



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 07:52 pm BST

PDB ID : 2JLW
Title : Dengue virus 4 NS3 helicase in complex with ssRNA2
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Deposited on : 2008-09-15
Resolution : 2.60 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 : 2.11
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) : 2.11

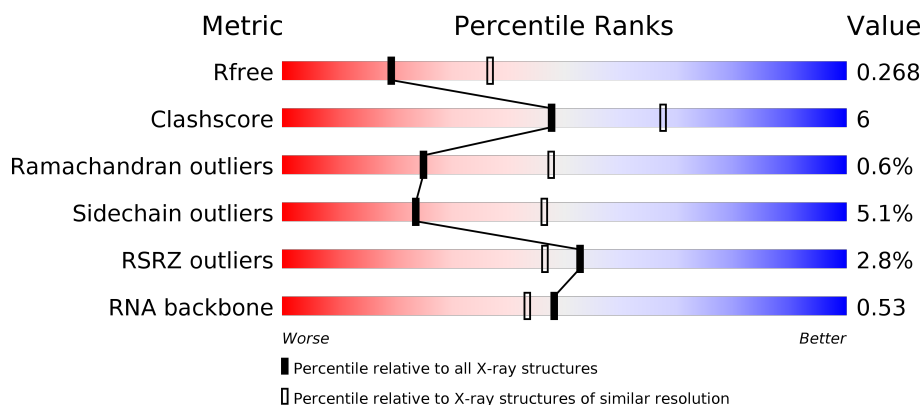
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)
RNA backbone	3102	1040 (2.90-2.30)

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 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	451	<div> <div>3%</div> <div>82%</div> <div>16%</div> <div>.</div> </div>
1	B	451	<div> <div>2%</div> <div>81%</div> <div>18%</div> <div>.</div> </div>
2	C	13	<div> <div>15%</div> <div>31%</div> <div>8%</div> <div>15%</div> <div>46%</div> </div>
2	D	13	<div> <div>15%</div> <div>23%</div> <div>15%</div> <div>8%</div> <div>8%</div> <div>46%</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7564 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 SERINE PROTEASE SUBUNIT NS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	0	0
			3591	2264	640	670	17			
1	B	449	Total	C	N	O	S	0	0	0
			3572	2254	634	667	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	250	ASP	GLU	conflict	UNP Q2YHF0
A	292	CYS	SER	conflict	UNP Q2YHF0
A	321	SER	THR	conflict	UNP Q2YHF0
A	322	ILE	THR	conflict	UNP Q2YHF0
A	381	ARG	LYS	conflict	UNP Q2YHF0
A	480	LYS	ARG	conflict	UNP Q2YHF0
B	250	ASP	GLU	conflict	UNP Q2YHF0
B	292	CYS	SER	conflict	UNP Q2YHF0
B	321	SER	THR	conflict	UNP Q2YHF0
B	322	ILE	THR	conflict	UNP Q2YHF0
B	381	ARG	LYS	conflict	UNP Q2YHF0
B	480	LYS	ARG	conflict	UNP Q2YHF0

- Molecule 2 is a RNA chain called 5'-R(*UP*AP*GP*AP*CP*UP*AP*AP*CP*AP*AP*CP*U)-3'.

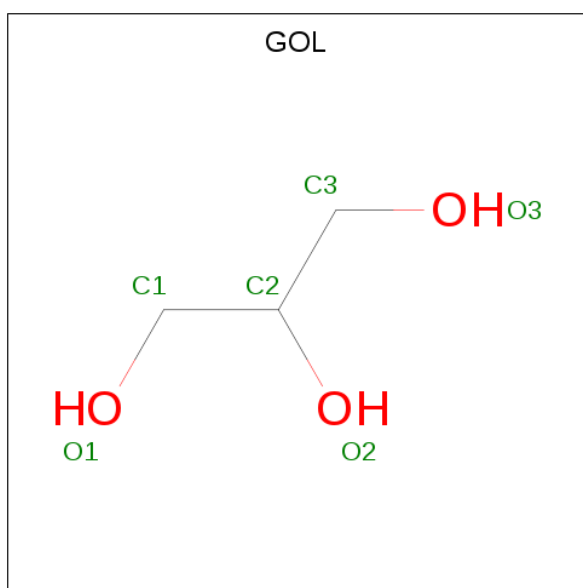
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	7	Total	C	N	O	P	0	0	0
			128	57	22	43	6			
2	D	7	Total	C	N	O	P	0	0	0
			128	57	22	43	6			

- 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	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

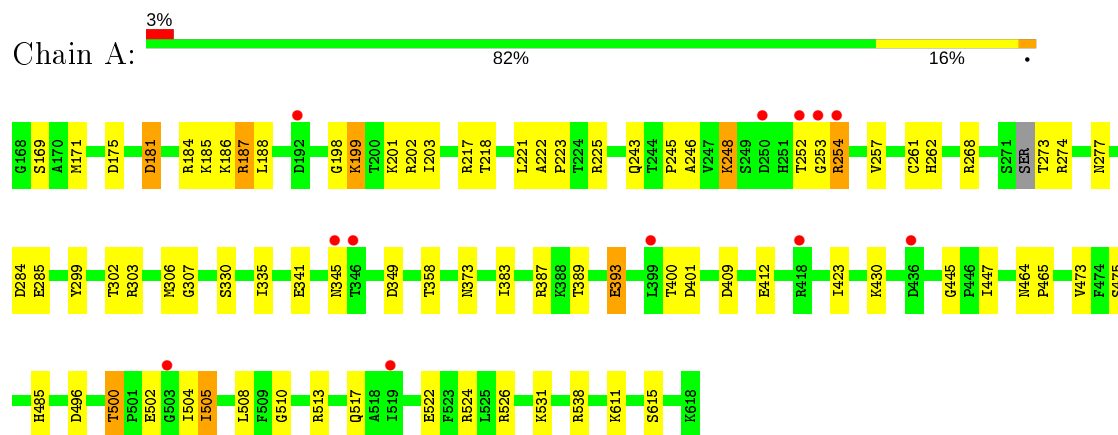
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	66	Total 66	O 66	0	0
5	B	62	Total 62	O 62	0	0
5	C	1	Total 1	O 1	0	0

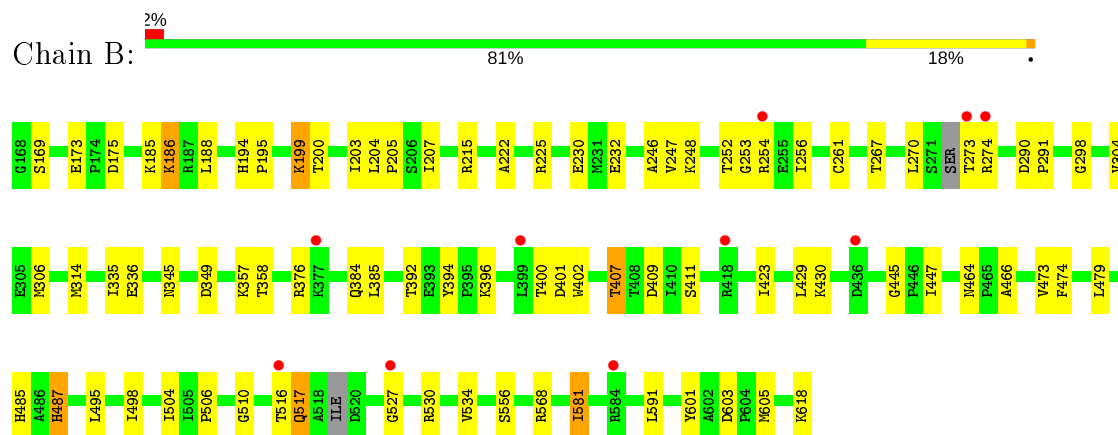
3 Residue-property plots [i](#)

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

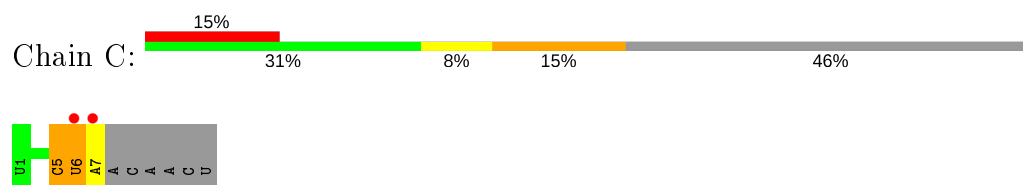
• Molecule 1: SERINE PROTEASE SUBUNIT NS3



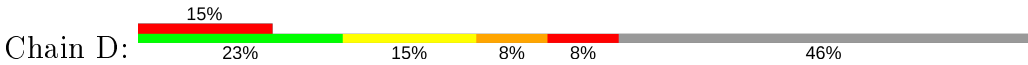
• Molecule 1: SERINE PROTEASE SUBUNIT NS3



• Molecule 2: 5'-R(*UP*AP*GP*AP*CP*UP*AP*AP*CP*AP*AP*CP*U)-3'



• Molecule 2: 5'-R(*UP*AP*GP*AP*CP*UP*AP*AP*CP*AP*AP*CP*U)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	132.43 Å 104.63 Å 72.32 Å 90.00° 117.67° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60 19.77 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.60) 100.0 (19.77-2.60)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.59 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.199 , 0.270 0.198 , 0.268	Depositor DCC
R_{free} test set	1348 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7564	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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: GOL, 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	0.41	0/3671	0.56	0/4969
1	B	0.40	0/3651	0.57	0/4941
2	C	0.81	0/142	2.26	4/220 (1.8%)
2	D	0.83	1/142 (0.7%)	2.18	3/220 (1.4%)
All	All	0.43	1/7606 (0.0%)	0.71	7/10350 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	7	A	P-O5'	-5.13	1.54	1.59

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	7	A	O5'-P-OP1	-25.27	80.38	110.70
2	D	7	A	O5'-P-OP2	-24.20	81.66	110.70
2	D	7	A	OP1-P-OP2	8.38	132.16	119.60
2	C	7	A	OP1-P-OP2	7.50	130.85	119.60
2	C	6	U	O3'-P-O5'	7.41	118.07	104.00
2	D	6	U	O3'-P-O5'	7.24	117.75	104.00
2	C	6	U	P-O3'-C3'	6.79	127.85	119.70

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	3591	0	3568	43	0
1	B	3572	0	3543	48	1
2	C	128	0	65	1	0
2	D	128	0	65	2	0
3	A	5	0	0	1	0
3	B	5	0	0	0	0
4	B	6	0	8	0	0
5	A	66	0	0	2	0
5	B	62	0	0	1	0
5	C	1	0	0	0	0
All	All	7564	0	7249	89	1

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 (89) 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:B:407:THR:HG21	1:B:411:SER:OG	1.70	0.91
1:B:430:LYS:HD3	1:B:447:ILE:HD11	1.61	0.80
1:A:252:THR:O	1:A:254:ARG:N	2.29	0.65
1:A:199:LYS:HE2	1:A:284:ASP:OD1	1.95	0.65
1:B:376:ARG:HH22	1:B:384:GLN:HE21	1.46	0.63
1:B:252:THR:O	1:B:254:ARG:N	2.31	0.62
1:B:199:LYS:HD2	1:B:200:THR:HG23	1.81	0.62
1:A:221:LEU:HB3	1:A:262:HIS:HD2	1.65	0.62
1:B:169:SER:HB3	1:B:175:ASP:HA	1.80	0.62
1:A:245:PRO:O	1:B:248:LYS:HG2	1.99	0.61
1:A:169:SER:HB3	1:A:175:ASP:HA	1.84	0.60
1:A:445:GLY:HA2	1:A:447:ILE:HG23	1.84	0.59
1:A:409:ASP:HB2	1:A:412:GLU:HB2	1.84	0.58
1:B:409:ASP:OD2	2:D:2:A:O2'	2.21	0.58
1:A:505:ILE:HG22	5:A:2047:HOH:O	2.03	0.58
1:B:396:LYS:O	1:B:400:THR:HB	2.04	0.58
1:A:500:THR:HG23	1:A:502:GLU:H	1.68	0.57
1:B:306:MET:SD	1:B:504:ILE:HD12	2.45	0.57
1:A:185:LYS:O	1:A:186:LYS:HB2	2.05	0.57
1:A:383:ILE:HG23	1:A:393:GLU:HG3	1.87	0.57
1:A:524:ARG:HD2	1:A:526:ARG:NH1	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:530:ARG:O	1:B:534:VAL:HG23	2.05	0.57
1:B:204:LEU:HB3	1:B:205:PRO:HD3	1.88	0.56
1:A:496:ASP:HA	1:A:505:ILE:HD11	1.88	0.55
1:B:298:GLY:HA3	1:B:495:LEU:HD21	1.86	0.55
1:B:581:ILE:HD13	1:B:591:LEU:HB2	1.88	0.55
1:A:243:GLN:OE1	1:A:268:ARG:NH2	2.39	0.55
1:A:430:LYS:HG3	1:A:485:HIS:CD2	2.42	0.55
1:A:223:PRO:HG3	1:A:285:GLU:HG3	1.88	0.54
1:B:385:LEU:HB2	1:B:407:THR:HG23	1.88	0.54
1:A:221:LEU:HB3	1:A:262:HIS:CD2	2.42	0.54
1:A:400:THR:HG22	1:A:401:ASP:N	2.23	0.53
2:D:6:U:H4'	2:D:7:A:OP1	2.08	0.52
1:B:222:ALA:O	1:B:261:CYS:HA	2.09	0.52
1:B:485:HIS:CE1	1:B:487:HIS:CD2	2.99	0.51
1:A:248:LYS:NZ	1:B:248:LYS:HG3	2.26	0.51
1:A:222:ALA:O	1:A:261:CYS:HA	2.11	0.50
1:A:500:THR:HG22	1:A:504:ILE:H	1.77	0.50
1:A:389:THR:O	1:A:393:GLU:HB2	2.12	0.50
1:B:203:ILE:O	1:B:207:ILE:HG12	2.12	0.49
1:A:225:ARG:HD3	2:C:5:C:OP2	2.13	0.49
1:B:194:HIS:HD2	1:B:195:PRO:O	1.96	0.49
1:B:407:THR:HG21	1:B:411:SER:HG	1.71	0.49
1:B:290:ASP:OD1	1:B:291:PRO:HD2	2.13	0.48
1:A:273:THR:HA	1:A:274:ARG:HA	1.59	0.48
1:B:306:MET:SD	1:B:504:ILE:CD1	3.02	0.48
1:B:267:THR:HA	1:B:270:LEU:HD12	1.96	0.47
1:B:445:GLY:HA2	1:B:447:ILE:HG23	1.97	0.47
1:B:199:LYS:HG2	1:B:314:MET:CE	2.44	0.46
1:A:268:ARG:HD2	5:A:2014:HOH:O	2.14	0.46
1:B:232:GLU:HG3	1:B:247:VAL:CG1	2.46	0.46
1:B:603:ASP:OD2	1:B:605:MET:HB2	2.15	0.46
1:B:358:THR:HG21	1:B:423:ILE:HD12	1.97	0.46
1:B:225:ARG:HB3	1:B:246:ALA:HB2	1.98	0.46
1:A:299:TYR:O	1:A:303:ARG:HG2	2.16	0.46
1:A:358:THR:HG21	1:A:423:ILE:HD12	1.97	0.45
1:A:302:THR:O	1:A:306:MET:HG3	2.16	0.45
1:B:429:LEU:HD23	1:B:445:GLY:HA3	1.99	0.45
1:A:181:ASP:HA	1:A:184:ARG:HD2	1.98	0.45
1:A:198:GLY:O	1:A:202:ARG:HB2	2.16	0.45
1:B:185:LYS:O	1:B:186:LYS:CB	2.65	0.44
1:A:199:LYS:HA	1:A:203:ILE:HD12	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:498:ILE:O	1:B:506:PRO:HD3	2.17	0.44
1:B:230:GLU:OE2	1:B:394:TYR:OH	2.35	0.44
1:A:252:THR:C	1:A:254:ARG:H	2.21	0.44
1:B:357:LYS:HB3	1:B:402:TRP:CZ3	2.53	0.44
1:A:225:ARG:HB3	1:A:246:ALA:HB2	1.99	0.43
1:B:464:ASN:HD21	1:B:466:ALA:HB3	1.82	0.43
1:A:508:LEU:O	1:A:513:ARG:NE	2.42	0.43
1:B:336:GLU:HA	1:B:474:PHE:O	2.18	0.43
1:A:248:LYS:HZ2	1:B:248:LYS:HG3	1.83	0.43
1:A:218:THR:HB	1:A:257:VAL:HG22	2.01	0.43
1:B:273:THR:HA	1:B:274:ARG:HA	1.54	0.43
1:B:430:LYS:HG2	1:B:485:HIS:CD2	2.54	0.43
1:A:277:ASN:OD1	1:A:303:ARG:NH2	2.52	0.42
1:B:517:GLN:H	1:B:517:GLN:CD	2.23	0.42
1:A:335:ILE:O	1:A:473:VAL:HA	2.19	0.42
1:A:500:THR:HG22	1:A:504:ILE:N	2.34	0.42
1:B:335:ILE:O	1:B:473:VAL:HA	2.20	0.42
1:B:252:THR:C	1:B:254:ARG:H	2.23	0.42
1:B:215:ARG:NH1	1:B:256:ILE:HD13	2.35	0.42
1:B:618:LYS:NZ	5:B:2062:HOH:O	2.53	0.42
1:A:464:ASN:HA	1:A:465:PRO:HD2	1.96	0.41
1:B:186:LYS:HA	1:B:304:VAL:O	2.20	0.41
1:A:184:ARG:O	1:A:187:ARG:HB3	2.21	0.41
1:B:385:LEU:CB	1:B:407:THR:HG23	2.50	0.41
1:A:199:LYS:HB3	3:A:1619:PO4:O4	2.20	0.41
1:A:186:LYS:HG3	1:A:307:GLY:HA2	2.03	0.40
1:B:400:THR:CG2	1:B:401:ASP:N	2.85	0.40

All (1) 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)
1:B:173:GLU:OE1	1:B:601:TYR:OH[4_454]	2.16	0.04

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	446/451 (99%)	425 (95%)	19 (4%)	2 (0%)	34	57
1	B	443/451 (98%)	420 (95%)	20 (4%)	3 (1%)	22	43
All	All	889/902 (99%)	845 (95%)	39 (4%)	5 (1%)	25	47

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	253	GLY
1	B	253	GLY
1	B	510	GLY
1	A	510	GLY
1	B	527	GLY

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	384/388 (99%)	359 (94%)	25 (6%)	17	34
1	B	381/388 (98%)	367 (96%)	14 (4%)	34	60
All	All	765/776 (99%)	726 (95%)	39 (5%)	24	46

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	171	MET
1	A	181	ASP
1	A	187	ARG
1	A	188	LEU
1	A	199	LYS
1	A	201	LYS
1	A	217	ARG

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Mol	Chain	Res	Type
1	A	248	LYS
1	A	254	ARG
1	A	330	SER
1	A	341	GLU
1	A	345	ASN
1	A	349	ASP
1	A	373	ASN
1	A	387	ARG
1	A	393	GLU
1	A	475	SER
1	A	500	THR
1	A	505	ILE
1	A	517	GLN
1	A	522	GLU
1	A	531	LYS
1	A	538	ARG
1	A	611	LYS
1	A	615	SER
1	B	186	LYS
1	B	188	LEU
1	B	199	LYS
1	B	345	ASN
1	B	349	ASP
1	B	392	THR
1	B	407	THR
1	B	479	LEU
1	B	487	HIS
1	B	516	THR
1	B	517	GLN
1	B	556	SER
1	B	568	ARG
1	B	581	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	HIS
1	A	251	HIS
1	A	529	GLN
1	B	194	HIS
1	B	384	GLN
1	B	464	ASN

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Mol	Chain	Res	Type
1	B	467	GLN
1	B	487	HIS
1	B	529	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	5/13 (38%)	2 (40%)	1 (20%)
2	D	5/13 (38%)	2 (40%)	1 (20%)
All	All	10/26 (38%)	4 (40%)	2 (20%)

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	5	C
2	C	6	U
2	D	5	C
2	D	6	U

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	6	U
2	D	6	U

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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
4	GOL	B	1619	-	5,5,5	0.37	0	5,5,5	0.34	0
3	PO4	B	1620	-	4,4,4	0.84	0	6,6,6	0.40	0
3	PO4	A	1619	-	4,4,4	0.88	0	6,6,6	0.65	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
4	GOL	B	1619	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1619	GOL	O1-C1-C2-O2
4	B	1619	GOL	O1-C1-C2-C3
4	B	1619	GOL	C1-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1619	PO4	1	0

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	450/451 (99%)	-0.12	12 (2%) 54 48	24, 33, 42, 46	4 (0%)
1	B	449/451 (99%)	-0.15	10 (2%) 62 56	24, 33, 41, 46	4 (0%)
2	C	7/13 (53%)	0.45	2 (28%) 0 0	36, 36, 67, 70	0
2	D	7/13 (53%)	0.72	2 (28%) 0 0	36, 36, 67, 70	0
All	All	913/928 (98%)	-0.12	26 (2%) 53 46	24, 33, 42, 70	8 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	7	A	4.6
2	D	6	U	3.7
2	C	7	A	3.6
1	B	399	LEU	3.3
1	B	273	THR	3.2
1	A	254	ARG	3.2
1	A	345	ASN	3.1
1	B	436	ASP	3.0
1	A	418	ARG	2.9
1	A	436	ASP	2.9
1	A	399	LEU	2.8
1	A	346	THR	2.7
2	C	6	U	2.7
1	B	254	ARG	2.7
1	A	252	THR	2.6
1	A	503	GLY	2.6
1	B	584	ARG	2.5
1	A	253	GLY	2.4
1	B	377	LYS	2.4
1	B	418	ARG	2.4
1	A	192	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	519	ILE	2.3
1	B	274	ARG	2.2
1	B	527	GLY	2.2
1	B	516	THR	2.1
1	A	250	ASP	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. 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	GOL	B	1619	6/6	0.93	0.22	61,61,61,62	0
3	PO4	B	1620	5/5	0.95	0.17	64,65,65,65	0
3	PO4	A	1619	5/5	0.96	0.12	60,60,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.