



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 13, 2017 – 08:44 AM EDT

PDB ID : 2YOJ
Title : HCV NS5B polymerase complexed with pyridonylindole compound
Authors : Chen, K.X.; Venkatraman, S.; Anilkumar, G.N.; Zeng, Q.; Lesburg, C.A.; Vibulbhan, B.; Yang, W.; Velazquez, F.; Chan, T.-Y.; Bennett, F.; Sannigrahi, M.; Jiang, Y.; Duca, J.S.; Pinto, P.; Gavalas, S.; Huang, Y.; Wu, W.; Selyutin, O.; Agrawal, S.; Feld, B.; Huang, H.-C.; Li, C.; Cheng, K.-C.; Shih, N.-Y.; Kozlowski, J.A.; Rosenblum, S.B.; Njoroge, F.G.
Deposited on : unknown
Resolution : 1.76 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

<http://wwpdb.org/validation/2016/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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20029824

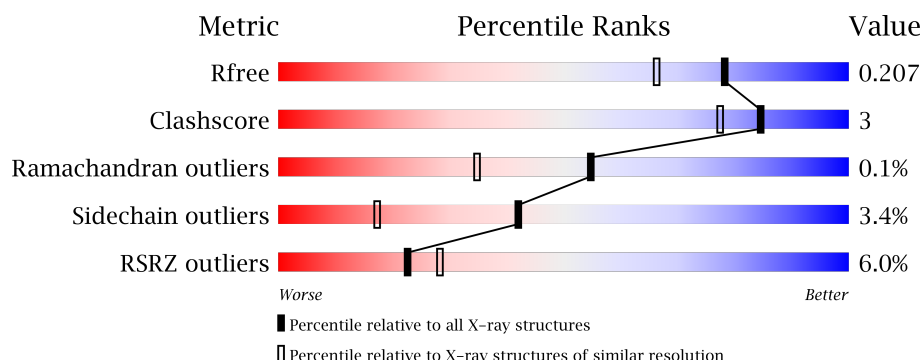
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.76 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1762 (1.76-1.76)
Clashscore	112137	1889 (1.76-1.76)
Ramachandran outliers	110173	1868 (1.76-1.76)
Sidechain outliers	110143	1868 (1.76-1.76)
RSRZ outliers	101464	1770 (1.76-1.76)

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 ≥ 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	585	<div> <div>3%</div> <div>90%</div> <div>6%</div> <div>•</div> </div>
1	B	585	<div> <div>8%</div> <div>86%</div> <div>8%</div> <div>• 5%</div> </div>

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	PO4	B	1566	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10444 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 RNA-DIRECTED RNA POLYMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	563	Total	C	N	O	S	0	4	0
			4414	2778	784	819	33			
1	B	558	Total	C	N	O	S	0	3	0
			4365	2748	772	812	33			

There are 34 discrepancies between the modelled and reference sequences:

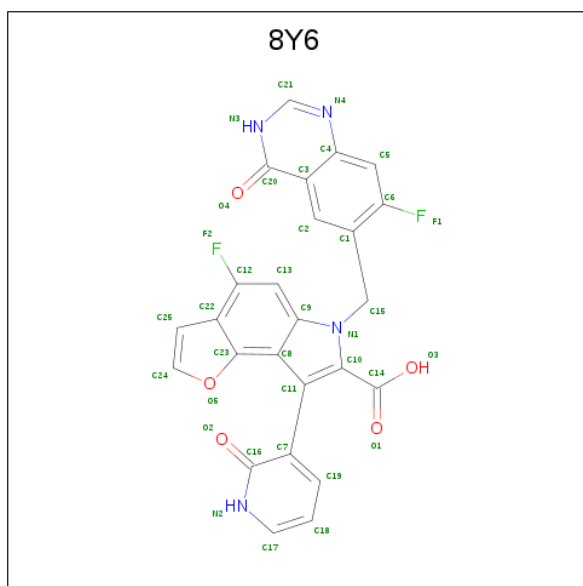
Chain	Residue	Modelled	Actual	Comment	Reference
A	571	GLU	-	expression tag	UNP O92972
A	572	ASN	-	expression tag	UNP O92972
A	573	LEU	-	expression tag	UNP O92972
A	574	TYR	-	expression tag	UNP O92972
A	575	PHE	-	expression tag	UNP O92972
A	576	GLN	-	expression tag	UNP O92972
A	577	GLY	-	expression tag	UNP O92972
A	578	LEU	-	expression tag	UNP O92972
A	579	GLU	-	expression tag	UNP O92972
A	580	HIS	-	expression tag	UNP O92972
A	581	HIS	-	expression tag	UNP O92972
A	582	HIS	-	expression tag	UNP O92972
A	583	HIS	-	expression tag	UNP O92972
A	584	HIS	-	expression tag	UNP O92972
A	585	HIS	-	expression tag	UNP O92972
A	440	GLY	GLU	conflict	UNP O92972
A	520	ILE	THR	conflict	UNP O92972
B	571	GLU	-	expression tag	UNP O92972
B	572	ASN	-	expression tag	UNP O92972
B	573	LEU	-	expression tag	UNP O92972
B	574	TYR	-	expression tag	UNP O92972
B	575	PHE	-	expression tag	UNP O92972
B	576	GLN	-	expression tag	UNP O92972
B	577	GLY	-	expression tag	UNP O92972
B	578	LEU	-	expression tag	UNP O92972

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Chain	Residue	Modelled	Actual	Comment	Reference
B	579	GLU	-	expression tag	UNP O92972
B	580	HIS	-	expression tag	UNP O92972
B	581	HIS	-	expression tag	UNP O92972
B	582	HIS	-	expression tag	UNP O92972
B	583	HIS	-	expression tag	UNP O92972
B	584	HIS	-	expression tag	UNP O92972
B	585	HIS	-	expression tag	UNP O92972
B	440	GLY	GLU	conflict	UNP O92972
B	520	ILE	THR	conflict	UNP O92972

- Molecule 2 is 4-fluoranyl-6-[(7-fluoranyl-4-oxidanylidene-3H-quinazolin-6-yl)methyl]-8-(2-oxidanylidene-1H-pyridin-3-yl)furo[2,3-e]indole-7-carboxylic acid (three-letter code: 8Y6) (formula: C₂₅H₁₄F₂N₄O₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			36	25	2	4	5		
2	B	1	Total	C	F	N	O	0	0
			36	25	2	4	5		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	875	Total	O	0	0
			875	875		
4	B	698	Total	O	0	0
			698	698		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.27Å 106.23Å 134.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.91 – 1.76 19.91 – 1.76	Depositor EDS
% Data completeness (in resolution range)	96.6 (19.91-1.76) 96.6 (19.91-1.76)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.90 (at 1.76Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.177 , 0.209 0.172 , 0.207	Depositor DCC
R_{free} test set	6195 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	21.3	Xtriage
Anisotropy	0.403	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10444	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, 8Y6

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	0/4508	0.60	0/6115
1	B	0.47	0/4459	0.63	0/6050
All	All	0.49	0/8967	0.61	0/12165

There are no bond length outliers.

There are no bond angle outliers.

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	4414	0	4434	18	0
1	B	4365	0	4377	28	0
2	A	36	0	13	1	0
2	B	36	0	13	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
4	A	875	0	0	6	0
4	B	698	0	0	3	0
All	All	10444	0	8837	47	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (47) 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:A:32:ARG:NH2	4:A:2101:HOH:O	2.20	0.73
1:B:24:ASN:HD21	1:B:27:SER:HB3	1.55	0.71
1:A:205:VAL:O	1:A:209:LYS:HG2	1.96	0.66
1:B:309:GLN:NE2	1:B:327:ALA:HB2	2.12	0.64
1:B:32:ARG:HD3	1:B:493:GLY:O	1.98	0.63
1:B:531:ARG:HG3	1:B:531:ARG:HH11	1.63	0.63
1:B:461:GLN:HB3	1:B:542:ALA:HA	1.86	0.57
1:B:201:VAL:HG23	1:B:370:VAL:HG22	1.86	0.57
1:B:461:GLN:HG2	1:B:541:ALA:HB3	1.90	0.53
1:B:309:GLN:HG2	1:B:325:GLU:HB3	1.92	0.52
1:B:23:ILE:HD12	1:B:23:ILE:H	1.74	0.52
1:A:83:LEU:HB2	1:A:173:MET:HA	1.91	0.51
1:B:31:LEU:HD11	1:B:492:LEU:HD12	1.92	0.51
1:A:498:ARG:NH1	4:A:2787:HOH:O	2.45	0.49
1:B:531:ARG:NH1	1:B:531:ARG:HG3	2.25	0.49
1:A:445:CYS:SG	1:A:454:ILE:HD12	2.54	0.48
1:B:268:ASN:HB3	1:B:274:CYS:SG	2.54	0.48
1:A:84:SER:OG	1:A:86:GLU:HG2	2.15	0.47
1:B:419:LEU:CD2	1:B:485:VAL:HG11	2.44	0.47
1:A:106:LYS:HD2	4:A:2144:HOH:O	2.14	0.47
1:A:361:GLU:HG2	1:A:370:VAL:O	2.15	0.47
1:B:309:GLN:HE22	1:B:327:ALA:HB2	1.80	0.46
1:B:309:GLN:O	1:B:324:CYS:HB2	2.15	0.46
1:A:488:CYS:HB2	4:A:2683:HOH:O	2.15	0.46
1:B:100:LYS:NZ	4:B:2255:HOH:O	2.47	0.45
1:B:528:TRP:HA	1:B:533:LYS:NZ	2.32	0.45
1:A:544:GLN:H	1:A:544:GLN:HG2	1.59	0.45
1:B:546:ASP:OD1	1:B:548:SER:HB3	2.16	0.45
1:B:466:LEU:HD22	1:B:551:PHE:HE2	1.82	0.45
1:A:359:ASP:HB3	1:A:362:LEU:HG	1.99	0.44
1:B:24:ASN:O	1:B:28:ASN:HB2	2.18	0.43
1:B:389:THR:HG23	1:B:492:LEU:HD21	2.00	0.43
1:A:372:VAL:HG22	1:A:382:TYR:CD1	2.54	0.43
1:A:309:GLN:O	1:A:324:CYS:HB2	2.19	0.42
1:B:32:ARG:NH2	1:B:495:PRO:HG3	2.34	0.42
1:B:100:LYS:HG3	1:B:101:PHE:CD1	2.53	0.42
1:B:221:THR:O	4:B:2426:HOH:O	2.22	0.42
1:B:419:LEU:HD22	1:B:485:VAL:HG21	2.01	0.42
1:A:172:LYS:HE3	1:A:560:ILE:HD13	2.01	0.42
1:B:33:HIS:HB3	1:B:36:MET:HG3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:GLU:HG2	4:B:2313:HOH:O	2.20	0.41
1:A:566:ARG:NH2	4:A:2865:HOH:O	2.52	0.41
1:B:219:TYR:HB3	1:B:320:LEU:HD23	2.02	0.41
2:A:1569:8Y6:H2	2:A:1569:8Y6:C10	2.50	0.41
1:B:90:LYS:HB3	1:B:90:LYS:HE2	1.96	0.41
1:A:154:ARG:HD2	4:A:2424:HOH:O	2.20	0.41
1:A:201:VAL:HG23	1:A:370:VAL:HG22	2.02	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	563/585 (96%)	554 (98%)	9 (2%)	0	100	100
1	B	557/585 (95%)	543 (98%)	13 (2%)	1 (0%)	51	31
All	All	1120/1170 (96%)	1097 (98%)	22 (2%)	1 (0%)	55	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	25	PRO

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	483/498 (97%)	472 (98%)	11 (2%)	56	31
1	B	478/498 (96%)	457 (96%)	21 (4%)	33	10
All	All	961/996 (96%)	929 (97%)	32 (3%)	42	18

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	A	57	LEU
1	A	100	LYS
1	A	184	GLN
1	A	209	LYS
1	A	336	LEU
1	A	377	SER
1	A	535	LYS
1	A	544	GLN
1	A	564	LEU
1	A	568	ARG
1	B	21	LEU
1	B	23	ILE
1	B	24	ASN
1	B	26	LEU
1	B	81	LYS
1	B	90	LYS
1	B	100	LYS
1	B	184	GLN
1	B	209	LYS
1	B	212	LYS
1	B	309	GLN
1	B	336	LEU
1	B	372	VAL
1	B	461	GLN
1	B	484	ARG
1	B	487	SER
1	B	499	THR
1	B	534	LEU
1	B	545	LEU
1	B	547	LEU
1	B	548	SER

Some 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 carbohydrates 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$
2	8Y6	A	1569	-	27,41,41	1.97	9 (33%)	36,62,62	2.23	11 (30%)
3	PO4	A	1570	-	4,4,4	2.75	1 (25%)	6,6,6	0.75	0
3	PO4	A	1571	-	4,4,4	2.85	1 (25%)	6,6,6	0.61	0
2	8Y6	B	1564	-	27,41,41	1.99	9 (33%)	36,62,62	2.27	11 (30%)
3	PO4	B	1565	-	4,4,4	2.65	1 (25%)	6,6,6	0.72	0
3	PO4	B	1566	-	4,4,4	2.70	1 (25%)	6,6,6	0.67	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	8Y6	A	1569	-	-	0/8/12/12	0/5/6/6
3	PO4	A	1570	-	-	0/0/0/0	0/0/0/0
3	PO4	A	1571	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	8Y6	B	1564	-	-	0/8/12/12	0/5/6/6
3	PO4	B	1565	-	-	0/0/0/0	0/0/0/0
3	PO4	B	1566	-	-	0/0/0/0	0/0/0/0

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1569	8Y6	C7-C11	-2.93	1.44	1.49
2	B	1564	8Y6	C7-C11	-2.84	1.44	1.49
2	B	1564	8Y6	C8-C9	-2.64	1.39	1.42
2	A	1569	8Y6	C21-N3	2.04	1.37	1.33
2	B	1564	8Y6	C4-N4	2.16	1.41	1.37
2	A	1569	8Y6	C2-C1	2.56	1.41	1.37
2	A	1569	8Y6	C4-N4	2.71	1.42	1.37
2	B	1564	8Y6	C20-N3	3.02	1.38	1.33
2	B	1564	8Y6	C2-C1	3.04	1.42	1.37
2	A	1569	8Y6	C5-C6	3.06	1.40	1.35
2	B	1564	8Y6	C18-C19	3.23	1.45	1.38
2	B	1564	8Y6	C16-N2	3.37	1.39	1.33
2	A	1569	8Y6	C16-N2	3.38	1.39	1.33
2	A	1569	8Y6	C18-C19	3.45	1.45	1.38
2	B	1564	8Y6	C5-C6	3.66	1.41	1.35
2	A	1569	8Y6	C20-N3	3.76	1.39	1.33
2	A	1569	8Y6	C16-C7	3.93	1.52	1.44
2	B	1564	8Y6	C16-C7	4.00	1.53	1.44
3	B	1565	PO4	P-O1	5.14	1.61	1.50
3	B	1566	PO4	P-O1	5.18	1.61	1.50
3	A	1570	PO4	P-O1	5.30	1.62	1.50
3	A	1571	PO4	P-O1	5.46	1.62	1.50

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1569	8Y6	C9-C8-C23	-5.25	118.71	123.90
2	A	1569	8Y6	N4-C21-N3	-4.84	124.64	128.86
2	B	1564	8Y6	C9-C8-C23	-4.77	119.18	123.90
2	B	1564	8Y6	N4-C21-N3	-4.74	124.73	128.86
2	B	1564	8Y6	C3-C20-N3	-4.53	114.62	123.48
2	A	1569	8Y6	C3-C20-N3	-4.52	114.64	123.48
2	A	1569	8Y6	C1-C15-N1	-3.76	107.08	113.31
2	B	1564	8Y6	C13-C12-C22	-3.69	122.02	123.96
2	B	1564	8Y6	C1-C15-N1	-3.51	107.49	113.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1569	8Y6	C13-C12-C22	-3.50	122.12	123.96
2	B	1564	8Y6	C18-C17-N2	-2.17	119.82	123.43
2	A	1569	8Y6	C21-N4-C4	2.03	117.40	115.27
2	B	1564	8Y6	C2-C3-C4	2.24	120.78	118.16
2	A	1569	8Y6	C2-C3-C4	2.27	120.82	118.16
2	A	1569	8Y6	F2-C12-C22	2.64	119.73	117.00
2	A	1569	8Y6	C7-C11-C8	3.02	130.97	124.32
2	B	1564	8Y6	C7-C11-C8	3.11	131.15	124.32
2	A	1569	8Y6	C17-N2-C16	3.27	124.03	116.60
2	B	1564	8Y6	C17-N2-C16	3.36	124.22	116.60
2	B	1564	8Y6	F2-C12-C22	3.54	120.66	117.00
2	A	1569	8Y6	C21-N3-C20	4.27	123.03	115.91
2	B	1564	8Y6	C21-N3-C20	4.34	123.15	115.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1569	8Y6	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	563/585 (96%)	-0.06	18 (3%)	48 55	14, 21, 42, 121	0
1	B	558/585 (95%)	0.34	49 (8%)	11 14	14, 25, 60, 105	0
All	All	1121/1170 (95%)	0.14	67 (5%)	23 28	14, 23, 50, 121	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	LEU	18.8
1	B	25	PRO	10.9
1	B	23	ILE	10.3
1	B	545	LEU	8.6
1	B	544	GLN	7.2
1	A	568	ARG	6.9
1	B	531	ARG	6.5
1	A	567	ALA	6.3
1	B	548	SER	6.2
1	B	563	SER	5.5
1	B	535	LYS	5.4
1	B	542	ALA	5.3
1	B	402	HIS	5.1
1	B	24	ASN	4.9
1	A	548	SER	4.8
1	B	14	CYS	4.7
1	B	29	SER	4.6
1	B	530	VAL	4.6
1	A	564	LEU	4.5
1	B	16	ALA	4.4
1	B	541	ALA	4.4
1	B	543	SER	4.2
1	B	30	LEU	4.1
1	B	532	THR	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	27	SER	3.9
1	B	502	HIS	3.8
1	B	377	SER	3.6
1	B	540	PRO	3.5
1	A	15	ALA	3.5
1	A	546	ASP	3.4
1	B	547	LEU	3.2
1	B	492	LEU	3.2
1	B	534	LEU	3.2
1	A	212	LYS	3.2
1	B	330	GLN	3.1
1	A	169	VAL	3.1
1	B	378	GLY	3.0
1	A	57	LEU	3.0
1	B	148	GLN	3.0
1	B	536	LEU	2.9
1	A	402	HIS	2.9
1	B	546	ASP	2.9
1	B	212	LYS	2.8
1	B	404	PRO	2.8
1	B	499	THR	2.8
1	A	16	ALA	2.7
1	B	533	LYS	2.7
1	B	34	HIS	2.7
1	B	435	ALA	2.7
1	A	545	LEU	2.7
1	B	21	LEU	2.7
1	B	22	PRO	2.6
1	A	544	GLN	2.6
1	A	352	ASP	2.6
1	B	537	THR	2.5
1	B	498	ARG	2.4
1	A	543	SER	2.4
1	A	566	ARG	2.4
1	B	496	PRO	2.2
1	B	483	ASN	2.2
1	B	31	LEU	2.1
1	B	465	ARG	2.1
1	A	14	CYS	2.1
1	A	547	LEU	2.1
1	B	501	ARG	2.1
1	B	362	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	529	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å ²)	Q<0.9
3	PO4	B	1566	5/5	0.89	0.19	3.48	47,47,49,52	0
3	PO4	A	1571	5/5	0.98	0.09	0.39	24,25,26,27	0
2	8Y6	A	1569	36/36	0.96	0.08	-0.01	13,17,23,24	0
2	8Y6	B	1564	36/36	0.96	0.07	-0.45	17,20,24,25	0
3	PO4	B	1565	5/5	0.96	0.11	-	35,38,39,43	0
3	PO4	A	1570	5/5	0.95	0.15	-	32,38,40,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.