



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:57 AM GMT

PDB ID : 3GNW
Title : HCV NS5B polymerase in complex with 1,5 benzodiazepine inhibitor 4c
Authors : Nyanguile, O.; De Bondt, H.
Deposited on : 2009-03-18
Resolution : 2.39 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

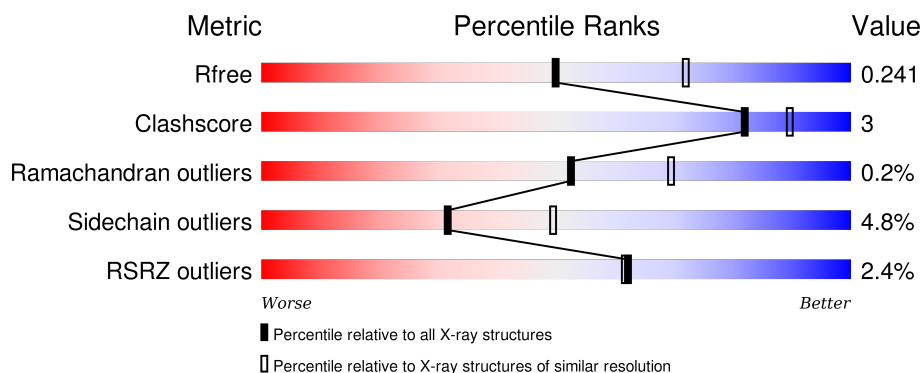
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	581	<div> <div></div> <div>85%</div> <div>11%</div> <div>• •</div> </div>
1	B	581	<div> <div>4%</div> <div>88%</div> <div>8%</div> <div>• •</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9386 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.

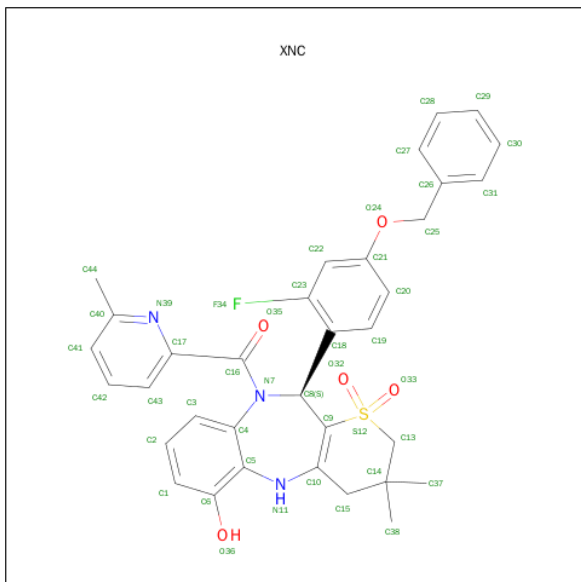
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	560	Total	C	N	O	S	0	2	0
			4368	2750	772	813	33			
1	B	568	Total	C	N	O	S	0	1	0
			4429	2788	788	821	32			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	EXPRESSION TAG	UNP O92972
A	-1	ALA	-	EXPRESSION TAG	UNP O92972
A	0	SER	-	EXPRESSION TAG	UNP O92972
A	571	LEU	-	EXPRESSION TAG	UNP O92972
A	572	GLU	-	EXPRESSION TAG	UNP O92972
A	573	HIS	-	EXPRESSION TAG	UNP O92972
A	574	HIS	-	EXPRESSION TAG	UNP O92972
A	575	HIS	-	EXPRESSION TAG	UNP O92972
A	576	HIS	-	EXPRESSION TAG	UNP O92972
A	577	HIS	-	EXPRESSION TAG	UNP O92972
A	578	HIS	-	EXPRESSION TAG	UNP O92972
B	-2	MET	-	EXPRESSION TAG	UNP O92972
B	-1	ALA	-	EXPRESSION TAG	UNP O92972
B	0	SER	-	EXPRESSION TAG	UNP O92972
B	571	LEU	-	EXPRESSION TAG	UNP O92972
B	572	GLU	-	EXPRESSION TAG	UNP O92972
B	573	HIS	-	EXPRESSION TAG	UNP O92972
B	574	HIS	-	EXPRESSION TAG	UNP O92972
B	575	HIS	-	EXPRESSION TAG	UNP O92972
B	576	HIS	-	EXPRESSION TAG	UNP O92972
B	577	HIS	-	EXPRESSION TAG	UNP O92972
B	578	HIS	-	EXPRESSION TAG	UNP O92972

- Molecule 2 is (11S)-11-[4-(BENZYLOXY)-2-FLUOROPHENYL]-3,3-DIMETHYL-10-[(6-METHYLPYRIDIN-2-YL)CARBONYL]-2,3,4,5,10,11-HEXAHYDROTHIOPYRAN

O[3,2-B][1,5]BENZODIAZEPIN-6-OL 1,1-DIOXIDE (three-letter code: XNC) (formula: $C_{34}H_{32}FN_3O_5S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			44	34	1	3	5	1		
2	B	1	Total	C	F	N	O	S	0	0
			44	34	1	3	5	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	2	Total	Cl	0	0
			2	2		

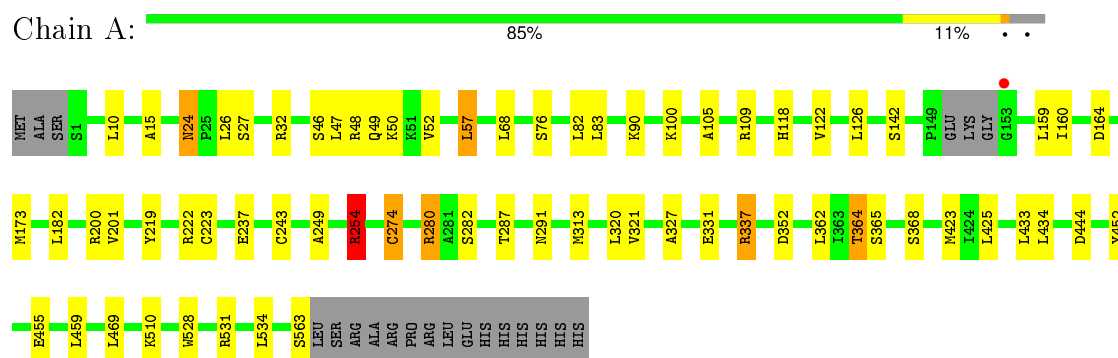
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	288	Total	O	0	0
			288	288		
4	B	210	Total	O	0	0
			210	210		

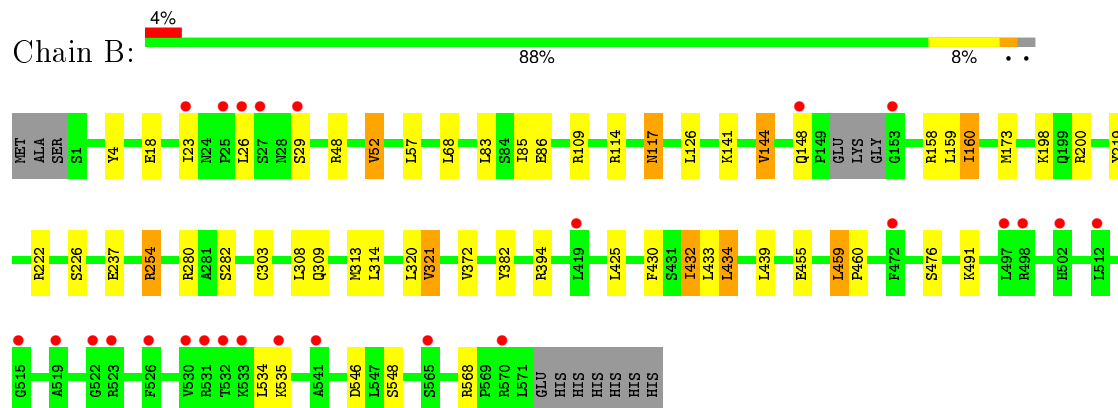
3 Residue-property plots [i](#)

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



• Molecule 1: RNA-directed RNA polymerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	106.53Å 107.57Å 133.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.59 – 2.39 75.70 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.5 (75.59-2.39) 98.5 (75.70-2.39)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.55 (at 2.40Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.167 , 0.217 0.197 , 0.241	Depositor DCC
R_{free} test set	3028 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.656	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.6	EDS
Estimated twinning fraction	0.022 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 60578 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9386	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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, 0.375 respectively for untwinned datasets, and 0.333, 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: XNC, CL

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	0.69	1/4463 (0.0%)	0.75	5/6057 (0.1%)
1	B	0.61	0/4525	0.75	5/6140 (0.1%)
All	All	0.65	1/8988 (0.0%)	0.75	10/12197 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	274	CYS	CB-SG	-9.10	1.66	1.82

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	254	ARG	NE-CZ-NH2	-9.24	115.68	120.30
1	B	280	ARG	NE-CZ-NH2	-8.61	115.99	120.30
1	B	280	ARG	NE-CZ-NH1	7.67	124.14	120.30
1	B	254	ARG	NE-CZ-NH1	7.33	123.96	120.30
1	A	254	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	B	109	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	A	280	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	254	ARG	NE-CZ-NH2	-5.07	117.77	120.30
1	A	200	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	A	337	ARG	NE-CZ-NH1	-5.03	117.79	120.30

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	4368	0	4377	28	0
1	B	4429	0	4451	29	0
2	A	44	0	31	0	0
2	B	44	0	31	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	288	0	0	3	0
4	B	210	0	0	6	0
All	All	9386	0	8890	57	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 (57) 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:76:SER:HB3	4:A:654:HOH:O	1.72	0.88
1:B:198:LYS:HD2	4:B:652:HOH:O	1.74	0.87
1:A:46:SER:HA	1:A:49:GLN:HE21	1.49	0.78
1:A:160:ILE:HD12	1:A:282:SER:OG	1.89	0.73
1:B:26:LEU:HD13	1:B:432:ILE:HD13	1.70	0.72
1:B:148:GLN:HE22	1:B:491:LYS:NZ	1.97	0.63
1:A:24:ASN:HD22	1:A:27:SER:H	1.48	0.60
1:B:237:GLU:OE2	1:B:254:ARG:HD2	2.03	0.59
1:B:200:ARG:NH1	4:B:645:HOH:O	2.36	0.58
1:A:48:ARG:HG2	1:A:159:LEU:HG	1.87	0.57
1:A:46:SER:HA	1:A:49:GLN:NE2	2.16	0.56
1:B:144:VAL:HG22	1:B:394:ARG:HG2	1.85	0.56
1:B:114:ARG:HA	1:B:117:ASN:HD22	1.72	0.55
1:B:303:CYS:SG	1:B:313:MET:CE	2.95	0.54
1:B:148:GLN:HE22	1:B:491:LYS:HZ2	1.54	0.53
1:A:280:ARG:HD2	1:A:291:ASN:OD1	2.09	0.53
1:A:100:LYS:NZ	1:A:164:ASP:OD1	2.41	0.53
1:B:455:GLU:HB3	4:B:656:HOH:O	2.10	0.52
1:B:303:CYS:SG	1:B:313:MET:HE3	2.50	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:LEU:HD13	1:A:249:ALA:HB2	1.94	0.49
1:A:337:ARG:NH2	4:A:623:HOH:O	2.45	0.49
1:A:57:LEU:HD12	1:A:57:LEU:C	2.33	0.49
1:B:86:GLU:HB2	4:B:789:HOH:O	2.13	0.48
1:B:546:ASP:OD1	1:B:548:SER:OG	2.31	0.48
1:A:83:LEU:HB2	1:A:173:MET:HA	1.96	0.48
1:B:198:LYS:HE2	4:B:632:HOH:O	2.14	0.47
1:B:219:TYR:HB3	1:B:320:LEU:HD23	1.96	0.47
1:A:423:MET:HA	1:A:528:TRP:CZ2	2.51	0.45
1:B:26:LEU:O	1:B:29:SER:HB2	2.17	0.45
1:B:83:LEU:HB2	1:B:173:MET:HA	1.98	0.45
1:A:455:GLU:HB3	4:A:620:HOH:O	2.17	0.45
1:A:237:GLU:OE1	1:A:254:ARG:NH1	2.50	0.45
1:A:24:ASN:HD21	1:A:26:LEU:HB2	1.82	0.45
1:A:364:THR:HA	1:A:368:SER:O	2.17	0.45
1:A:118:HIS:O	1:A:122:VAL:HG23	2.18	0.44
1:B:52:VAL:HG22	1:B:226:SER:OG	2.17	0.44
1:A:105:ALA:O	1:A:109:ARG:HG3	2.18	0.44
1:B:459:LEU:N	1:B:460:PRO:CD	2.82	0.43
1:B:309:GLN:HA	1:B:309:GLN:HE21	1.82	0.43
1:A:182:LEU:HD12	1:A:243:CYS:SG	2.58	0.43
1:A:327:ALA:O	1:A:331:GLU:HG3	2.18	0.43
1:A:52:VAL:HG12	1:A:223[B]:CYS:SG	2.59	0.43
1:B:4:TYR:CE2	1:B:52:VAL:HG13	2.53	0.43
1:B:430:PHE:O	1:B:434:LEU:HB2	2.19	0.43
1:B:372:VAL:HG22	1:B:382:TYR:CD1	2.54	0.42
1:A:321:VAL:HG21	1:A:365:SER:HB2	2.00	0.42
1:B:160:ILE:HA	1:B:282:SER:OG	2.19	0.42
1:B:303:CYS:SG	1:B:313:MET:HE1	2.58	0.42
1:A:444:ASP:HA	1:A:452:TYR:O	2.20	0.42
1:A:280:ARG:HD3	1:A:287:THR:HG23	2.01	0.41
1:B:48:ARG:HG2	1:B:159:LEU:HG	2.03	0.41
1:B:314:LEU:HB3	1:B:321:VAL:HG13	2.03	0.41
1:A:219:TYR:HB3	1:A:320:LEU:HD23	2.02	0.41
1:A:32:ARG:HD2	1:A:32:ARG:N	2.36	0.41
1:B:198:LYS:CD	4:B:652:HOH:O	2.49	0.40
1:B:141:LYS:HE3	1:B:158:ARG:NH2	2.37	0.40
1:A:434:LEU:HD21	1:A:510:LYS:HB2	2.02	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	557/581 (96%)	548 (98%)	8 (1%)	1 (0%)	52	69
1	B	564/581 (97%)	551 (98%)	12 (2%)	1 (0%)	52	69
All	All	1121/1162 (96%)	1099 (98%)	20 (2%)	2 (0%)	52	69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	ALA
1	B	23	ILE

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	478/495 (97%)	454 (95%)	24 (5%)	30	48
1	B	484/495 (98%)	462 (96%)	22 (4%)	34	52
All	All	962/990 (97%)	916 (95%)	46 (5%)	31	49

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	24	ASN
1	A	47	LEU
1	A	50	LYS

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Mol	Chain	Res	Type
1	A	57	LEU
1	A	68	LEU
1	A	90	LYS
1	A	126	LEU
1	A	142	SER
1	A	201	VAL
1	A	222	ARG
1	A	254	ARG
1	A	274	CYS
1	A	313	MET
1	A	352	ASP
1	A	362	LEU
1	A	364	THR
1	A	425	LEU
1	A	433	LEU
1	A	459	LEU
1	A	469	LEU
1	A	531	ARG
1	A	534	LEU
1	A	563	SER
1	B	18	GLU
1	B	52	VAL
1	B	57	LEU
1	B	68	LEU
1	B	85	ILE
1	B	117	ASN
1	B	126	LEU
1	B	144	VAL
1	B	160	ILE
1	B	222	ARG
1	B	308	LEU
1	B	321	VAL
1	B	425	LEU
1	B	432	ILE
1	B	433	LEU
1	B	434	LEU
1	B	439	LEU
1	B	459	LEU
1	B	476	SER
1	B	534	LEU
1	B	535	LYS
1	B	568	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	34	HIS
1	A	35	ASN
1	A	49	GLN
1	A	148	GLN
1	A	206	ASN
1	A	273	ASN
1	A	309	GLN
1	A	483	ASN
1	A	502	HIS
1	A	514	GLN
1	A	562	HIS
1	B	117	ASN
1	B	148	GLN
1	B	194	GLN
1	B	206	ASN
1	B	273	ASN
1	B	309	GLN
1	B	483	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

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	XNC	A	579	-	44,49,49	1.16	4 (9%)	61,74,74	1.78	8 (13%)
2	XNC	B	579	-	44,49,49	1.09	5 (11%)	61,74,74	1.53	9 (14%)

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	XNC	A	579	-	-	0/17/54/54	0/4/6/6
2	XNC	B	579	-	-	0/17/54/54	0/4/6/6

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	579	XNC	C16-N7	2.03	1.40	1.37
2	A	579	XNC	O32-S12	2.13	1.46	1.44
2	B	579	XNC	C40-N39	2.35	1.39	1.34
2	B	579	XNC	C17-N39	2.35	1.38	1.34
2	A	579	XNC	C18-C8	2.47	1.55	1.52
2	A	579	XNC	C16-N7	2.58	1.41	1.37
2	B	579	XNC	C18-C8	2.70	1.55	1.52
2	B	579	XNC	C15-C10	2.89	1.53	1.50
2	A	579	XNC	C17-N39	3.69	1.40	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	579	XNC	O33-S12-O32	-9.72	107.66	117.11
2	B	579	XNC	O33-S12-O32	-6.17	111.11	117.11
2	B	579	XNC	C43-C17-N39	-3.02	119.30	122.90
2	A	579	XNC	C22-C23-C18	-2.63	120.98	123.78
2	A	579	XNC	C41-C40-N39	-2.27	118.27	121.44
2	A	579	XNC	C15-C14-C13	-2.20	105.59	110.53
2	B	579	XNC	C22-C23-C18	-2.17	121.47	123.78
2	A	579	XNC	C19-C18-C23	2.18	119.10	116.14
2	B	579	XNC	C19-C18-C23	2.22	119.16	116.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	579	XNC	C25-O24-C21	2.30	123.88	117.70
2	B	579	XNC	O32-S12-C13	2.58	111.80	109.11
2	B	579	XNC	O33-S12-C13	2.67	111.90	109.11
2	B	579	XNC	C17-N39-C40	2.97	120.64	118.13
2	A	579	XNC	O32-S12-C13	3.19	112.44	109.11
2	A	579	XNC	C14-C15-C10	3.49	117.75	113.39
2	B	579	XNC	C14-C15-C10	3.51	117.78	113.39
2	A	579	XNC	C17-N39-C40	3.77	121.31	118.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	560/581 (96%)	0.14	1 (0%) 95 95	19, 29, 41, 54	0
1	B	568/581 (97%)	0.43	26 (4%) 36 37	20, 29, 44, 77	0
All	All	1128/1162 (97%)	0.29	27 (2%) 62 61	19, 29, 42, 77	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	LEU	5.3
1	B	25	PRO	4.7
1	B	531	ARG	4.6
1	B	27	SER	4.5
1	B	532	THR	3.9
1	B	153	GLY	3.9
1	B	535	LYS	3.4
1	B	23	ILE	3.4
1	B	526	PHE	3.1
1	B	502	HIS	3.0
1	B	148	GLN	2.9
1	B	523	ARG	2.8
1	B	29	SER	2.7
1	A	153	GLY	2.7
1	B	512	LEU	2.6
1	B	533	LYS	2.4
1	B	498	ARG	2.3
1	B	419	LEU	2.3
1	B	570	ARG	2.3
1	B	565	SER	2.3
1	B	515	GLY	2.2
1	B	541	ALA	2.2
1	B	497	LEU	2.2
1	B	519	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	472	PHE	2.2
1	B	522	GLY	2.0
1	B	530	VAL	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
2	XNC	B	579	44/44	0.97	0.11	-0.69	16,21,24,32	0
2	XNC	A	579	44/44	0.98	0.12	-1.42	13,15,18,22	0
3	CL	B	580	1/1	0.98	0.07	-2.83	36,36,36,36	0
3	CL	A	581	1/1	0.98	0.06	-10.57	24,24,24,24	0
3	CL	A	580	1/1	0.99	0.07	-	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.