



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:05 PM UTC

PDB ID : 3GNW / pdb_00003gnw
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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

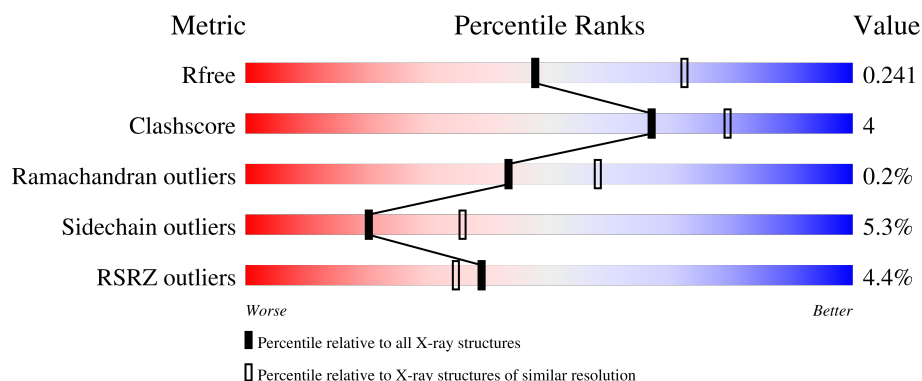
1 Overall quality at a glance

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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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	581	
1	B	581	

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.

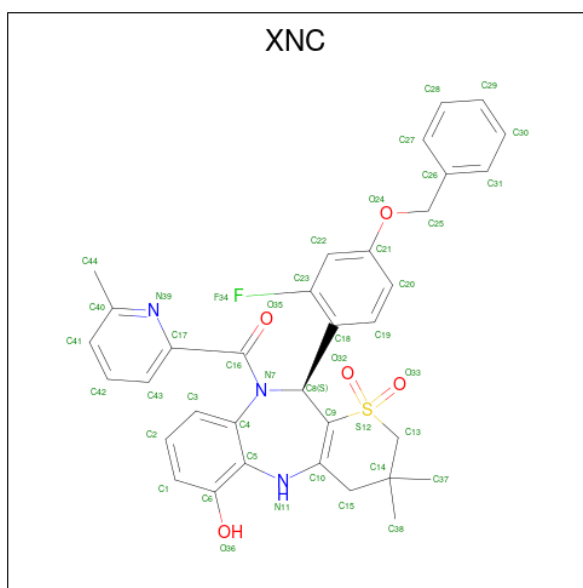
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	560	Total	C	N	O	S	0	2	0
			4368	2750	772	813	33			
1	B	568	Total	C	N	O	S	0	1	0
			4429	2788	788	821	32			

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-hexahydrothiopyrano[3,2-b][1,5]benzodiazepin-6-ol 1,1-dioxide

(CCD ID: 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 (CCD ID: CL) (formula: Cl).

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

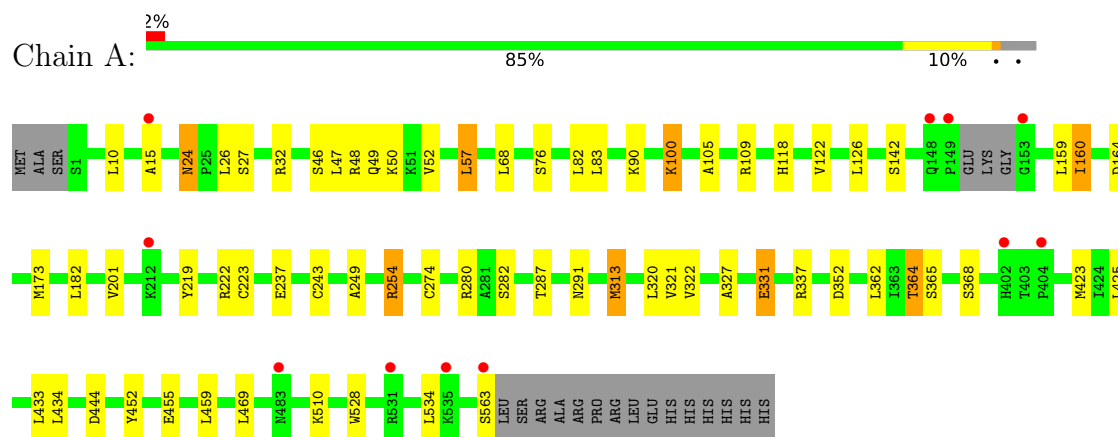
- Molecule 4 is water.

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

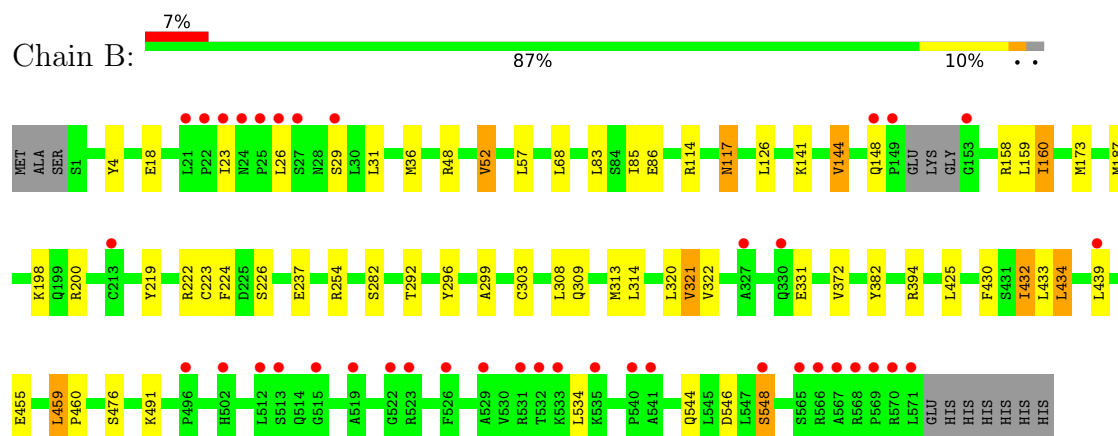
3 Residue-property plots

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: 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 75.59 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.5 (75.59-2.39) 98.5 (75.59-2.39)	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	3029 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.656	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.022 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	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, 0.333 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: CL, XNC

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	0.82	0/4463	0.90	0/6057
1	B	0.75	0/4525	0.90	0/6140
All	All	0.78	0/8988	0.90	0/12197

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	4368	0	4377	31	0
1	B	4429	0	4451	34	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	65	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 4.

All (65) 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:B:303:CYS:SG	1:B:313:MET:HE3	2.22	0.79
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:187:MET:HE1	1:B:292:THR:HG22	1.90	0.54
1:A:57:LEU:HD12	1:A:57:LEU:C	2.33	0.53
1:A:280:ARG:HD2	1:A:291:ASN:OD1	2.09	0.53
1:B:148:GLN:HE22	1:B:491:LYS:HZ2	1.54	0.53
1:A:100:LYS:NZ	1:A:164:ASP:OD1	2.41	0.53
1:B:303:CYS:SG	1:B:313:MET:CE	2.95	0.52
1:B:455:GLU:HB3	4:B:656:HOH:O	2.10	0.52
1:B:299:ALA:HB1	1:B:313:MET:HE1	1.93	0.51
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: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:A:423:MET:HE3	1:A:528:TRP:CZ3	2.52	0.45
1:B:187:MET:HE3	1:B:296:TYR:CD2	2.50	0.45
1:B:26:LEU:O	1:B:29:SER:HB2	2.17	0.45
1:A:313:MET:HE3	1:A:313:MET:HB2	1.75	0.45
1:B:83:LEU:HB2	1:B:173:MET:HA	1.98	0.45
1:B:313:MET:HE2	1:B:322:VAL:HG22	1.96	0.45
1:A:237:GLU:OE1	1:A:254:ARG:NH1	2.50	0.45
1:A:455:GLU:HB3	4:A:620:HOH:O	2.17	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:309:GLN:HE21	1:B:309:GLN:HA	1.82	0.43
1:B:459:LEU:N	1:B:460:PRO:CD	2.82	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:313:MET:HE2	1:B:322:VAL:CG2	2.49	0.42
1:B:160:ILE:HA	1:B:282:SER:OG	2.19	0.42
1:A:444:ASP:HA	1:A:452:TYR:O	2.20	0.42
1:A:313:MET:HE2	1:A:322:VAL:HG22	2.02	0.41
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:B:223:CYS:O	1:B:224:PHE:C	2.64	0.41
1:A:219:TYR:HB3	1:A:320:LEU:HD23	2.02	0.41
1:A:32:ARG:N	1:A:32:ARG:HD2	2.36	0.41
1:A:434:LEU:HD21	1:A:510:LYS:HB2	2.02	0.40
1:B:141:LYS:HE3	1:B:158:ARG:NH2	2.37	0.40
1:B:198:LYS:CD	4:B:652:HOH:O	2.49	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%)	43 58

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	564/581 (97%)	551 (98%)	12 (2%)	1 (0%)	43 58
All	All	1121/1162 (96%)	1099 (98%)	20 (2%)	2 (0%)	43 58

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%)	452 (95%)	26 (5%)	20 35
1	B	484/495 (98%)	459 (95%)	25 (5%)	21 36
All	All	962/990 (97%)	911 (95%)	51 (5%)	20 36

All (51) 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
1	A	57	LEU
1	A	68	LEU
1	A	90	LYS
1	A	100	LYS
1	A	126	LEU
1	A	142	SER
1	A	160	ILE
1	A	201	VAL
1	A	222	ARG
1	A	254	ARG

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Mol	Chain	Res	Type
1	A	274	CYS
1	A	313	MET
1	A	331	GLU
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	534	LEU
1	A	563	SER
1	B	18	GLU
1	B	31	LEU
1	B	36	MET
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	331	GLU
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	544	GLN
1	B	548	SER

Sometimes 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

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Mol	Chain	Res	Type
1	A	35	ASN
1	A	49	GLN
1	A	206	ASN
1	A	273	ASN
1	A	411	ASN
1	A	483	ASN
1	A	514	GLN
1	A	562	HIS
1	B	35	ASN
1	B	117	ASN
1	B	148	GLN
1	B	194	GLN
1	B	206	ASN
1	B	309	GLN
1	B	411	ASN
1	B	446	GLN
1	B	461	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 oligosaccharides 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	-	45,49,49	1.27	7 (15%)	64,74,74	1.78	8 (12%)
2	XNC	B	579	-	45,49,49	1.29	7 (15%)	64,74,74	1.63	11 (17%)

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	-	-	1/17/54/54	0/4/6/6
2	XNC	B	579	-	-	2/17/54/54	0/4/6/6

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	579	XNC	C13-S12	4.33	1.82	1.76
2	A	579	XNC	C17-N39	3.72	1.40	1.34
2	B	579	XNC	C15-C10	3.18	1.53	1.50
2	A	579	XNC	C13-S12	2.96	1.80	1.76
2	B	579	XNC	C40-N39	2.70	1.39	1.34
2	A	579	XNC	C16-N7	2.64	1.41	1.37
2	B	579	XNC	C18-C8	2.55	1.55	1.52
2	A	579	XNC	O32-S12	2.40	1.46	1.43
2	B	579	XNC	C17-N39	2.35	1.38	1.34
2	A	579	XNC	C18-C8	2.31	1.55	1.52
2	A	579	XNC	C15-C10	2.11	1.52	1.50
2	B	579	XNC	O32-S12	2.08	1.46	1.43
2	A	579	XNC	C10-N11	2.07	1.39	1.35
2	B	579	XNC	C16-N7	2.06	1.40	1.37

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	579	XNC	O33-S12-O32	-9.80	107.66	117.14
2	B	579	XNC	O33-S12-O32	-6.24	111.11	117.14
2	B	579	XNC	C9-C10-N11	-4.03	123.71	126.58
2	B	579	XNC	C14-C15-C10	4.00	117.78	113.40
2	A	579	XNC	C14-C15-C10	3.98	117.75	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	579	XNC	C17-N39-C40	3.82	121.31	118.24
2	B	579	XNC	C43-C17-N39	-3.10	119.30	122.92
2	A	579	XNC	O32-S12-C13	3.00	112.44	109.14
2	B	579	XNC	C17-N39-C40	2.99	120.64	118.24
2	A	579	XNC	C22-C23-C18	-2.66	120.98	123.93
2	B	579	XNC	C25-O24-C21	2.63	123.88	117.62
2	A	579	XNC	C41-C40-N39	-2.55	118.27	121.59
2	B	579	XNC	O33-S12-C13	2.51	111.90	109.14
2	B	579	XNC	O32-S12-C13	2.42	111.80	109.14
2	B	579	XNC	C19-C18-C23	2.33	119.16	116.22
2	A	579	XNC	C19-C18-C23	2.29	119.10	116.22
2	A	579	XNC	C25-O24-C21	2.24	122.97	117.62
2	B	579	XNC	C22-C23-C18	-2.22	121.47	123.93
2	B	579	XNC	O35-C16-N7	-2.00	118.94	121.51

There are no chirality outliers.

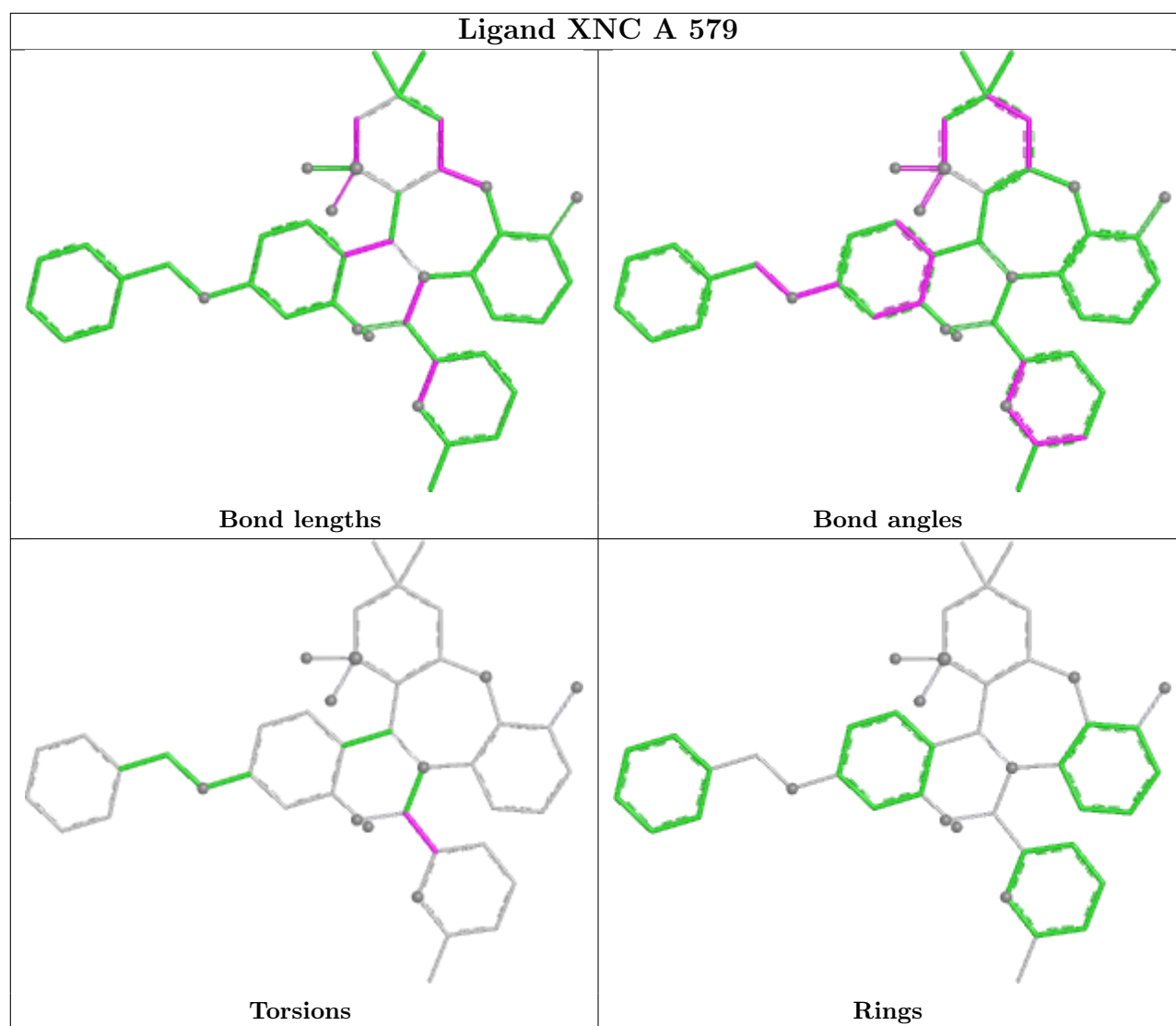
All (3) torsion outliers are listed below:

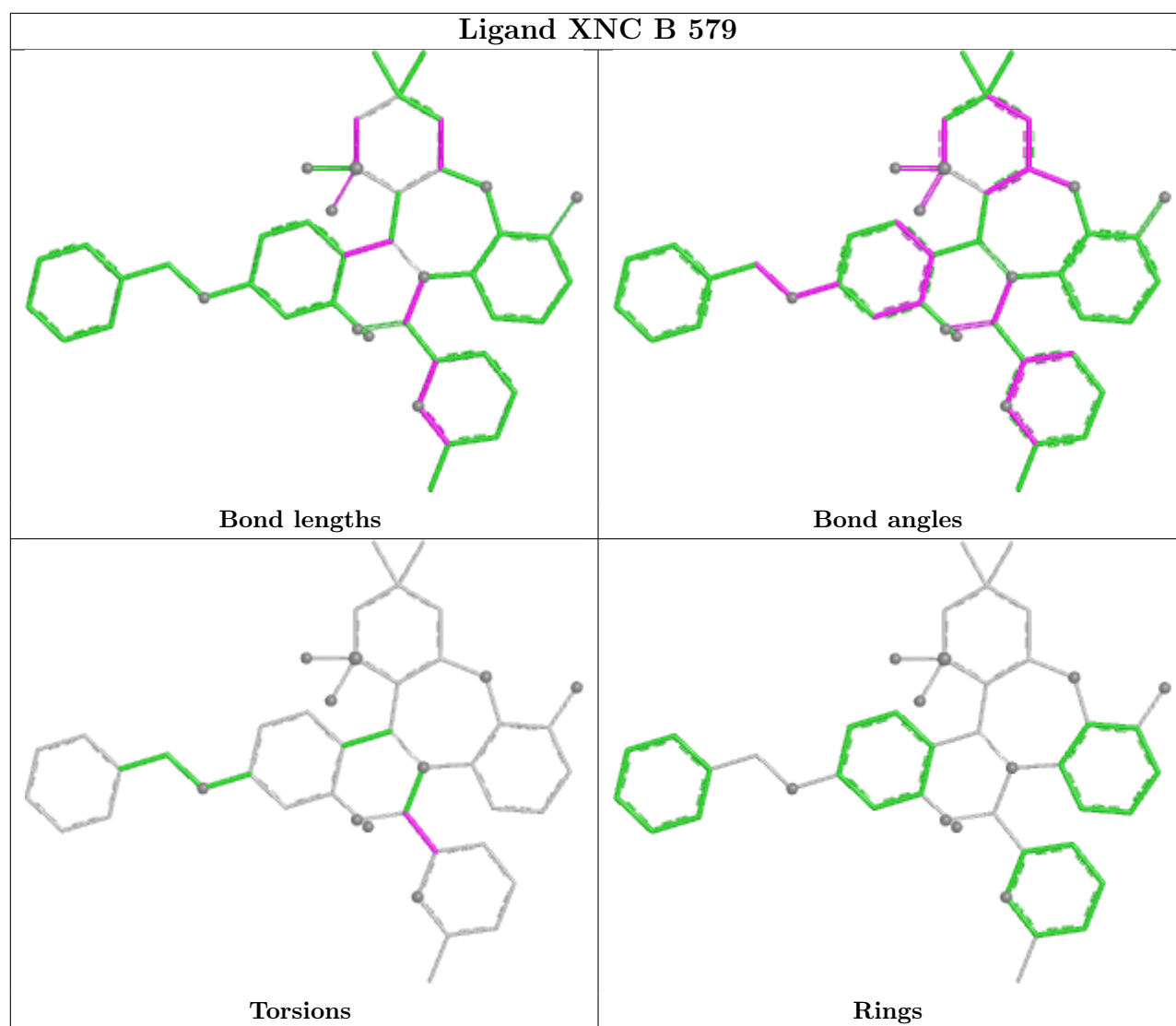
Mol	Chain	Res	Type	Atoms
2	B	579	XNC	N7-C16-C17-N39
2	A	579	XNC	N7-C16-C17-N39
2	B	579	XNC	O35-C16-C17-N39

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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%)	0.20	11 (1%) 65 60	16, 29, 41, 54	1 (0%)
1	B	568/581 (97%)	0.56	39 (6%) 23 19	20, 29, 44, 77	0
All	All	1128/1162 (97%)	0.38	50 (4%) 39 34	16, 29, 42, 77	1 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	149	PRO	6.3
1	B	571	LEU	5.5
1	A	153	GLY	5.3
1	B	23	ILE	5.0
1	B	25	PRO	4.9
1	B	149	PRO	4.6
1	B	26	LEU	4.2
1	B	27	SER	4.0
1	B	153	GLY	3.8
1	B	502	HIS	3.6
1	B	531	ARG	3.3
1	B	148	GLN	3.1
1	B	535	LYS	3.1
1	B	519	ALA	3.1
1	B	532	THR	3.0
1	B	569	PRO	3.0
1	B	567	ALA	3.0
1	B	568	ARG	2.8
1	B	496	PRO	2.8
1	B	541	ALA	2.7
1	B	523	ARG	2.7
1	B	330	GLN	2.6
1	B	21	LEU	2.5
1	B	327	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	15	ALA	2.4
1	B	22	PRO	2.4
1	B	515	GLY	2.4
1	A	563	SER	2.4
1	B	540	PRO	2.4
1	B	533	LYS	2.4
1	B	29	SER	2.4
1	B	512	LEU	2.4
1	A	148	GLN	2.4
1	A	535	LYS	2.3
1	B	566	ARG	2.3
1	B	522	GLY	2.3
1	A	402	HIS	2.3
1	B	570	ARG	2.2
1	B	526	PHE	2.2
1	B	565	SER	2.2
1	B	529	ALA	2.1
1	A	483	ASN	2.1
1	A	404	PRO	2.1
1	B	548	SER	2.1
1	B	513	SER	2.0
1	B	439	LEU	2.0
1	A	212	LYS	2.0
1	B	213	CYS	2.0
1	B	24	ASN	2.0
1	A	531	ARG	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](#)

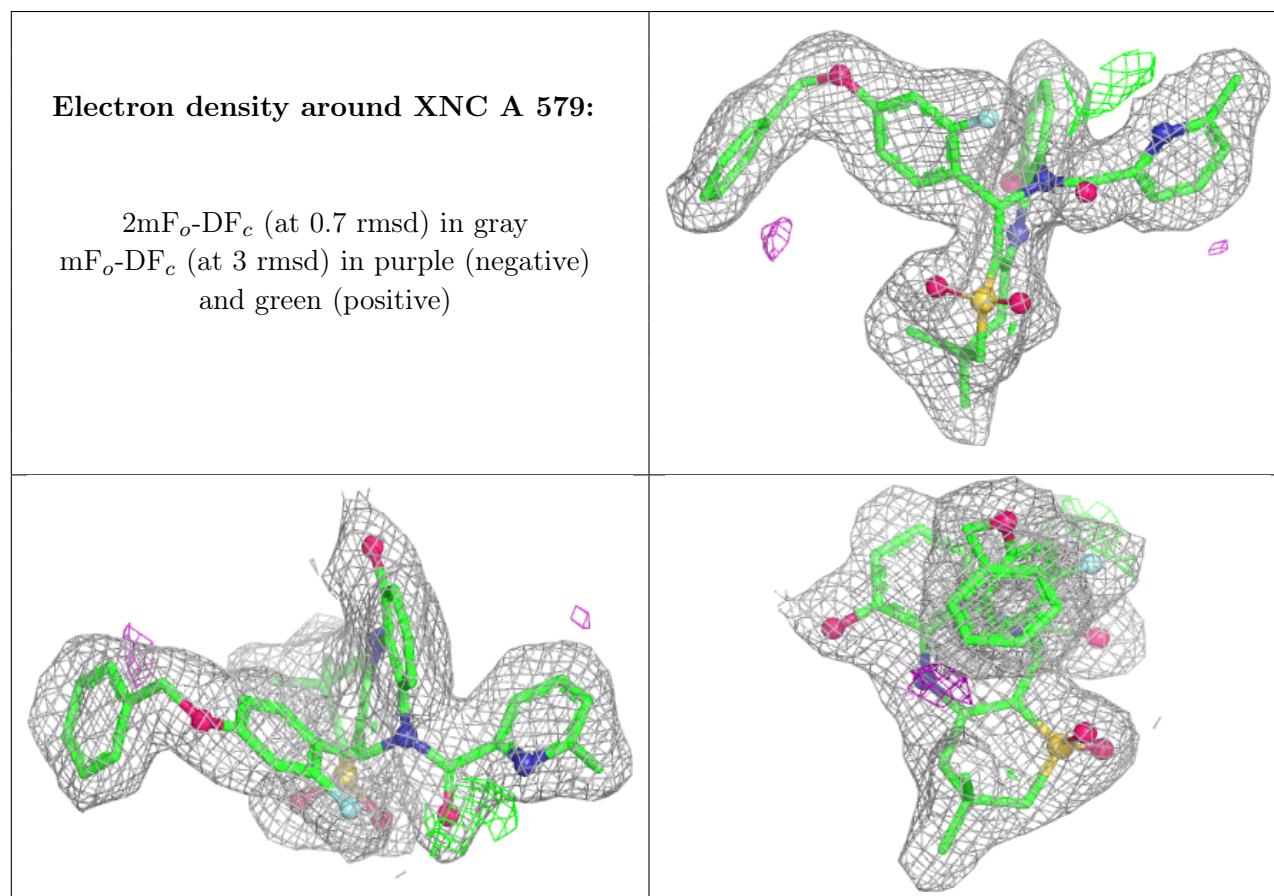
There are no oligosaccharides in this entry.

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.

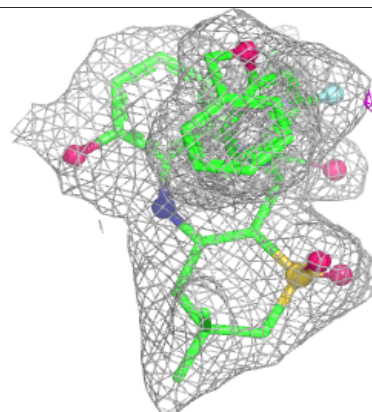
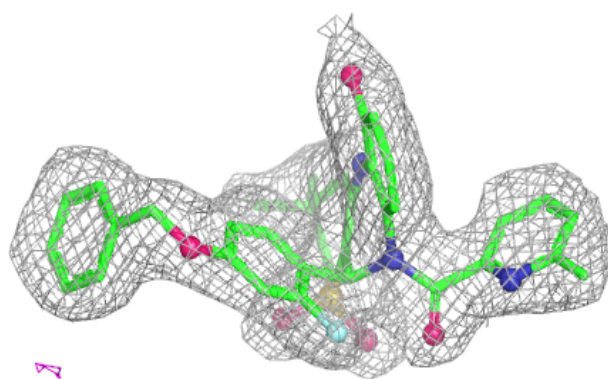
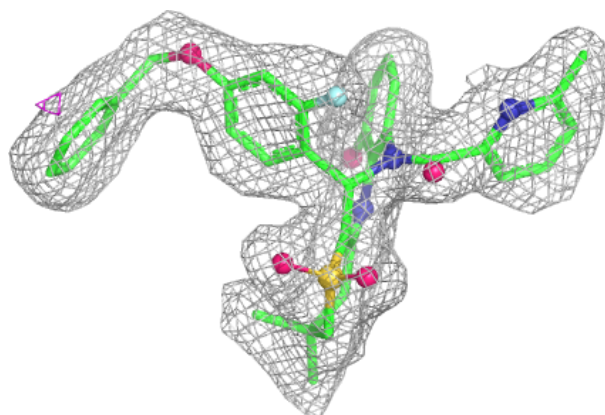
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	XNC	A	579	44/44	0.97	0.06	13,15,18,22	0
2	XNC	B	579	44/44	0.97	0.07	16,21,24,32	0
3	CL	B	580	1/1	0.97	0.09	36,36,36,36	0
3	CL	A	581	1/1	0.98	0.07	24,24,24,24	0
3	CL	A	580	1/1	0.99	0.10	34,34,34,34	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around XNC B 579:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.