



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:41 PM UTC

PDB ID : 4EAW / pdb_00004eaw
Title : HCV NS5B in complex with IDX375
Authors : Dousson, C.B.; Papparin, J.-L.; Surleraux, D.; Augustin, M.; Blaesse, M.;
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Deposited on : 2012-03-22
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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)
Xtrriage (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

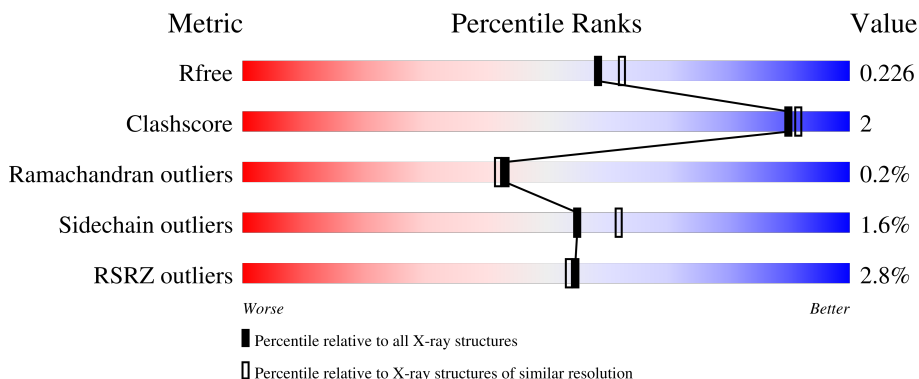
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	562	
1	B	562	

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
4	PEG	A	609	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9719 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.

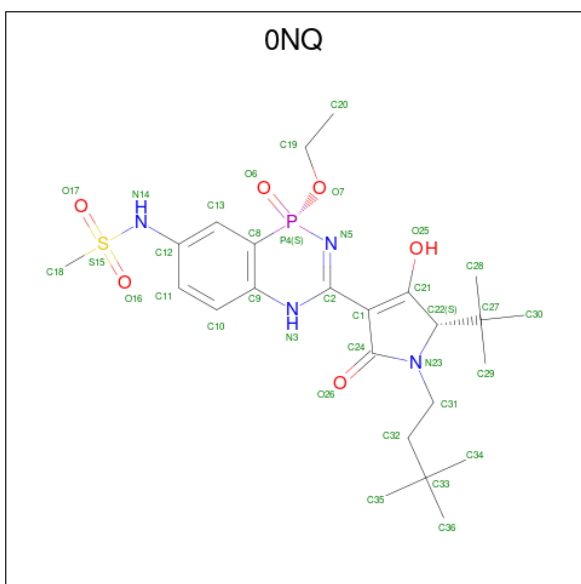
- Molecule 1 is a protein called RNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	559	4439	2793	789	823	34	38	9	0
1	B	559	4405	2773	781	818	33	59	5	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	316	TYR	ASN	conflict	UNP O92972
B	316	TYR	ASN	conflict	UNP O92972

- Molecule 2 is N-{(1S)-3-[(5S)-5-tert-butyl-1-(3,3-dimethylbutyl)-4-hydroxy-2-oxo-2,5-dihydro-1H-pyrrol-3-yl]-1-ethoxy-1-oxido-1,4-dihydro-2,4,1-benzodiazaphosphinin-7-yl}methanesulfonamide (CCD ID: 0NQ) (formula: C₂₄H₃₇N₄O₆PS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	A	1	Total	C	N	O	P	S	0	0
			36	24	4	6	1	1		
2	B	1	Total	C	N	O	P	S	0	0
			36	24	4	6	1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



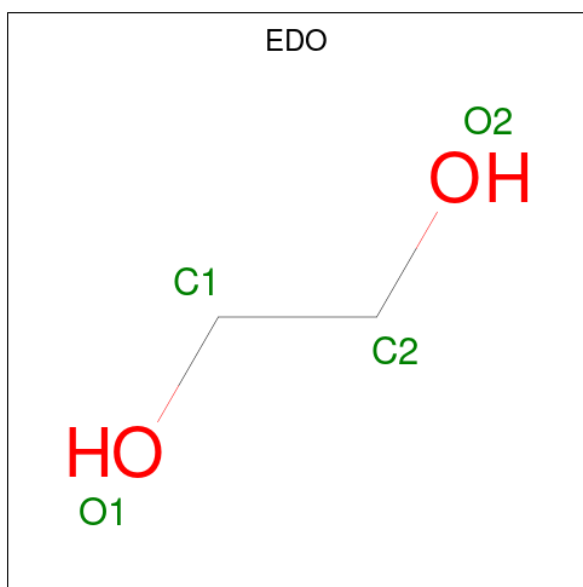
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

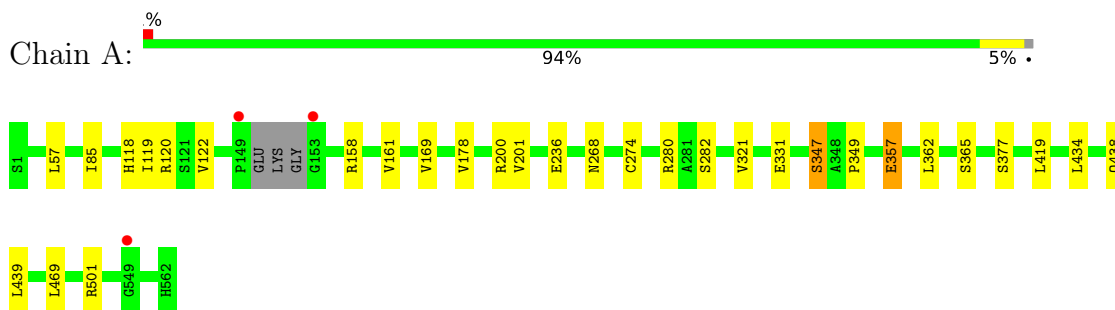
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	431	Total	O	0	2
			433	433		
6	B	299	Total	O	0	2
			301	301		

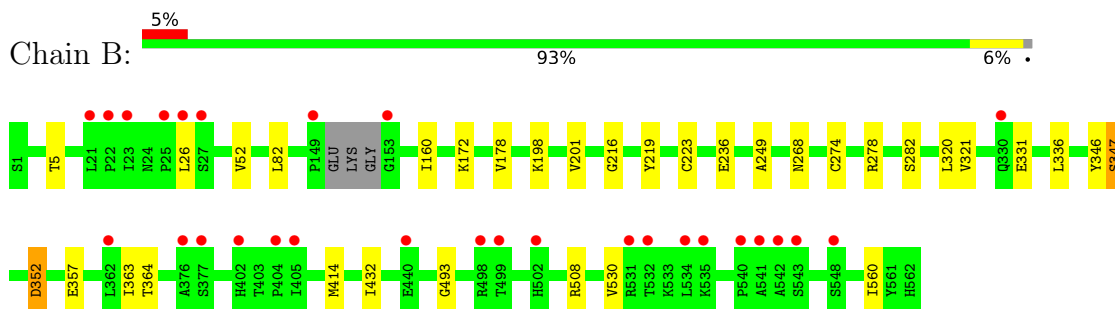
3 Residue-property plots [i](#)

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	106.31Å 108.03Å 133.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.59 – 2.00 41.59 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (41.59-2.00) 99.7 (41.59-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.00Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.173 , 0.219 0.186 , 0.226	Depositor DCC
R_{free} test set	1041 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtrriage
Anisotropy	0.519	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.018 for k,h,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9719	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% 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: 0NQ, SO4, PEG, EDO

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	1.00	2/4536 (0.0%)	0.95	0/6153
1	B	0.93	0/4501	0.96	2/6107 (0.0%)
All	All	0.97	2/9037 (0.0%)	0.96	2/12260 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	158	ARG	NE-CZ	5.87	1.39	1.33
1	A	501	ARG	NE-CZ	5.79	1.39	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	352	ASP	CA-C-N	-5.59	114.62	120.38
1	B	352	ASP	C-N-CA	-5.59	114.62	120.38

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	4439	0	4439	16	0
1	B	4405	0	4410	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	36	0	36	0	0
2	B	36	0	36	1	0
3	A	30	0	0	0	0
3	B	10	0	0	0	0
4	A	14	0	20	4	0
4	B	7	0	10	0	0
5	A	8	0	12	0	0
6	A	433	0	0	2	0
6	B	301	0	0	2	0
All	All	9719	0	8963	31	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 2.

All (31) 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:5:THR:HG23	1:B:278[A]:ARG:HH22	1.51	0.75
1:A:200:ARG:NH2	1:A:365:SER:OG	2.24	0.70
1:A:161:VAL:HG23	4:A:609:PEG:H11	1.77	0.67
1:B:82:LEU:HD13	1:B:249:ALA:HB2	1.76	0.65
1:A:434:LEU:HD13	1:A:439:LEU:HD11	1.79	0.64
1:B:172:LYS:HE3	1:B:560:ILE:HD13	1.91	0.52
1:A:161:VAL:CG2	4:A:609:PEG:H11	2.40	0.52
1:A:282:SER:HB3	4:A:609:PEG:C2	2.40	0.52
1:B:52:VAL:HG12	1:B:223[B]:CYS:SG	2.50	0.51
1:A:268:ASN:HB3	1:A:274[B]:CYS:SG	2.52	0.50
1:B:198:LYS:HA	1:B:201:VAL:HG12	1.93	0.50
1:A:321:VAL:HG21	1:A:365:SER:HB2	1.94	0.50
1:A:282:SER:HB3	4:A:609:PEG:H21	1.93	0.49
1:B:160:ILE:HD12	1:B:282:SER:OG	2.12	0.49
1:A:118[A]:HIS:O	1:A:122:VAL:HG23	2.12	0.49
1:A:119:ILE:HD13	1:A:169:VAL:HG11	1.93	0.49
1:A:236[B]:GLU:OE2	1:A:280:ARG:NH2	2.38	0.47
1:B:414:MET:HG3	2:B:601:ONQ:H22	1.97	0.47
1:A:85:ILE:HD12	1:A:120:ARG:NH2	2.30	0.46
1:B:268:ASN:HB3	1:B:274:CYS:SG	2.55	0.46
1:B:219:TYR:HB3	1:B:320:LEU:HD23	1.99	0.44
1:B:26:LEU:HD22	1:B:432:ILE:HG12	2.00	0.44
1:A:178:VAL:HG23	6:A:927:HOH:O	2.17	0.43
1:B:493:GLY:HA2	6:B:892:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:GLY:HA3	1:B:363:ILE:HD11	2.01	0.42
1:B:178:VAL:HG23	6:B:759:HOH:O	2.20	0.42
1:B:508:ARG:CZ	1:B:530:VAL:HG11	2.50	0.41
1:A:118[B]:HIS:HE1	6:A:706:HOH:O	2.03	0.41
1:B:346:TYR:O	1:B:347:SER:HB3	2.20	0.41
1:A:347:SER:O	1:A:349:PRO:HD3	2.22	0.40
1:A:357:GLU:HG3	1:A:362:LEU:HB2	2.03	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	564/562 (100%)	553 (98%)	10 (2%)	1 (0%)	43 42
1	B	560/562 (100%)	549 (98%)	10 (2%)	1 (0%)	43 42
All	All	1124/1124 (100%)	1102 (98%)	20 (2%)	2 (0%)	43 42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	347	SER
1	B	347	SER

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	485/478 (102%)	477 (98%)	8 (2%)	55	62
1	B	481/478 (101%)	473 (98%)	8 (2%)	53	60
All	All	966/956 (101%)	950 (98%)	16 (2%)	55	60

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LEU
1	A	201	VAL
1	A	331	GLU
1	A	357	GLU
1	A	377	SER
1	A	419	LEU
1	A	438	GLN
1	A	469	LEU
1	B	236[A]	GLU
1	B	236[B]	GLU
1	B	321	VAL
1	B	331	GLU
1	B	336	LEU
1	B	352	ASP
1	B	357	GLU
1	B	364	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	330	GLN
1	A	374	HIS
1	A	406	ASN
1	A	438	GLN
1	A	446	GLN
1	A	562	HIS
1	B	33	HIS
1	B	184	GLN
1	B	273	ASN
1	B	446	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](#)

15 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
3	SO4	A	604	-	4,4,4	0.25	0	6,6,6	0.30	0
4	PEG	A	609	-	6,6,6	0.56	0	5,5,5	1.15	0
3	SO4	B	602	-	4,4,4	0.30	0	6,6,6	0.36	0
4	PEG	B	604	-	6,6,6	0.59	0	5,5,5	0.65	0
3	SO4	A	607	-	4,4,4	0.42	0	6,6,6	0.60	0
5	EDO	A	611	-	3,3,3	0.75	0	2,2,2	0.23	0
2	ONQ	A	601	-	34,38,38	1.40	5 (14%)	41,61,61	2.10	8 (19%)
4	PEG	A	608	-	6,6,6	0.72	0	5,5,5	0.72	0
2	ONQ	B	601	-	34,38,38	1.17	3 (8%)	41,61,61	2.14	6 (14%)
3	SO4	A	603	-	4,4,4	0.28	0	6,6,6	0.59	0
3	SO4	B	603	-	4,4,4	0.21	0	6,6,6	0.28	0
3	SO4	A	605	-	4,4,4	0.33	0	6,6,6	0.60	0
3	SO4	A	602	-	4,4,4	0.28	0	6,6,6	1.07	0
5	EDO	A	610	-	3,3,3	0.39	0	2,2,2	0.16	0
3	SO4	A	606	-	4,4,4	0.24	0	6,6,6	0.49	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PEG	A	609	-	-	3/4/4/4	-
5	EDO	A	611	-	-	1/1/1/1	-
4	PEG	B	604	-	-	4/4/4/4	-
2	ONQ	A	601	-	-	2/18/60/60	0/2/3/3
4	PEG	A	608	-	-	3/4/4/4	-
2	ONQ	B	601	-	-	2/18/60/60	0/2/3/3
5	EDO	A	610	-	-	1/1/1/1	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	ONQ	C2-N3	4.24	1.41	1.36
2	A	601	ONQ	C9-C8	-3.83	1.38	1.41
2	B	601	ONQ	C2-N3	3.66	1.40	1.36
2	B	601	ONQ	C22-N23	3.24	1.49	1.46
2	A	601	ONQ	C18-S15	2.44	1.81	1.75
2	B	601	ONQ	C18-S15	2.37	1.80	1.75
2	A	601	ONQ	C22-N23	2.24	1.48	1.46
2	A	601	ONQ	C9-N3	2.12	1.43	1.39

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	ONQ	C22-N23-C24	-8.82	105.62	113.85
2	A	601	ONQ	C22-N23-C24	-7.96	106.42	113.85
2	B	601	ONQ	C18-S15-N14	6.33	113.69	106.56
2	A	601	ONQ	C18-S15-N14	5.48	112.73	106.56
2	A	601	ONQ	O16-S15-O17	-4.74	112.52	118.87
2	B	601	ONQ	O7-P4-C8	3.97	116.89	105.78
2	A	601	ONQ	C31-N23-C24	-3.30	120.06	123.71
2	B	601	ONQ	O26-C24-C1	-2.63	124.18	130.65
2	A	601	ONQ	O7-P4-C8	2.61	113.08	105.78
2	A	601	ONQ	O6-P4-C8	-2.47	108.49	113.75
2	B	601	ONQ	O16-S15-O17	-2.27	115.83	118.87
2	B	601	ONQ	C1-C24-N23	2.19	112.23	108.33
2	A	601	ONQ	C28-C27-C30	-2.10	104.72	108.80
2	A	601	ONQ	O26-C24-C1	-2.07	125.55	130.65

There are no chirality outliers.

All (16) torsion outliers are listed below:

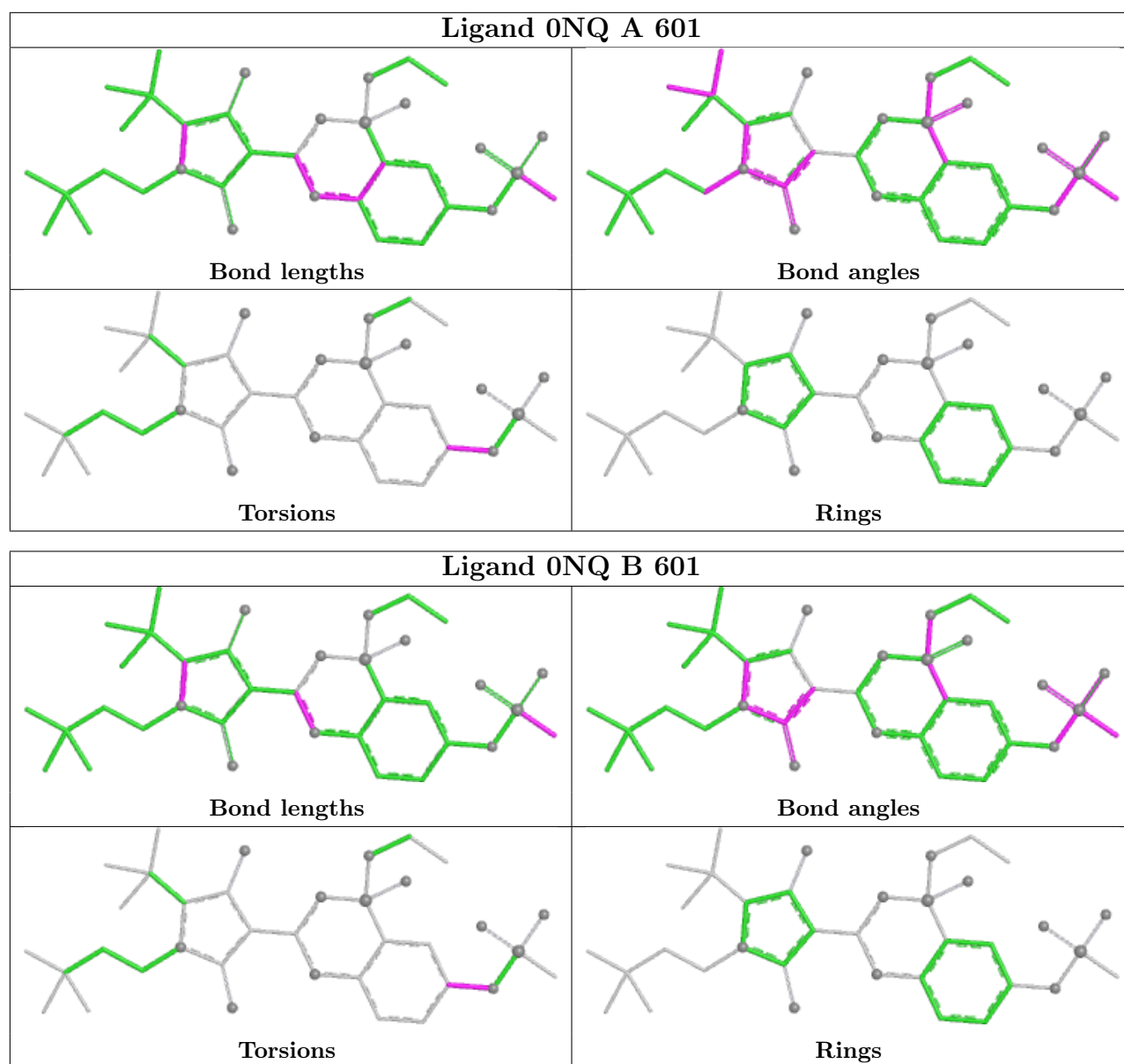
Mol	Chain	Res	Type	Atoms
4	A	608	PEG	O2-C3-C4-O4
4	A	608	PEG	C4-C3-O2-C2
4	B	604	PEG	C4-C3-O2-C2
4	A	609	PEG	O1-C1-C2-O2
4	B	604	PEG	O2-C3-C4-O4
5	A	611	EDO	O1-C1-C2-O2
4	A	609	PEG	O2-C3-C4-O4
4	B	604	PEG	C1-C2-O2-C3
4	A	608	PEG	C1-C2-O2-C3
2	A	601	0NQ	C11-C12-N14-S15
2	B	601	0NQ	C11-C12-N14-S15
2	A	601	0NQ	C13-C12-N14-S15
4	B	604	PEG	O1-C1-C2-O2
4	A	609	PEG	C1-C2-O2-C3
2	B	601	0NQ	C13-C12-N14-S15
5	A	610	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	609	PEG	4	0
2	B	601	0NQ	1	0

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	559/562 (99%)	-0.50	3 (0%) 87 87	7, 21, 44, 66	21 (3%)
1	B	559/562 (99%)	-0.02	28 (5%) 34 33	8, 32, 72, 94	22 (3%)
All	All	1118/1124 (99%)	-0.26	31 (2%) 55 54	7, 25, 62, 94	43 (3%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	23	ILE	6.5
1	B	149	PRO	4.8
1	A	149	PRO	4.7
1	B	26	LEU	4.5
1	B	25	PRO	4.4
1	B	22	PRO	4.0
1	B	543	SER	3.6
1	B	153	GLY	3.4
1	B	27	SER	3.3
1	B	531	ARG	3.2
1	B	542	ALA	3.2
1	B	405	ILE	3.1
1	B	404	PRO	3.1
1	B	541	ALA	2.9
1	B	532	THR	2.8
1	B	498	ARG	2.8
1	B	499	THR	2.8
1	B	377	SER	2.6
1	B	21	LEU	2.6
1	B	540	PRO	2.5
1	A	549	GLY	2.4
1	B	330	GLN	2.3
1	B	440	GLU	2.3
1	A	153	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	362	LEU	2.2
1	B	376	ALA	2.2
1	B	534	LEU	2.2
1	B	502	HIS	2.1
1	B	548	SER	2.1
1	B	402	HIS	2.1
1	B	535	LYS	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](#)

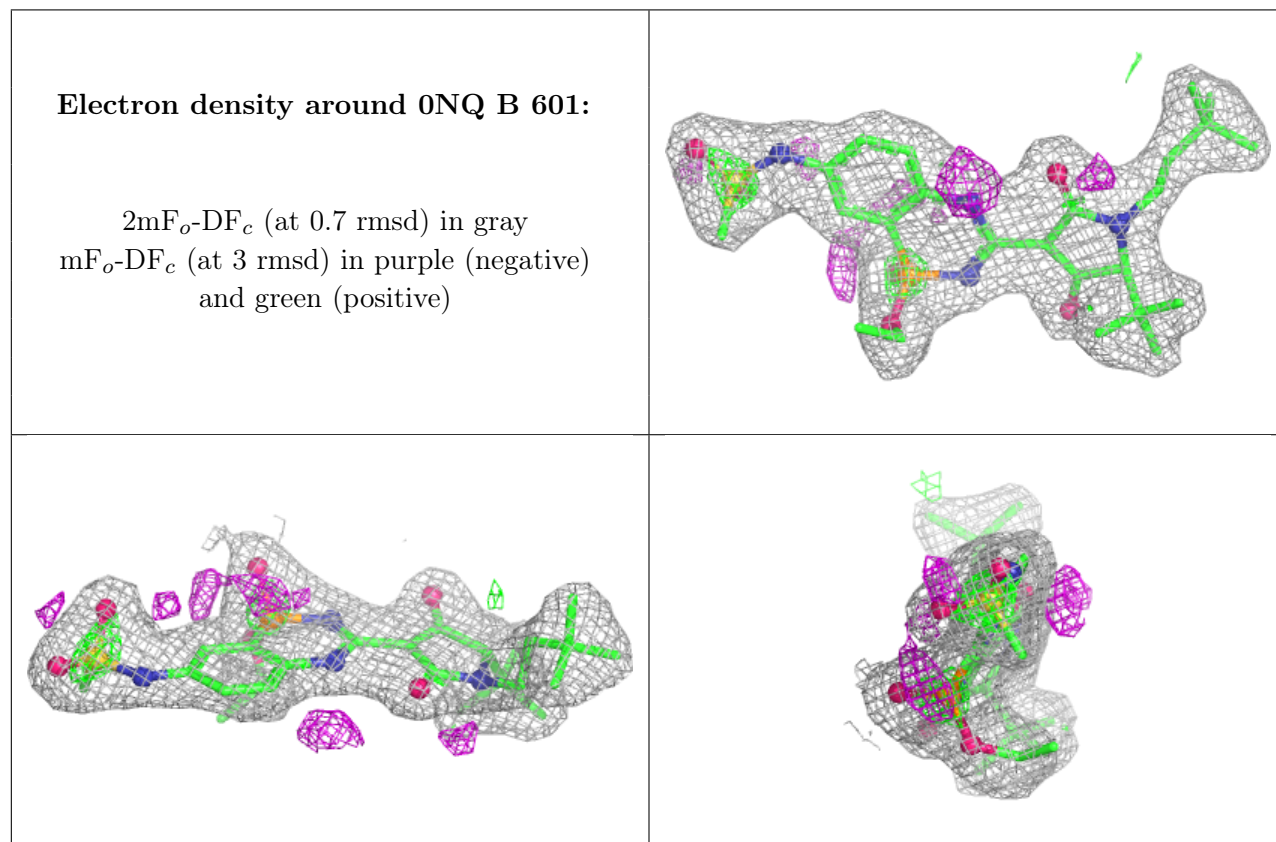
There are no oligosaccharides in this entry.

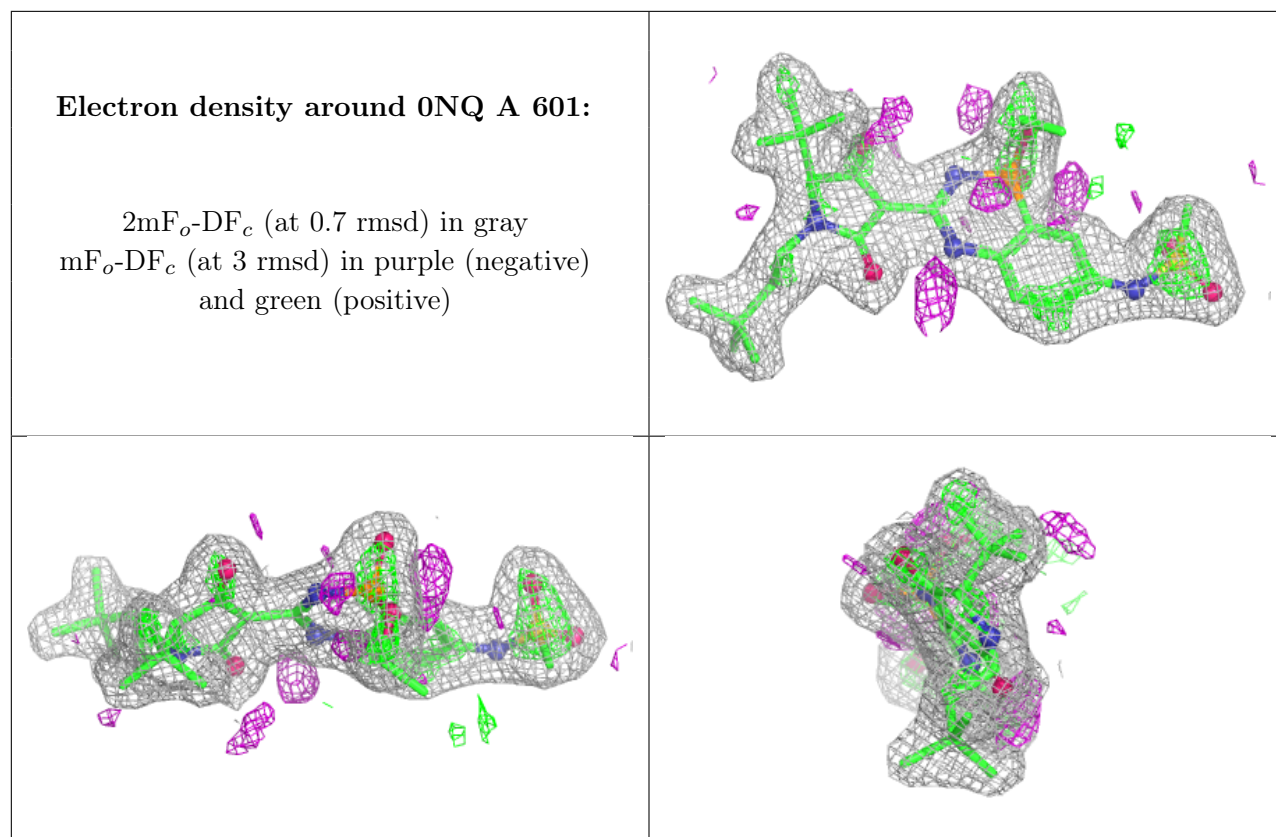
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
4	PEG	A	609	7/7	0.71	0.37	73,81,86,87	0
3	SO4	A	607	5/5	0.73	0.20	81,82,96,96	0
3	SO4	A	605	5/5	0.77	0.14	74,75,84,85	0
4	PEG	B	604	7/7	0.78	0.31	53,67,69,71	0
3	SO4	A	603	5/5	0.82	0.12	70,78,81,89	0
5	EDO	A	611	4/4	0.83	0.23	56,61,63,63	0
3	SO4	B	603	5/5	0.84	0.11	61,77,83,88	0
4	PEG	A	608	7/7	0.85	0.26	52,63,63,65	0
3	SO4	A	604	5/5	0.86	0.11	89,90,98,99	0
3	SO4	A	606	5/5	0.88	0.19	92,96,101,107	0
3	SO4	A	602	5/5	0.93	0.15	51,58,65,67	0
5	EDO	A	610	4/4	0.94	0.23	72,73,74,74	0
2	ONQ	B	601	36/36	0.95	0.14	34,47,54,64	0
3	SO4	B	602	5/5	0.95	0.10	53,54,62,63	0
2	ONQ	A	601	36/36	0.96	0.14	31,39,49,66	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.