



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:44 PM UTC

PDB ID : 4NCG / pdb_00004ncg
Title : Discovery of Doravirine, an orally bioavailable non-nucleoside reverse transcriptase inhibitor potent against a wide range of resistant mutant HIV viruses
Authors : Yan, Y.
Deposited on : 2013-10-24
Resolution : 2.58 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

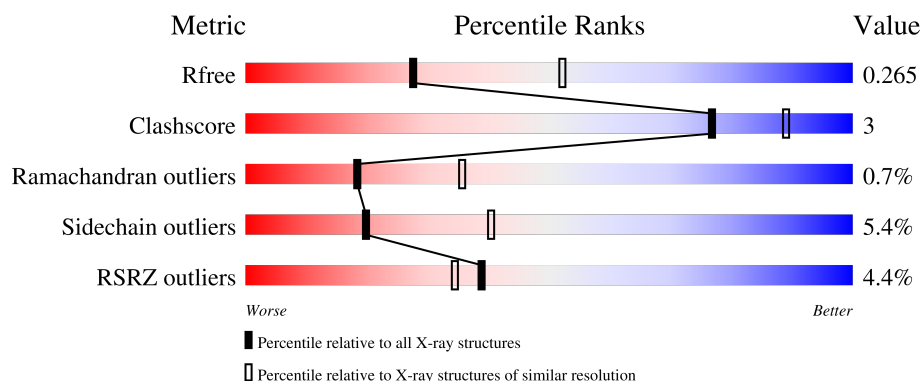
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	563	4% 82% 16% ..
2	B	443	5% 79% 11% • 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7913 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 Reverse transcriptase/ribonuclease H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	553	Total	C	N	O	S	0	0	0
			4491	2903	749	831	8			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	expression tag	UNP P04585
A	-1	ASN	-	expression tag	UNP P04585
A	0	SER	-	expression tag	UNP P04585

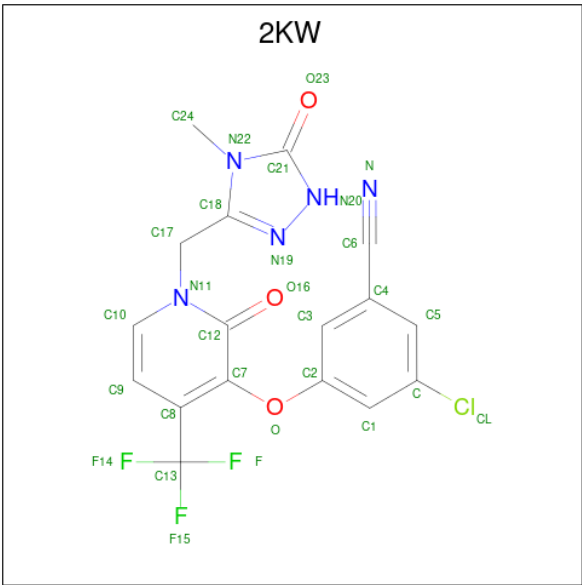
- Molecule 2 is a protein called p51 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	402	Total	C	N	O	S	0	0	0
			3307	2153	546	602	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	MET	LEU	engineered mutation	UNP P04585
B	0	SER	PHE	engineered mutation	UNP P04585

- Molecule 3 is 3-chloro-5-({1-[(4-methyl-5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2-oxo-4-(trifluoromethyl)-1,2-dihydropyridin-3-yl}oxy)benzonitrile (CCD ID: 2KW) (formula: C₁₇H₁₁ClF₃N₅O₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	0	0
			29	17	1	3	5	3		

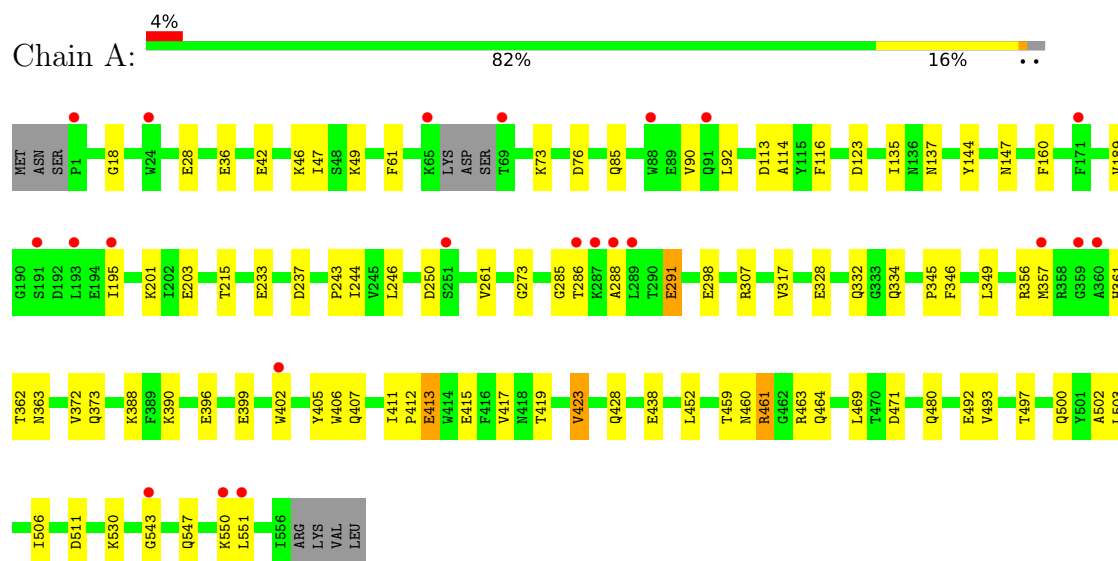
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	41	Total	O	0	0
			41	41		
4	B	45	Total	O	0	0
			45	45		

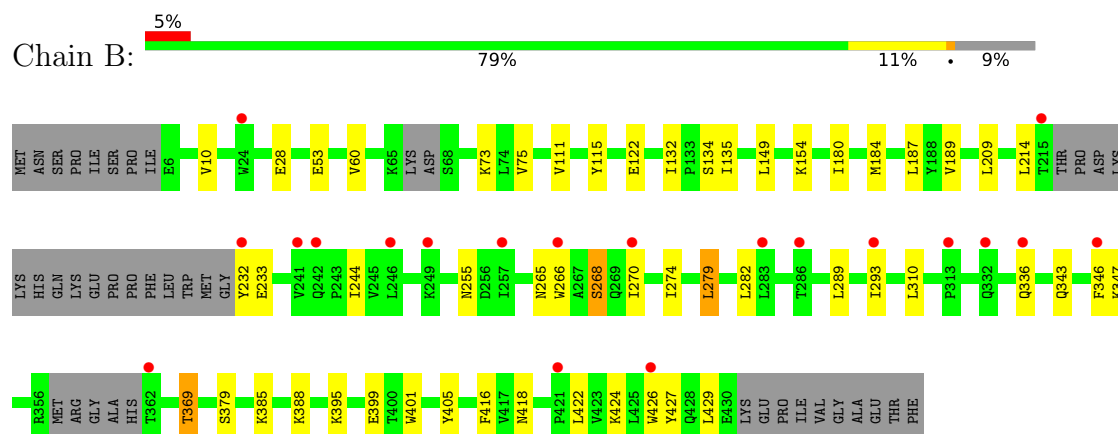
3 Residue-property plots [i](#)

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: Reverse transcriptase/ribonuclease H



- Molecule 2: p51 RT



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.39Å 153.62Å 154.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.01 – 2.58 26.01 – 2.58	Depositor EDS
% Data completeness (in resolution range)	96.3 (26.01-2.58) 96.2 (26.01-2.58)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 2.57Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.217 , 0.260 0.219 , 0.265	Depositor DCC
R_{free} test set	2163 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	69.5	Xtriage
Anisotropy	0.253	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7913	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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: 2KW

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	0.86	0/4607	1.28	21/6262 (0.3%)
2	B	0.84	0/3398	1.26	9/4620 (0.2%)
All	All	0.85	0/8005	1.27	30/10882 (0.3%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	423	VAL	N-CA-CB	7.49	118.89	110.72
2	B	416	PHE	CA-CB-CG	-6.43	107.37	113.80
1	A	285	GLY	CA-C-N	5.74	132.49	121.54
1	A	285	GLY	C-N-CA	5.74	132.49	121.54
2	B	343	GLN	N-CA-C	-5.70	106.98	114.04
1	A	388	LYS	N-CA-C	-5.59	100.12	109.24
1	A	113	ASP	CA-C-N	5.57	130.05	121.19
1	A	113	ASP	C-N-CA	5.57	130.05	121.19
1	A	250	ASP	CA-CB-CG	5.56	118.16	112.60
2	B	401	TRP	N-CA-C	5.51	121.03	113.97
2	B	347	LYS	CA-C-N	5.35	128.94	121.02
2	B	347	LYS	C-N-CA	5.35	128.94	121.02
2	B	293	ILE	CB-CA-C	5.33	115.30	110.13
1	A	147	ASN	N-CA-C	-5.32	107.34	113.88
1	A	349	LEU	N-CA-C	-5.29	104.94	111.40
1	A	261	VAL	CA-C-N	5.26	125.78	119.94
1	A	261	VAL	C-N-CA	5.26	125.78	119.94
1	A	250	ASP	N-CA-C	-5.23	105.25	111.69
1	A	237	ASP	CA-CB-CG	5.19	117.79	112.60
1	A	480	GLN	CA-C-N	5.14	127.16	120.28
1	A	480	GLN	C-N-CA	5.14	127.16	120.28
1	A	503	LEU	CA-C-N	5.12	125.77	119.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	503	LEU	C-N-CA	5.12	125.77	119.99
1	A	460	ASN	CA-C-N	5.08	130.78	122.65
1	A	460	ASN	C-N-CA	5.08	130.78	122.65
2	B	424	LYS	CA-C-N	5.07	127.03	120.44
2	B	424	LYS	C-N-CA	5.07	127.03	120.44
2	B	132	ILE	N-CA-C	-5.06	102.54	107.55
1	A	550	LYS	CA-C-N	5.05	127.46	120.29
1	A	550	LYS	C-N-CA	5.05	127.46	120.29

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	4491	0	4523	26	0
2	B	3307	0	3319	22	0
3	A	29	0	11	0	0
4	A	41	0	0	0	0
4	B	45	0	0	0	0
All	All	7913	0	7853	45	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 (45) 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:328:GLU:HG2	1:A:390:LYS:HB2	1.76	0.67
1:A:406:TRP:CH2	2:B:418:ASN:HA	2.33	0.64
1:A:46:LYS:HD2	1:A:116:PHE:HB3	1.81	0.62
1:A:405:TYR:CE2	1:A:407:GLN:HB3	2.35	0.61
2:B:154:LYS:HE2	2:B:184:MET:HE2	1.85	0.58
2:B:60:VAL:HG23	2:B:75:VAL:HG22	1.89	0.55
2:B:255:ASN:HB2	2:B:289:LEU:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:209:LEU:HD22	2:B:214:LEU:HD22	1.90	0.53
1:A:288:ALA:O	1:A:291:GLU:HB3	2.09	0.52
1:A:502:ALA:O	1:A:506:ILE:HG12	2.11	0.50
2:B:369:THR:HG21	2:B:405:TYR:HB2	1.94	0.50
2:B:266:TRP:HZ2	2:B:427:TYR:CZ	2.30	0.50
2:B:244:ILE:HB	2:B:310:LEU:HD22	1.94	0.49
1:A:500:GLN:CD	2:B:422:LEU:HD12	2.37	0.49
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.96	0.48
1:A:233:GLU:OE1	1:A:243:PRO:HD3	2.14	0.48
2:B:268:SER:HB3	2:B:274:ILE:HB	1.97	0.47
1:A:396:GLU:HA	1:A:399:GLU:HG2	1.96	0.46
1:A:28:GLU:HG2	1:A:135:ILE:HG12	1.97	0.46
2:B:111:VAL:HG11	2:B:187:LEU:HD22	1.98	0.46
1:A:438:GLU:HG3	1:A:461:ARG:HB2	1.98	0.46
2:B:266:TRP:CZ3	2:B:426:TRP:HB3	2.51	0.46
2:B:180:ILE:HG12	2:B:189:VAL:HG13	1.96	0.45
1:A:492:GLU:HG2	1:A:530:LYS:HB2	1.98	0.45
1:A:273:GLY:HA2	1:A:332:GLN:NE2	2.32	0.45
1:A:246:LEU:HD12	1:A:307:ARG:HD3	1.98	0.44
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.18	0.43
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.53	0.43
1:A:372:VAL:HG11	1:A:411:ILE:HG23	2.00	0.43
1:A:345:PRO:HA	1:A:346:PHE:HA	1.74	0.43
2:B:279:LEU:HA	2:B:282:LEU:HD12	1.99	0.43
1:A:412:PRO:O	1:A:413:GLU:C	2.61	0.43
1:A:61:PHE:HE1	1:A:76:ASP:HB2	1.84	0.43
1:A:201:LYS:HD3	1:A:201:LYS:HA	1.88	0.43
2:B:28:GLU:HA	2:B:135:ILE:HD11	2.01	0.43
2:B:209:LEU:HB3	2:B:214:LEU:HB2	2.01	0.42
1:A:417:VAL:HG22	1:A:419:THR:HG23	2.02	0.42
1:A:49:LYS:HG2	1:A:144:TYR:CE1	2.55	0.41
1:A:399:GLU:HA	1:A:402:TRP:CD1	2.55	0.41
2:B:270:ILE:HG12	2:B:346:PHE:O	2.20	0.41
2:B:232:TYR:N	2:B:233:GLU:HA	2.35	0.41
1:A:85:GLN:HE22	2:B:53:GLU:HA	1.86	0.41
2:B:154:LYS:CE	2:B:184:MET:HE2	2.49	0.41
1:A:42:GLU:HA	1:A:47:ILE:O	2.22	0.40
2:B:266:TRP:CZ2	2:B:427:TYR:CZ	3.09	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	549/563 (98%)	520 (95%)	22 (4%)	7 (1%)	9	19
2	B	394/443 (89%)	377 (96%)	17 (4%)	0	100	100
All	All	943/1006 (94%)	897 (95%)	39 (4%)	7 (1%)	18	36

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	334	GLN
1	A	543	GLY
1	A	286	THR
1	A	90	VAL
1	A	413	GLU
1	A	195	ILE
1	A	18	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	490/503 (97%)	459 (94%)	31 (6%)	16	34
2	B	362/403 (90%)	347 (96%)	15 (4%)	27	51
All	All	852/906 (94%)	806 (95%)	46 (5%)	20	40

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

Mol	Chain	Res	Type
1	A	36	GLU
1	A	73	LYS
1	A	92	LEU
1	A	123	ASP
1	A	137	ASN
1	A	189	VAL
1	A	203	GLU
1	A	215	THR
1	A	244	ILE
1	A	291	GLU
1	A	298	GLU
1	A	317	VAL
1	A	356	ARG
1	A	357	MET
1	A	361	HIS
1	A	362	THR
1	A	373	GLN
1	A	415	GLU
1	A	423	VAL
1	A	428	GLN
1	A	452	LEU
1	A	459	THR
1	A	461	ARG
1	A	463	ARG
1	A	464	GLN
1	A	469	LEU
1	A	471	ASP
1	A	493	VAL
1	A	497	THR
1	A	547	GLN
1	A	551	LEU
2	B	10	VAL
2	B	73	LYS
2	B	122	GLU
2	B	134	SER
2	B	265	ASN
2	B	268	SER
2	B	279	LEU
2	B	336	GLN
2	B	369	THR
2	B	379	SER
2	B	385	LYS
2	B	388	LYS

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Mol	Chain	Res	Type
2	B	395	LYS
2	B	399	GLU
2	B	429	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	147	ASN
1	A	332	GLN
1	A	334	GLN
1	A	487	GLN
1	A	520	GLN
2	B	91	GLN
2	B	147	ASN
2	B	269	GLN
2	B	278	GLN
2	B	394	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	2KW	A	601	-	31,31,31	1.19	4 (12%)	40,46,46	2.06	9 (22%)

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
3	2KW	A	601	-	-	0/16/16/16	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	2KW	C7-C12	2.92	1.51	1.43
3	A	601	2KW	C21-N20	2.77	1.37	1.35
3	A	601	2KW	C7-C8	-2.61	1.34	1.37
3	A	601	2KW	C17-C18	2.51	1.51	1.49

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	2KW	C9-C8-C7	7.60	121.25	114.94
3	A	601	2KW	C18-C17-N11	-6.51	102.89	111.63
3	A	601	2KW	F15-C13-C8	3.07	118.61	112.68
3	A	601	2KW	C17-N11-C12	2.97	119.56	117.38
3	A	601	2KW	C24-N22-C18	2.50	129.81	126.96
3	A	601	2KW	C12-C7-C8	-2.39	119.64	122.82
3	A	601	2KW	O23-C21-N22	2.31	128.43	126.04
3	A	601	2KW	O16-C12-N11	2.06	123.28	120.53
3	A	601	2KW	N22-C18-N19	2.00	112.16	111.51

There are no chirality outliers.

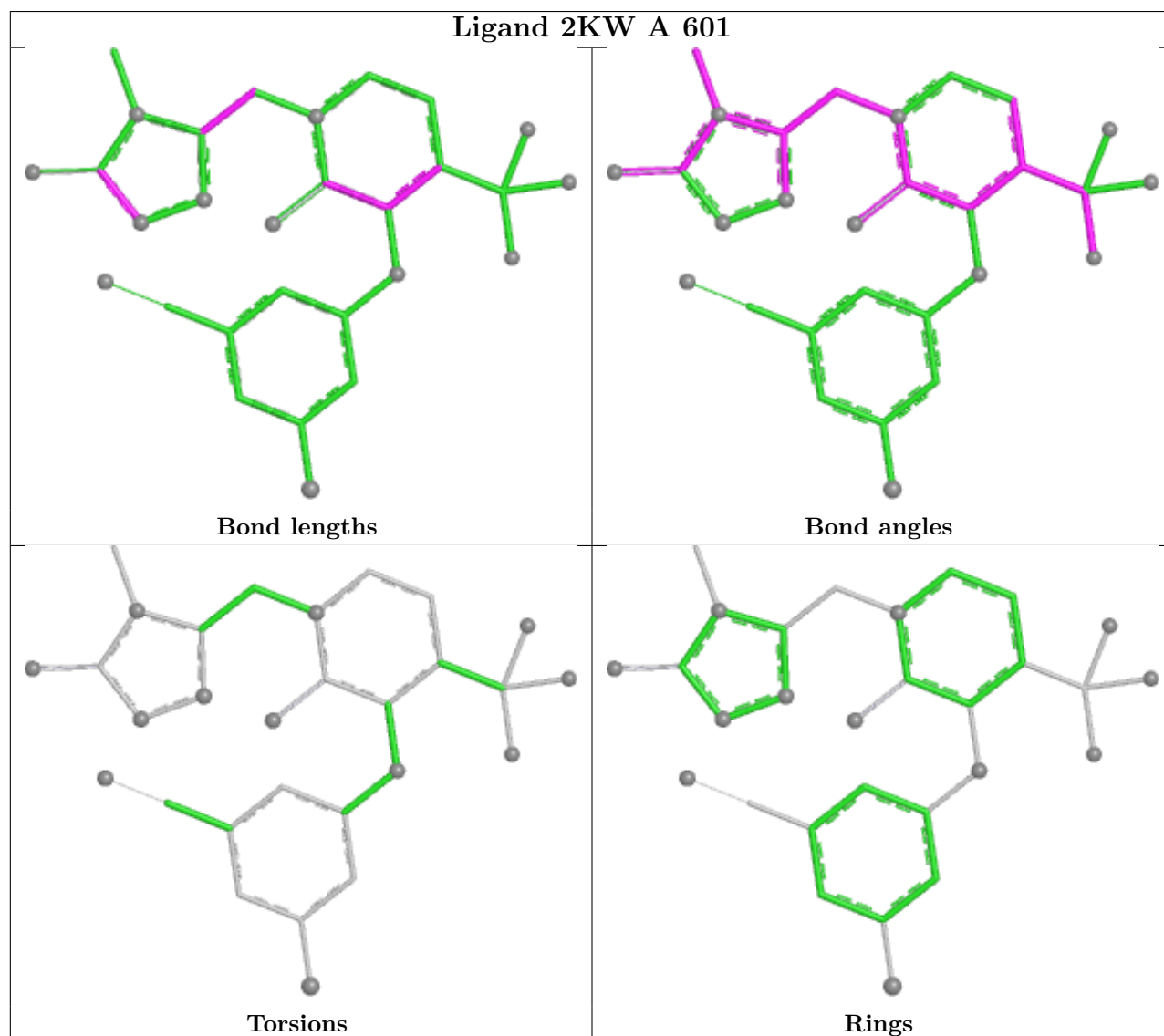
There are no torsion outliers.

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 ⓘ

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	553/563 (98%)	0.38	22 (3%)	42	38	47, 76, 110, 156	0
2	B	402/443 (90%)	0.33	20 (4%)	34	29	48, 71, 117, 140	0
All	All	955/1006 (94%)	0.36	42 (4%)	39	34	47, 75, 114, 156	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	215	THR	4.5
2	B	362	THR	3.9
2	B	24	TRP	3.6
2	B	286	THR	3.2
2	B	426	TRP	3.2
2	B	313	PRO	3.1
1	A	1	PRO	3.1
2	B	266	TRP	3.1
1	A	251	SER	3.0
1	A	193	LEU	3.0
1	A	286	THR	3.0
1	A	357	MET	3.0
1	A	91	GLN	2.9
2	B	241	VAL	2.9
2	B	232	TYR	2.7
1	A	88	TRP	2.5
1	A	69	THR	2.5
1	A	543	GLY	2.5
1	A	551	LEU	2.5
2	B	332	GLN	2.5
2	B	257	ILE	2.4
1	A	65	LYS	2.4
2	B	293	ILE	2.4
2	B	346	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	360	ALA	2.3
2	B	246	LEU	2.3
1	A	24	TRP	2.3
2	B	421	PRO	2.3
1	A	550	LYS	2.2
1	A	359	GLY	2.2
2	B	242	GLN	2.2
2	B	249	LYS	2.2
2	B	336	GLN	2.1
1	A	171	PHE	2.1
1	A	289	LEU	2.1
1	A	195	ILE	2.1
1	A	287	LYS	2.1
1	A	191	SER	2.1
1	A	288	ALA	2.1
2	B	283	LEU	2.1
1	A	402	TRP	2.0
2	B	270	ILE	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](#)

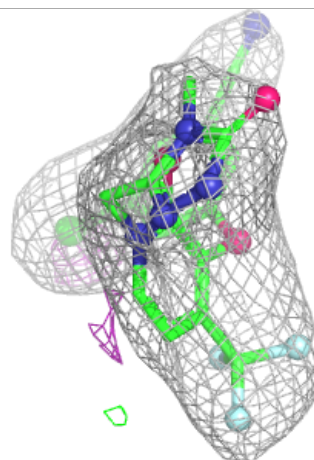
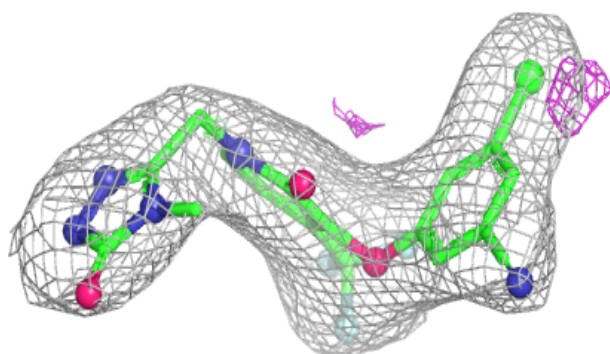
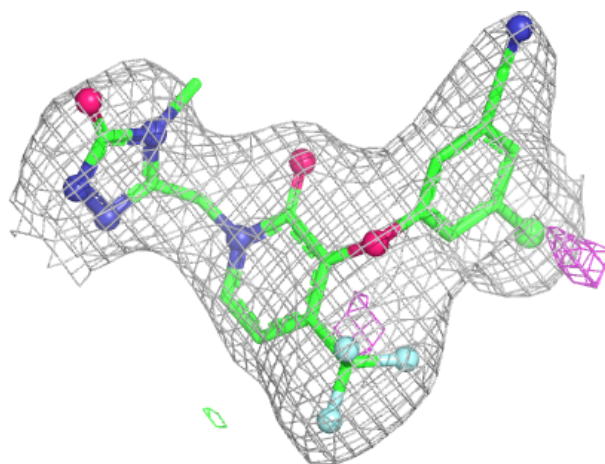
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	2KW	A	601	29/29	0.95	0.07	54,66,80,83	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 2KW A 601:

$2mF_o-DF_c$ (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.