Full wwPDB X-ray Structure Validation Report (i)

Sep 26, 2023 – 06:20 PM EDT

PDB ID	:	6D7U
Title	:	The crystal structure of hemagglutinin from A/Guangdong/17SF003/2016 H7N9 influenza virus
Authors	:	Yang, H.; Stevens, J.
Deposited on	:	2018-04-25
Resolution	:	2.70 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

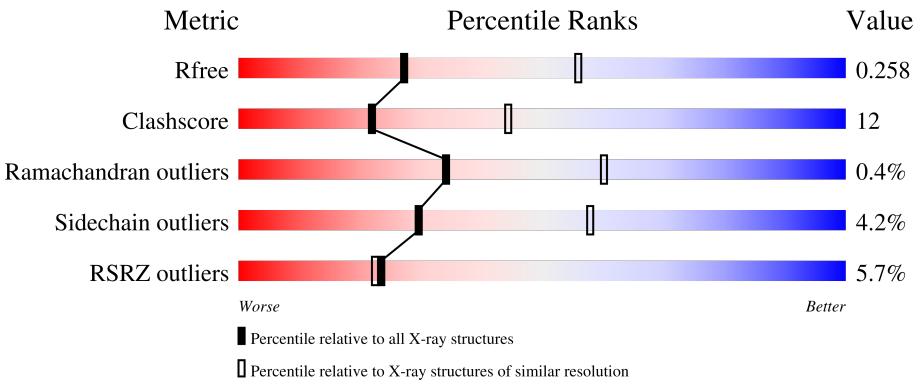
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	316	5% 75% 24% •
1	C	316	4% 76% 22% •
1	E	316	3% 75% 22% •
2	B	171	9% 75% 19% 6%
2	D	171	8% 66% 27% 5% •

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Mol	Chain	Length	Quality of chain
2	F	171	8% 71% 24% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 11559 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 HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	316	2423	1502	440	467	14	1	0	0
1	C	316	2423	1502	440	467	14	1	0	0
1	E	316	2423	1502	440	467	14	1	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	125	ALA	VAL	conflict	UNP A0A1S6R2B6
A	261	ARG	GLY	conflict	UNP A0A1S6R2B6
C	125	ALA	VAL	conflict	UNP A0A1S6R2B6
C	261	ARG	GLY	conflict	UNP A0A1S6R2B6
E	125	ALA	VAL	conflict	UNP A0A1S6R2B6
E	261	ARG	GLY	conflict	UNP A0A1S6R2B6

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	171	1388	859	246	276	7	0	0	0
2	D	171	1388	859	246	276	7	0	0	0
2	F	171	1388	859	246	276	7	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	2	ILE	LEU	conflict	UNP A0A2I7YV20
D	2	ILE	LEU	conflict	UNP A0A2I7YV20

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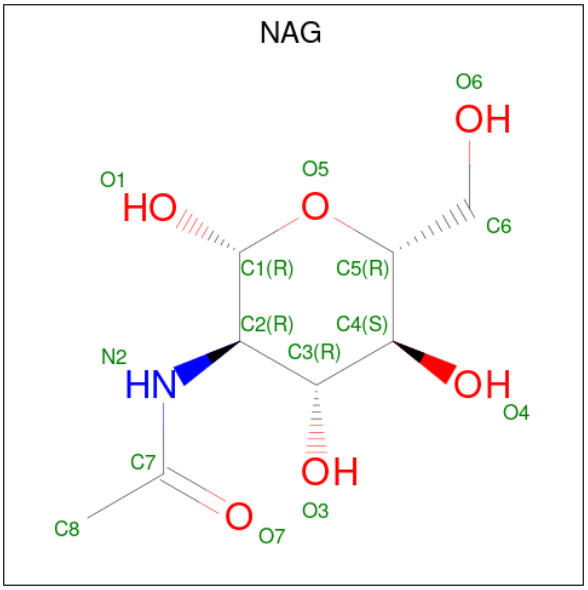
Chain	Residue	Modelled	Actual	Comment	Reference
F	2	ILE	LEU	conflict	UNP A0A2I7YV20

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



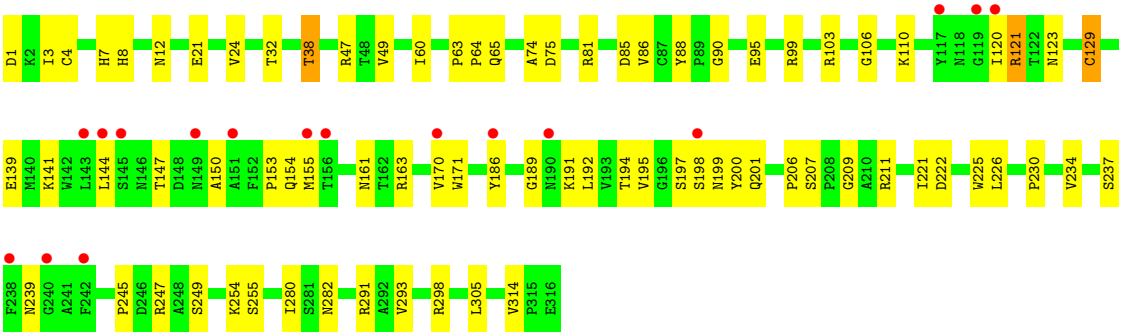
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	14	8	1	5	0	0
			Total	C	N	O		
3	A	1	14	8	1	5	0	0
			Total	C	N	O		
3	B	1	14	8	1	5	0	0
			Total	C	N	O		
3	C	1	14	8	1	5	0	0
			Total	C	N	O		
3	C	1	14	8	1	5	0	0
			Total	C	N	O		
3	D	1	14	8	1	5	0	0
			Total	C	N	O		
3	E	1	14	8	1	5	0	0
			Total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

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

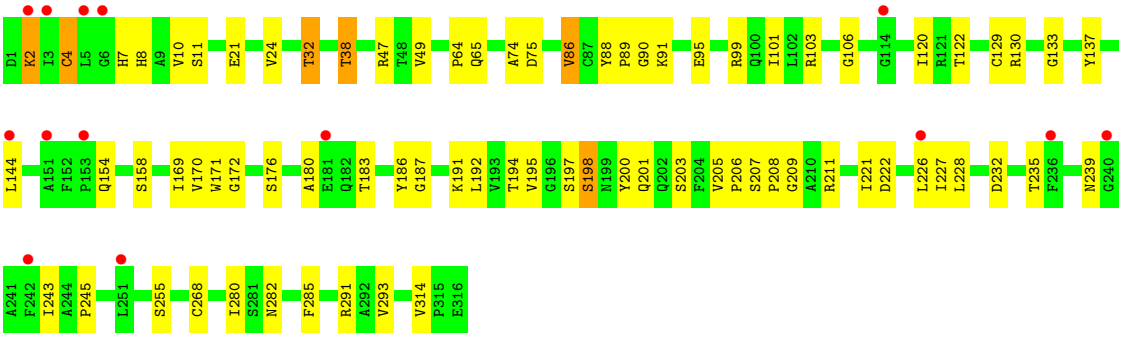
• Molecule 1: Hemagglutinin HA1 chain

Chain A:



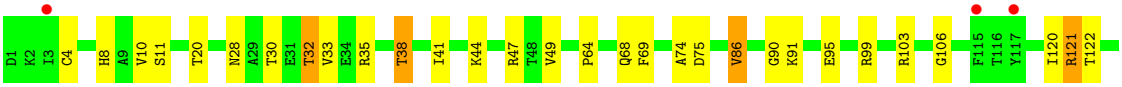
• Molecule 1: Hemagglutinin HA1 chain

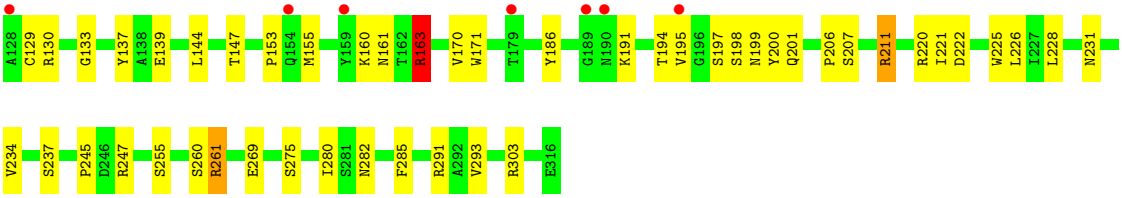
Chain C:



• Molecule 1: Hemagglutinin HA1 chain

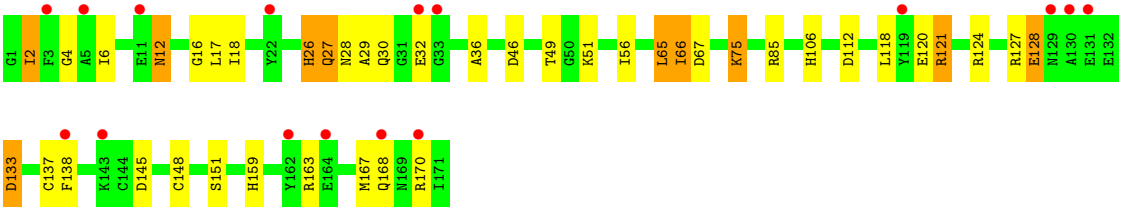
Chain E:





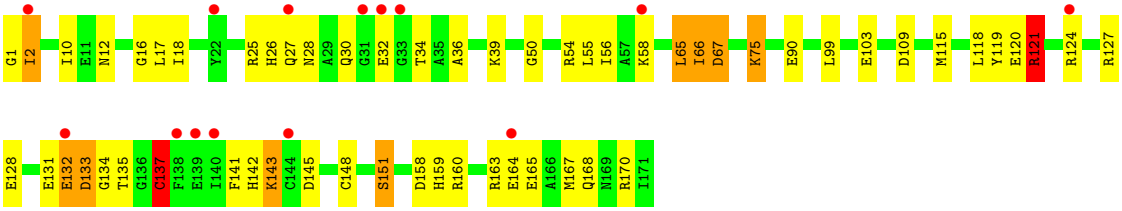
• Molecule 2: Hemagglutinin HA2 chain

Chain B:



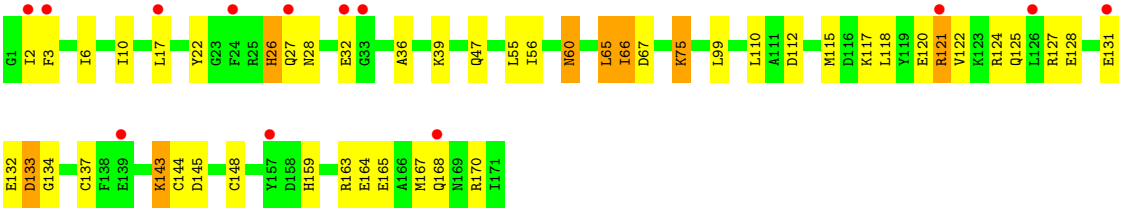
• Molecule 2: Hemagglutinin HA2 chain

Chain D:



• Molecule 2: Hemagglutinin HA2 chain

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	203.12Å 117.05Å 119.97Å 90.00° 124.46° 90.00°	Depositor
Resolution (Å)	38.00 – 2.70 38.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.7 (38.00-2.70) 98.1 (38.00-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.24 (at 2.69Å)	Xtriage
Refinement program	PHENIX (dev_2733: ???)	Depositor
R, R_{free}	0.221 , 0.258 0.221 , 0.258	Depositor DCC
R_{free} test set	3266 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å^2)	61.3	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 36.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.010 for -1/2*h-3/2*k-l,-1/2*h+1/2*k-l,1/2*h+1/2*k 0.003 for 1/2*h+1/2*k+2*l,1/2*h+1/2*k,-1/2*h+1/2*k-l 0.000 for -h-k-l,l,k 0.019 for -h+k-l,-l,-k 0.018 for 1/2*h-1/2*k+2*l,-1/2*h+1/2*k,-1/2*h-1/2*k-l 0.010 for -1/2*h+3/2*k-l,1/2*h+1/2*k+l,1/2*h-1/2*k 0.460 for 1/2*h-3/2*k,-1/2*h-1/2*k,-1/2*h+1/2*k-l 0.467 for 1/2*h+3/2*k,1/2*h-1/2*k,-1/2*h-1/2*k-l 0.005 for -1/2*h-1/2*k+l,-1/2*h-1/2*k-l,1/2*h-1/2*k 0.018 for -1/2*h+1/2*k+l,1/2*h-1/2*k+l,1/2*h+1/2*k 0.014 for -h-2*l,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11559	wwPDB-VP
Average B, all atoms (Å^2)	75.0	wwPDB-VP

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% of the height of the origin peak. No significant pseudotranslation is detected.*

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: 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	0.33	0/2470	0.61	0/3341
1	C	0.38	1/2470 (0.0%)	0.65	2/3341 (0.1%)
1	E	0.34	0/2470	0.71	8/3341 (0.2%)
2	B	0.34	0/1412	0.73	4/1901 (0.2%)
2	D	0.44	1/1412 (0.1%)	0.80	7/1901 (0.4%)
2	F	0.39	0/1412	0.75	5/1901 (0.3%)
All	All	0.37	2/11646 (0.0%)	0.70	26/15726 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
1	E	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	4	CYS	CB-SG	7.86	1.95	1.82
2	D	132	GLU	CB-CG	-5.15	1.42	1.52

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	163	ARG	NE-CZ-NH1	9.88	125.24	120.30
2	F	39	LYS	CA-CB-CG	9.77	134.90	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	163	ARG	NE-CZ-NH2	-9.68	115.46	120.30
2	D	121	ARG	CG-CD-NE	-8.40	94.16	111.80
2	D	137	CYS	CA-CB-SG	7.55	127.58	114.00
2	B	133	ASP	CB-CG-OD2	7.36	124.92	118.30
2	D	121	ARG	NE-CZ-NH2	-7.28	116.66	120.30
2	D	121	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	C	198	SER	CB-CA-C	6.90	123.20	110.10
1	E	91	LYS	CB-CG-CD	-6.77	94.00	111.60
2	F	144	CYS	CA-CB-SG	6.68	126.02	114.00
2	F	39	LYS	N-CA-CB	6.53	122.35	110.60
2	B	128	GLU	CA-CB-CG	6.52	127.75	113.40
1	C	4	CYS	CA-CB-SG	6.36	125.44	114.00
2	B	168	GLN	CA-CB-CG	6.23	127.11	113.40
2	F	39	LYS	CB-CA-C	-6.01	98.38	110.40
1	E	91	LYS	CD-CE-NZ	-5.95	98.02	111.70
2	D	143	LYS	CB-CG-CD	-5.91	96.25	111.60
2	D	121	ARG	CA-CB-CG	5.81	126.18	113.40
2	F	39	LYS	CD-CE-NZ	5.54	124.44	111.70
2	D	121	ARG	CB-CG-CD	5.41	125.66	111.60
2	B	128	GLU	CG-CD-OE1	5.30	128.90	118.30
1	E	261	ARG	NE-CZ-NH1	5.25	122.92	120.30
1	E	261	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	E	163	ARG	CD-NE-CZ	5.11	130.76	123.60
1	E	261	ARG	CG-CD-NE	5.05	122.41	111.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	85	ASP	Peptide
1	C	122	THR	Peptide
1	E	122	THR	Peptide

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	2423	0	2373	56	0
1	C	2423	0	2373	57	0
1	E	2423	0	2373	60	0
2	B	1388	0	1305	41	0
2	D	1388	0	1305	55	0
2	F	1388	0	1305	39	0
3	A	28	0	26	0	0
3	B	14	0	13	1	0
3	C	28	0	26	0	0
3	D	14	0	13	0	0
3	E	28	0	26	5	0
3	F	14	0	13	1	0
All	All	11559	0	11151	267	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 12.

All (267) 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:28:ASN:HD21	3:E:401:NAG:C1	1.31	1.40
2:B:124:ARG:HA	2:D:132:GLU:OE2	1.30	1.27
1:E:28:ASN:ND2	3:E:401:NAG:C1	2.00	1.25
1:E:121:ARG:NH2	1:E:147:THR:HA	1.56	1.20
1:C:2:LYS:NZ	2:D:137:CYS:SG	2.29	1.05
1:E:121:ARG:HH22	1:E:147:THR:HA	1.21	0.98
1:C:4:CYS:HA	2:D:137:CYS:HB3	1.47	0.97
2:B:133:ASP:OD2	2:B:137:CYS:HB2	1.68	0.92
1:A:194:THR:HG23	1:A:237:SER:HB3	1.54	0.89
2:D:75:LYS:H	2:D:75:LYS:HD2	1.37	0.88
2:B:75:LYS:H	2:B:75:LYS:HD2	1.39	0.85
1:A:209:GLY:HA3	1:E:194:THR:HG21	1.59	0.84
1:E:28:ASN:CG	3:E:401:NAG:C1	2.46	0.83
2:F:75:LYS:H	2:F:75:LYS:HD2	1.43	0.82
2:B:124:ARG:HB3	2:D:132:GLU:OE1	1.80	0.82
2:B:124:ARG:CA	2:D:132:GLU:OE2	2.23	0.82
1:E:121:ARG:HG3	1:E:144:LEU:HB2	1.60	0.82
2:F:164:GLU:O	2:F:168:GLN:OE1	1.98	0.81
2:D:1:GLY:O	2:D:2:ILE:HG22	1.82	0.80
2:B:67:ASP:OD2	2:B:85:ARG:NH2	2.13	0.78
1:E:28:ASN:OD1	3:E:401:NAG:C1	2.30	0.78
1:E:121:ARG:NH2	1:E:147:THR:CA	2.40	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:17:LEU:HD11	2:D:36:ALA:HB2	1.66	0.77
2:B:124:ARG:NH2	2:D:119:TYR:OH	2.18	0.77
1:E:194:THR:HG23	1:E:237:SER:HB2	1.65	0.77
1:A:121:ARG:HG2	1:A:123:ASN:OD1	1.85	0.76
1:A:194:THR:HG21	1:C:209:GLY:HA3	1.67	0.76
1:A:192:LEU:H	1:A:239:ASN:HD21	1.32	0.75
2:D:25:ARG:HG3	2:D:34:THR:HG22	1.69	0.75
2:B:133:ASP:OD2	2:B:137:CYS:CB	2.37	0.72
2:F:143:LYS:NZ	2:F:165:GLU:OE2	2.21	0.71
2:B:17:LEU:HD11	2:B:36:ALA:HB2	1.73	0.71
2:F:60:ASN:ND2	2:F:60:ASN:O	2.21	0.70
1:E:38:THR:HG21	1:E:280:ILE:HD12	1.73	0.70
1:C:201:GLN:HE21	1:E:222:ASP:HB3	1.57	0.69
2:D:124:ARG:HE	2:F:134:GLY:HA2	1.56	0.69
1:A:38:THR:HG21	1:A:280:ILE:HD12	1.75	0.68
2:F:122:VAL:HA	2:F:125:GLN:HE21	1.58	0.68
2:B:4:GLY:HA2	2:F:117:LYS:HD3	1.75	0.68
1:A:170:VAL:HG12	1:A:225:TRP:HB3	1.75	0.68
2:B:30:GLN:OE1	2:B:30:GLN:N	2.26	0.68
1:A:222:ASP:HB3	1:E:201:GLN:HE21	1.58	0.67
1:E:35:ARG:HD3	1:E:303:ARG:HG3	1.75	0.67
1:C:38:THR:HG21	1:C:280:ILE:HD12	1.77	0.67
1:C:4:CYS:CA	2:D:137:CYS:HB3	2.09	0.67
1:A:154:GLN:OE1	1:A:239:ASN:HB3	1.96	0.66
1:A:201:GLN:HE21	1:C:222:ASP:HB3	1.60	0.66
2:D:120:GLU:O	2:D:124:ARG:HG3	1.98	0.64
2:F:6:ILE:N	2:F:112:ASP:OD1	2.30	0.64
2:D:27:GLN:OE1	2:D:32:GLU:HB2	1.97	0.64
1:E:30:THR:HB	3:E:401:NAG:H62	1.80	0.64
1:E:293:VAL:HG11	2:F:65:LEU:HD13	1.80	0.64
1:A:121:ARG:HH12	1:A:144:LEU:HD12	1.62	0.63
1:A:121:ARG:NH1	1:A:144:LEU:HD12	2.14	0.63
1:A:293:VAL:HG11	2:B:65:LEU:HD13	1.80	0.63
1:C:192:LEU:HD11	1:C:194:THR:HG23	1.79	0.63
2:F:17:LEU:HD11	2:F:36:ALA:HB2	1.79	0.63
2:B:127:ARG:HG3	2:B:159:HIS:CG	2.34	0.62
2:D:168:GLN:OE1	2:D:168:GLN:N	2.26	0.62
1:E:68:GLN:HG2	1:E:69:PHE:CE2	2.34	0.62
1:A:291:ARG:HH21	2:B:67:ASP:HB3	1.65	0.62
1:A:314:VAL:HG13	2:B:12:ASN:HA	1.82	0.62
2:D:58:LYS:HG3	2:D:58:LYS:O	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:55:LEU:HD22	2:D:99:LEU:HD21	1.82	0.61
1:E:291:ARG:HH21	2:F:67:ASP:HB3	1.65	0.61
2:B:124:ARG:HA	2:D:132:GLU:CD	2.17	0.61
2:F:145:ASP:O	2:F:148:CYS:N	2.34	0.60
2:D:133:ASP:OD1	2:D:134:GLY:N	2.34	0.60
1:A:21:GLU:OE1	1:A:24:VAL:HB	2.00	0.60
1:A:195:VAL:HG13	1:A:200:TYR:HE1	1.67	0.60
2:F:75:LYS:HD2	2:F:75:LYS:N	2.15	0.60
3:F:201:NAG:O7	3:F:201:NAG:O3	2.20	0.60
2:F:55:LEU:HD22	2:F:99:LEU:HD21	1.82	0.59
1:A:192:LEU:H	1:A:239:ASN:ND2	2.01	0.59
2:F:122:VAL:HA	2:F:125:GLN:NE2	2.17	0.59
2:B:75:LYS:HD2	2:B:75:LYS:N	2.15	0.59
2:D:127:ARG:HG3	2:D:159:HIS:CG	2.37	0.58
1:A:7:HIS:HE1	1:A:314:VAL:HG12	1.67	0.58
1:A:161:ASN:HD21	1:A:230:PRO:HA	1.69	0.58
1:C:88:TYR:CD1	1:C:221:ILE:HD12	2.38	0.58
1:C:192:LEU:CD1	1:C:194:THR:HG23	2.34	0.57
2:B:128:GLU:O	2:B:128:GLU:HG2	2.04	0.57
2:F:163:ARG:O	2:F:167:MET:HG3	2.04	0.57
1:C:195:VAL:HG23	1:C:200:TYR:HE1	1.68	0.57
2:D:132:GLU:HG3	2:D:132:GLU:O	2.02	0.57
2:B:4:GLY:CA	2:F:117:LYS:HD3	2.34	0.57
2:D:10:ILE:HD12	2:D:115:MET:SD	2.45	0.57
1:C:291:ARG:NH1	2:D:67:ASP:HB3	2.19	0.57
1:E:160:LYS:HE2	1:E:231:ASN:HA	1.85	0.57
1:C:291:ARG:HH12	2:D:67:ASP:HB3	1.70	0.57
1:A:207:SER:OG	1:A:207:SER:O	2.22	0.56
1:C:7:HIS:HE1	1:C:314:VAL:HG12	1.68	0.56
1:C:198:SER:OG	1:C:232:ASP:OD2	2.23	0.56
1:A:60:ILE:HD12	1:A:170:VAL:HG11	1.86	0.56
1:A:121:ARG:HH12	1:A:144:LEU:CD1	2.19	0.56
1:C:171:TRP:CZ2	1:C:195:VAL:HG21	2.41	0.55
1:C:194:THR:HG22	1:C:203:SER:HB3	1.87	0.55
1:C:228:LEU:HD21	1:C:232:ASP:O	2.07	0.55
1:E:195:VAL:HG13	1:E:200:TYR:HE1	1.72	0.55
2:D:2:ILE:HG21	2:D:109:ASP:HB3	1.88	0.55
2:D:118:LEU:HD12	2:D:121:ARG:HD2	1.89	0.55
2:F:3:PHE:HB2	2:F:112:ASP:OD2	2.07	0.55
1:C:95:GLU:OE1	1:C:99:ARG:NH1	2.40	0.54
2:D:54:ARG:NH2	2:D:103:GLU:OE2	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:LYS:O	1:C:206:PRO:HD2	2.07	0.54
2:D:75:LYS:HD2	2:D:75:LYS:N	2.15	0.54
2:D:145:ASP:O	2:D:148:CYS:N	2.40	0.54
1:E:228:LEU:HD13	1:E:234:VAL:HG23	1.89	0.54
2:F:118:LEU:O	2:F:121:ARG:HG3	2.08	0.53
1:A:110:LYS:HE3	1:A:249:SER:HB3	1.90	0.53
1:C:120:ILE:HB	1:C:144:LEU:O	2.07	0.53
2:B:27:GLN:NE2	2:B:32:GLU:HB2	2.23	0.53
1:E:139:GLU:OE1	1:E:247:ARG:HD3	2.09	0.53
1:A:139:GLU:OE1	1:A:247:ARG:HD3	2.08	0.53
1:C:282:ASN:HB3	2:D:56:ILE:HG23	1.91	0.53
1:E:194:THR:CG2	1:E:237:SER:HB2	2.35	0.53
1:A:63:PRO:HB2	1:A:65:GLN:OE1	2.09	0.53
2:F:127:ARG:HG3	2:F:159:HIS:CG	2.44	0.53
1:A:298:ARG:NH2	2:D:90:GLU:OE1	2.42	0.52
1:E:197:SER:OG	1:E:198:SER:N	2.43	0.52
1:C:201:GLN:NE2	1:E:222:ASP:HB3	2.25	0.52
2:D:66:ILE:HD13	2:D:66:ILE:H	1.73	0.52
1:E:170:VAL:HG22	1:E:225:TRP:HB3	1.91	0.52
1:A:95:GLU:OE1	1:A:99:ARG:NH1	2.40	0.52
1:C:171:TRP:CZ3	1:C:226:LEU:HD22	2.45	0.52
2:B:124:ARG:CB	2:D:132:GLU:OE1	2.55	0.51
2:B:148:CYS:O	2:B:151:SER:HB3	2.11	0.51
1:C:32:THR:HG23	1:C:285:PHE:HD2	1.75	0.51
1:C:293:VAL:HG11	2:D:65:LEU:HD13	1.92	0.51
2:D:39:LYS:HE2	2:D:121:ARG:NH2	2.25	0.51
2:D:158:ASP:OD2	2:D:160:ARG:NH2	2.43	0.51
1:E:170:VAL:O	1:E:245:PRO:HB3	2.10	0.51
1:E:207:SER:O	1:E:207:SER:OG	2.26	0.51
1:E:106:GLY:HA2	1:E:255:SER:HB3	1.93	0.51
2:D:133:ASP:OD2	2:D:135:THR:HG23	2.10	0.50
1:A:222:ASP:HB3	1:E:201:GLN:NE2	2.25	0.50
2:D:148:CYS:O	2:D:151:SER:HB3	2.10	0.50
2:B:66:ILE:HD13	2:B:66:ILE:H	1.75	0.50
1:A:64:PRO:HD3	1:A:129:CYS:SG	2.51	0.50
1:C:171:TRP:HZ3	1:C:226:LEU:HD22	1.77	0.50
2:D:50:GLY:HA3	1:E:20:THR:O	2.11	0.50
2:B:16:GLY:O	2:B:18:ILE:HG12	2.10	0.50
1:E:282:ASN:HB3	2:F:56:ILE:HG23	1.93	0.50
1:A:4:CYS:HA	2:B:137:CYS:HA	1.94	0.50
1:A:153:PRO:O	1:A:155:MET:HG3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:130:ARG:HH12	1:E:133:GLY:HA2	1.76	0.50
1:C:192:LEU:HD11	1:C:194:THR:CG2	2.42	0.50
1:E:191:LYS:O	1:E:206:PRO:HD2	2.12	0.50
1:A:60:ILE:CD1	1:A:170:VAL:HG11	2.42	0.49
1:E:68:GLN:HG2	1:E:69:PHE:CD2	2.47	0.49
1:A:86:VAL:HG22	1:A:88:TYR:O	2.12	0.49
1:A:106:GLY:HA2	1:A:255:SER:HB3	1.93	0.49
2:F:120:GLU:O	2:F:124:ARG:HG3	2.13	0.49
1:E:120:ILE:HB	1:E:144:LEU:O	2.13	0.49
1:C:169:ILE:HB	1:C:226:LEU:HD23	1.95	0.49
1:A:90:GLY:HA3	1:A:221:ILE:O	2.13	0.48
1:C:4:CYS:HA	2:D:137:CYS:CB	2.32	0.48
2:D:124:ARG:HA	2:F:132:GLU:OE1	2.13	0.48
1:E:64:PRO:HD3	1:E:129:CYS:SG	2.53	0.48
1:C:99:ARG:O	1:C:103:ARG:HG3	2.13	0.48
1:C:106:GLY:HA2	1:C:255:SER:HB3	1.95	0.48
2:D:16:GLY:O	2:D:18:ILE:HG13	2.14	0.48
2:F:66:ILE:HD13	2:F:66:ILE:H	1.78	0.48
2:B:163:ARG:O	2:B:167:MET:HG3	2.14	0.48
2:D:25:ARG:HE	2:D:34:THR:CG2	2.27	0.48
1:C:88:TYR:CD1	1:C:89:PRO:HD2	2.49	0.48
2:B:120:GLU:O	2:B:124:ARG:HG3	2.14	0.47
1:C:21:GLU:CD	1:C:24:VAL:HG11	2.35	0.47
2:F:27:GLN:NE2	2:F:32:GLU:HB2	2.29	0.47
1:A:99:ARG:O	1:A:103:ARG:HG3	2.14	0.47
1:A:170:VAL:HG12	1:A:225:TRP:CB	2.42	0.47
2:B:6:ILE:N	2:B:112:ASP:OD1	2.47	0.47
1:E:226:LEU:HD11	1:E:234:VAL:HG21	1.97	0.47
1:C:90:GLY:HA3	1:C:221:ILE:O	2.14	0.47
1:C:86:VAL:HG22	1:C:137:TYR:OH	2.14	0.47
1:A:197:SER:HB3	1:A:234:VAL:HG23	1.96	0.47
1:E:99:ARG:O	1:E:103:ARG:HG3	2.15	0.47
1:A:12:ASN:OD1	1:A:12:ASN:O	2.33	0.46
1:C:21:GLU:HB2	1:C:24:VAL:CG1	2.45	0.46
1:C:170:VAL:O	1:C:245:PRO:HB3	2.15	0.46
1:E:32:THR:HG23	1:E:285:PHE:HD2	1.79	0.46
1:A:3:ILE:HG22	2:B:138:PHE:HB2	1.98	0.46
2:D:142:HIS:HB2	2:D:165:GLU:OE2	2.15	0.46
2:D:163:ARG:NH2	2:F:131:GLU:OE2	2.48	0.46
1:A:254:LYS:HB2	1:A:254:LYS:NZ	2.31	0.46
2:F:10:ILE:HD12	2:F:115:MET:SD	2.56	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:28:ASN:HD22	2:F:145:ASP:HA	1.80	0.46
1:A:191:LYS:O	1:A:206:PRO:HD2	2.16	0.46
1:E:86:VAL:HG22	1:E:137:TYR:OH	2.15	0.46
1:E:163:ARG:HD2	1:E:163:ARG:HA	1.51	0.45
1:C:64:PRO:HD3	1:C:129:CYS:SG	2.56	0.45
2:F:121:ARG:HD3	2:F:125:GLN:NE2	2.31	0.45
1:C:180:ALA:O	1:C:183:THR:HB	2.17	0.45
2:B:46:ASP:HA	2:B:49:THR:HB	1.98	0.45
1:C:154:GLN:HG2	1:C:239:ASN:HB3	1.99	0.45
1:E:171:TRP:CH2	1:E:226:LEU:HD23	2.52	0.45
2:F:56:ILE:O	2:F:56:ILE:HG22	2.16	0.45
1:A:282:ASN:HB3	2:B:56:ILE:HG23	1.97	0.45
1:C:130:ARG:NH1	1:C:133:GLY:HA2	2.31	0.45
2:D:56:ILE:O	2:D:56:ILE:HG22	2.16	0.45
1:E:95:GLU:OE1	1:E:99:ARG:NH1	2.46	0.45
2:B:118:LEU:O	2:B:121:ARG:HG3	2.17	0.44
1:E:130:ARG:NH1	1:E:133:GLY:HA2	2.32	0.44
1:E:261:ARG:HG2	1:E:275:SER:O	2.17	0.44
2:F:22:TYR:HD1	2:F:115:MET:HE3	1.82	0.44
1:A:161:ASN:ND2	1:A:230:PRO:HA	2.32	0.44
3:B:201:NAG:O7	3:B:201:NAG:O3	2.32	0.44
1:A:170:VAL:O	1:A:245:PRO:HB3	2.17	0.44
2:B:51:LYS:HZ1	2:B:106:HIS:HB3	1.81	0.44
1:E:10:VAL:HG22	1:E:11:SER:N	2.32	0.44
1:C:7:HIS:HB3	2:D:115:MET:HE3	1.99	0.44
1:A:32:THR:HG22	1:A:305:LEU:HB2	1.99	0.44
1:A:201:GLN:NE2	1:C:222:ASP:HB3	2.30	0.44
1:C:101:ILE:HD13	1:C:227:ILE:HD11	2.00	0.44
2:F:133:ASP:OD1	2:F:137:CYS:N	2.47	0.44
1:C:49:VAL:HG23	1:C:74:ALA:HB2	1.99	0.44
2:B:26:HIS:O	2:B:32:GLU:HA	2.16	0.43
2:B:56:ILE:HG22	2:B:56:ILE:O	2.18	0.43
1:C:10:VAL:HG22	1:C:11:SER:N	2.33	0.43
1:E:90:GLY:HA3	1:E:221:ILE:O	2.18	0.43
2:D:131:GLU:OE1	2:D:170:ARG:NE	2.51	0.43
1:E:171:TRP:CZ2	1:E:195:VAL:HG11	2.54	0.43
2:F:47:GLN:NE2	2:F:110:LEU:HD11	2.34	0.43
1:A:32:THR:HG22	1:A:305:LEU:O	2.19	0.43
1:E:41:ILE:HD11	1:E:260:SER:OG	2.19	0.43
1:A:7:HIS:CE1	1:A:314:VAL:HG12	2.50	0.42
1:E:4:CYS:HA	2:F:137:CYS:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:260:SER:HA	1:E:275:SER:HA	2.02	0.42
2:D:163:ARG:O	2:D:167:MET:HG3	2.19	0.42
1:C:21:GLU:OE2	1:C:24:VAL:HG11	2.20	0.42
2:D:128:GLU:OE1	2:F:170:ARG:NH2	2.53	0.42
1:E:153:PRO:O	1:E:155:MET:HG3	2.19	0.42
2:B:128:GLU:O	2:B:170:ARG:NH1	2.53	0.42
1:C:197:SER:OG	1:C:200:TYR:HB3	2.20	0.42
1:E:32:THR:HG22	1:E:33:VAL:HG23	2.02	0.42
1:E:44:LYS:HE2	1:E:269:GLU:HB2	2.01	0.42
1:A:171:TRP:CH2	1:A:226:LEU:HD23	2.54	0.42
2:F:27:GLN:NE2	2:F:32:GLU:OE1	2.52	0.42
1:A:47:ARG:NH1	1:A:75:ASP:OD1	2.53	0.41
2:D:30:GLN:C	2:D:30:GLN:OE1	2.58	0.41
2:B:29:ALA:HB3	2:B:30:GLN:OE1	2.20	0.41
1:C:176:SER:O	1:C:208:PRO:HA	2.20	0.41
1:E:47:ARG:NH1	1:E:75:ASP:OD1	2.53	0.41
1:E:160:LYS:HG2	1:E:161:ASN:N	2.33	0.41
1:A:147:THR:HG22	1:A:150:ALA:HB2	2.03	0.41
1:A:189:GLY:O	1:A:191:LYS:HG2	2.20	0.41
2:B:2:ILE:H	2:F:117:LYS:HE2	1.85	0.41
1:C:47:ARG:NH1	1:C:75:ASP:OD2	2.52	0.41
1:C:91:LYS:O	1:C:222:ASP:HA	2.21	0.41
1:A:120:ILE:HB	1:A:144:LEU:O	2.19	0.41
1:E:211:ARG:HD2	1:E:220:ARG:CG	2.51	0.41
2:F:26:HIS:C	2:F:26:HIS:HD1	2.24	0.41
2:B:127:ARG:HG3	2:B:159:HIS:CD2	2.56	0.41
1:C:172:GLY:HA3	1:C:243:ILE:HB	2.03	0.41
1:E:49:VAL:HG23	1:E:74:ALA:HB2	2.03	0.41
2:B:170:ARG:NH2	2:F:128:GLU:OE1	2.54	0.41
2:D:28:ASN:HD22	2:D:145:ASP:HA	1.86	0.41
1:C:158:SER:OG	1:C:235:THR:HG23	2.20	0.40
2:D:141:PHE:CD2	2:D:170:ARG:HG2	2.56	0.40
2:D:164:GLU:O	2:D:168:GLN:OE1	2.39	0.40
1:C:191:LYS:O	1:C:205:VAL:HG13	2.21	0.40
1:A:49:VAL:HG23	1:A:74:ALA:HB2	2.01	0.40
2:B:28:ASN:HD22	2:B:145:ASP:HA	1.87	0.40
1:C:183:THR:HA	1:C:187:GLY:O	2.21	0.40
2:D:25:ARG:HE	2:D:34:THR:HG21	1.86	0.40
1:E:199:ASN:C	1:E:199:ASN:OD1	2.59	0.40
1:C:207:SER:O	1:C:207:SER:OG	2.25	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	314/316 (99%)	294 (94%)	19 (6%)	1 (0%)	41	66
1	C	314/316 (99%)	295 (94%)	19 (6%)	0	100	100
1	E	314/316 (99%)	295 (94%)	19 (6%)	0	100	100
2	B	169/171 (99%)	152 (90%)	15 (9%)	2 (1%)	13	32
2	D	169/171 (99%)	152 (90%)	15 (9%)	2 (1%)	13	32
2	F	169/171 (99%)	153 (90%)	15 (9%)	1 (1%)	25	50
All	All	1449/1461 (99%)	1341 (92%)	102 (7%)	6 (0%)	34	60

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	2	ILE
2	D	12	ASN
2	B	12	ASN
2	D	2	ILE
1	A	198	SER
2	F	2	ILE

5.3.2 Protein sidechains (i)

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	267/267 (100%)	256 (96%)	11 (4%)	30	59
1	C	267/267 (100%)	258 (97%)	9 (3%)	37	66

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	267/267 (100%)	259 (97%)	8 (3%)	41	70
2	B	145/145 (100%)	139 (96%)	6 (4%)	30	59
2	D	145/145 (100%)	135 (93%)	10 (7%)	15	35
2	F	145/145 (100%)	137 (94%)	8 (6%)	21	46
All	All	1236/1236 (100%)	1184 (96%)	52 (4%)	30	58

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ASP
1	A	8	HIS
1	A	38	THR
1	A	81	ARG
1	A	121	ARG
1	A	129	CYS
1	A	141	LYS
1	A	163	ARG
1	A	186	TYR
1	A	199	ASN
1	A	211	ARG
2	B	26	HIS
2	B	27	GLN
2	B	65	LEU
2	B	66	ILE
2	B	75	LYS
2	B	121	ARG
1	C	2	LYS
1	C	8	HIS
1	C	32	THR
1	C	38	THR
1	C	65	GLN
1	C	86	VAL
1	C	186	TYR
1	C	211	ARG
1	C	268	CYS
2	D	26	HIS
2	D	65	LEU
2	D	66	ILE
2	D	67	ASP
2	D	75	LYS
2	D	121	ARG

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Mol	Chain	Res	Type
2	D	133	ASP
2	D	137	CYS
2	D	143	LYS
2	D	151	SER
1	E	8	HIS
1	E	32	THR
1	E	38	THR
1	E	86	VAL
1	E	121	ARG
1	E	163	ARG
1	E	186	TYR
1	E	211	ARG
2	F	26	HIS
2	F	60	ASN
2	F	65	LEU
2	F	66	ILE
2	F	75	LYS
2	F	121	ARG
2	F	133	ASP
2	F	143	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
2	B	27	GLN
2	B	61	GLN
2	D	62	GLN
2	D	71	ASN
1	E	65	GLN
2	F	27	GLN
2	F	47	GLN
2	F	125	GLN
2	F	169	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 monosaccharides 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	$\#\|Z\| > 2$
3	NAG	E	401	-	14,14,15	0.49	0	17,19,21	1.03	0
3	NAG	C	402	1	14,14,15	0.60	0	17,19,21	1.95	4 (23%)
3	NAG	D	201	2	14,14,15	0.34	0	17,19,21	0.50	0
3	NAG	B	201	2	14,14,15	0.38	0	17,19,21	0.53	0
3	NAG	A	402	1	14,14,15	0.48	0	17,19,21	1.52	3 (17%)
3	NAG	C	401	1	14,14,15	0.19	0	17,19,21	0.81	1 (5%)
3	NAG	E	402	1	14,14,15	0.37	0	17,19,21	1.20	1 (5%)
3	NAG	A	401	1	14,14,15	0.25	0	17,19,21	1.27	1 (5%)
3	NAG	F	201	2	14,14,15	0.44	0	17,19,21	0.71	1 (5%)

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	E	401	-	-	0/6/23/26	0/1/1/1
3	NAG	C	402	1	-	3/6/23/26	0/1/1/1
3	NAG	D	201	2	-	0/6/23/26	0/1/1/1
3	NAG	B	201	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	402	1	-	3/6/23/26	0/1/1/1
3	NAG	C	401	1	-	1/6/23/26	0/1/1/1
3	NAG	E	402	1	-	3/6/23/26	0/1/1/1
3	NAG	A	401	1	-	2/6/23/26	0/1/1/1
3	NAG	F	201	2	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	402	NAG	C1-O5-C5	5.96	120.27	112.19
3	A	401	NAG	C1-O5-C5	4.43	118.19	112.19
3	A	402	NAG	C1-O5-C5	4.25	117.95	112.19
3	E	402	NAG	C1-O5-C5	3.50	116.94	112.19
3	C	401	NAG	C1-O5-C5	2.62	115.74	112.19
3	C	402	NAG	C2-N2-C7	2.44	126.38	122.90
3	C	402	NAG	C1-C2-N2	2.42	114.63	110.49
3	A	402	NAG	C2-N2-C7	2.28	126.15	122.90
3	A	402	NAG	C1-C2-N2	2.24	114.31	110.49
3	F	201	NAG	C2-N2-C7	2.21	126.04	122.90
3	C	402	NAG	C3-C4-C5	2.14	114.05	110.24

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	NAG	O5-C5-C6-O6
3	C	402	NAG	O5-C5-C6-O6
3	E	402	NAG	O5-C5-C6-O6
3	A	402	NAG	C4-C5-C6-O6
3	B	201	NAG	C1-C2-N2-C7
3	E	402	NAG	C4-C5-C6-O6
3	F	201	NAG	O5-C5-C6-O6
3	C	402	NAG	C1-C2-N2-C7
3	C	402	NAG	C4-C5-C6-O6
3	F	201	NAG	C4-C5-C6-O6
3	A	402	NAG	C1-C2-N2-C7
3	F	201	NAG	C1-C2-N2-C7
3	E	402	NAG	C1-C2-N2-C7
3	F	201	NAG	C3-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
3	A	401	NAG	C4-C5-C6-O6
3	A	401	NAG	O5-C5-C6-O6
3	B	201	NAG	C3-C2-N2-C7
3	C	401	NAG	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	401	NAG	5	0
3	B	201	NAG	1	0
3	F	201	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 (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	316/316 (100%)	0.37	17 (5%)	25	24	33, 67, 105, 135	2 (0%)
1	C	316/316 (100%)	0.35	14 (4%)	34	33	33, 68, 103, 133	2 (0%)
1	E	316/316 (100%)	0.32	10 (3%)	47	48	32, 67, 103, 130	2 (0%)
2	B	171/171 (100%)	0.53	16 (9%)	8	6	25, 91, 130, 143	0
2	D	171/171 (100%)	0.54	14 (8%)	11	9	24, 91, 130, 140	0
2	F	171/171 (100%)	0.54	13 (7%)	13	12	22, 90, 128, 139	0
All	All	1461/1461 (100%)	0.41	84 (5%)	23	22	22, 73, 122, 143	6 (0%)

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	119	GLY	9.2
2	F	168	GLN	7.2
1	A	149	ASN	5.4
2	D	139	GLU	5.2
2	B	22	TYR	4.4
2	D	2	ILE	4.3
1	A	242	PHE	4.1
2	D	33	GLY	4.1
1	A	145	SER	4.0
1	E	195	VAL	3.9
1	A	240	GLY	3.7
2	D	32	GLU	3.7
2	D	140	ILE	3.6
2	D	58	LYS	3.5
1	E	189	GLY	3.5
2	B	33	GLY	3.4
1	A	156	THR	3.4
1	C	144	LEU	3.4
1	A	120	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	186	TYR	3.4
1	A	190	ASN	3.3
2	B	32	GLU	3.3
1	A	151	ALA	3.2
1	C	114	GLY	3.2
2	D	138	PHE	3.2
2	D	164	GLU	3.1
1	E	3	ILE	3.1
1	C	181	GLU	2.9
2	D	27	GLN	2.9
1	A	198	SER	2.9
1	C	3	ILE	2.9
2	B	119	TYR	2.9
2	B	131	GLU	2.8
2	F	3	PHE	2.8
2	D	31	GLY	2.8
2	B	138	PHE	2.7
2	F	32	GLU	2.7
1	C	6	GLY	2.7
2	F	157	TYR	2.7
2	F	2	ILE	2.7
1	A	143	LEU	2.6
1	E	128	ALA	2.6
1	E	190	ASN	2.6
2	D	22	TYR	2.6
2	B	130	ALA	2.6
1	E	179	THR	2.6
2	B	3	PHE	2.6
2	F	17	LEU	2.6
1	C	151	ALA	2.5
1	E	117	TYR	2.5
1	C	2	LYS	2.5
2	B	5	ALA	2.5
1	C	251	LEU	2.5
2	F	33	GLY	2.5
1	C	236	PHE	2.4
2	F	27	GLN	2.4
2	B	11	GLU	2.4
1	A	170	VAL	2.3
1	E	159	TYR	2.3
2	B	168	GLN	2.3
1	C	240	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	154	GLN	2.3
1	C	5	LEU	2.3
2	F	139	GLU	2.3
1	C	153	PRO	2.3
2	B	164	GLU	2.3
2	F	24	PHE	2.3
1	A	238	PHE	2.2
2	B	129	ASN	2.2
1	A	155	MET	2.2
2	F	126	LEU	2.2
2	D	124	ARG	2.2
2	F	131	GLU	2.2
2	F	121	ARG	2.2
2	B	170	ARG	2.1
2	D	144	CYS	2.1
1	C	226	LEU	2.1
2	B	143	LYS	2.1
2	D	132	GLU	2.1
1	A	117	TYR	2.0
2	B	162	TYR	2.0
1	C	242	PHE	2.0
1	E	115	PHE	2.0
1	A	144	LEU	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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	NAG	E	402	14/15	0.71	0.22	101,119,128,129	0
3	NAG	A	402	14/15	0.75	0.23	102,114,122,123	0
3	NAG	E	401	14/15	0.80	0.26	85,111,118,118	0
3	NAG	A	401	14/15	0.82	0.25	86,110,117,120	0
3	NAG	C	402	14/15	0.82	0.13	87,112,124,128	0
3	NAG	F	201	14/15	0.82	0.17	67,86,104,114	0
3	NAG	C	401	14/15	0.83	0.20	87,111,118,119	0
3	NAG	D	201	14/15	0.86	0.21	64,85,99,108	0
3	NAG	B	201	14/15	0.89	0.17	62,79,104,119	0

6.5 Other polymers (i)

There are no such residues in this entry.