



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:22 PM UTC

PDB ID : 5TF5 / pdb_00005tf5
Title : CRYSTAL STRUCTURE OF HUMAN KAT-2 IN COMPLEX WITH A REVERSIBLE INHIBITOR
Authors : Nematollahi, A.; Sun, G.; Church, W.B.
Deposited on : 2016-09-24
Resolution : 1.81 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

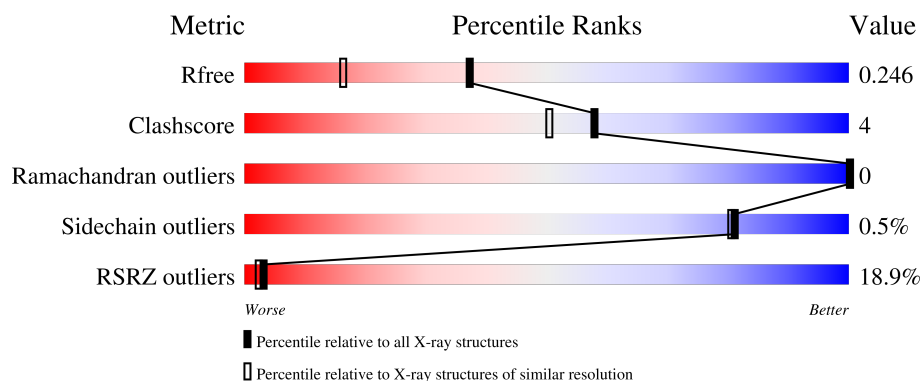
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	425	21% 92% 8%
1	B	425	16% 92% 8%

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
1	LLP	B	263	-	-	X	-
2	7AR	B	501	-	-	X	-

2 Entry composition [i](#)

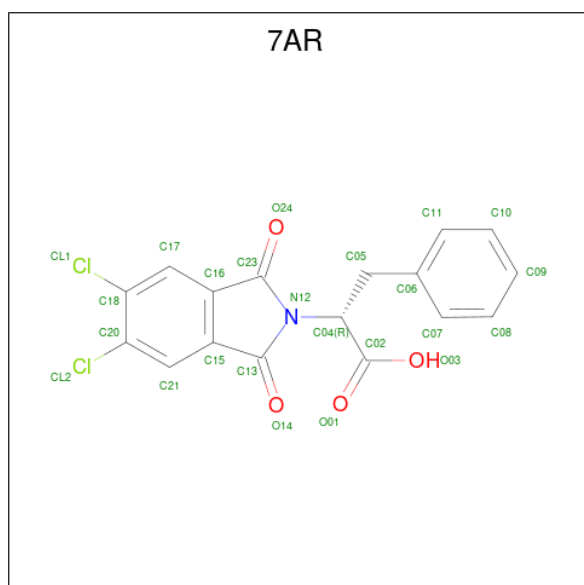
There are 3 unique types of molecules in this entry. The entry contains 14080 atoms, of which 6680 are hydrogens and 0 are deuteriums.

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 Kynurenine/ α -aminoadipate aminotransferase, mitochondrial.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	425	Total	C	H	N	O	P	S	0	0	0
			6687	2147	3340	560	621	1	18			
1	B	425	Total	C	H	N	O	P	S	0	0	0
			6687	2147	3340	560	621	1	18			

- Molecule 2 is (2R)-2-(5,6-dichloro-1,3-dioxo-1,3-dihydro-2H-isoindol-2-yl)-3-phenylpropanoic acid (CCD ID: 7AR) (formula: $C_{17}H_{11}Cl_2NO_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	Cl	N	O	0	0
			24	17	2	1	4		

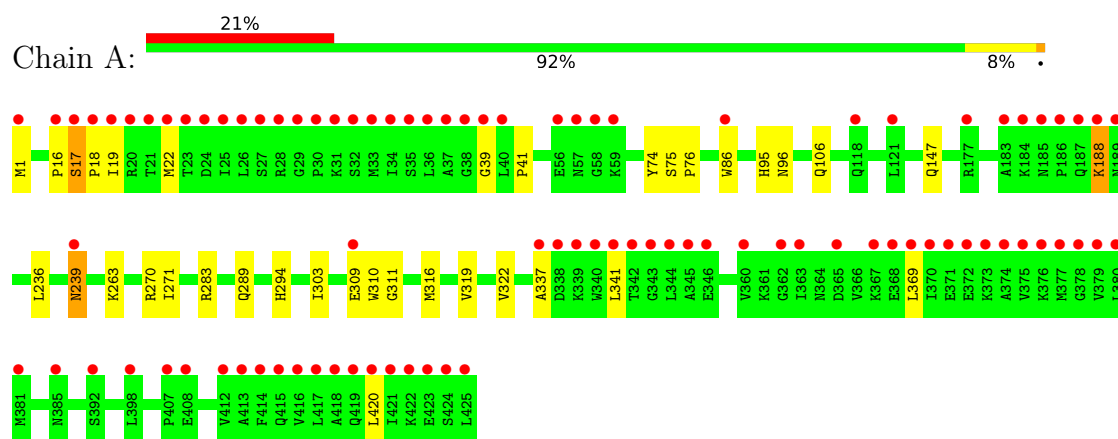
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	326	Total 326	O 326	0	0
3	B	356	Total 356	O 356	0	0

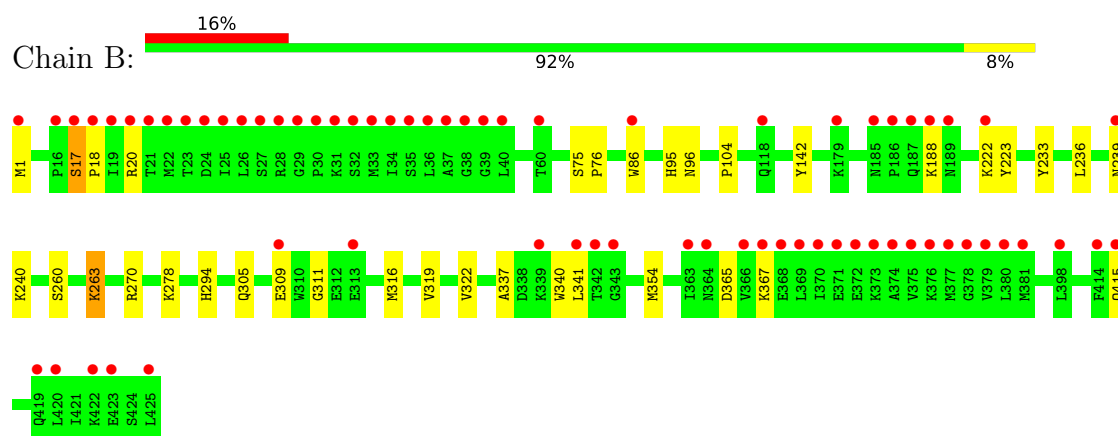
3 Residue-property plots [i](#)

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: Kynurenine/alpha-aminoadipate aminotransferase, mitochondrial



- Molecule 1: Kynurenine/alpha-aminoadipate aminotransferase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.88Å 99.28Å 107.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.13 – 1.81 46.13 – 1.81	Depositor EDS
% Data completeness (in resolution range)	98.3 (46.13-1.81) 89.6 (46.13-1.81)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.67 (at 1.81Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: 2015)	Depositor
R, R_{free}	0.219 , 0.245 0.220 , 0.246	Depositor DCC
R_{free} test set	1995 reflections (2.27%)	wwPDB-VP
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.688	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14080	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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: LLP, 7AR

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	0.23	0/3404	0.42	0/4620
1	B	0.26	0/3404	0.45	0/4620
All	All	0.25	0/6808	0.44	0/9240

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	17	SER	Peptide
1	A	188	LYS	Peptide
1	B	17	SER	Peptide
1	B	188	LYS	Peptide

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	3347	3340	3358	26	1
1	B	3347	3340	3358	36	1
2	B	24	0	0	20	0
3	A	326	0	0	2	4
3	B	356	0	0	1	4
All	All	7400	6680	6716	60	5

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 (60) 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:263:LLP:H4'1	2:B:501:7AR:C04	1.79	1.12
1:B:263:LLP:NZ	2:B:501:7AR:C06	2.53	0.72
1:B:263:LLP:NZ	2:B:501:7AR:C11	2.58	0.67
1:B:263:LLP:C4'	2:B:501:7AR:C07	2.75	0.65
1:B:263:LLP:H4'1	2:B:501:7AR:C07	2.29	0.63
1:B:263:LLP:C4'	2:B:501:7AR:C06	2.77	0.63
1:B:263:LLP:HD3	2:B:501:7AR:C10	2.31	0.61
1:A:294:HIS:CE1	1:B:270:ARG:HD3	2.38	0.58
1:A:239:ASN:OD1	1:A:239:ASN:N	2.35	0.58
1:A:39:GLY:O	2:B:501:7AR:C06	2.52	0.57
1:A:16:PRO:O	3:A:501:HOH:O	2.17	0.56
1:A:270:ARG:HD3	1:B:294:HIS:CE1	2.42	0.55
1:B:263:LLP:H4'1	2:B:501:7AR:C06	2.38	0.54
1:A:17:SER:HB3	1:A:18:PRO:CD	2.37	0.54
1:B:17:SER:HB3	1:B:18:PRO:HD3	1.90	0.53
1:B:86:TRP:C	1:B:86:TRP:CD1	2.86	0.53
1:B:263:LLP:C4'	2:B:501:7AR:C04	2.70	0.53
1:A:19:ILE:HG12	2:B:501:7AR:O24	2.10	0.52
1:A:86:TRP:HH2	1:A:311:GLY:C	2.18	0.51
1:B:236:LEU:HD21	1:B:322:VAL:HG12	1.93	0.51
1:B:260:SER:OG	1:B:263:LLP:OP1	2.28	0.51
1:B:142:TYR:CE2	2:B:501:7AR:O14	2.64	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:TRP:C	1:A:86:TRP:CD1	2.89	0.50
1:B:354:MET:SD	2:B:501:7AR:C10	3.00	0.49
1:B:17:SER:HB3	1:B:18:PRO:CD	2.42	0.49
1:A:337:ALA:HA	1:A:341:LEU:HD23	1.95	0.48
1:A:147:GLN:OE1	3:A:502:HOH:O	2.20	0.47
1:A:316:MET:O	1:A:319:VAL:HG22	2.14	0.47
1:A:39:GLY:O	2:B:501:7AR:C11	2.62	0.47
1:B:17:SER:C	1:B:20:ARG:HE	2.22	0.47
1:B:233:TYR:OH	2:B:501:7AR:C08	2.63	0.47
1:A:41:PRO:HD3	2:B:501:7AR:C11	2.45	0.46
1:B:316:MET:O	1:B:319:VAL:HG22	2.16	0.46
1:B:365:ASP:OD1	1:B:367:LYS:HG2	2.16	0.45
1:B:104:PRO:HB3	1:B:278:LYS:HD3	1.98	0.45
1:A:17:SER:HB3	1:A:18:PRO:HD3	1.98	0.45
1:B:305:GLN:O	1:B:309:GLU:HG2	2.18	0.44
1:B:222:LYS:HD3	1:B:223:TYR:CE2	2.52	0.44
1:A:236:LEU:HD21	1:A:322:VAL:HG12	2.00	0.44
1:B:270:ARG:NH2	3:B:614:HOH:O	2.49	0.44
1:A:271:ILE:HG13	1:A:303:ILE:HD12	1.98	0.44
1:B:95:HIS:O	1:B:96:ASN:C	2.62	0.43
1:B:86:TRP:HH2	1:B:311:GLY:C	2.26	0.43
1:A:309:GLU:HG3	1:A:310:TRP:N	2.34	0.43
1:A:75:SER:HB2	1:A:76:PRO:CD	2.48	0.43
1:B:340:TRP:CE2	1:B:415:GLN:HG3	2.54	0.43
1:A:22:MET:HE3	2:B:501:7AR:CL1	2.56	0.42
1:B:263:LLP:H4'1	2:B:501:7AR:C05	2.45	0.42
1:A:283:ARG:HD2	1:A:283:ARG:N	2.35	0.42
1:B:75:SER:HB2	1:B:76:PRO:CD	2.49	0.42
1:B:340:TRP:CZ2	1:B:415:GLN:CG	3.02	0.42
1:B:354:MET:SD	2:B:501:7AR:C09	3.08	0.41
1:A:95:HIS:O	1:A:96:ASN:C	2.64	0.41
1:A:17:SER:HB2	1:A:289:GLN:HB3	2.03	0.41
1:B:17:SER:CB	1:B:18:PRO:CD	2.99	0.41
1:A:74:TYR:CE2	2:B:501:7AR:C05	3.03	0.40
1:A:188:LYS:O	1:A:188:LYS:CG	2.69	0.40
1:B:337:ALA:HA	1:B:341:LEU:HD23	2.02	0.40
1:B:340:TRP:CZ2	1:B:415:GLN:HG3	2.56	0.40
1:A:369:LEU:HD11	1:A:420:LEU:HB3	2.02	0.40

All (5) 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)
3:A:786:HOH:O	3:B:922:HOH:O[1_655]	2.00	0.20
3:A:738:HOH:O	3:B:920:HOH:O[3_745]	2.04	0.16
1:A:106:GLN:OE1	1:B:240:LYS:HZ2[2_655]	1.48	0.12
3:A:531:HOH:O	3:B:879:HOH:O[3_745]	2.09	0.11
3:A:810:HOH:O	3:B:951:HOH:O[2_655]	2.15	0.05

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	422/425 (99%)	404 (96%)	18 (4%)	0	100	100
1	B	422/425 (99%)	400 (95%)	22 (5%)	0	100	100
All	All	844/850 (99%)	804 (95%)	40 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	369/369 (100%)	367 (100%)	2 (0%)	81	80
1	B	369/369 (100%)	367 (100%)	2 (0%)	81	80
All	All	738/738 (100%)	734 (100%)	4 (0%)	81	80

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	239	ASN
1	B	1	MET
1	B	239	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	199	ASN
1	B	108	GLN
1	B	199	ASN
1	B	237	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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$
1	LLP	A	263	1	23,24,25	3.86	7 (30%)	25,32,34	2.00	7 (28%)
1	LLP	B	263	1	23,24,25	3.93	8 (34%)	25,32,34	1.94	7 (28%)

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
1	LLP	A	263	1	-	6/16/17/19	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	LLP	B	263	1	-	6/16/17/19	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	263	LLP	C3-C2	10.98	1.52	1.41
1	A	263	LLP	C3-C2	10.37	1.51	1.41
1	B	263	LLP	C4'-NZ	8.45	1.55	1.27
1	A	263	LLP	C4'-NZ	8.42	1.55	1.27
1	A	263	LLP	C4-C5	7.87	1.52	1.42
1	B	263	LLP	C4-C3	7.62	1.53	1.41
1	B	263	LLP	C4-C5	7.45	1.52	1.42
1	A	263	LLP	C4-C3	7.27	1.53	1.41
1	A	263	LLP	C4-C4'	4.38	1.56	1.46
1	B	263	LLP	C4-C4'	3.66	1.54	1.46
1	B	263	LLP	C2-N1	3.08	1.39	1.33
1	A	263	LLP	C2-N1	3.00	1.39	1.33
1	A	263	LLP	C6-C5	2.69	1.43	1.37
1	B	263	LLP	C6-C5	2.60	1.42	1.37
1	B	263	LLP	C6-N1	2.13	1.38	1.34

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	LLP	C3-C4-C5	-5.30	114.02	118.28
1	B	263	LLP	C3-C4-C5	-5.05	114.22	118.28
1	A	263	LLP	OP3-P-OP2	3.74	121.82	107.80
1	B	263	LLP	CD-CE-NZ	3.48	120.05	110.83
1	B	263	LLP	OP4-C5'-C5	3.40	115.74	109.36
1	A	263	LLP	CD-CE-NZ	3.35	119.70	110.83
1	A	263	LLP	OP4-C5'-C5	3.02	115.01	109.36
1	A	263	LLP	CE-NZ-C4'	2.77	127.59	118.72
1	B	263	LLP	OP3-P-OP1	2.67	121.23	110.83
1	B	263	LLP	CE-NZ-C4'	2.61	127.08	118.72
1	A	263	LLP	C2'-C2-C3	-2.61	117.74	120.80
1	B	263	LLP	C2'-C2-N1	2.57	122.49	117.64
1	A	263	LLP	C2'-C2-N1	2.53	122.41	117.64
1	B	263	LLP	C2'-C2-C3	-2.26	118.15	120.80

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	263	LLP	C4-C4'-NZ-CE
1	A	263	LLP	C5'-OP4-P-OP1
1	A	263	LLP	C5'-OP4-P-OP2
1	A	263	LLP	C5'-OP4-P-OP3
1	B	263	LLP	C4-C4'-NZ-CE
1	B	263	LLP	C5'-OP4-P-OP1
1	B	263	LLP	C5'-OP4-P-OP2
1	B	263	LLP	C5'-OP4-P-OP3
1	A	263	LLP	C3-C4-C4'-NZ
1	B	263	LLP	C3-C4-C4'-NZ
1	B	263	LLP	CD-CE-NZ-C4'
1	A	263	LLP	CD-CE-NZ-C4'

There are no ring outliers.

1 monomer is involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	263	LLP	11	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
2	7AR	B	501	-	26,26,26	3.14	8 (30%)	38,38,38	1.42	6 (15%)

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	7AR	B	501	-	-	6/12/28/28	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	7AR	O14-C13	8.83	1.40	1.22
2	B	501	7AR	O24-C23	8.81	1.40	1.22
2	B	501	7AR	C15-C13	-5.21	1.40	1.48
2	B	501	7AR	C16-C23	-4.45	1.41	1.48
2	B	501	7AR	C13-N12	-4.38	1.31	1.40
2	B	501	7AR	C23-N12	-3.81	1.32	1.40
2	B	501	7AR	C18-CL1	2.42	1.79	1.73
2	B	501	7AR	C20-CL2	2.22	1.78	1.73

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	7AR	C06-C05-C04	-3.44	105.33	113.31
2	B	501	7AR	C17-C16-C23	2.80	134.19	129.39
2	B	501	7AR	C16-C15-C13	-2.51	106.06	108.26
2	B	501	7AR	C15-C13-N12	2.48	108.70	105.96
2	B	501	7AR	O03-C02-C04	2.17	120.42	113.34
2	B	501	7AR	C21-C15-C13	2.12	133.03	129.39

There are no chirality outliers.

All (6) torsion outliers are listed below:

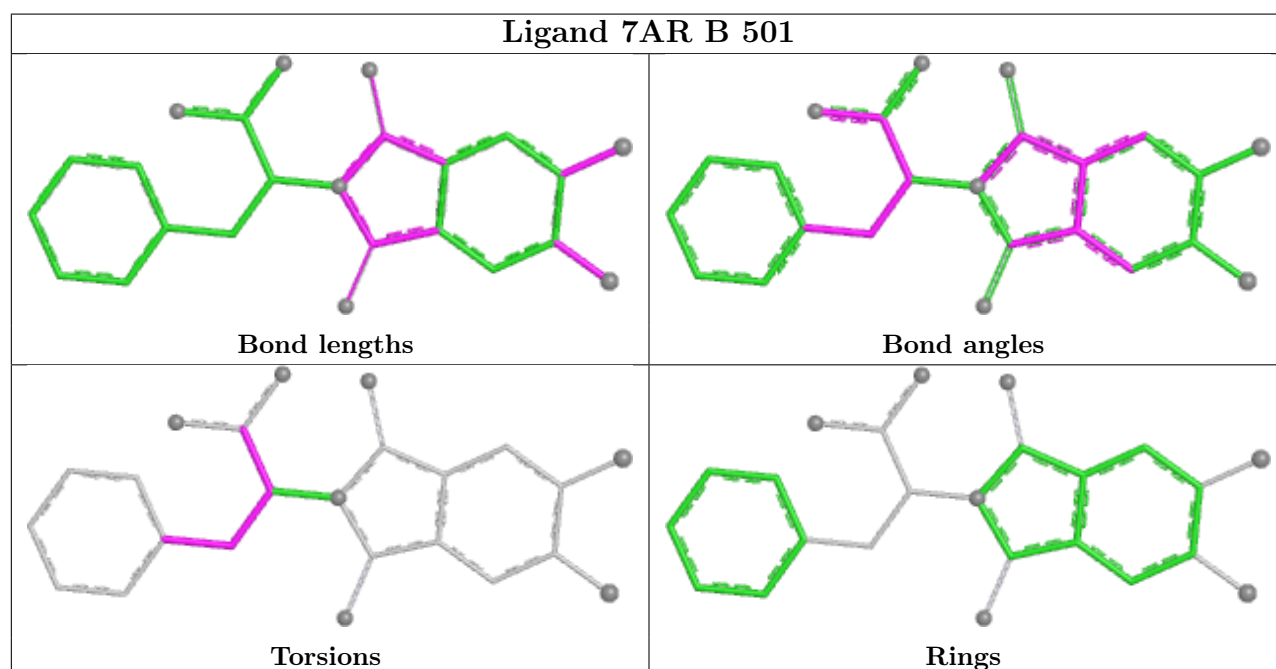
Mol	Chain	Res	Type	Atoms
2	B	501	7AR	N12-C04-C05-C06
2	B	501	7AR	C02-C04-C05-C06
2	B	501	7AR	C04-C05-C06-C11
2	B	501	7AR	C04-C05-C06-C07
2	B	501	7AR	O01-C02-C04-N12
2	B	501	7AR	O03-C02-C04-N12

There are no ring outliers.

1 monomer is involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	7AR	20	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	424/425 (99%)	1.07	91 (21%) 2 2	14, 27, 97, 167	0
1	B	424/425 (99%)	0.83	69 (16%) 4 3	14, 24, 96, 168	0
All	All	848/850 (99%)	0.95	160 (18%) 3 2	14, 26, 98, 168	0

All (160) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	18	PRO	12.0
1	A	342	THR	10.9
1	A	18	PRO	10.3
1	B	17	SER	9.4
1	B	20	ARG	9.1
1	B	342	THR	9.0
1	B	21	THR	8.8
1	B	19	ILE	8.6
1	B	186	PRO	8.5
1	A	19	ILE	8.5
1	B	379	VAL	8.2
1	B	25	ILE	7.8
1	A	21	THR	7.7
1	B	370	ILE	7.7
1	B	26	LEU	7.5
1	A	26	LEU	7.4
1	A	34	ILE	7.2
1	A	25	ILE	7.1
1	B	188	LYS	7.1
1	A	30	PRO	6.7
1	B	222	LYS	6.6
1	A	341	LEU	6.5
1	A	23	THR	6.3
1	B	23	THR	6.3

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Mol	Chain	Res	Type	RSRZ
1	B	16	PRO	6.0
1	B	375	VAL	5.9
1	B	30	PRO	5.9
1	A	379	VAL	5.8
1	B	380	LEU	5.8
1	B	374	ALA	5.8
1	A	188	LYS	5.7
1	A	20	ARG	5.7
1	A	186	PRO	5.7
1	A	425	LEU	5.6
1	A	35	SER	5.5
1	A	17	SER	5.5
1	A	29	GLY	5.4
1	A	57	ASN	5.4
1	A	58	GLY	5.3
1	A	22	MET	5.3
1	A	59	LYS	5.3
1	A	380	LEU	5.2
1	A	370	ILE	5.2
1	A	36	LEU	5.2
1	A	378	GLY	5.1
1	A	375	VAL	5.0
1	B	34	ILE	4.9
1	A	340	TRP	4.9
1	A	343	GLY	4.9
1	B	377	MET	4.8
1	A	32	SER	4.7
1	A	37	ALA	4.7
1	B	22	MET	4.7
1	B	27	SER	4.6
1	A	367	LYS	4.5
1	A	376	LYS	4.5
1	B	339	LYS	4.5
1	A	377	MET	4.4
1	A	33	MET	4.4
1	A	369	LEU	4.4
1	A	368	GLU	4.3
1	B	369	LEU	4.3
1	B	341	LEU	4.2
1	A	31	LYS	4.2
1	A	363	ILE	4.2
1	B	425	LEU	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	38	GLY	4.2
1	A	24	ASP	4.1
1	B	24	ASP	4.1
1	A	424	SER	4.1
1	B	32	SER	4.1
1	B	33	MET	4.0
1	A	374	ALA	4.0
1	A	344	LEU	4.0
1	A	420	LEU	4.0
1	A	27	SER	3.9
1	B	367	LYS	3.9
1	B	378	GLY	3.9
1	A	28	ARG	3.8
1	B	420	LEU	3.7
1	A	373	LYS	3.7
1	B	373	LYS	3.7
1	B	29	GLY	3.7
1	B	31	LYS	3.7
1	B	86	TRP	3.7
1	A	339	LYS	3.6
1	B	28	ARG	3.6
1	A	40	LEU	3.6
1	B	372	GLU	3.6
1	B	37	ALA	3.5
1	B	187	GLN	3.5
1	B	363	ILE	3.5
1	B	376	LYS	3.5
1	A	421	ILE	3.5
1	B	368	GLU	3.4
1	B	35	SER	3.4
1	A	86	TRP	3.4
1	A	337	ALA	3.3
1	B	343	GLY	3.3
1	B	38	GLY	3.3
1	A	16	PRO	3.3
1	A	422	LYS	3.2
1	A	414	PHE	3.2
1	A	419	GLN	3.2
1	A	423	GLU	3.2
1	B	366	VAL	3.2
1	A	56	GLU	3.2
1	B	415	GLN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	417	LEU	3.1
1	A	39	GLY	3.1
1	B	381	MET	3.1
1	A	416	VAL	3.0
1	A	187	GLN	3.0
1	A	185	ASN	3.0
1	B	423	GLU	3.0
1	B	60	THR	2.9
1	A	398	LEU	2.9
1	A	372	GLU	2.9
1	B	39	GLY	2.8
1	B	118	GLN	2.8
1	A	309	GLU	2.8
1	B	419	GLN	2.8
1	A	189	ASN	2.7
1	A	184	LYS	2.7
1	B	36	LEU	2.7
1	B	309	GLU	2.7
1	B	371	GLU	2.7
1	B	1	MET	2.6
1	B	414	PHE	2.6
1	B	364	ASN	2.6
1	A	365	ASP	2.5
1	A	239	ASN	2.5
1	B	239	ASN	2.5
1	A	371	GLU	2.5
1	B	40	LEU	2.5
1	A	392	SER	2.5
1	A	183	ALA	2.5
1	A	412	VAL	2.5
1	B	189	ASN	2.4
1	A	121	LEU	2.4
1	A	415	GLN	2.4
1	A	1	MET	2.4
1	B	398	LEU	2.3
1	A	338	ASP	2.3
1	B	179	LYS	2.3
1	B	313	GLU	2.2
1	A	118	GLN	2.2
1	A	381	MET	2.2
1	A	407	PRO	2.2
1	A	360	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	177	ARG	2.2
1	A	418	ALA	2.1
1	B	185	ASN	2.1
1	A	362	GLY	2.1
1	B	422	LYS	2.1
1	A	385	ASN	2.1
1	A	345	ALA	2.1
1	A	413	ALA	2.0
1	A	346	GLU	2.0
1	A	408	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	LLP	A	263	24/25	0.85	0.25	8,20,81,90	0
1	LLP	B	263	24/25	0.86	0.27	4,19,79,86	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

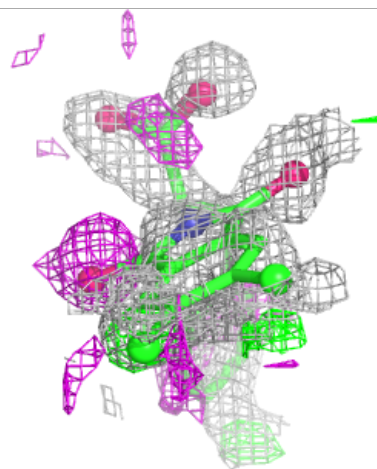
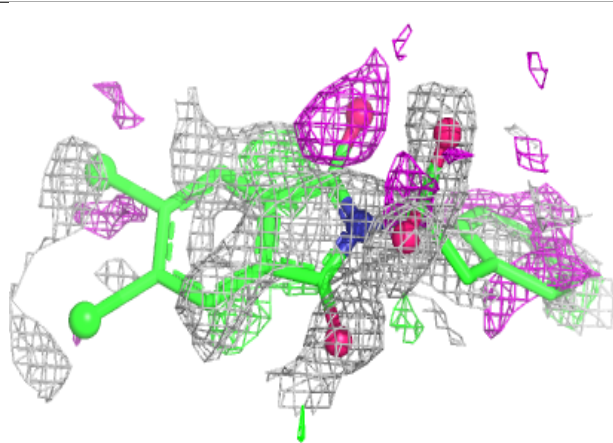
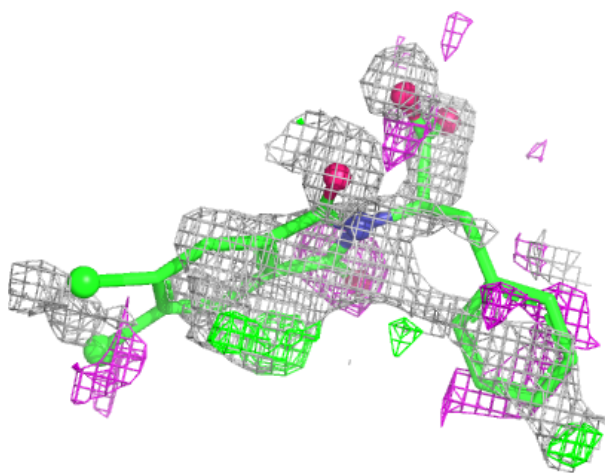
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	7AR	B	501	24/24	0.50	0.29	36,60,97,117	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 7AR B 501:

$2mF_o-DF_c$ (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.