	584	ASP
1	F	1	GLN
1	F	29	GLU
1	F	46	ASN
1	F	101	ASP
1	F	104	PHE
1	F	143	ARG
1	F	154	SER
1	F	202	SER
1	F	203	ASP
1	F	252	HIS
1	F	346	THR
1	F	412	LEU
1	F	442	CYS
1	F	450	THR
1	F	460	ILE
1	F	493	ASP
1	F	541	LEU
1	F	550	ASP
1	F	584	ASP
2	G	43	THR
2	H	43	THR
2	I	32	SER

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Mol	Chain	Res	Type
2	I	43	THR
2	J	43	THR
2	K	43	THR
2	L	43	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (54) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	GLN
1	A	126	GLN
1	A	127	ASN
1	A	137	HIS
1	A	165	HIS
1	A	179	HIS
1	A	206	HIS
1	A	458	GLN
1	B	68	GLN
1	B	126	GLN
1	B	127	ASN
1	B	137	HIS
1	B	165	HIS
1	B	179	HIS
1	B	206	HIS
1	B	458	GLN
1	C	126	GLN
1	C	127	ASN
1	C	137	HIS
1	C	165	HIS
1	C	179	HIS
1	C	206	HIS
1	C	458	GLN
1	D	68	GLN
1	D	119	ASN
1	D	126	GLN
1	D	137	HIS
1	D	144	ASN
1	D	165	HIS
1	D	179	HIS
1	D	206	HIS
1	D	458	GLN
1	D	606	HIS
1	E	68	GLN

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Mol	Chain	Res	Type
1	E	126	GLN
1	E	137	HIS
1	E	144	ASN
1	E	165	HIS
1	E	179	HIS
1	E	206	HIS
1	E	421	GLN
1	E	458	GLN
1	E	606	HIS
1	F	68	GLN
1	F	126	GLN
1	F	127	ASN
1	F	137	HIS
1	F	165	HIS
1	F	179	HIS
1	F	206	HIS
1	F	458	GLN
1	F	533	GLN
2	I	18	ASN
2	K	18	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

26 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	1001	1	14,14,15	1.20	1 (7%)	15,19,21	2.72	8 (53%)
3	NAG	A	1002	1	14,14,15	1.55	1 (7%)	15,19,21	2.57	6 (40%)
3	NAG	A	1003	1	14,14,15	1.20	1 (7%)	15,19,21	1.72	4 (26%)
3	NAG	A	1004	1	14,14,15	2.19	3 (21%)	15,19,21	4.44	10 (66%)
3	NAG	B	1001	1	14,14,15	1.27	1 (7%)	15,19,21	3.06	7 (46%)
3	NAG	B	1002	1	14,14,15	1.49	1 (7%)	15,19,21	2.42	6 (40%)
3	NAG	B	1003	1	14,14,15	1.61	2 (14%)	15,19,21	2.26	7 (46%)
3	NAG	B	1004	1	14,14,15	3.51	8 (57%)	15,19,21	3.99	7 (46%)
3	NAG	B	1005	1	14,14,15	2.65	5 (35%)	15,19,21	2.66	7 (46%)
3	NAG	B	1006	1	14,14,15	1.81	4 (28%)	15,19,21	2.28	4 (26%)
3	NAG	C	1001	1	14,14,15	1.15	0	15,19,21	3.19	7 (46%)
3	NAG	C	1002	1	14,14,15	1.56	2 (14%)	15,19,21	1.99	6 (40%)
3	NAG	D	1001	1	14,14,15	1.46	3 (21%)	15,19,21	2.81	6 (40%)
3	NAG	D	1002	1	14,14,15	1.66	4 (28%)	15,19,21	2.18	4 (26%)
3	NAG	D	1003	1	14,14,15	1.20	1 (7%)	15,19,21	2.16	5 (33%)
3	NAG	D	1004	1	14,14,15	1.64	3 (21%)	15,19,21	2.20	3 (20%)
3	NAG	D	1005	1	14,14,15	1.39	2 (14%)	15,19,21	3.33	8 (53%)
3	NAG	E	1001	1	14,14,15	1.27	2 (14%)	15,19,21	3.01	8 (53%)
3	NAG	E	1002	1	14,14,15	1.24	1 (7%)	15,19,21	2.11	5 (33%)
3	NAG	E	1003	1	14,14,15	2.22	4 (28%)	15,19,21	2.12	5 (33%)
3	NAG	F	1001	1	14,14,15	1.60	3 (21%)	15,19,21	3.35	9 (60%)
3	NAG	F	1002	1	14,14,15	1.21	1 (7%)	15,19,21	2.42	6 (40%)
3	NAG	F	1003	1	14,14,15	1.05	1 (7%)	15,19,21	2.00	4 (26%)
3	NAG	F	1004	1	14,14,15	2.22	4 (28%)	15,19,21	2.24	6 (40%)
3	NAG	F	1005	1	14,14,15	1.73	3 (21%)	15,19,21	2.02	5 (33%)
3	NAG	F	1006	1	14,14,15	2.15	4 (28%)	15,19,21	1.09	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
3	NAG	A	1001	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1002	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1003	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1004	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1001	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1002	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1003	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1004	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1005	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1006	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1001	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1002	1	-	0/6/23/26	0/1/1/1
3	NAG	D	1001	1	-	0/6/23/26	0/1/1/1
3	NAG	D	1002	1	-	0/6/23/26	0/1/1/1
3	NAG	D	1003	1	-	0/6/23/26	0/1/1/1
3	NAG	D	1004	1	-	0/6/23/26	0/1/1/1
3	NAG	D	1005	1	-	0/6/23/26	0/1/1/1
3	NAG	E	1001	1	-	0/6/23/26	0/1/1/1
3	NAG	E	1002	1	-	0/6/23/26	0/1/1/1
3	NAG	E	1003	1	-	0/6/23/26	0/1/1/1
3	NAG	F	1001	1	-	0/6/23/26	0/1/1/1
3	NAG	F	1002	1	-	0/6/23/26	0/1/1/1
3	NAG	F	1003	1	-	0/6/23/26	0/1/1/1
3	NAG	F	1004	1	-	0/6/23/26	0/1/1/1
3	NAG	F	1005	1	-	0/6/23/26	0/1/1/1
3	NAG	F	1006	1	-	0/6/23/26	0/1/1/1

All (65) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	1001	NAG	C2-N2	-3.73	1.39	1.46
3	B	1004	NAG	O5-C1	-3.15	1.38	1.43
3	E	1001	NAG	C2-N2	-2.76	1.41	1.46
3	D	1001	NAG	C2-N2	-2.55	1.41	1.46
3	F	1004	NAG	O5-C5	-2.47	1.38	1.43
3	B	1004	NAG	O5-C5	-2.45	1.38	1.43
3	F	1004	NAG	O5-C1	-2.42	1.39	1.43
3	B	1005	NAG	C8-C7	2.00	1.54	1.50
3	F	1003	NAG	C4-C3	2.01	1.57	1.52
3	A	1001	NAG	O3-C3	2.05	1.47	1.43
3	E	1003	NAG	C2-N2	2.06	1.50	1.46
3	F	1004	NAG	C7-N2	2.08	1.42	1.34
3	D	1001	NAG	O5-C5	2.11	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1005	NAG	C3-C2	2.12	1.57	1.52
3	F	1002	NAG	O5-C1	2.14	1.47	1.43
3	F	1006	NAG	O5-C5	2.15	1.48	1.43
3	D	1004	NAG	O5-C5	2.15	1.48	1.43
3	F	1005	NAG	C4-C3	2.15	1.58	1.52
3	A	1004	NAG	C2-N2	2.19	1.50	1.46
3	B	1003	NAG	C3-C2	2.20	1.57	1.52
3	D	1002	NAG	O5-C1	2.20	1.47	1.43
3	E	1001	NAG	O3-C3	2.24	1.48	1.43
3	C	1002	NAG	O5-C1	2.25	1.47	1.43
3	D	1005	NAG	C6-C5	2.26	1.60	1.51
3	D	1002	NAG	C3-C2	2.26	1.57	1.52
3	B	1001	NAG	O5-C5	2.27	1.48	1.43
3	D	1002	NAG	C4-C3	2.31	1.58	1.52
3	B	1006	NAG	C4-C5	2.34	1.58	1.53
3	F	1001	NAG	C4-C3	2.40	1.58	1.52
3	E	1002	NAG	C1-C2	2.44	1.55	1.52
3	F	1001	NAG	C8-C7	2.49	1.55	1.50
3	A	1003	NAG	C1-C2	2.51	1.55	1.52
3	B	1006	NAG	C3-C2	2.52	1.58	1.52
3	A	1004	NAG	C3-C2	2.53	1.58	1.52
3	B	1002	NAG	C1-C2	2.54	1.56	1.52
3	D	1004	NAG	C4-C3	2.58	1.59	1.52
3	B	1006	NAG	C4-C3	2.75	1.59	1.52
3	B	1004	NAG	C4-C5	2.79	1.59	1.53
3	B	1004	NAG	O7-C7	2.80	1.29	1.23
3	E	1003	NAG	C4-C3	2.81	1.59	1.52
3	D	1001	NAG	O3-C3	2.81	1.49	1.43
3	F	1005	NAG	C4-C5	2.83	1.59	1.53
3	F	1006	NAG	C4-C5	2.95	1.59	1.53
3	D	1003	NAG	C1-C2	2.96	1.56	1.52
3	A	1002	NAG	C1-C2	2.99	1.56	1.52
3	B	1005	NAG	C4-C3	3.08	1.60	1.52
3	D	1005	NAG	C1-C2	3.18	1.56	1.52
3	B	1006	NAG	C1-C2	3.30	1.57	1.52
3	D	1004	NAG	C1-C2	3.35	1.57	1.52
3	F	1006	NAG	C3-C2	3.36	1.60	1.52
3	C	1002	NAG	C1-C2	3.38	1.57	1.52
3	B	1004	NAG	C8-C7	3.53	1.57	1.50
3	F	1005	NAG	C1-C2	3.86	1.57	1.52
3	D	1002	NAG	C1-C2	4.03	1.58	1.52
3	B	1004	NAG	C2-N2	4.07	1.53	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	1003	NAG	C3-C2	4.08	1.61	1.52
3	B	1003	NAG	C1-C2	4.26	1.58	1.52
3	B	1004	NAG	C7-N2	4.34	1.51	1.34
3	F	1006	NAG	C1-C2	4.97	1.59	1.52
3	B	1005	NAG	C4-C5	4.99	1.63	1.53
3	E	1003	NAG	C1-C2	5.36	1.59	1.52
3	F	1004	NAG	C1-C2	6.02	1.60	1.52
3	B	1005	NAG	C1-C2	6.22	1.61	1.52
3	A	1004	NAG	C1-C2	6.42	1.61	1.52
3	B	1004	NAG	C1-C2	9.29	1.65	1.52

All (153) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1004	NAG	C1-O5-C5	-9.77	99.85	112.25
3	D	1005	NAG	C3-C2-N2	-8.04	91.31	110.56
3	A	1004	NAG	O4-C4-C3	-6.84	94.95	110.34
3	D	1001	NAG	C3-C2-N2	-6.78	94.33	110.56
3	C	1001	NAG	C3-C2-N2	-6.32	95.41	110.56
3	B	1001	NAG	C3-C2-N2	-6.21	95.69	110.56
3	F	1001	NAG	C3-C2-N2	-6.01	96.17	110.56
3	A	1004	NAG	C3-C2-N2	-5.94	96.33	110.56
3	E	1001	NAG	C3-C2-N2	-5.91	96.41	110.56
3	A	1001	NAG	C3-C2-N2	-5.85	96.55	110.56
3	B	1005	NAG	C1-O5-C5	-4.40	106.66	112.25
3	B	1003	NAG	C3-C4-C5	-4.22	102.85	110.20
3	F	1003	NAG	C8-C7-N2	-3.93	108.59	116.11
3	B	1004	NAG	O7-C7-C8	-3.92	114.87	122.06
3	A	1002	NAG	O3-C3-C2	-3.88	101.43	109.11
3	B	1004	NAG	O5-C5-C6	-3.86	98.99	107.35
3	D	1002	NAG	O3-C3-C2	-3.75	101.69	109.11
3	D	1003	NAG	C8-C7-N2	-3.71	109.00	116.11
3	D	1005	NAG	C3-C4-C5	-3.68	103.78	110.20
3	B	1003	NAG	C1-O5-C5	-3.46	107.86	112.25
3	F	1004	NAG	O5-C5-C6	-3.41	99.97	107.35
3	D	1005	NAG	C1-O5-C5	-3.33	108.02	112.25
3	A	1003	NAG	C8-C7-N2	-3.30	109.79	116.11
3	B	1002	NAG	C8-C7-N2	-3.26	109.87	116.11
3	B	1005	NAG	O5-C5-C6	-3.19	100.44	107.35
3	B	1002	NAG	C2-N2-C7	-3.17	118.97	123.04
3	B	1004	NAG	O4-C4-C3	-3.10	103.35	110.34
3	F	1001	NAG	O4-C4-C5	-3.04	101.18	109.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1001	NAG	C8-C7-N2	-2.94	110.48	116.11
3	F	1005	NAG	O5-C5-C6	-2.61	101.69	107.35
3	C	1002	NAG	O3-C3-C2	-2.59	103.98	109.11
3	F	1005	NAG	O4-C4-C3	-2.57	104.55	110.34
3	E	1002	NAG	O3-C3-C2	-2.57	104.02	109.11
3	F	1001	NAG	C8-C7-N2	-2.57	111.19	116.11
3	D	1002	NAG	C2-N2-C7	-2.53	119.79	123.04
3	A	1001	NAG	O4-C4-C5	-2.52	102.57	109.24
3	C	1001	NAG	C8-C7-N2	-2.52	111.29	116.11
3	C	1001	NAG	O4-C4-C5	-2.51	102.59	109.24
3	A	1004	NAG	C8-C7-N2	-2.49	111.34	116.11
3	E	1002	NAG	O3-C3-C4	-2.47	104.77	110.34
3	A	1002	NAG	C8-C7-N2	-2.44	111.44	116.11
3	E	1001	NAG	C6-C5-C4	-2.42	107.04	113.02
3	A	1001	NAG	C8-C7-N2	-2.41	111.49	116.11
3	A	1003	NAG	C3-C2-N2	-2.41	104.79	110.56
3	A	1002	NAG	C2-N2-C7	-2.41	119.94	123.04
3	B	1005	NAG	C8-C7-N2	-2.37	111.56	116.11
3	C	1002	NAG	O3-C3-C4	-2.16	105.47	110.34
3	F	1002	NAG	O5-C5-C6	-2.14	102.71	107.35
3	E	1001	NAG	C8-C7-N2	-2.14	112.01	116.11
3	C	1002	NAG	C8-C7-N2	-2.12	112.06	116.11
3	F	1002	NAG	C8-C7-N2	-2.10	112.09	116.11
3	B	1006	NAG	O4-C4-C5	-2.02	103.89	109.24
3	F	1001	NAG	C6-C5-C4	-2.02	108.04	113.02
3	A	1004	NAG	O7-C7-C8	2.01	125.75	122.06
3	F	1003	NAG	O5-C5-C6	2.03	111.74	107.35
3	E	1003	NAG	C6-C5-C4	2.04	118.04	113.02
3	E	1003	NAG	O7-C7-N2	2.05	126.04	121.86
3	A	1003	NAG	C2-N2-C7	2.08	125.71	123.04
3	D	1004	NAG	C2-N2-C7	2.14	125.79	123.04
3	F	1005	NAG	O3-C3-C4	2.15	115.18	110.34
3	A	1001	NAG	O5-C5-C6	2.15	112.01	107.35
3	B	1005	NAG	C3-C4-C5	2.16	113.97	110.20
3	F	1002	NAG	O7-C7-N2	2.17	126.30	121.86
3	D	1002	NAG	C3-C2-N2	2.19	115.81	110.56
3	D	1005	NAG	O3-C3-C4	2.21	115.31	110.34
3	A	1001	NAG	O3-C3-C4	2.24	115.39	110.34
3	D	1003	NAG	C2-N2-C7	2.30	126.00	123.04
3	D	1005	NAG	O7-C7-N2	2.33	126.61	121.86
3	B	1003	NAG	O3-C3-C2	2.36	113.78	109.11
3	F	1001	NAG	C3-C4-C5	2.38	114.35	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1002	NAG	C3-C2-N2	2.41	116.33	110.56
3	B	1002	NAG	C3-C2-N2	2.44	116.40	110.56
3	B	1001	NAG	C3-C4-C5	2.45	114.48	110.20
3	F	1002	NAG	C3-C2-N2	2.46	116.44	110.56
3	B	1003	NAG	C6-C5-C4	2.49	119.15	113.02
3	E	1002	NAG	C3-C2-N2	2.52	116.60	110.56
3	F	1004	NAG	O7-C7-N2	2.53	127.03	121.86
3	A	1002	NAG	C3-C2-N2	2.54	116.65	110.56
3	B	1003	NAG	O3-C3-C4	2.64	116.27	110.34
3	B	1002	NAG	O7-C7-N2	2.65	127.27	121.86
3	D	1003	NAG	O3-C3-C4	2.67	116.36	110.34
3	D	1001	NAG	O7-C7-C8	2.70	127.02	122.06
3	E	1001	NAG	O3-C3-C4	2.71	116.43	110.34
3	E	1001	NAG	C3-C4-C5	2.74	114.98	110.20
3	D	1003	NAG	O7-C7-C8	2.77	127.15	122.06
3	F	1001	NAG	O3-C3-C4	2.78	116.59	110.34
3	E	1003	NAG	C2-N2-C7	2.78	126.61	123.04
3	A	1001	NAG	O7-C7-C8	2.97	127.51	122.06
3	D	1001	NAG	C3-C4-C5	2.97	115.38	110.20
3	F	1004	NAG	C2-N2-C7	2.98	126.87	123.04
3	D	1001	NAG	O3-C3-C4	3.00	117.09	110.34
3	D	1005	NAG	O5-C5-C6	3.03	113.91	107.35
3	F	1003	NAG	O3-C3-C4	3.03	117.16	110.34
3	A	1001	NAG	C1-O5-C5	3.04	116.11	112.25
3	B	1005	NAG	C4-C3-C2	3.12	116.08	111.23
3	B	1006	NAG	C3-C4-C5	3.15	115.69	110.20
3	C	1002	NAG	C4-C3-C2	3.17	116.15	111.23
3	B	1002	NAG	C4-C3-C2	3.22	116.24	111.23
3	F	1004	NAG	C6-C5-C4	3.24	121.02	113.02
3	B	1001	NAG	O3-C3-C4	3.28	117.72	110.34
3	B	1003	NAG	O5-C5-C6	3.32	114.54	107.35
3	D	1005	NAG	C4-C3-C2	3.33	116.41	111.23
3	F	1004	NAG	O3-C3-C4	3.35	117.87	110.34
3	E	1002	NAG	C4-C3-C2	3.39	116.50	111.23
3	A	1003	NAG	C4-C3-C2	3.40	116.52	111.23
3	D	1004	NAG	O3-C3-C4	3.46	118.13	110.34
3	E	1001	NAG	O7-C7-C8	3.48	128.45	122.06
3	B	1006	NAG	C4-C3-C2	3.60	116.83	111.23
3	B	1003	NAG	C2-N2-C7	3.62	127.69	123.04
3	C	1001	NAG	O3-C3-C4	3.62	118.49	110.34
3	F	1004	NAG	C4-C3-C2	3.71	117.00	111.23
3	A	1004	NAG	C2-N2-C7	3.80	127.92	123.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	1002	NAG	C4-C3-C2	3.84	117.19	111.23
3	B	1004	NAG	C3-C4-C5	3.87	116.94	110.20
3	D	1001	NAG	C4-C3-C2	4.00	117.44	111.23
3	F	1005	NAG	C4-C3-C2	4.01	117.47	111.23
3	E	1003	NAG	O5-C5-C6	4.02	116.06	107.35
3	F	1003	NAG	C4-C3-C2	4.04	117.50	111.23
3	B	1004	NAG	C6-C5-C4	4.04	122.98	113.02
3	A	1004	NAG	C6-C5-C4	4.06	123.02	113.02
3	F	1001	NAG	O7-C7-C8	4.18	129.74	122.06
3	A	1002	NAG	C4-C3-C2	4.21	117.78	111.23
3	B	1004	NAG	O7-C7-N2	4.22	130.47	121.86
3	B	1001	NAG	C1-O5-C5	4.22	117.61	112.25
3	C	1001	NAG	O7-C7-C8	4.31	129.97	122.06
3	F	1005	NAG	C1-O5-C5	4.43	117.87	112.25
3	B	1001	NAG	O7-C7-C8	4.48	130.28	122.06
3	B	1005	NAG	C6-C5-C4	4.50	124.12	113.02
3	A	1004	NAG	C4-C3-C2	4.58	118.36	111.23
3	B	1005	NAG	C2-N2-C7	4.69	129.07	123.04
3	A	1004	NAG	O3-C3-C2	4.72	118.46	109.11
3	E	1001	NAG	C4-C3-C2	4.81	118.71	111.23
3	D	1003	NAG	C4-C3-C2	4.82	118.72	111.23
3	D	1001	NAG	C1-O5-C5	4.82	118.37	112.25
3	C	1002	NAG	C1-O5-C5	4.88	118.44	112.25
3	E	1003	NAG	C4-C3-C2	4.91	118.87	111.23
3	F	1001	NAG	C4-C3-C2	5.07	119.11	111.23
3	B	1001	NAG	C4-C3-C2	5.09	119.15	111.23
3	C	1001	NAG	C4-C3-C2	5.26	119.41	111.23
3	A	1001	NAG	C4-C3-C2	5.40	119.62	111.23
3	E	1002	NAG	C1-O5-C5	5.41	119.12	112.25
3	C	1001	NAG	C1-O5-C5	5.61	119.36	112.25
3	A	1004	NAG	O5-C5-C6	5.61	119.48	107.35
3	D	1002	NAG	C1-O5-C5	5.70	119.48	112.25
3	E	1001	NAG	C1-O5-C5	5.83	119.65	112.25
3	B	1006	NAG	C1-O5-C5	5.87	119.70	112.25
3	B	1002	NAG	C1-O5-C5	5.88	119.71	112.25
3	A	1002	NAG	C1-O5-C5	5.93	119.78	112.25
3	D	1005	NAG	C6-C5-C4	6.16	128.21	113.02
3	F	1002	NAG	C1-O5-C5	6.49	120.49	112.25
3	D	1004	NAG	C1-O5-C5	6.51	120.51	112.25
3	F	1001	NAG	C1-O5-C5	6.80	120.87	112.25
3	B	1004	NAG	C2-N2-C7	11.99	138.45	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1005	NAG	2	0
3	D	1004	NAG	2	0
3	D	1005	NAG	4	0
3	F	1004	NAG	2	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	550/615 (89%)	-0.10	25 (4%) 37 15	47, 85, 249, 280	0
1	B	601/615 (97%)	-0.31	0 100 100	45, 69, 115, 151	0
1	C	598/615 (97%)	-0.09	18 (3%) 54 25	50, 93, 173, 229	0
1	D	600/615 (97%)	-0.30	2 (0%) 94 84	45, 77, 117, 172	0
1	E	593/615 (96%)	-0.09	11 (1%) 70 41	45, 93, 164, 228	0
1	F	591/615 (96%)	-0.14	16 (2%) 58 28	43, 72, 155, 278	0
2	G	50/55 (90%)	-0.37	0 100 100	68, 98, 142, 151	0
2	H	51/55 (92%)	-0.21	0 100 100	60, 85, 130, 140	0
2	I	50/55 (90%)	-0.26	0 100 100	85, 104, 143, 147	0
2	J	51/55 (92%)	-0.25	0 100 100	75, 100, 136, 140	0
2	K	49/55 (89%)	-0.12	2 (4%) 41 17	87, 110, 143, 158	0
2	L	51/55 (92%)	-0.31	1 (1%) 68 39	62, 86, 126, 143	0
All	All	3835/4020 (95%)	-0.18	75 (1%) 68 39	43, 83, 156, 280	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	614	THR	6.4
1	F	610	TYR	6.1
1	A	545	HIS	4.7
1	A	547	PRO	4.5
1	A	513	LEU	4.3
1	F	606	HIS	4.2
1	F	605	SER	3.9
1	A	530	CYS	3.9
1	A	528	VAL	3.9
1	F	609	ILE	3.9
1	C	524	ASP	3.8

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Mol	Chain	Res	Type	RSRZ
1	F	604	THR	3.7
1	A	546	GLY	3.5
1	C	509	GLU	3.5
1	A	543	THR	3.4
1	F	589	CYS	3.4
1	F	600	CYS	3.4
1	F	567	LYS	3.3
1	C	547	PRO	3.3
1	F	569	PRO	3.2
1	C	307	LYS	3.2
1	E	583	ALA	3.1
1	E	597	THR	3.0
1	E	589	CYS	3.0
1	F	595	ASN	3.0
1	A	523	ASN	3.0
1	C	510	SER	3.0
1	E	605	SER	3.0
1	C	517	GLU	3.0
1	D	613	TRP	2.9
1	F	558	PHE	2.9
1	E	599	GLY	2.9
1	A	525	SER	2.9
1	C	587	ARG	2.8
1	A	152	LEU	2.8
1	C	525	SER	2.8
1	F	598	GLN	2.7
1	A	526	ILE	2.7
1	C	508	ILE	2.7
1	C	598	GLN	2.6
1	F	601	ASN	2.6
1	A	537	MET	2.6
1	A	544	CYS	2.6
1	A	522	GLU	2.6
1	C	601	ASN	2.5
1	C	152	LEU	2.5
1	C	518	PHE	2.5
1	A	524	ASP	2.5
1	C	500	ARG	2.5
1	C	570	ASP	2.5
1	F	302	THR	2.5
1	E	580	PHE	2.5
1	A	99	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	539	ASP	2.4
1	C	501	PHE	2.4
1	F	591	PRO	2.4
2	L	54	SER	2.3
1	A	554	LYS	2.3
1	E	601	ASN	2.2
1	F	599	GLY	2.2
1	A	551	ASN	2.2
1	A	521	PHE	2.2
1	E	600	CYS	2.2
1	A	548	GLY	2.2
1	E	536	LYS	2.1
1	C	560	ASP	2.1
1	E	541	LEU	2.1
2	K	26	LYS	2.1
1	A	148	SER	2.1
1	A	550	ASP	2.0
1	E	560	ASP	2.0
2	K	53	ALA	2.0
1	A	536	LYS	2.0
1	A	542	LEU	2.0
1	C	589	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. 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( $\text{\AA}^2$ )	Q<0.9
3	NAG	D	1005	14/15	0.83	0.34	6.68	108,111,116,118	0
3	NAG	E	1003	14/15	0.78	0.28	3.85	147,149,154,158	0
3	NAG	B	1004	14/15	0.86	0.27	3.77	89,93,97,97	0
3	NAG	D	1004	14/15	0.85	0.29	2.96	125,129,134,135	0
3	NAG	D	1002	14/15	0.93	0.23	2.12	73,77,80,81	0
3	NAG	A	1004	14/15	0.86	0.27	0.79	117,121,125,127	0
3	NAG	F	1002	14/15	0.93	0.20	0.72	70,74,77,78	0
3	NAG	C	1002	14/15	0.90	0.23	0.64	72,77,81,82	0
3	NAG	E	1002	14/15	0.93	0.18	0.10	77,82,85,87	0
3	NAG	F	1004	14/15	0.91	0.21	-0.26	77,81,84,85	0
3	NAG	A	1002	14/15	0.92	0.17	-0.36	71,76,78,80	0
3	NAG	B	1002	14/15	0.88	0.18	-0.42	81,86,88,89	0
3	NAG	F	1003	14/15	0.85	0.25	-	118,123,128,130	0
3	NAG	A	1003	14/15	0.84	0.29	-	145,151,158,160	0
3	NAG	B	1005	14/15	0.88	0.24	-	100,105,108,111	0
3	NAG	F	1006	14/15	0.79	0.22	-	127,129,135,136	0
3	NAG	D	1001	14/15	0.95	0.17	-	66,70,74,74	0
3	NAG	B	1006	14/15	0.87	0.26	-	118,121,125,125	0
3	NAG	F	1005	14/15	0.89	0.17	-	126,130,135,136	0
3	NAG	A	1001	14/15	0.95	0.18	-	69,72,77,78	0
3	NAG	B	1003	14/15	0.85	0.25	-	140,145,150,152	0
3	NAG	C	1001	14/15	0.95	0.13	-	72,75,80,81	0
3	NAG	D	1003	14/15	0.91	0.19	-	126,131,134,137	0
3	NAG	B	1001	14/15	0.97	0.15	-	60,64,67,68	0
3	NAG	F	1001	14/15	0.97	0.13	-	55,59,63,63	0
3	NAG	E	1001	14/15	0.95	0.17	-	66,70,74,74	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.