



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 05:08 AM GMT

PDB ID : 4MZ4
Title : Discovery of an Irreversible HCV NS5B Polymerase Inhibitor
Authors : Zeng, Q.; Anilkumar, G.N.; Rosenblum, S.B.; Huang, H.-C.; Lesburg, C.A.; Jiang, Y.; Selyutin, O.; Chan, T.-Y.; Bennett, F.; Chen, K.X.; Venkatraman, S.; Sannigrahi, M.; Velazquez, F.; Duca, J.S.; Gavalas, S.; Huang, Y.; Pu, H.; Wang, L.; Pinto, P.; Vibulbhan, B.; Agrawal, S.; Ferrari, E.; Jiang, C.-K.; Li, C.; Hesk, D.; Gesell, J.; Sorota, S.; Shih, N.-Y.; Njoroge, F.G.; Kozlowski, J.A.
Deposited on : 2013-09-29
Resolution : 1.63 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

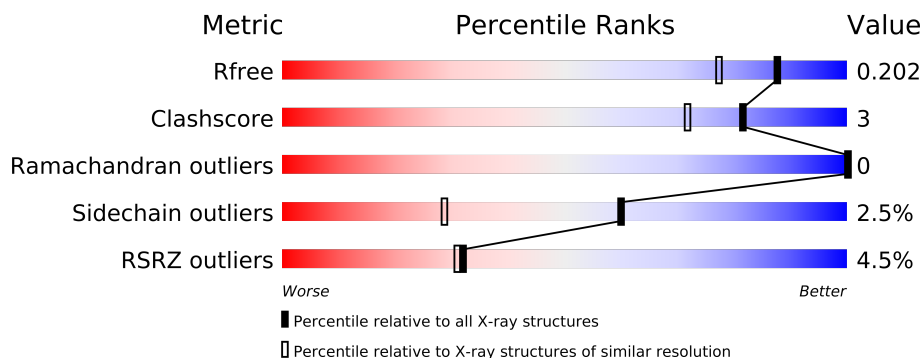
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.63 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}	66092	1326 (1.66-1.62)
Clashscore	79885	1525 (1.66-1.62)
Ramachandran outliers	78287	1490 (1.66-1.62)
Sidechain outliers	78261	1490 (1.66-1.62)
RSRZ outliers	66119	1326 (1.66-1.62)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	576	
1	B	576	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	PO4	A	603	-	X
3	PO4	A	604	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10615 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

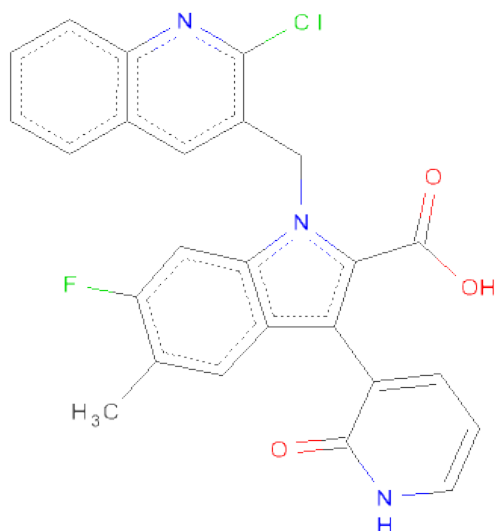
- Molecule 1 is a protein called RNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	563	Total	C	N	O	S	0	12	0
			4468	2808	792	832	36			
1	B	560	Total	C	N	O	S	0	8	0
			4412	2777	779	821	35			

There are 16 discrepancies between the modelled and reference sequences:

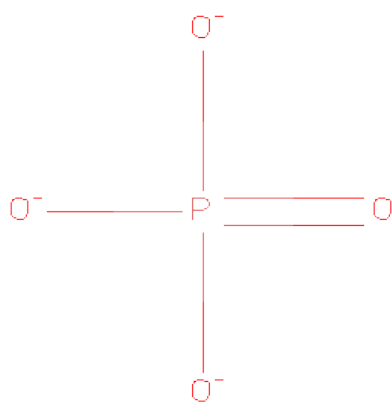
Chain	Residue	Modelled	Actual	Comment	Reference
A	440	GLY	GLU	CONFLICT	UNP O92972
A	520	ILE	THR	CONFLICT	UNP O92972
A	571	GLU	-	EXPRESSION TAG	UNP O92972
A	572	ASN	-	EXPRESSION TAG	UNP O92972
A	573	LEU	-	EXPRESSION TAG	UNP O92972
A	574	TYR	-	EXPRESSION TAG	UNP O92972
A	575	PHE	-	EXPRESSION TAG	UNP O92972
A	576	GLN	-	EXPRESSION TAG	UNP O92972
B	440	GLY	GLU	CONFLICT	UNP O92972
B	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

- Molecule 2 is 1-[(2-CHLOROQUINOLIN-3-YL)METHYL]-6-FLUORO-5-METHYL-3-(2-OXO-1,2-DIHYDROPYRIDIN-3-YL)-1H-INDOLE-2-CARBOXYLICACID (three-letter code: 2F3) (formula: C₂₅H₁₇ClFN₃O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			32	25	1	3	3		
2	B	1	Total	C	F	N	O	0	0
			32	25	1	3	3		

- 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		

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

- Molecule 4 is water.

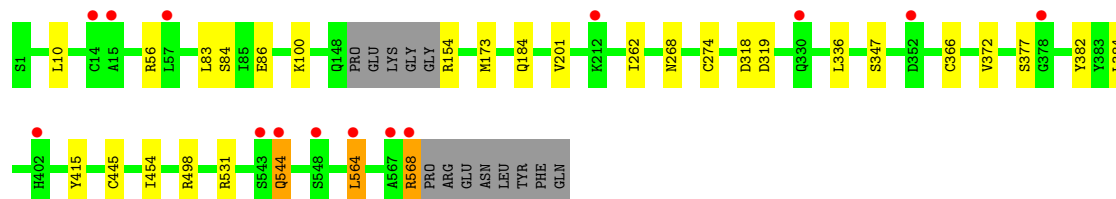
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	902	Total	O	0	0
			902	902		
4	B	754	Total	O	0	0
			754	754		

3 Residue-property plots i

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

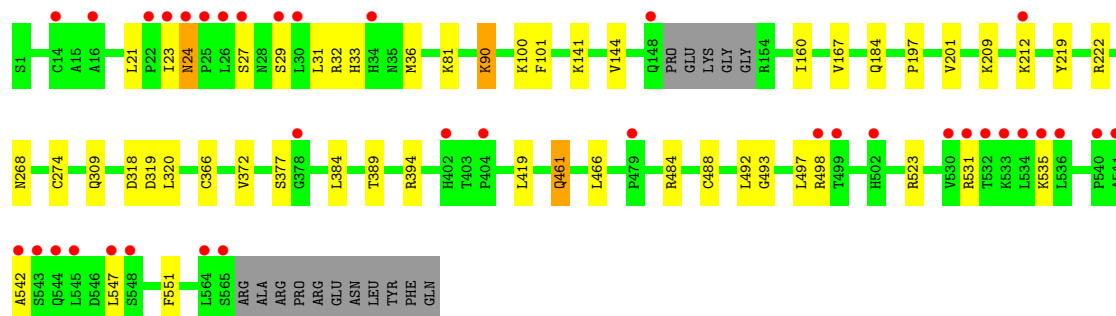
- Molecule 1: RNA-directed RNA polymerase

Chain A:



- Molecule 1: RNA-directed RNA polymerase

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.99Å 106.60Å 134.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 1.63 19.82 – 1.63	Depositor EDS
% Data completeness (in resolution range)	99.0 (19.92-1.63) 99.3 (19.82-1.63)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.91 (at 1.63Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.175 , 0.204 0.172 , 0.202	Depositor DCC
R_{free} test set	8018 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	19.0	Xtriage
Anisotropy	0.404	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 159667 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10615	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.04% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 2F3, PO4

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	0.53	0/4562	0.62	0/6189
1	B	0.50	0/4506	0.64	0/6114
All	All	0.52	0/9068	0.63	0/12303

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4468	0	4474	24	0
1	B	4412	0	4421	32	0
2	A	32	0	16	9	0
2	B	32	0	16	5	0
3	A	15	0	0	0	0
4	A	902	0	0	6	0
4	B	754	0	0	8	0
All	All	10615	0	8927	56	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including

hydrogens) of the entry. The overall clashscore for this entry is 3.

All (56) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:366[A]:CYS:SG	2:A:601:2F3:C18	2.50	0.98
1:B:24:ASN:HD21	1:B:27:SER:HB3	1.35	0.92
1:A:366[A]:CYS:SG	2:A:601:2F3:H12	2.13	0.88
1:A:366[B]:CYS:HG	2:A:601:2F3:C9	1.77	0.76
1:B:201[A]:VAL:HG12	1:B:384:LEU:HG	1.69	0.74
1:B:201[A]:VAL:CG1	1:B:384:LEU:HG	2.19	0.73
1:B:201[B]:VAL:HG22	1:B:384:LEU:HG	1.71	0.71
1:B:366[A]:CYS:SG	2:B:601:2F3:C18	2.80	0.69
1:A:366[A]:CYS:SG	2:A:601:2F3:C14	2.84	0.65
1:A:201[B]:VAL:HG22	1:A:384:LEU:HG	1.80	0.62
1:B:523:ARG:NH2	1:B:535:LYS:HB3	2.16	0.61
1:B:523:ARG:HH21	1:B:535:LYS:HB3	1.67	0.59
1:B:461:GLN:HB3	1:B:542:ALA:HA	1.84	0.58
1:B:366[A]:CYS:SG	2:B:601:2F3:H12	2.43	0.57
1:A:262:ILE:HD11	4:A:1249:HOH:O	2.04	0.56
1:A:268:ASN:HB3	1:A:274[B]:CYS:SG	2.44	0.56
1:A:498:ARG:NH1	4:A:1027:HOH:O	2.38	0.55
1:B:498:ARG:NH1	4:B:1007:HOH:O	2.40	0.54
1:B:141:LYS:HD2	1:B:160:ILE:HD11	1.90	0.54
1:A:347[B]:SER:OG	1:A:347[B]:SER:O	2.24	0.54
1:A:531:ARG:HD2	4:A:1139:HOH:O	2.09	0.53
1:B:389:THR:HG23	1:B:492:LEU:HD21	1.89	0.53
1:B:488:CYS:HB2	4:B:1248:HOH:O	2.08	0.53
1:A:366[B]:CYS:SG	2:A:601:2F3:N1	2.66	0.52
1:A:366[B]:CYS:SG	2:A:601:2F3:C7	2.88	0.52
1:B:319:ASP:HB2	2:B:601:2F3:H11	1.91	0.52
1:B:309:GLN:HG2	4:B:1211:HOH:O	2.10	0.51
1:B:219:TYR:HB3	1:B:320:LEU:HD23	1.93	0.51
1:A:201[B]:VAL:CG2	1:A:384:LEU:HG	2.42	0.50
1:A:445:CYS:SG	1:A:454:ILE:HD12	2.52	0.49
1:A:544:GLN:HG3	4:A:1404:HOH:O	2.12	0.49
1:A:318:ASP:OD1	2:A:601:2F3:H10	2.13	0.48
1:B:100:LYS:HG3	1:B:101:PHE:CD1	2.48	0.48
1:B:100:LYS:NZ	4:B:1187:HOH:O	2.46	0.48
1:A:568:ARG:HG3	4:A:1178:HOH:O	2.14	0.48
1:A:319:ASP:HB2	2:A:601:2F3:C6	2.43	0.48
1:A:84:SER:OG	1:A:86:GLU:HG2	2.14	0.48
1:B:24:ASN:ND2	1:B:27:SER:HB3	2.17	0.48
1:B:466:LEU:HD22	1:B:551:PHE:HE2	1.78	0.48
1:B:318:ASP:OD1	2:B:601:2F3:H10	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:144:VAL:HB	1:B:394:ARG:HG2	1.98	0.46
1:B:531:ARG:NH2	4:B:912:HOH:O	2.49	0.45
1:B:268:ASN:HB3	1:B:274:CYS:SG	2.57	0.45
1:B:419:LEU:HD21	1:B:497:LEU:CD1	2.47	0.45
1:A:154:ARG:N	4:A:1470:HOH:O	2.50	0.44
1:A:415:TYR:CE1	2:A:601:2F3:H18	2.53	0.43
1:B:222:ARG:HG2	4:B:1199:HOH:O	2.17	0.43
1:A:83:LEU:HB2	1:A:173:MET:HA	2.00	0.43
1:A:564:LEU:HA	1:A:564:LEU:HD22	1.87	0.43
1:B:167:VAL:HG23	4:B:1381:HOH:O	2.19	0.42
1:B:90:LYS:HE3	4:B:1364:HOH:O	2.20	0.42
1:B:197:PRO:O	1:B:201[B]:VAL:HG23	2.19	0.42
1:B:366[A]:CYS:SG	2:B:601:2F3:C14	3.07	0.42
1:A:372:VAL:HG22	1:A:382:TYR:CD1	2.56	0.40
1:B:33:HIS:HB3	1:B:36:MET:HG3	2.04	0.40
1:B:32:ARG:HD3	1:B:493:GLY:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	571/576 (99%)	563 (99%)	8 (1%)	0	100	100
1	B	564/576 (98%)	551 (98%)	13 (2%)	0	100	100
All	All	1135/1152 (98%)	1114 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	491/490 (100%)	482 (98%)	9 (2%)	71	46
1	B	485/490 (99%)	470 (97%)	15 (3%)	52	19
All	All	976/980 (100%)	952 (98%)	24 (2%)	60	29

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	56	ARG
1	A	100	LYS
1	A	184	GLN
1	A	336	LEU
1	A	377	SER
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	29	SER
1	B	31	LEU
1	B	81	LYS
1	B	90	LYS
1	B	184	GLN
1	B	209	LYS
1	B	212	LYS
1	B	372	VAL
1	B	377	SER
1	B	461	GLN
1	B	484	ARG
1	B	547	LEU

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 chains 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 ⓘ

5 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	2F3	A	601	1	35,36,37	20.46	10 (28%)	47,53,55	1.58	8 (17%)
3	PO4	A	602	-	4,4,4	1.61	1 (25%)	6,6,6	0.31	0
3	PO4	A	603	-	4,4,4	1.02	0	6,6,6	0.30	0
3	PO4	A	604	-	4,4,4	0.45	0	6,6,6	0.30	0
2	2F3	B	601	1	35,36,37	33.15	9 (25%)	47,53,55	1.46	6 (12%)

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	2F3	A	601	1	-	0/12/12/12	0/1/5/5
3	PO4	A	602	-	-	0/0/0/0	0/0/0/0
3	PO4	A	603	-	-	0/0/0/0	0/0/0/0
3	PO4	A	604	-	-	0/0/0/0	0/0/0/0
2	2F3	B	601	1	-	0/12/12/12	0/1/5/5

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	2F3	C9-C7	195.80	1.40	1.39
2	A	601	2F3	C9-C7	120.58	1.39	1.39
2	B	601	2F3	C10-C16	-4.47	1.43	1.50
2	B	601	2F3	C18-C17	4.30	1.40	1.35
2	A	601	2F3	C10-C16	-4.21	1.44	1.50
2	A	601	2F3	O3-C19	3.91	1.35	1.23
2	A	601	2F3	C18-C17	3.80	1.39	1.35
2	B	601	2F3	O3-C19	3.69	1.34	1.23
2	A	601	2F3	C21-C10	3.37	1.52	1.43
2	B	601	2F3	C21-C10	3.25	1.52	1.43
2	A	601	2F3	C23-C24	3.17	1.46	1.39
2	B	601	2F3	C12-C17	3.04	1.40	1.37
2	A	601	2F3	C20-C7	-2.93	1.48	1.51
2	B	601	2F3	C20-C7	-2.72	1.48	1.51
3	A	602	PO4	P-O4	2.57	1.63	1.52
2	A	601	2F3	O1-C19	-2.55	1.22	1.30
2	A	601	2F3	C12-C17	2.48	1.40	1.37
2	B	601	2F3	O1-C19	-2.47	1.22	1.30
2	B	601	2F3	C23-C24	2.45	1.44	1.39
2	A	601	2F3	C22-N3	2.02	1.37	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	2F3	C11-C12-C17	4.40	119.11	116.12
2	B	601	2F3	C11-C12-C17	4.14	118.94	116.12
2	A	601	2F3	C15-N2-C14	3.91	110.79	106.78
2	B	601	2F3	C15-N2-C14	3.48	110.36	106.78
2	A	601	2F3	C25-C12-C17	-3.31	119.43	121.68
2	A	601	2F3	C7-C20-N2	2.75	117.15	112.81
2	B	601	2F3	C24-C10-C16	2.68	124.12	119.01
2	A	601	2F3	C16-C15-C19	-2.54	125.91	130.71
2	B	601	2F3	C16-C15-C19	-2.43	126.11	130.71
2	B	601	2F3	C7-C20-N2	2.35	116.52	112.81
2	A	601	2F3	C24-C10-C16	2.24	123.28	119.01
2	B	601	2F3	F1-C17-C18	-2.20	117.65	120.07
2	A	601	2F3	C16-C15-N2	-2.19	106.08	108.24
2	A	601	2F3	F1-C17-C18	-2.09	117.77	120.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	563/576 (97%)	-0.22	14 (2%) 54 55	13, 19, 39, 83	0
1	B	560/576 (97%)	0.12	37 (6%) 18 16	13, 24, 57, 95	0
All	All	1123/1152 (97%)	-0.05	51 (4%) 32 31	13, 21, 48, 95	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	LEU	15.8
1	B	545	LEU	10.5
1	B	23	ILE	9.3
1	B	25	PRO	8.8
1	B	565	SER	7.5
1	A	568	ARG	7.5
1	B	531	ARG	6.6
1	B	544	GLN	6.2
1	A	567	ALA	6.0
1	B	535	LYS	5.7
1	B	548	SER	5.2
1	B	541	ALA	4.7
1	B	502	HIS	4.7
1	B	534	LEU	4.5
1	B	542	ALA	4.4
1	B	24	ASN	4.4
1	B	532	THR	4.2
1	B	543	SER	4.1
1	A	564	LEU	4.0
1	B	14	CYS	3.9
1	A	548	SER	3.8
1	B	29	SER	3.8
1	B	30	LEU	3.7
1	B	536	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	564	LEU	3.6
1	B	530	VAL	3.5
1	B	16	ALA	3.5
1	A	212	LYS	3.4
1	B	402	HIS	3.3
1	B	540	PRO	3.3
1	B	27	SER	3.3
1	B	547	LEU	3.2
1	B	404	PRO	3.1
1	B	499	THR	3.1
1	B	479	PRO	3.1
1	A	15	ALA	2.9
1	A	402	HIS	2.8
1	A	544	GLN	2.8
1	A	57	LEU	2.8
1	B	22	PRO	2.7
1	A	352	ASP	2.7
1	A	14	CYS	2.7
1	B	533	LYS	2.6
1	B	498	ARG	2.5
1	B	212	LYS	2.4
1	B	34	HIS	2.4
1	A	543	SER	2.4
1	B	148	GLN	2.4
1	B	378	GLY	2.2
1	A	330	GLN	2.0
1	A	378	GLY	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 carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	PO4	A	603	5/5	0.20	5.60	36,39,44,46	0
3	PO4	A	604	5/5	0.14	2.74	66,68,70,71	0
2	2F3	A	601	32/33	0.10	1.34	15,19,43,45	0
2	2F3	B	601	32/33	0.09	1.23	18,22,34,35	0
3	PO4	A	602	5/5	0.09	0.51	21,22,22,24	0

6.5 Other polymers ⓘ

There are no such residues in this entry.