



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 04:35 pm BST

PDB ID : 4JA8
Title : Complex of Mitochondrial Isocitrate Dehydrogenase R140Q Mutant with AGI-6780 Inhibitor
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Deposited on : 2013-02-18
Resolution : 1.55 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

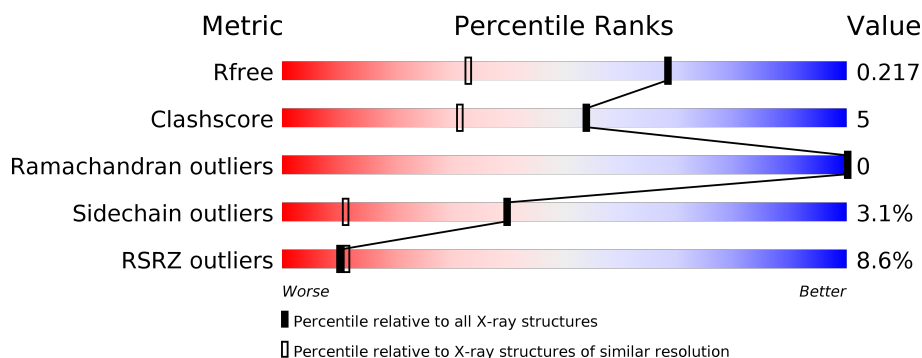
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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}	130704	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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 on 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	416	<div> <div></div> <div> <div></div> <div>84%</div> <div>16%</div> </div> </div>
1	B	416	<div> <div>16%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div>••</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7629 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 Isocitrate dehydrogenase [NADP], mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	13	0
			3428	2175	602	633	18			
1	B	406	Total	C	N	O	S	0	6	0
			3269	2081	562	607	19			

There are 10 discrepancies between the modelled and reference sequences:

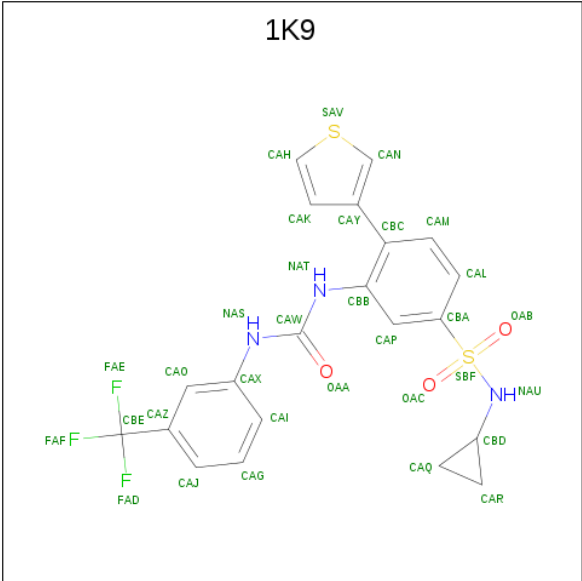
Chain	Residue	Modelled	Actual	Comment	Reference
A	140	GLN	ARG	ENGINEERED MUTATION	UNP P48735
A	453	HIS	-	EXPRESSION TAG	UNP P48735
A	454	HIS	-	EXPRESSION TAG	UNP P48735
A	455	HIS	-	EXPRESSION TAG	UNP P48735
A	456	HIS	-	EXPRESSION TAG	UNP P48735
B	140	GLN	ARG	ENGINEERED MUTATION	UNP P48735
B	453	HIS	-	EXPRESSION TAG	UNP P48735
B	454	HIS	-	EXPRESSION TAG	UNP P48735
B	455	HIS	-	EXPRESSION TAG	UNP P48735
B	456	HIS	-	EXPRESSION TAG	UNP P48735

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is 1-[5-(cyclopropylsulfamoyl)-2-thiophen-3-yl-phenyl]-3-[3-(trifluoromethyl)phenyl]urea (three-letter code: 1K9) (formula: C₂₁H₁₈F₃N₃O₃S₂).

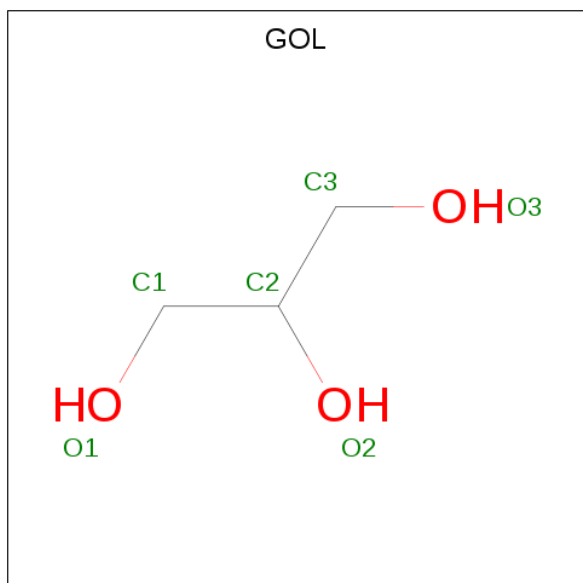


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	S	0	1
			64	42	6	6	6	4		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ca 1 1	0	0
4	A	1	Total Ca 1 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0

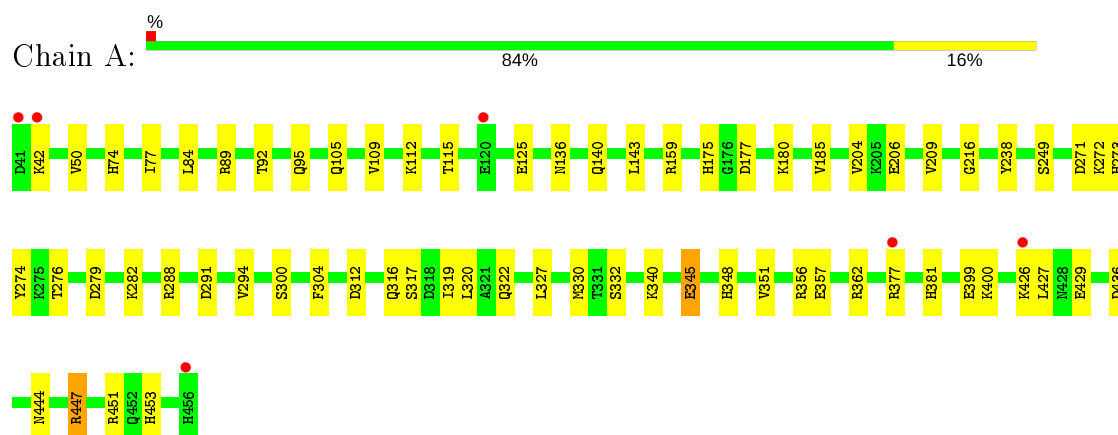
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	469	Total O 469 469	0	0
6	B	277	Total O 277 277	0	0

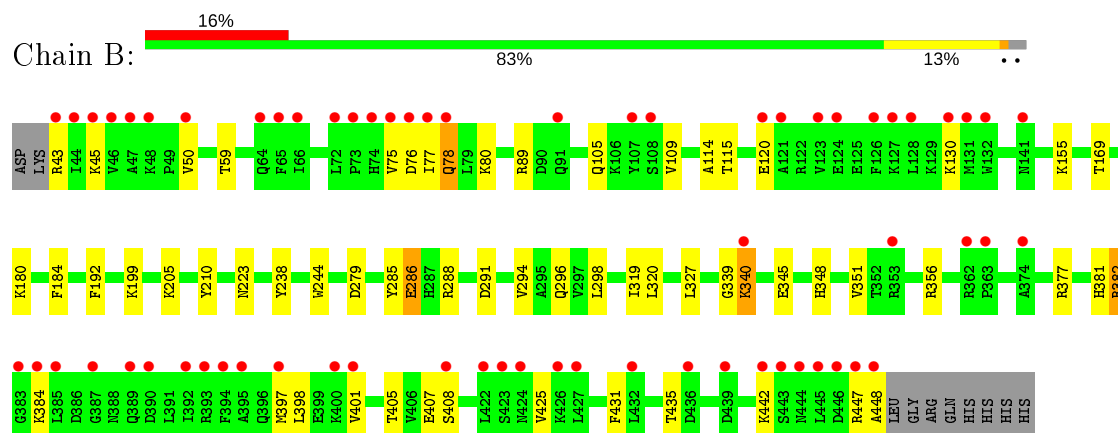
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Isocitrate dehydrogenase [NADP], mitochondrial



- Molecule 1: Isocitrate dehydrogenase [NADP], mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.66Å 119.66Å 125.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.90 – 1.55 31.90 – 1.55	Depositor EDS
% Data completeness (in resolution range)	99.5 (31.90-1.55) 99.4 (31.90-1.55)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.182 , 0.218 0.180 , 0.217	Depositor DCC
R_{free} test set	6429 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	14.6	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7629	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: GOL, CA, 1K9, NDP

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	1.35	15/3507 (0.4%)	1.15	13/4733 (0.3%)
1	B	1.23	9/3343 (0.3%)	1.03	3/4516 (0.1%)
All	All	1.29	24/6850 (0.4%)	1.09	16/9249 (0.2%)

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	185	VAL	CB-CG1	6.72	1.67	1.52
1	B	345	GLU	CD-OE2	6.72	1.33	1.25
1	A	216	GLY	N-CA	6.46	1.55	1.46
1	A	209	VAL	CB-CG1	6.28	1.66	1.52
1	B	285	TYR	CE1-CZ	6.14	1.46	1.38
1	A	238	TYR	CD2-CE2	6.11	1.48	1.39
1	B	244	TRP	CG-CD1	5.70	1.44	1.36
1	A	447	ARG	CZ-NH1	5.57	1.40	1.33
1	B	114	ALA	CA-CB	5.55	1.64	1.52
1	A	357	GLU	CG-CD	5.53	1.60	1.51
1	B	286	GLU	CD-OE1	5.47	1.31	1.25
1	A	317	SER	CA-CB	-5.37	1.44	1.52
1	A	400	LYS	CG-CD	5.32	1.70	1.52
1	A	206	GLU	CG-CD	5.24	1.59	1.51
1	B	192	PHE	CE1-CZ	5.20	1.47	1.37
1	A	274	TYR	CD1-CE1	5.19	1.47	1.39
1	B	210	TYR	CD1-CE1	5.18	1.47	1.39
1	A	345[A]	GLU	CD-OE2	5.17	1.31	1.25
1	A	345[B]	GLU	CD-OE2	5.17	1.31	1.25
1	B	184	PHE	CD1-CE1	5.15	1.49	1.39
1	A	125	GLU	CG-CD	5.15	1.59	1.51
1	A	282	LYS	CD-CE	5.13	1.64	1.51
1	A	140	GLN	CG-CD	5.09	1.62	1.51
1	B	192	PHE	CG-CD2	5.01	1.46	1.38

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	362	ARG	NE-CZ-NH1	9.45	125.02	120.30
1	B	89	ARG	NE-CZ-NH2	-7.32	116.64	120.30
1	A	362	ARG	NE-CZ-NH2	-6.53	117.04	120.30
1	A	447	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	A	356	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	A	271	ASP	CB-CG-OD2	-6.42	112.52	118.30
1	A	177	ASP	CB-CG-OD1	6.26	123.93	118.30
1	A	436	ASP	CB-CG-OD2	-5.92	112.97	118.30
1	A	238	TYR	CB-CG-CD1	-5.75	117.55	121.00
1	B	279	ASP	CB-CG-OD1	5.58	123.32	118.30
1	A	204	VAL	CG1-CB-CG2	5.57	119.82	110.90
1	A	84	LEU	CB-CG-CD2	5.51	120.36	111.00
1	A	89	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	451	ARG	NE-CZ-NH1	-5.40	117.60	120.30
1	B	356	ARG	NE-CZ-NH2	-5.35	117.63	120.30
1	A	279	ASP	CB-CG-OD1	5.30	123.07	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3428	0	3382	36	0
1	B	3269	0	3241	36	0
2	A	48	0	26	1	0
2	B	48	0	26	1	0
3	A	64	0	36	3	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	24	0	30	4	0
6	A	469	0	0	14	0
6	B	277	0	0	9	0
All	All	7629	0	6741	72	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 5.

All (72) 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:320[B]:LEU:CD2	6:A:1068:HOH:O	1.85	1.19
1:B:320[A]:LEU:CD2	6:B:686:HOH:O	1.92	1.14
1:B:320[A]:LEU:HD23	6:B:686:HOH:O	1.53	1.07
1:A:330:MET:CE	6:A:993:HOH:O	2.04	1.05
1:A:330:MET:HE2	6:A:993:HOH:O	1.56	1.02
1:A:320[B]:LEU:HD22	6:A:1068:HOH:O	1.52	1.00
1:B:288:ARG:HE	1:B:296:GLN:HE22	1.04	0.99
1:B:286:GLU:OE2	1:B:288:ARG:NH1	2.06	0.89
1:A:320[B]:LEU:HD23	6:A:1068:HOH:O	1.55	0.86
1:B:320[A]:LEU:HD22	6:B:686:HOH:O	1.72	0.75
1:A:294:VAL:HB	1:B:319[B]:ILE:HD11	1.69	0.75
1:B:109:VAL:HG22	1:B:381:HIS:HD2	1.50	0.74
5:A:506:GOL:H32	1:B:205:LYS:HE3	1.72	0.71
1:A:105[A]:GLN:HG3	1:A:143:LEU:CD2	2.23	0.68
1:B:75:VAL:HG12	1:B:77:ILE:HG22	1.77	0.66
1:A:322:GLN:HG3	6:A:855:HOH:O	1.96	0.65
1:A:319[A]:ILE:HD11	1:B:294:VAL:HB	1.81	0.63
5:A:506:GOL:H12	1:B:205:LYS:HE3	1.79	0.62
1:B:43:ARG:HD3	1:B:76:ASP:HB2	1.80	0.62
1:B:288:ARG:HE	1:B:296:GLN:NE2	1.89	0.60
1:B:425:VAL:HG22	1:B:431:PHE:HD2	1.66	0.59
1:B:320[A]:LEU:HB3	6:B:686:HOH:O	2.02	0.58
1:B:50:VAL:CG2	1:B:382:ARG:HG3	2.34	0.58
1:B:288:ARG:NE	1:B:296:GLN:HE22	1.88	0.57
1:A:273:HIS:HE1	6:A:894:HOH:O	1.86	0.56
6:A:946:HOH:O	1:B:199:LYS:HD2	2.03	0.56
1:A:74:HIS:HE1	6:A:818:HOH:O	1.89	0.56
1:B:327:LEU:HB3	1:B:348:HIS:HB3	1.88	0.56
1:B:320[A]:LEU:CB	6:B:686:HOH:O	2.55	0.54
1:B:77:ILE:HD12	6:B:718:HOH:O	2.07	0.54
1:A:105[A]:GLN:CG	1:A:143:LEU:CD2	2.86	0.54
1:A:327:LEU:HB3	1:A:348:HIS:HB3	1.90	0.53
1:B:340:LYS:HD3	6:B:868:HOH:O	2.10	0.51
1:A:330:MET:HB2	6:A:993:HOH:O	2.11	0.50
1:A:105[A]:GLN:HG3	1:A:143:LEU:HD21	1.92	0.50
1:B:115:THR:O	2:B:501:NDP:H2N	2.12	0.49
1:A:42:LYS:HD3	1:A:453[B]:HIS:CD2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:444:ASN:OD1	1:A:447:ARG:NH2	2.46	0.49
1:A:377[B]:ARG:HD3	1:A:399:GLU:OE2	2.13	0.49
5:A:506:GOL:H32	1:B:205:LYS:CE	2.41	0.48
1:A:74:HIS:HD2	6:A:939:HOH:O	1.97	0.48
1:A:115:THR:O	2:A:501:NDP:H2N	2.14	0.48
1:A:300[B]:SER:HG	1:A:304:PHE:HE2	1.62	0.47
1:B:155:LYS:HG2	1:B:407:GLU:OE2	2.14	0.47
1:B:120:GLU:HG2	1:B:120:GLU:H	1.55	0.47
1:B:398:LEU:HA	1:B:401:VAL:HG22	1.97	0.46
1:A:249:SER:HA	1:A:288:ARG:O	2.15	0.46
1:B:50:VAL:HG23	1:B:382:ARG:HG3	1.96	0.46
1:B:169:THR:HG21	1:B:238:TYR:CE2	2.51	0.46
3:A:502[A]:1K9:H7	3:A:502[A]:1K9:OAA	2.15	0.45
1:B:296:GLN:CG	6:B:874:HOH:O	2.64	0.45
1:B:75:VAL:HG12	1:B:77:ILE:CG2	2.46	0.45
1:A:332:SER:HB3	1:A:345[A]:GLU:HG2	1.98	0.44
1:A:109:VAL:HG22	1:A:381:HIS:HD2	1.82	0.44
1:A:319[B]:ILE:HG23	1:B:298:LEU:HD11	2.00	0.43
1:A:92:THR:O	1:A:95[B]:GLN:HG3	2.19	0.42
1:A:112:LYS:NZ	1:A:136:ASN:HD21	2.17	0.42
3:A:502[B]:1K9:H7	3:A:502[B]:1K9:OAA	2.19	0.42
1:A:272[B]:LYS:HD2	6:A:1041:HOH:O	2.19	0.42
1:B:405:THR:O	1:B:408:SER:HB2	2.18	0.42
1:A:345[B]:GLU:HG3	6:A:604:HOH:O	2.20	0.42
1:A:175:HIS:CD2	1:A:180:LYS:HE2	2.55	0.41
1:A:276:THR:OG1	5:A:504:GOL:H12	2.21	0.41
1:A:312:ASP:O	1:A:316[B]:GLN:HG2	2.20	0.41
1:A:42:LYS:CD	1:A:453[B]:HIS:CD2	3.04	0.41
1:B:78:GLN:HG3	1:B:78:GLN:O	2.20	0.41
3:A:502[B]:1K9:H1	3:A:502[B]:1K9:OAA	2.20	0.41
1:B:397:MET:SD	1:B:448:ALA:HB1	2.60	0.41
1:A:320[B]:LEU:HB3	6:A:1068:HOH:O	2.20	0.41
1:A:426:LYS:HB3	1:A:429:GLU:HB3	2.03	0.40
1:B:339:GLY:HA2	6:B:825:HOH:O	2.21	0.40
1:A:50:VAL:HG23	1:A:77:ILE:HD11	2.04	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	427/416 (103%)	413 (97%)	14 (3%)	0	100	100
1	B	410/416 (99%)	391 (95%)	19 (5%)	0	100	100
All	All	837/832 (101%)	804 (96%)	33 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	370/357 (104%)	365 (99%)	5 (1%)	67	41
1	B	354/357 (99%)	337 (95%)	17 (5%)	25	3
All	All	724/714 (101%)	702 (97%)	22 (3%)	40	12

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	159	ARG
1	A	291	ASP
1	A	340	LYS
1	A	351	VAL
1	A	427	LEU
1	B	45	LYS
1	B	59	THR
1	B	78	GLN

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Mol	Chain	Res	Type
1	B	80	LYS
1	B	105	GLN
1	B	130	LYS
1	B	180	LYS
1	B	223	ASN
1	B	291	ASP
1	B	340	LYS
1	B	351	VAL
1	B	377	ARG
1	B	382	ARG
1	B	384	LYS
1	B	435	THR
1	B	442	LYS
1	B	447	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	HIS
1	A	136	ASN
1	A	273	HIS
1	B	105	GLN
1	B	211	ASN
1	B	241	GLN
1	B	296	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 carbohydrates in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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	NDP	B	501	-	45,52,52	2.01	11 (24%)	53,80,80	1.69	13 (24%)
2	NDP	A	501	-	45,52,52	1.92	11 (24%)	53,80,80	1.63	12 (22%)
5	GOL	A	507	-	5,5,5	1.12	0	5,5,5	0.82	0
5	GOL	A	504	-	5,5,5	1.31	1 (20%)	5,5,5	0.69	0
3	1K9	A	502[B]	-	35,35,35	2.17	9 (25%)	47,52,52	2.68	8 (17%)
5	GOL	A	505	-	5,5,5	0.48	0	5,5,5	0.63	0
3	1K9	A	502[A]	-	35,35,35	2.64	15 (42%)	47,52,52	1.94	8 (17%)
5	GOL	A	506	-	5,5,5	0.82	0	5,5,5	0.90	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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NDP	B	501	-	-	6/30/77/77	0/5/5/5
2	NDP	A	501	-	-	8/30/77/77	0/5/5/5
5	GOL	A	507	-	-	2/4/4/4	-
5	GOL	A	504	-	-	0/4/4/4	-
3	1K9	A	502[B]	-	-	1/27/31/31	0/4/4/4
5	GOL	A	505	-	-	0/4/4/4	-
3	1K9	A	502[A]	-	-	1/27/31/31	0/4/4/4
5	GOL	A	506	-	-	4/4/4/4	-

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502[A]	1K9	OAC-SBF	9.48	1.54	1.43
3	A	502[B]	1K9	OAC-SBF	7.07	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	NDP	C4N-C3N	5.28	1.60	1.49
3	A	502[B]	1K9	CBC-CAY	-5.17	1.40	1.49
2	A	501	NDP	C2A-N3A	5.01	1.40	1.32
2	A	501	NDP	O4B-C1B	4.87	1.47	1.41
2	B	501	NDP	C6N-C5N	4.71	1.41	1.33
3	A	502[A]	1K9	CAN-CAY	-4.68	1.34	1.37
2	B	501	NDP	O4B-C1B	4.46	1.47	1.41
3	A	502[A]	1K9	CBE-CAZ	-4.33	1.40	1.49
2	B	501	NDP	O4D-C1D	4.13	1.51	1.42
2	B	501	NDP	P2B-O2B	4.12	1.67	1.59
2	A	501	NDP	C6N-C5N	4.10	1.40	1.33
3	A	502[A]	1K9	CBC-CAY	-3.97	1.42	1.49
3	A	502[B]	1K9	CAR-CBD	3.93	1.57	1.48
3	A	502[B]	1K9	CBE-CAZ	-3.90	1.41	1.49
3	A	502[A]	1K9	CBA-SBF	3.89	1.82	1.76
2	A	501	NDP	C7N-C3N	3.64	1.56	1.48
2	B	501	NDP	C4A-N3A	3.55	1.40	1.35
3	A	502[B]	1K9	CAN-CAY	-3.23	1.35	1.37
2	A	501	NDP	C4N-C5N	-3.06	1.40	1.48
2	A	501	NDP	C8A-N7A	-2.96	1.29	1.34
3	A	502[A]	1K9	CAX-NAS	-2.80	1.36	1.41
3	A	502[A]	1K9	CAP-CBA	2.76	1.44	1.39
2	A	501	NDP	PN-O5D	-2.75	1.48	1.59
2	B	501	NDP	C4N-C5N	-2.74	1.41	1.48
3	A	502[A]	1K9	CAO-CAX	2.71	1.43	1.39
3	A	502[A]	1K9	CAO-CAZ	2.70	1.43	1.39
3	A	502[A]	1K9	OAB-SBF	-2.57	1.40	1.43
3	A	502[A]	1K9	CBB-NAT	-2.55	1.36	1.41
3	A	502[B]	1K9	CAK-CAY	-2.51	1.38	1.42
2	B	501	NDP	O3B-C3B	2.48	1.48	1.43
2	A	501	NDP	PA-O2A	-2.48	1.43	1.55
3	A	502[B]	1K9	CAN-SAV	2.44	1.74	1.70
2	A	501	NDP	O4D-C4D	2.44	1.50	1.45
2	A	501	NDP	C4N-C3N	2.41	1.54	1.49
3	A	502[A]	1K9	CAK-CAY	-2.37	1.38	1.42
5	A	504	GOL	C1-C2	-2.33	1.42	1.51
2	A	501	NDP	PN-O2N	-2.24	1.44	1.55
3	A	502[B]	1K9	CAG-CAI	2.22	1.43	1.38
3	A	502[A]	1K9	FAD-CBE	-2.19	1.24	1.32
3	A	502[A]	1K9	CAH-SAV	-2.19	1.60	1.70
2	B	501	NDP	C2A-N1A	-2.17	1.29	1.33
3	A	502[B]	1K9	CAX-NAS	-2.12	1.37	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	NDP	C5D-C4D	2.09	1.58	1.51
3	A	502[A]	1K9	OAA-CAW	-2.08	1.19	1.23
2	B	501	NDP	PN-O2N	-2.05	1.45	1.55

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502[B]	1K9	CAY-CAN-SAV	-11.97	104.57	112.29
3	A	502[A]	1K9	CAY-CAN-SAV	-8.75	106.65	112.29
3	A	502[B]	1K9	CBD-NAU-SBF	-7.93	108.36	122.08
3	A	502[B]	1K9	CAQ-CBD-NAU	-6.24	109.33	118.51
2	B	501	NDP	O4D-C1D-N1N	-5.36	97.58	108.06
2	A	501	NDP	N3A-C2A-N1A	-4.56	121.54	128.68
3	A	502[B]	1K9	CAH-SAV-CAN	4.36	101.26	92.37
2	B	501	NDP	C4A-C5A-N7A	-3.91	105.32	109.40
2	A	501	NDP	C1D-N1N-C2N	-3.91	114.60	121.11
3	A	502[A]	1K9	CAQ-CBD-NAU	-3.87	112.81	118.51
3	A	502[B]	1K9	CAR-CAQ-CBD	3.69	62.93	59.84
3	A	502[A]	1K9	CBD-NAU-SBF	-3.57	115.90	122.08
2	A	501	NDP	C4A-C5A-N7A	-3.53	105.72	109.40
2	A	501	NDP	O4D-C1D-N1N	-3.33	101.54	108.06
2	B	501	NDP	O4D-C4D-C5D	-3.14	99.03	109.37
2	B	501	NDP	C3D-C2D-C1D	3.07	107.25	101.43
2	B	501	NDP	O3X-P2B-O2X	3.06	119.32	107.64
3	A	502[A]	1K9	CAY-CBC-CBB	2.97	125.54	122.32
2	B	501	NDP	C1D-N1N-C2N	-2.80	116.44	121.11
3	A	502[B]	1K9	CAY-CBC-CBB	2.79	125.34	122.32
3	A	502[A]	1K9	OAC-SBF-CBA	-2.77	104.56	107.97
2	B	501	NDP	O4D-C4D-C3D	2.71	110.48	105.11
2	A	501	NDP	O2A-PA-O1A	2.70	125.59	112.24
2	A	501	NDP	C1D-N1N-C6N	2.66	126.56	120.83
3	A	502[A]	1K9	CBC-CBB-NAT	-2.60	115.09	119.32
3	A	502[B]	1K9	CAQ-CAR-CBD	-2.49	57.75	59.84
2	A	501	NDP	O2N-PN-O1N	2.49	124.54	112.24
2	A	501	NDP	O5D-PN-O1N	-2.44	99.53	109.07
3	A	502[A]	1K9	CAM-CAL-CBA	-2.36	117.00	119.45
2	A	501	NDP	O5B-C5B-C4B	2.36	117.11	108.99
2	A	501	NDP	O2N-PN-O5D	2.34	118.60	107.75
2	B	501	NDP	C3B-C2B-C1B	-2.25	98.66	102.89
3	A	502[B]	1K9	CBC-CBB-NAT	-2.21	115.71	119.32
2	B	501	NDP	C3N-C2N-N1N	2.17	126.20	123.10
2	B	501	NDP	O4D-C1D-C2D	-2.14	101.97	106.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NDP	O3D-C3D-C2D	-2.13	104.93	111.82
2	B	501	NDP	C4D-O4D-C1D	-2.12	104.79	109.47
3	A	502[A]	1K9	CAI-CAX-NAS	-2.12	113.28	120.40
2	A	501	NDP	C2A-N1A-C6A	2.04	122.24	118.75
2	B	501	NDP	C5A-C6A-N6A	2.03	123.43	120.35
2	A	501	NDP	O2B-C2B-C1B	-2.03	102.80	110.10

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	506	GOL	O1-C1-C2-C3
5	A	506	GOL	C1-C2-C3-O3
2	B	501	NDP	C2D-C1D-N1N-C6N
3	A	502[B]	1K9	CAN-CAY-CBC-CAM
5	A	507	GOL	C1-C2-C3-O3
5	A	506	GOL	O1-C1-C2-O2
3	A	502[A]	1K9	CAN-CAY-CBC-CAM
5	A	506	GOL	O2-C2-C3-O3
2	A	501	NDP	C5D-O5D-PN-O3
5	A	507	GOL	O1-C1-C2-O2
2	A	501	NDP	C5D-O5D-PN-O2N
2	B	501	NDP	C3B-C4B-C5B-O5B
2	B	501	NDP	C2D-C1D-N1N-C2N
2	A	501	NDP	C2D-C1D-N1N-C2N
2	A	501	NDP	C2D-C1D-N1N-C6N
2	B	501	NDP	O4D-C1D-N1N-C6N
2	B	501	NDP	O4D-C1D-N1N-C2N
2	A	501	NDP	O4D-C1D-N1N-C2N
2	B	501	NDP	O4B-C4B-C5B-O5B
2	A	501	NDP	O4D-C1D-N1N-C6N
2	A	501	NDP	C5B-O5B-PA-O1A
2	A	501	NDP	C3B-C4B-C5B-O5B

There are no ring outliers.

6 monomers are involved in 9 short contacts:

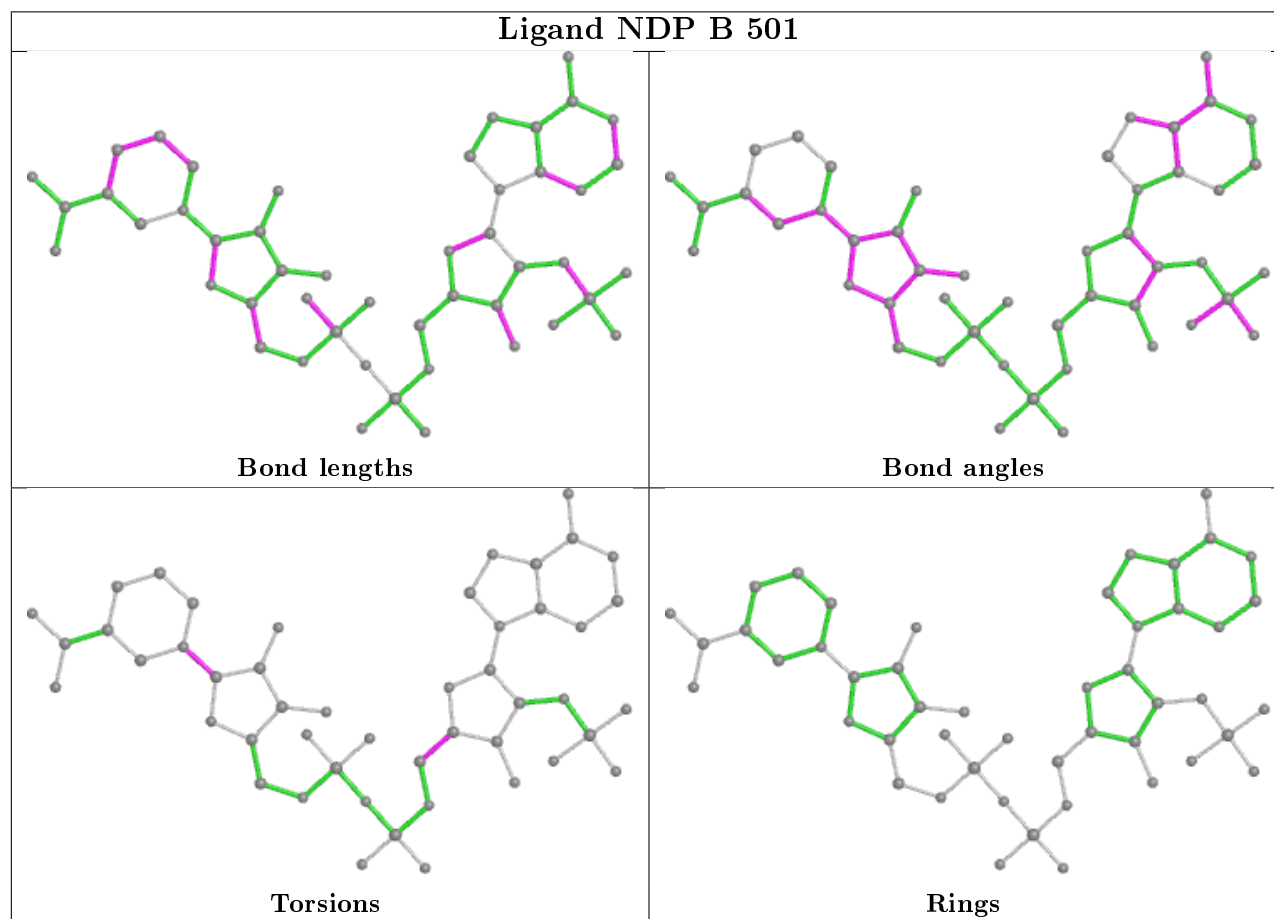
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	NDP	1	0
2	A	501	NDP	1	0
5	A	504	GOL	1	0

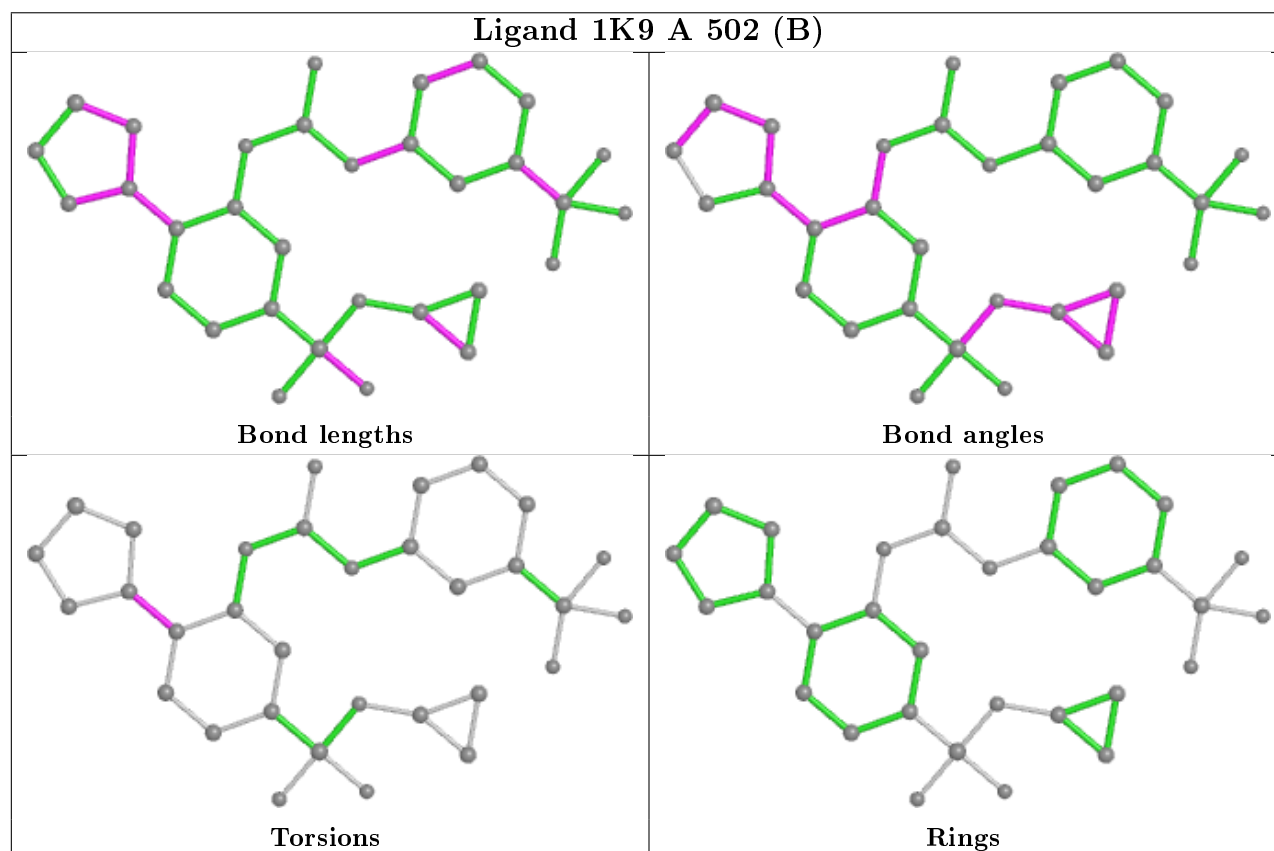
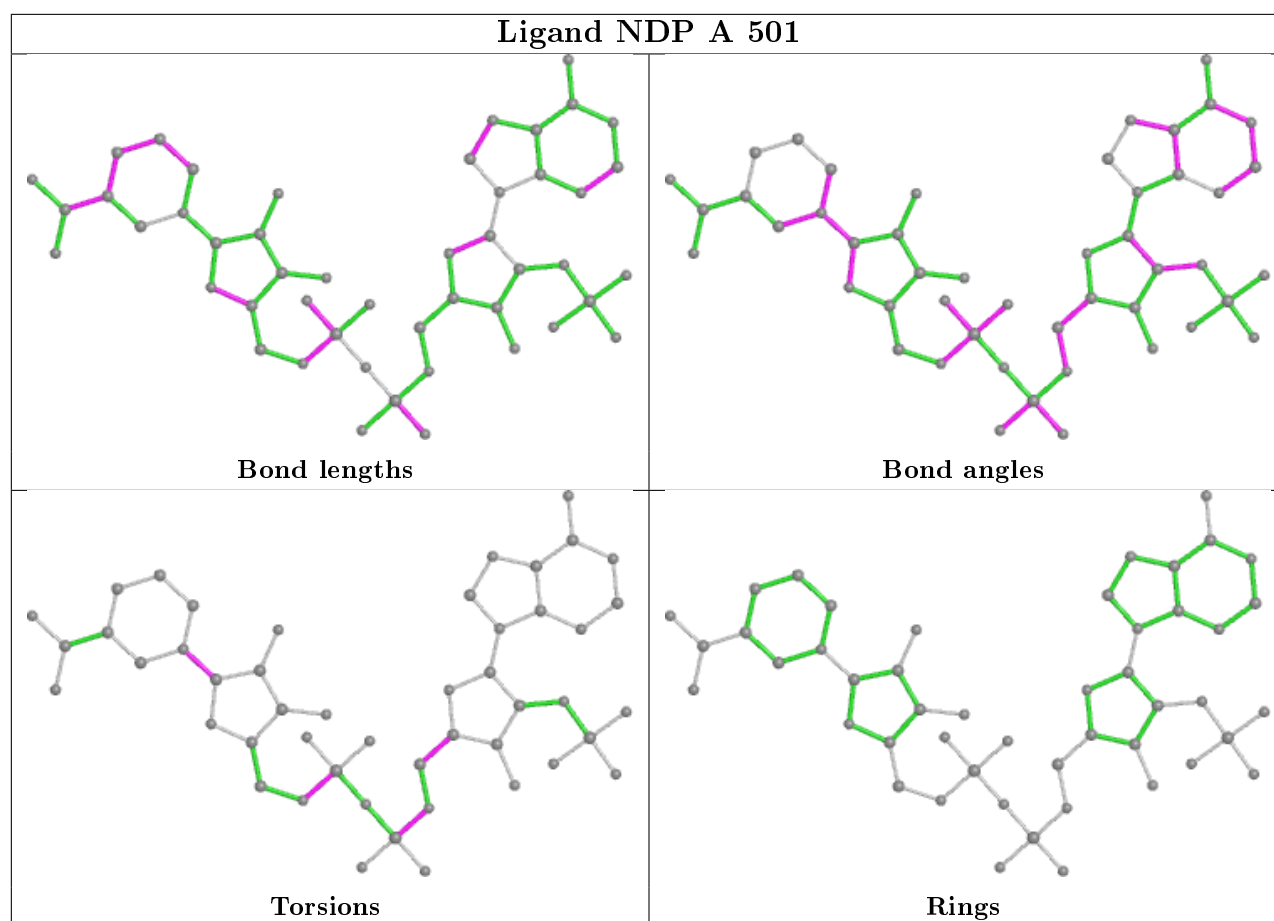
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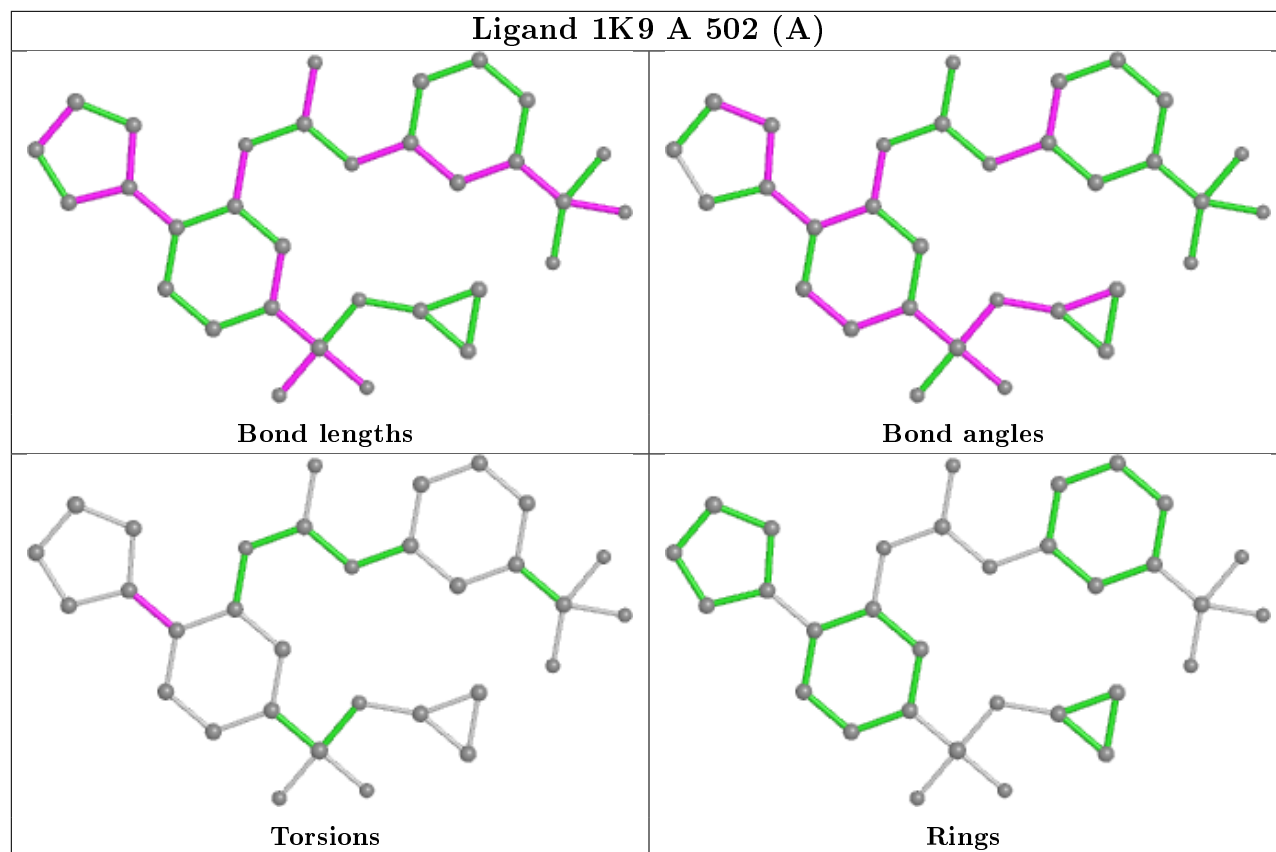
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502[B]	1K9	2	0
3	A	502[A]	1K9	1	0
5	A	506	GOL	3	0

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	416/416 (100%)	-0.06	6 (1%) 75 80	7, 13, 28, 36	0
1	B	406/416 (97%)	0.86	65 (16%) 1 1	8, 25, 54, 71	0
All	All	822/832 (98%)	0.39	71 (8%) 10 11	7, 17, 47, 71	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	44	ILE	7.7
1	B	77	ILE	5.6
1	B	75	VAL	5.5
1	B	397	MET	5.3
1	B	447	ARG	5.2
1	B	383	GLY	5.0
1	B	123	VAL	4.8
1	B	132	TRP	4.8
1	B	46	VAL	4.8
1	B	448	ALA	4.8
1	B	128	LEU	4.3
1	B	73	PRO	4.3
1	B	120	GLU	4.2
1	B	43	ARG	4.2
1	B	47	ALA	4.2
1	B	76	ASP	4.0
1	B	72	LEU	3.9
1	B	126	PHE	3.9
1	B	48	LYS	3.8
1	B	423	SER	3.8
1	B	443	SER	3.8
1	B	390	ASP	3.7
1	B	78	GLN	3.7
1	A	456	HIS	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	45	LYS	3.6
1	B	107	TYR	3.6
1	B	50	VAL	3.5
1	B	130	LYS	3.5
1	A	41	ASP	3.5
1	B	74	HIS	3.4
1	B	124	GLU	3.4
1	B	446	ASP	3.4
1	B	340	LYS	3.3
1	B	394	PHE	3.1
1	B	127	LYS	3.1
1	B	424	ASN	3.1
1	B	353	ARG	3.0
1	B	387	GLY	3.0
1	B	392	ILE	3.0
1	B	432	LEU	3.0
1	A	42	LYS	2.9
1	B	389	GLN	2.9
1	B	393	ARG	2.9
1	B	384	LYS	2.9
1	B	444	ASN	2.8
1	B	426	LYS	2.8
1	B	65	PHE	2.8
1	B	445	LEU	2.7
1	B	395	ALA	2.6
1	B	427	LEU	2.6
1	B	362	ARG	2.6
1	B	91	GLN	2.6
1	B	131	MET	2.5
1	A	377[A]	ARG	2.5
1	B	408	SER	2.5
1	B	401	VAL	2.4
1	B	66	ILE	2.4
1	B	385	LEU	2.4
1	B	422	LEU	2.4
1	B	363	PRO	2.3
1	B	400	LYS	2.3
1	A	120	GLU	2.3
1	B	141	ASN	2.2
1	B	436	ASP	2.2
1	B	439	ASP	2.2
1	B	121	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	442	LYS	2.1
1	B	64	GLN	2.1
1	B	374	ALA	2.1
1	A	426	LYS	2.0
1	B	108	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

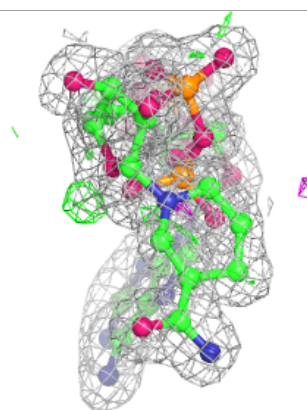
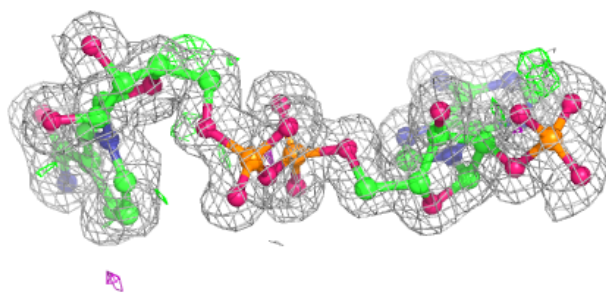
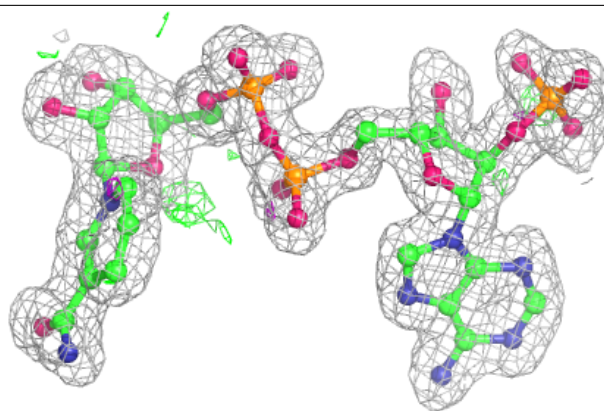
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GOL	A	507	6/6	0.83	0.15	21,24,28,33	0
5	GOL	A	506	6/6	0.89	0.16	28,30,33,35	0
5	GOL	A	505	6/6	0.90	0.19	30,36,41,44	0
5	GOL	A	504	6/6	0.92	0.18	17,17,32,34	0
4	CA	A	503	1/1	0.94	0.08	21,21,21,21	0
2	NDP	B	501	48/48	0.95	0.07	15,19,26,28	0
3	1K9	A	502[B]	32/32	0.97	0.11	6,12,15,17	32
3	1K9	A	502[A]	32/32	0.97	0.11	4,7,11,12	32
4	CA	B	502	1/1	0.97	0.06	20,20,20,20	0
2	NDP	A	501	48/48	0.98	0.06	7,10,14,18	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.

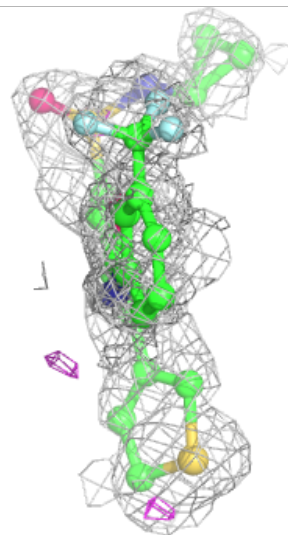
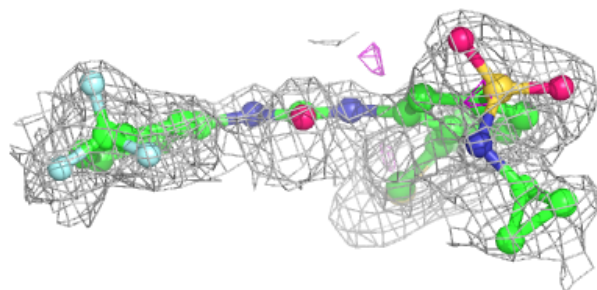
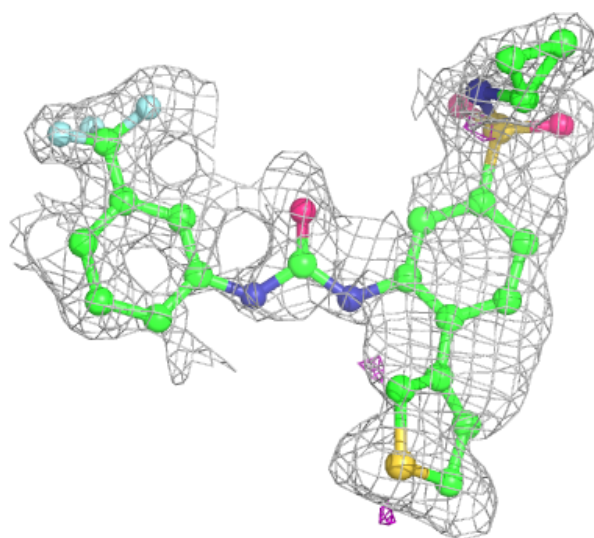
Electron density around NDP B 501:

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



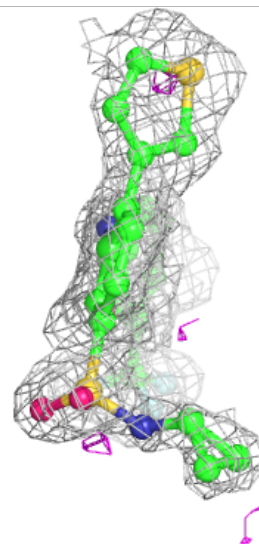
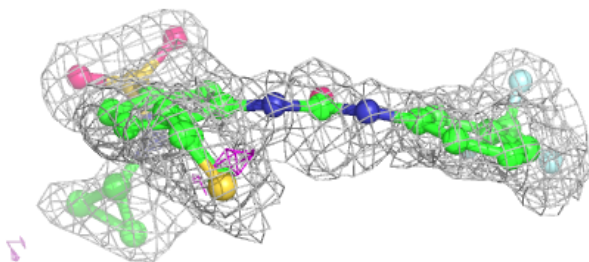
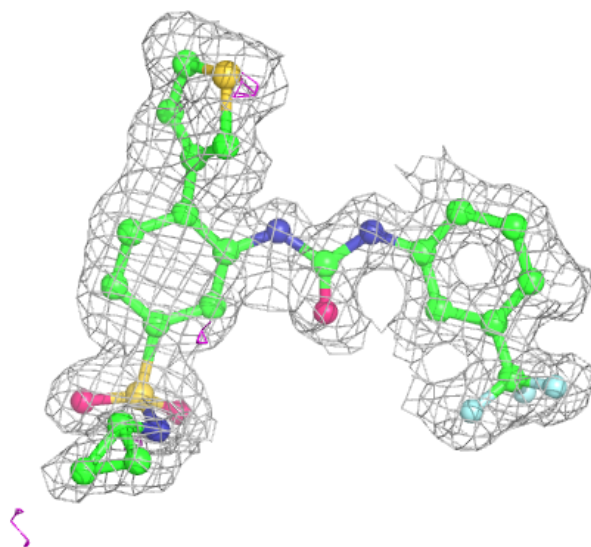
Electron density around 1K9 A 502 (B):

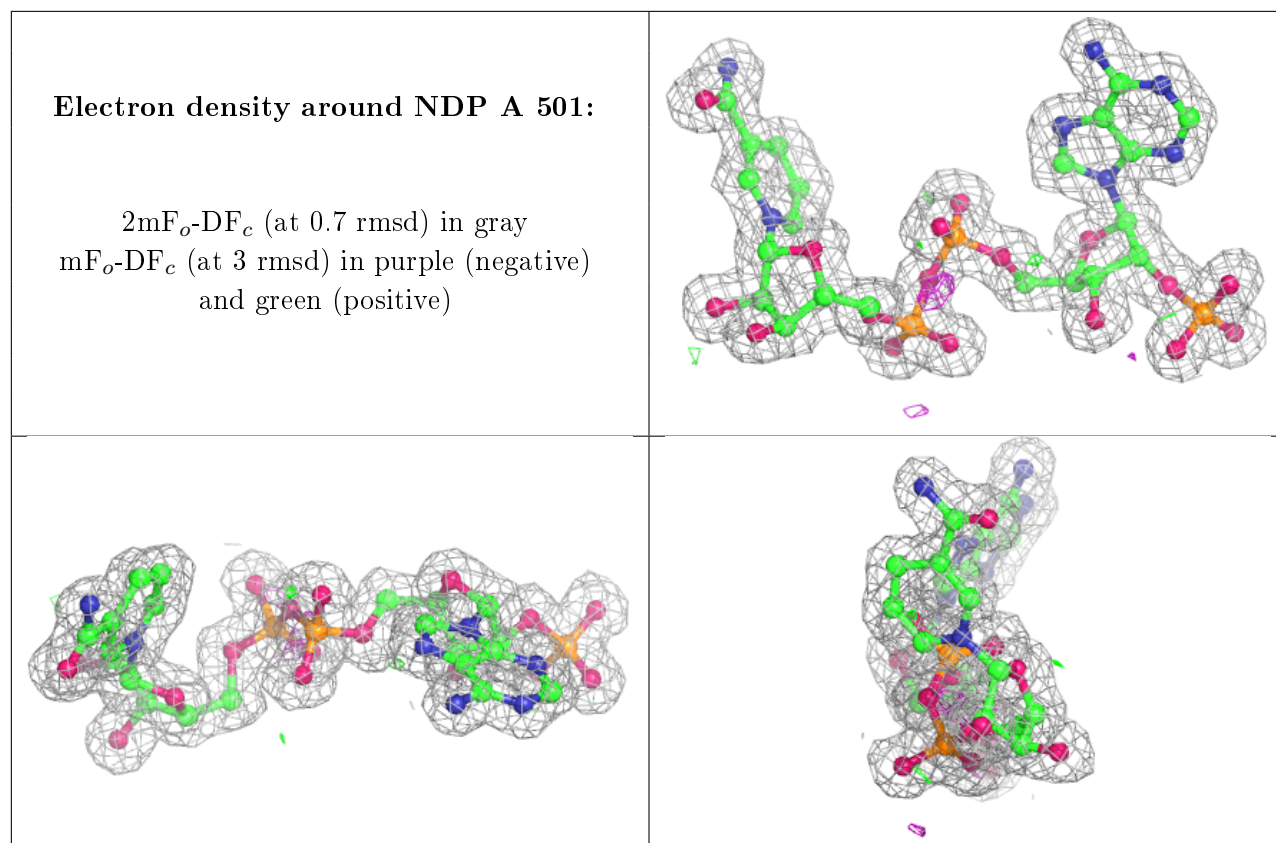
2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 1K9 A 502 (A):

$2mF_o-DF_c$ (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.