



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 26, 2021 – 08:04 PM EST

PDB ID : 7KJH
Title : Plasmodium falciparum protein Pf12p bound to nanobody B9
Authors : Dietrich, M.H.; Tham, W.H.
Deposited on : 2020-10-26
Resolution : 2.00 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.16
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) : 2.16

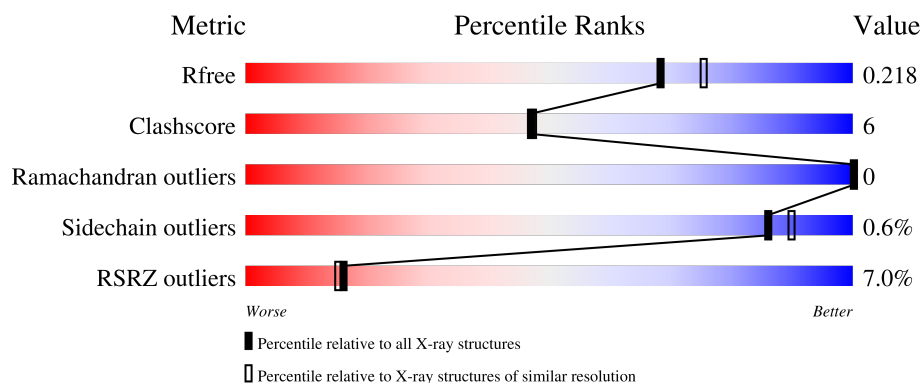
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 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 $\leq 5\%$. 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	126	<div> <div>4%</div> <div>83%</div> <div>13%</div> <div>5%</div> </div>
1	B	126	<div> <div>7%</div> <div>92%</div> <div>8%</div> </div>
2	C	321	<div> <div>6%</div> <div>66%</div> <div>15%</div> <div>19%</div> </div>
2	D	321	<div> <div>7%</div> <div>70%</div> <div>11%</div> <div>19%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6790 atoms, of which 48 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 Nanobody B9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	120	Total	C	N	O	S	0	1	0
			903	572	157	170	4			
1	B	126	Total	C	N	O	S	0	1	0
			978	619	174	181	4			

- Molecule 2 is a protein called Surface protein P12p.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	261	Total	C	N	O	S	0	3	0
			2103	1357	343	385	18			
2	D	260	Total	C	N	O	S	0	4	0
			2085	1348	337	381	19			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	21	GLY	-	expression tag	UNP C6KSX1
C	22	ALA	-	expression tag	UNP C6KSX1
C	23	SER	-	expression tag	UNP C6KSX1
D	21	GLY	-	expression tag	UNP C6KSX1
D	22	ALA	-	expression tag	UNP C6KSX1
D	23	SER	-	expression tag	UNP C6KSX1

- Molecule 3 is CITRATE ANION (three-letter code: FLC) (formula: C₆H₅O₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			18	6	5	7		
3	A	1	Total	C	H	O	0	0
			18	6	5	7		
3	B	1	Total	C	H	O	0	0
			18	6	5	7		
3	B	1	Total	C	H	O	0	0
			18	6	5	7		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
4	D	1	Total	C	H	N	O	0	0
			28	8	14	1	5		

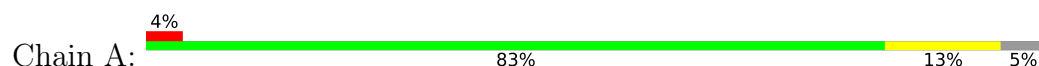
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	95	Total	O	0	0
			95	95		
5	C	220	Total	O	0	0
			220	220		
5	D	171	Total	O	0	0
			171	171		
5	B	107	Total	O	0	0
			107	107		

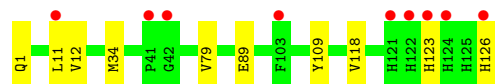
3 Residue-property plots [i](#)

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.

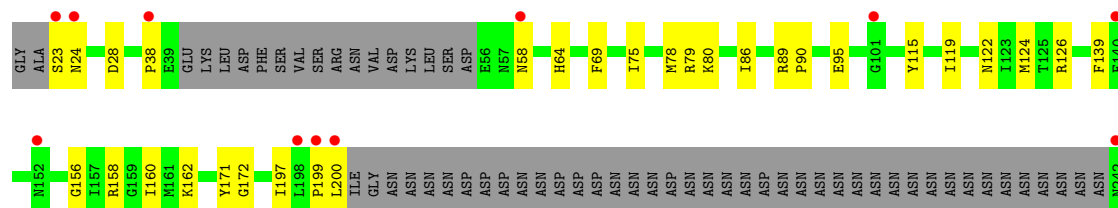
- Molecule 1: Nanobody B9



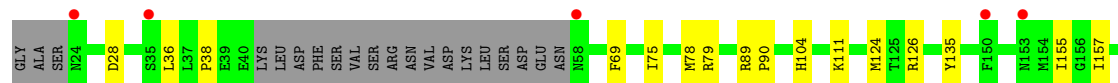
- Molecule 1: Nanobody B9

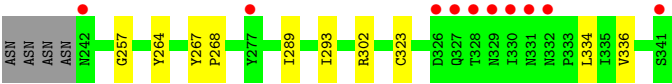
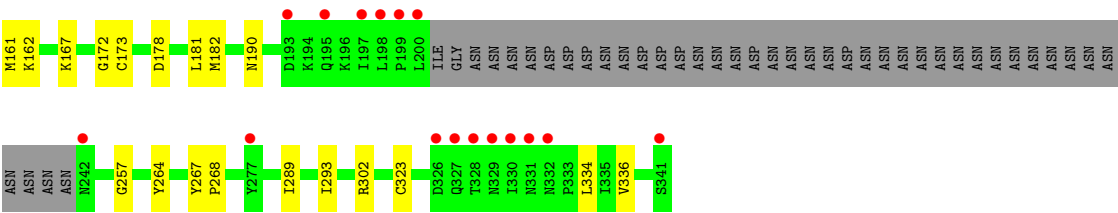


- Molecule 2: Surface protein P12p



- Molecule 2: Surface protein P12p





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.99Å 107.04Å 114.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.34 – 2.00 48.45 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.34-2.00) 100.0 (48.45-2.00)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.15_3459	Depositor
R, R_{free}	0.180 , 0.217 0.181 , 0.218	Depositor DCC
R_{free} test set	2126 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6790	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FLC, 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	0.47	1/927 (0.1%)	0.62	0/1257
1	B	0.40	0/1008	0.54	0/1366
2	C	0.42	0/2155	0.59	0/2916
2	D	0.38	0/2147	0.57	0/2907
All	All	0.41	1/6237 (0.0%)	0.58	0/8446

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	96	CYS	CB-SG	-5.35	1.73	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	903	0	838	9	0
1	B	978	0	894	6	0
2	C	2103	0	1999	35	0
2	D	2085	0	1976	24	0
3	A	26	10	10	1	0
3	B	26	10	10	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	14	14	13	0	0
4	D	14	14	13	0	0
5	A	95	0	0	2	1
5	B	107	0	0	1	0
5	C	220	0	0	5	1
5	D	171	0	0	5	2
All	All	6742	48	5753	74	2

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 6.

All (74) 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)
3:B:201:FLC:OB2	5:B:301:HOH:O	1.95	0.84
1:A:27:ARG:NH2	5:A:301:HOH:O	1.97	0.82
2:C:58:ASN:O	5:C:501:HOH:O	2.02	0.77
2:C:78:MET:HE2	2:C:79:ARG:H	1.49	0.76
2:C:122[B]:ASN:ND2	5:C:504:HOH:O	2.20	0.73
2:D:78[A]:MET:HE3	2:D:79:ARG:H	1.56	0.69
2:C:78:MET:HE3	2:C:78:MET:HA	1.74	0.69
2:D:182:MET:HE3	5:D:562:HOH:O	1.91	0.69
2:D:78[A]:MET:CE	2:D:79:ARG:H	2.07	0.67
1:A:106:ASN:HB3	1:B:126:HIS:HB2	1.78	0.66
2:C:23:SER:N	2:C:64:HIS:HD1	1.95	0.65
2:C:171:TYR:OH	5:C:502:HOH:O	2.15	0.64
2:C:139:PHE:CD1	2:C:162:LYS:HE3	2.35	0.61
2:C:78:MET:HE1	5:C:714:HOH:O	1.99	0.61
2:C:78:MET:CE	2:C:79:ARG:H	2.13	0.61
2:C:280:ASN:OD1	2:C:281:ILE:HG12	2.01	0.60
2:C:139:PHE:CE1	2:C:162:LYS:HE3	2.37	0.59
2:C:283:ILE:HD11	2:C:288:ILE:HD13	1.83	0.59
2:D:124:MET:HE2	2:D:126:ARG:HH21	1.70	0.57
2:D:124:MET:CE	2:D:126:ARG:HH21	2.18	0.56
2:D:323:CYS:HB2	2:D:334:LEU:HB2	1.88	0.56
2:D:78[B]:MET:HG3	2:D:79:ARG:N	2.22	0.56
2:D:36:LEU:HA	2:D:157:ILE:HB	1.88	0.55
2:D:172:GLY:HA3	2:D:257:GLY:O	2.07	0.55
2:C:124:MET:HE1	2:C:126:ARG:HH11	1.73	0.53
2:C:199:PRO:O	2:C:200:LEU:HB2	2.08	0.53
2:D:267:TYR:HA	2:D:268:PRO:C	2.30	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:158:ARG:NH1	2:C:160:ILE:HD11	2.24	0.52
2:D:104:HIS:HE1	5:D:573:HOH:O	1.92	0.52
2:D:135:TYR:O	2:D:167:LYS:HE3	2.11	0.51
2:D:38:PRO:HA	2:D:155:ILE:O	2.11	0.50
3:A:202:FLC:OG1	1:B:123:HIS:HA	2.12	0.49
2:C:248:THR:CG2	2:C:341:SER:HB3	2.42	0.49
2:C:280:ASN:O	2:C:281:ILE:HD13	2.13	0.49
2:C:80:LYS:NZ	2:C:122[B]:ASN:OD1	2.45	0.48
1:A:74:ASN:HA	5:A:367:HOH:O	2.13	0.48
2:C:79:ARG:HG2	2:C:86:ILE:HG22	1.95	0.48
2:D:178:ASP:CG	2:D:181:LEU:HG	2.34	0.48
2:C:172:GLY:HA3	2:C:257:GLY:O	2.14	0.47
2:C:124:MET:CE	2:C:126:ARG:HH11	2.27	0.47
1:B:89:GLU:CD	1:B:89:GLU:H	2.18	0.46
2:C:38:PRO:HA	2:C:156:GLY:HA2	1.97	0.46
1:B:34:MET:HG3	1:B:79:VAL:HG21	1.97	0.46
1:B:12:VAL:CG1	1:B:118:VAL:HG22	2.46	0.46
1:A:34:MET:HG3	1:A:79:VAL:HG21	1.97	0.45
2:C:267:TYR:HA	2:C:268:PRO:C	2.37	0.45
2:C:259:ILE:HD13	2:C:303:ILE:HG23	1.98	0.45
2:D:264:TYR:O	2:D:302:ARG:HG2	2.16	0.45
2:C:197:ILE:O	2:C:200:LEU:N	2.50	0.45
2:C:264:TYR:O	2:C:302:ARG:HG2	2.17	0.44
2:D:124:MET:HE2	5:D:549:HOH:O	2.16	0.44
2:C:89:ARG:HA	2:C:90:PRO:C	2.37	0.44
2:D:28:ASP:HA	2:D:75:ILE:HB	2.00	0.44
1:A:13:GLN:HA	1:A:119:SER:HB2	1.98	0.44
1:A:67:ARG:CZ	1:A:87:LYS:HE2	2.48	0.44
2:C:95:GLU:HG3	2:C:115:TYR:CE2	2.53	0.44
2:C:316:PRO:HD2	2:C:340:PHE:CG	2.53	0.43
1:A:68:PHE:HA	1:A:82:GLN:O	2.19	0.43
1:A:14:ALA:HB2	1:A:119:SER:O	2.19	0.43
2:D:173:CYS:SG	2:D:336:VAL:HG13	2.59	0.43
2:D:162:LYS:HE3	5:D:559:HOH:O	2.19	0.43
2:C:78:MET:SD	2:C:119:ILE:HD11	2.59	0.43
1:A:48:LEU:HD22	1:A:64:VAL:HG11	2.00	0.42
2:C:260:CYS:O	2:C:302:ARG:NH2	2.52	0.42
2:D:289:ILE:HB	2:D:293:ILE:HD11	2.01	0.42
2:D:89:ARG:HA	2:D:90:PRO:C	2.40	0.42
2:C:24:ASN:N	2:C:24:ASN:OD1	2.51	0.41
2:D:111:LYS:HE2	5:D:520:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:GLN:HA	1:B:109:TYR:CZ	2.55	0.41
2:C:28:ASP:HA	2:C:75:ILE:HB	2.03	0.41
2:C:78:MET:CE	5:C:714:HOH:O	2.64	0.41
2:D:173:CYS:HA	2:D:190:ASN:O	2.21	0.41
2:D:161:MET:HE2	2:D:161:MET:HB3	1.98	0.41
2:C:323:CYS:HB2	2:C:334:LEU:HB2	2.04	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:392:HOH:O	5:D:648:HOH:O[2_455]	2.14	0.06
5:C:695:HOH:O	5:D:637:HOH:O[1_655]	2.18	0.02

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	119/126 (94%)	113 (95%)	6 (5%)	0	100	100
1	B	125/126 (99%)	121 (97%)	4 (3%)	0	100	100
2	C	256/321 (80%)	249 (97%)	7 (3%)	0	100	100
2	D	258/321 (80%)	254 (98%)	4 (2%)	0	100	100
All	All	758/894 (85%)	737 (97%)	21 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	87/101 (86%)	86 (99%)	1 (1%)	73	78
1	B	95/101 (94%)	94 (99%)	1 (1%)	73	78
2	C	232/304 (76%)	231 (100%)	1 (0%)	91	93
2	D	228/304 (75%)	227 (100%)	1 (0%)	91	93
All	All	642/810 (79%)	638 (99%)	4 (1%)	86	90

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	ASP
2	C	69	PHE
2	D	69	PHE
1	B	11	LEU

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 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	FLC	A	202	-	3,12,12	1.04	0	3,17,17	1.85	1 (33%)
4	NAG	D	401	2	14,14,15	1.90	4 (28%)	17,19,21	1.36	3 (17%)
3	FLC	A	201	-	3,12,12	1.25	0	3,17,17	2.08	1 (33%)
3	FLC	B	201	-	3,12,12	1.32	0	3,17,17	2.98	2 (66%)
4	NAG	C	401	2	14,14,15	2.06	4 (28%)	17,19,21	1.55	4 (23%)
3	FLC	B	202	-	3,12,12	0.85	0	3,17,17	0.85	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	FLC	A	202	-	-	0/6/16/16	-
4	NAG	D	401	2	-	0/6/23/26	0/1/1/1
3	FLC	A	201	-	-	2/6/16/16	-
3	FLC	B	201	-	-	3/6/16/16	-
4	NAG	C	401	2	-	0/6/23/26	0/1/1/1
3	FLC	B	202	-	-	1/6/16/16	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	401	NAG	O5-C1	4.43	1.50	1.43
4	D	401	NAG	O5-C1	4.40	1.50	1.43
4	C	401	NAG	C7-N2	3.46	1.46	1.34
4	D	401	NAG	C7-N2	3.25	1.45	1.34
4	C	401	NAG	C2-N2	3.17	1.51	1.46
4	C	401	NAG	O5-C5	2.50	1.48	1.43
4	D	401	NAG	C2-N2	2.13	1.49	1.46
4	D	401	NAG	O5-C5	2.05	1.47	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	201	FLC	CB-CG-CGC	-4.61	107.60	114.98
4	C	401	NAG	C4-C3-C2	3.46	116.08	111.02
3	A	201	FLC	CB-CA-CAC	-3.33	109.65	114.98
4	D	401	NAG	C6-C5-C4	-2.93	106.14	113.00
4	C	401	NAG	C1-C2-N2	-2.80	105.70	110.49
3	A	202	FLC	CB-CG-CGC	-2.62	110.79	114.98
4	D	401	NAG	C3-C4-C5	2.37	114.47	110.24
4	C	401	NAG	C8-C7-N2	2.34	120.06	116.10
4	C	401	NAG	C1-O5-C5	2.27	115.26	112.19
4	D	401	NAG	C8-C7-N2	2.24	119.89	116.10
3	B	201	FLC	CB-CA-CAC	-2.22	111.42	114.98

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	201	FLC	CAC-CA-CB-CBC
3	B	201	FLC	CAC-CA-CB-CG
3	B	201	FLC	CAC-CA-CB-OHB
3	A	201	FLC	OHB-CB-CG-CGC
3	B	202	FLC	CAC-CA-CB-OHB
3	A	201	FLC	CA-CB-CG-CGC

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	202	FLC	1	0
3	B	201	FLC	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, 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	120/126 (95%)	0.19	5 (4%)	36 35	19, 34, 58, 71	0
1	B	126/126 (100%)	0.17	9 (7%)	16 15	22, 32, 54, 76	0
2	C	261/321 (81%)	0.41	19 (7%)	15 14	19, 29, 60, 84	0
2	D	260/321 (80%)	0.29	21 (8%)	12 11	21, 34, 65, 82	0
All	All	767/894 (85%)	0.30	54 (7%)	16 15	19, 32, 62, 84	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	199	PRO	7.0
2	D	328	THR	6.0
2	C	330	ILE	5.6
2	D	198	LEU	5.3
2	D	197	ILE	5.1
2	C	199	PRO	4.7
2	C	313	ASN	4.5
1	B	126	HIS	4.4
2	D	329	ASN	4.2
2	D	200	LEU	4.1
1	A	11	LEU	3.9
2	D	58	ASN	3.9
2	C	200	LEU	3.8
2	D	330	ILE	3.8
1	B	124	HIS	3.7
1	B	123	HIS	3.6
2	C	312	ILE	3.3
2	C	329	ASN	3.3
1	B	11	LEU	3.0
2	C	198	LEU	3.0
2	C	24	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	119	SER	2.9
2	C	152	ASN	2.9
2	C	23	SER	2.8
1	B	103[A]	PHE	2.8
2	D	150	PHE	2.8
2	C	328	THR	2.7
2	D	24	ASN	2.7
1	B	122	HIS	2.7
2	C	58	ASN	2.7
2	D	326	ASP	2.6
2	C	38	PRO	2.5
2	C	331	ASN	2.5
2	D	331	ASN	2.4
1	B	121	HIS	2.4
1	A	10	GLY	2.4
2	D	35	SER	2.3
1	B	41	PRO	2.3
2	D	193	ASP	2.3
2	D	277	TYR	2.3
1	A	40	ALA	2.3
1	A	41	PRO	2.3
2	D	332	ASN	2.2
2	C	242	ASN	2.2
2	C	279	ASN	2.2
2	D	242	ASN	2.1
1	B	42	GLY	2.1
2	D	153	ASN	2.1
2	D	195	GLN	2.1
2	D	327	GLN	2.1
2	D	341	SER	2.1
2	C	101	GLY	2.1
2	C	140	PHE	2.0
2	C	339	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands [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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	NAG	C	401	14/15	0.70	0.31	43,83,106,112	0
4	NAG	D	401	14/15	0.76	0.40	84,101,115,123	0
3	FLC	B	202	13/13	0.82	0.28	30,51,61,72	0
3	FLC	B	201	13/13	0.83	0.20	38,48,61,61	0
3	FLC	A	201	13/13	0.92	0.12	29,41,50,51	0
3	FLC	A	202	13/13	0.95	0.21	27,43,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.