



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 09:44 pm BST

PDB ID : 3AU8
Title : Crystal structure of the ternary complex of an isomerase
Authors : Umeda, T.; Tanaka, N.; Kusakabe, Y.; Nakanishi, M.; Kitade, Y.; Nakamura, K.T.
Deposited on : 2011-02-01
Resolution : 1.86 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

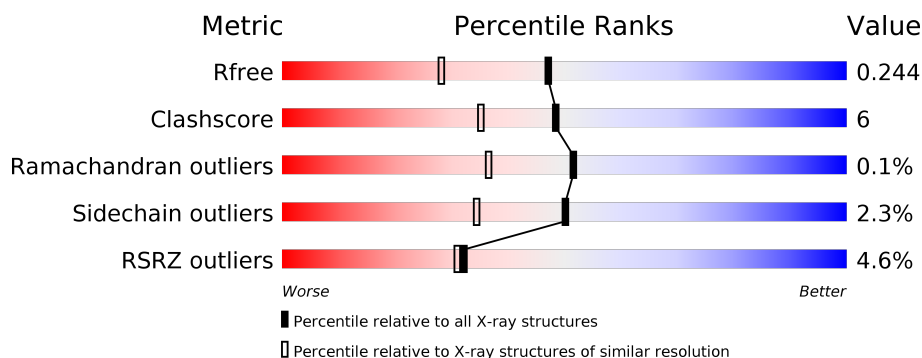
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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	488	<div> <div>5%</div> <div> <div></div> <div>68%</div> <div>13%</div> <div>18%</div> </div> </div>
1	B	488	<div> <div>3%</div> <div> <div></div> <div>73%</div> <div>9%</div> <div>17%</div> </div> </div>

2 Entry composition [i](#)

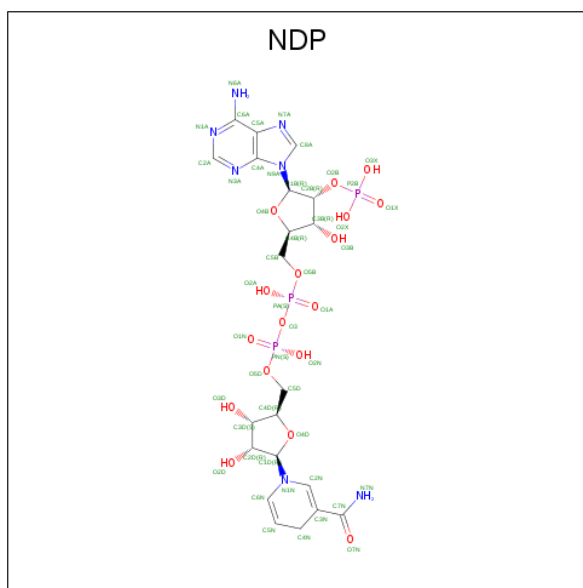
There are 4 unique types of molecules in this entry. The entry contains 6885 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 1-deoxy-D-xylulose 5-phosphate reductoisomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	0	0	0
			3180	2040	517	604	19			
1	B	405	Total	C	N	O	S	0	0	0
			3237	2080	528	610	19			

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



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 MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

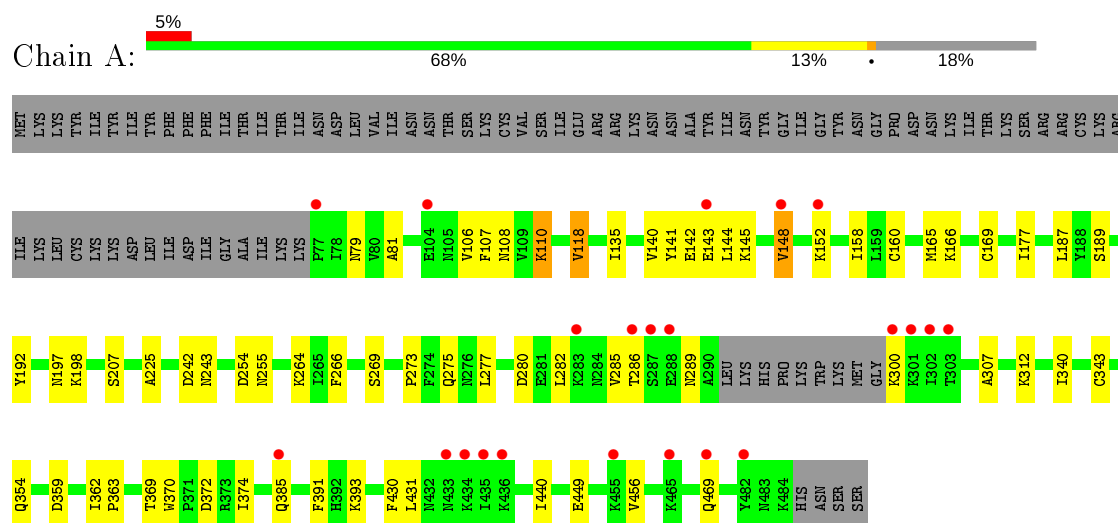
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	183	Total 183	O 183	0	0
4	B	187	Total 187	O 187	0	0

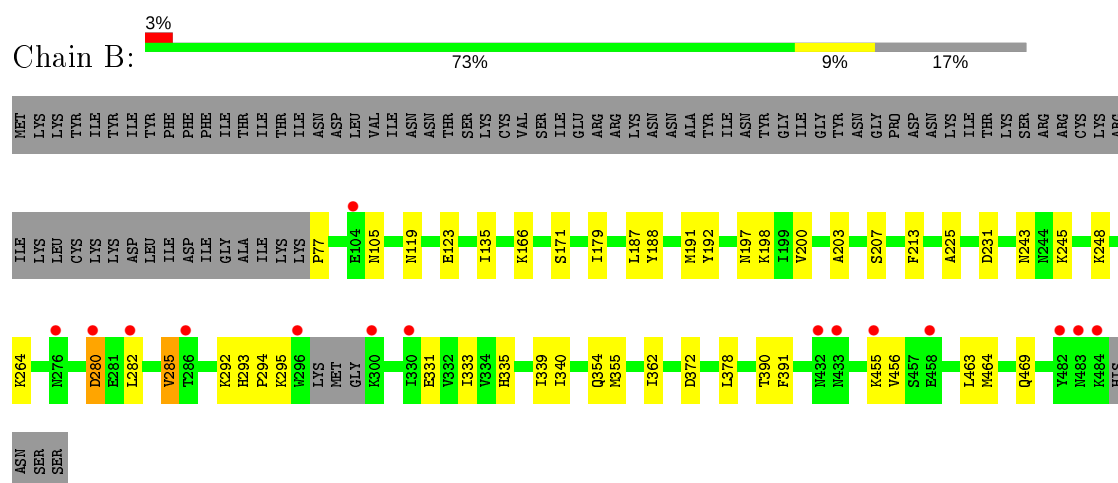
3 Residue-property plots [i](#)

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: 1-deoxy-D-xylulose 5-phosphate reductoisomerase



- Molecule 1: 1-deoxy-D-xylulose 5-phosphate reductoisomerase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	168.38Å 58.53Å 86.07Å 90.00° 117.74° 90.00°	Depositor
Resolution (Å)	50.00 – 1.86 37.87 – 1.86	Depositor EDS
% Data completeness (in resolution range)	95.2 (50.00-1.86) 95.2 (37.87-1.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.200 , 0.246 0.201 , 0.244	Depositor DCC
R_{free} test set	3017 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.024 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6885	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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: NDP, MN

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	0.50	0/3237	0.56	0/4371
1	B	0.48	0/3298	0.56	0/4454
All	All	0.49	0/6535	0.56	0/8825

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3180	0	3222	41	0
1	B	3237	0	3283	38	0
2	A	48	0	26	0	0
2	B	48	0	26	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	183	0	0	7	0
4	B	187	0	0	7	0
All	All	6885	0	6557	79	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 6.

All (79) 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:165:MET:CE	1:A:189:SER:HA	1.73	1.19
1:A:165:MET:HE3	1:A:189:SER:HA	0.98	0.96
1:A:165:MET:HE3	1:A:189:SER:CA	1.95	0.90
1:A:148:VAL:HG11	1:A:158:ILE:HD11	1.69	0.75
1:B:231:ASP:OD1	2:B:502:NDP:H5N	1.87	0.73
1:B:213:PHE:HD2	4:B:517:HOH:O	1.76	0.69
1:A:165:MET:CE	1:A:189:SER:CA	2.63	0.66
1:B:293:HIS:HE1	1:B:295:LYS:HE2	1.64	0.63
1:B:282:LEU:HA	1:B:285:VAL:HG22	1.80	0.63
4:A:669:HOH:O	1:B:372:ASP:HB3	1.99	0.62
1:A:108:ASN:HB2	4:A:655:HOH:O	1.99	0.62
1:B:340:ILE:HD12	1:B:391:PHE:HZ	1.67	0.60
1:B:335:HIS:CE1	1:B:339:ILE:H	2.20	0.59
1:B:335:HIS:CE1	1:B:340:ILE:H	2.22	0.58
4:A:582:HOH:O	1:B:372:ASP:HB3	2.05	0.57
1:B:362:ILE:HD11	1:B:378:LEU:HD12	1.86	0.57
1:B:245:LYS:HD3	1:B:248:LYS:HE2	1.89	0.55
1:A:243:ASN:ND2	1:A:370:TRP:HE1	2.05	0.55
1:B:335:HIS:HE1	1:B:339:ILE:H	1.54	0.54
1:B:166:LYS:HE2	1:B:192:TYR:CZ	2.43	0.54
1:A:456:VAL:HG22	4:A:562:HOH:O	2.09	0.53
1:B:191:MET:HE1	1:B:463:LEU:HD23	1.89	0.53
1:A:118:VAL:HG11	1:A:140:VAL:HB	1.89	0.52
1:B:264:LYS:HE3	1:B:331:GLU:HB3	1.91	0.52
1:A:141:TYR:CD1	1:A:160:CYS:HB2	2.44	0.52
1:B:455:LYS:HB3	4:B:639:HOH:O	2.09	0.51
2:B:502:NDP:N7N	2:B:502:NDP:O1N	2.42	0.50
1:A:169:CYS:O	1:A:198:LYS:HE2	2.11	0.50
1:B:455:LYS:HG3	4:B:673:HOH:O	2.10	0.50
1:B:335:HIS:HE1	1:B:340:ILE:H	1.58	0.50
1:B:245:LYS:NZ	4:B:603:HOH:O	2.45	0.50
1:A:145:LYS:O	1:A:148:VAL:HG12	2.12	0.49
1:B:213:PHE:CE1	1:B:456:VAL:HG11	2.48	0.49
1:A:372:ASP:HB3	4:B:603:HOH:O	2.11	0.49
1:B:171:SER:O	1:B:198:LYS:NZ	2.45	0.49
1:B:335:HIS:HD2	1:B:390:THR:O	1.96	0.49
1:A:187:LEU:HD13	1:A:207:SER:HA	1.95	0.49
1:B:187:LEU:HD13	1:B:207:SER:HA	1.94	0.49
1:A:282:LEU:O	1:A:285:VAL:HG22	2.13	0.48
1:B:77:PRO:N	4:B:651:HOH:O	2.45	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:ASN:HB3	1:A:110:LYS:HG3	1.95	0.48
1:B:188:TYR:CD1	1:B:464:MET:HG3	2.49	0.47
1:A:143:GLU:C	4:A:607:HOH:O	2.53	0.47
1:A:166:LYS:HE2	1:A:192:TYR:CZ	2.49	0.47
1:A:275:GLN:HB2	4:A:659:HOH:O	2.13	0.47
1:A:118:VAL:HG13	1:A:135:ILE:HD13	1.95	0.47
1:B:213:PHE:CE1	1:B:456:VAL:CG1	2.97	0.47
1:B:293:HIS:CE1	1:B:295:LYS:HE2	2.49	0.46
1:B:231:ASP:CG	2:B:502:NDP:H5N	2.36	0.46
1:A:118:VAL:CG1	1:A:140:VAL:HB	2.45	0.46
1:A:369:THR:HG21	1:A:374:ILE:HG12	1.98	0.45
1:B:340:ILE:HG12	1:B:355:MET:HG2	1.99	0.44
1:A:300:LYS:HD2	1:A:431:LEU:HD11	2.00	0.44
1:B:203:ALA:O	2:B:502:NDP:H1D	2.18	0.44
1:B:119:ASN:O	1:B:123:GLU:HG2	2.18	0.43
1:A:340:ILE:HD12	1:A:391:PHE:HZ	1.82	0.43
1:A:393:LYS:HD2	4:B:557:HOH:O	2.19	0.43
1:B:333:ILE:HD11	1:B:391:PHE:HB3	2.00	0.43
1:A:118:VAL:HG12	1:A:144:LEU:HD23	2.01	0.43
1:A:286:THR:H	1:A:289:ASN:ND2	2.17	0.43
1:A:266:PHE:HB2	1:A:343:CYS:HB2	2.00	0.42
1:B:280:ASP:OD1	1:B:280:ASP:N	2.52	0.42
1:A:362:ILE:HB	1:A:363:PRO:HD3	2.01	0.42
1:B:179:ILE:HG13	1:B:200:VAL:HG13	2.00	0.42
1:A:106:VAL:HG13	1:A:107:PHE:CD2	2.55	0.42
1:A:197:ASN:HA	1:A:225:ALA:HB2	2.01	0.42
1:B:293:HIS:HA	1:B:294:PRO:HD3	1.90	0.42
1:A:273:PRO:HD2	1:A:307:ALA:HA	2.01	0.42
1:A:277:LEU:HD13	1:A:285:VAL:HG11	2.02	0.42
1:A:269:SER:HB2	1:A:312:LYS:HE3	2.03	0.41
1:A:242:ASP:OD1	1:B:243:ASN:HB2	2.20	0.41
1:A:430:PHE:CG	1:A:440:ILE:HD11	2.55	0.41
1:A:148:VAL:HG11	1:A:158:ILE:CD1	2.46	0.41
1:A:254:ASP:O	1:A:255:ASN:HB2	2.20	0.41
1:B:264:LYS:NZ	1:B:331:GLU:OE1	2.40	0.41
1:A:145:LYS:HA	1:A:158:ILE:HD13	2.03	0.41
1:A:81:ALA:HB3	1:A:177:ILE:HG12	2.03	0.40
1:A:359:ASP:HB3	4:A:664:HOH:O	2.20	0.40
1:B:197:ASN:HA	1:B:225:ALA:HB2	2.04	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	395/488 (81%)	384 (97%)	11 (3%)	0	100	100
1	B	401/488 (82%)	385 (96%)	15 (4%)	1 (0%)	47	33
All	All	796/976 (82%)	769 (97%)	26 (3%)	1 (0%)	51	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	105	ASN

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	367/449 (82%)	356 (97%)	11 (3%)	41	24
1	B	373/449 (83%)	367 (98%)	6 (2%)	62	49
All	All	740/898 (82%)	723 (98%)	17 (2%)	50	34

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	LYS
1	A	118	VAL
1	A	142	GLU
1	A	148	VAL
1	A	152	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	264	LYS
1	A	280	ASP
1	A	354	GLN
1	A	385	GLN
1	A	449	GLU
1	A	469	GLN
1	B	135	ILE
1	B	280	ASP
1	B	285	VAL
1	B	292	LYS
1	B	354	GLN
1	B	469	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	ASN
1	A	243	ASN
1	A	253	GLN
1	A	289	ASN
1	A	328	ASN
1	A	354	GLN
1	A	459	ASN
1	B	95	ASN
1	B	220	ASN
1	B	253	GLN
1	B	328	ASN
1	B	335	HIS
1	B	354	GLN
1	B	361	GLN
1	B	407	GLN
1	B	469	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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NDP	A	501	-	45,52,52	1.63	5 (11%)	53,80,80	1.30	5 (9%)
2	NDP	B	502	-	45,52,52	1.66	5 (11%)	53,80,80	1.29	5 (9%)

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	A	501	-	-	7/30/77/77	0/5/5/5
2	NDP	B	502	-	-	8/30/77/77	0/5/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NDP	O7N-C7N	7.27	1.41	1.24
2	B	502	NDP	O7N-C7N	7.13	1.41	1.24
2	B	502	NDP	C2A-N3A	4.13	1.38	1.32
2	A	501	NDP	C2A-N3A	3.85	1.38	1.32
2	A	501	NDP	C6N-C5N	3.62	1.39	1.33
2	B	502	NDP	C6N-C5N	3.41	1.39	1.33
2	B	502	NDP	C2A-N1A	2.91	1.39	1.33
2	A	501	NDP	C2A-N1A	2.63	1.38	1.33
2	B	502	NDP	P2B-O2B	2.41	1.63	1.59
2	A	501	NDP	C2N-C3N	2.06	1.40	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NDP	N3A-C2A-N1A	-6.04	119.24	128.68
2	B	502	NDP	N3A-C2A-N1A	-5.34	120.34	128.68
2	A	501	NDP	PN-O3-PA	-3.62	120.39	132.83
2	B	502	NDP	PN-O3-PA	-3.55	120.65	132.83
2	A	501	NDP	O4D-C1D-N1N	2.72	113.37	108.06
2	B	502	NDP	O4D-C1D-N1N	2.66	113.27	108.06
2	B	502	NDP	C1B-N9A-C4A	-2.24	122.70	126.64
2	A	501	NDP	C1B-N9A-C4A	-2.17	122.83	126.64
2	B	502	NDP	C3B-C2B-C1B	-2.15	98.85	102.89
2	A	501	NDP	C1D-N1N-C6N	-2.13	116.23	120.83

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	502	NDP	C5D-O5D-PN-O3
2	B	502	NDP	C5D-O5D-PN-O1N
2	B	502	NDP	C5D-O5D-PN-O2N
2	B	502	NDP	O4D-C4D-C5D-O5D
2	B	502	NDP	C3D-C4D-C5D-O5D
2	A	501	NDP	O4D-C4D-C5D-O5D
2	B	502	NDP	O4D-C1D-N1N-C6N
2	A	501	NDP	C3D-C4D-C5D-O5D
2	A	501	NDP	O4D-C1D-N1N-C6N
2	A	501	NDP	C2D-C1D-N1N-C6N
2	A	501	NDP	C2B-O2B-P2B-O2X
2	A	501	NDP	C2N-C3N-C7N-N7N
2	B	502	NDP	C2N-C3N-C7N-N7N
2	A	501	NDP	O4B-C4B-C5B-O5B
2	B	502	NDP	O4B-C4B-C5B-O5B

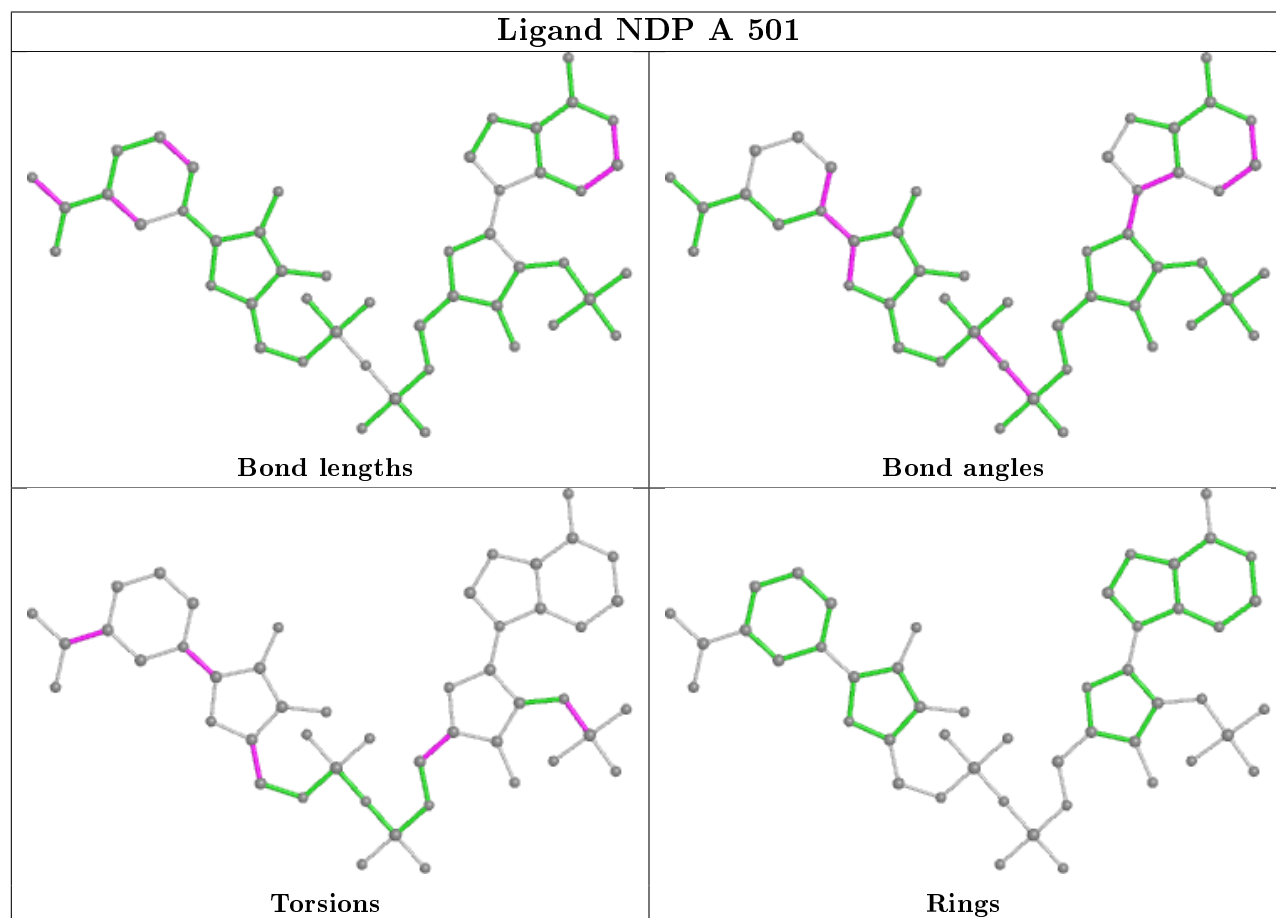
There are no ring outliers.

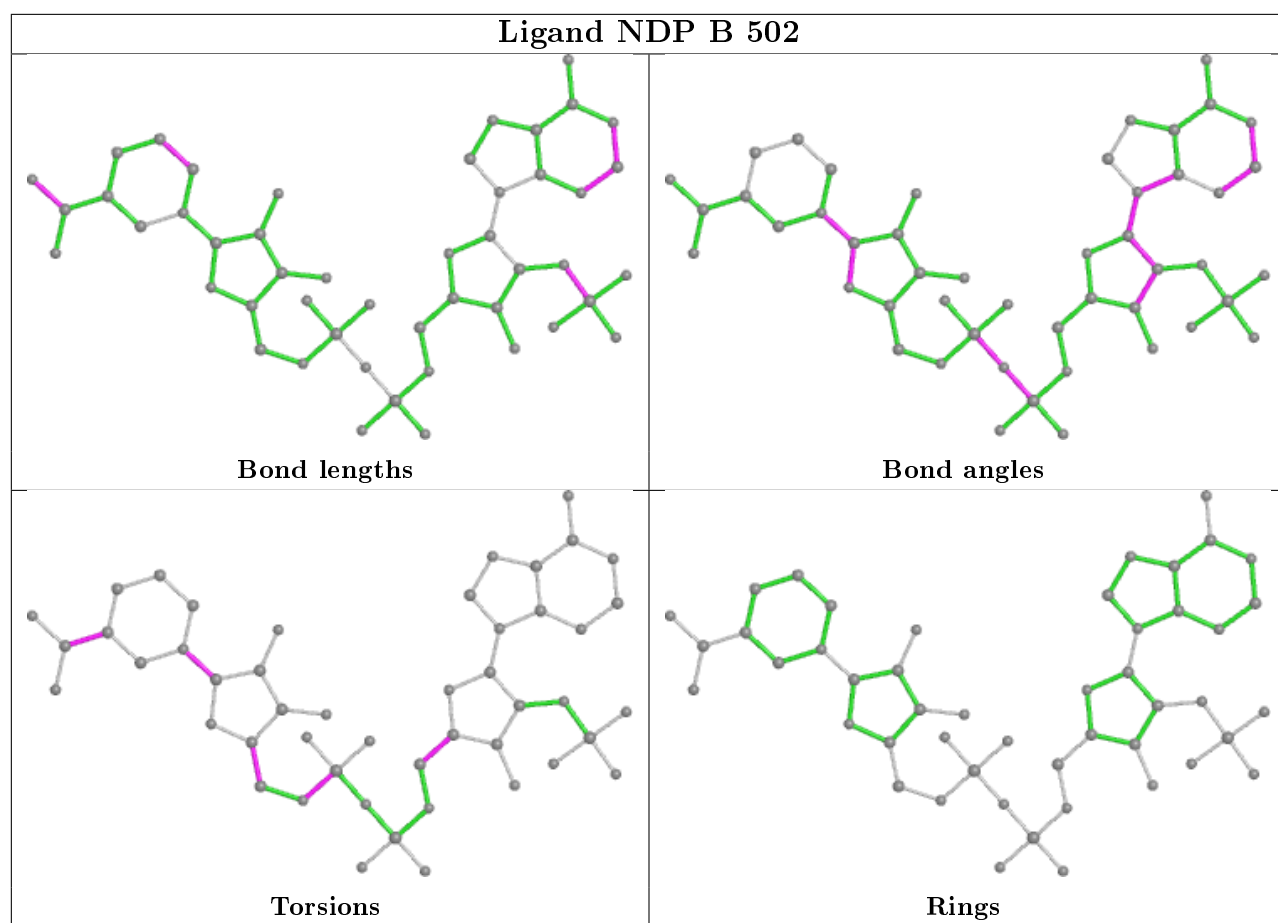
1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	NDP	4	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	399/488 (81%)	0.13	22 (5%) 25 24	16, 28, 49, 62	0
1	B	405/488 (82%)	0.19	15 (3%) 41 39	16, 30, 48, 68	0
All	All	804/976 (82%)	0.16	37 (4%) 32 31	16, 29, 49, 68	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	287	SER	4.2
1	B	483	ASN	3.9
1	A	288	GLU	3.8
1	A	300	LYS	3.3
1	B	482	TYR	3.1
1	A	435	ILE	2.9
1	A	465	LYS	2.9
1	B	276	ASN	2.8
1	B	432	ASN	2.7
1	B	433	ASN	2.7
1	A	436	LYS	2.7
1	B	484	LYS	2.7
1	A	104	GLU	2.7
1	B	458	GLU	2.6
1	A	433	ASN	2.6
1	B	280	ASP	2.6
1	A	77	PRO	2.6
1	A	302	ILE	2.5
1	A	152	LYS	2.5
1	A	434	LYS	2.4
1	B	282	LEU	2.3
1	A	286	THR	2.3
1	B	296	TRP	2.3
1	A	482	TYR	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	301	LYS	2.2
1	B	286	THR	2.2
1	A	148	VAL	2.2
1	A	303	THR	2.2
1	A	469	GLN	2.1
1	A	455	LYS	2.1
1	B	455	LYS	2.1
1	A	283	LYS	2.0
1	B	300	LYS	2.0
1	A	385	GLN	2.0
1	B	104	GLU	2.0
1	B	330	ILE	2.0
1	A	143	GLU	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](#)

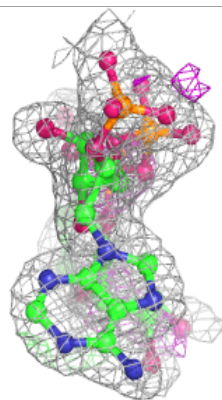
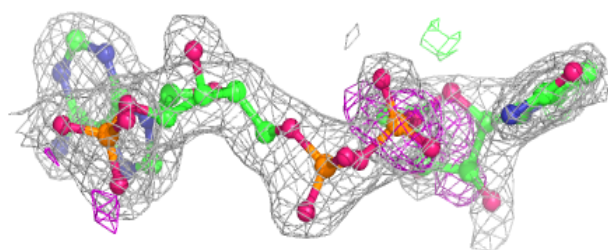
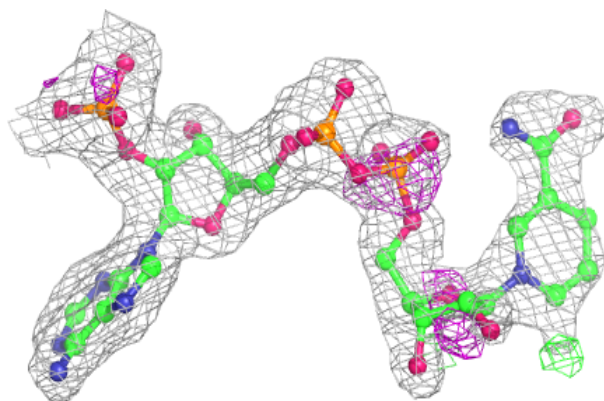
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NDP	A	501	48/48	0.92	0.11	27,38,50,50	0
2	NDP	B	502	48/48	0.93	0.12	22,39,56,56	0
3	MN	A	601	1/1	0.99	0.05	25,25,25,25	0
3	MN	B	602	1/1	0.99	0.04	27,27,27,27	0

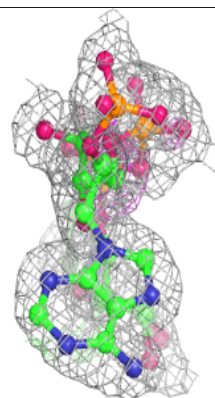
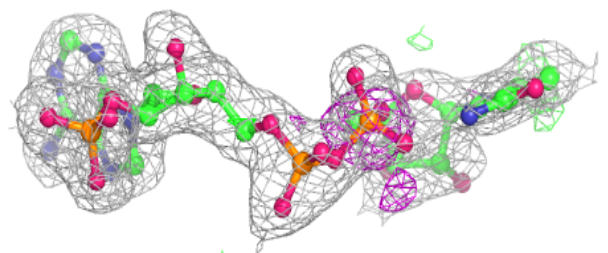
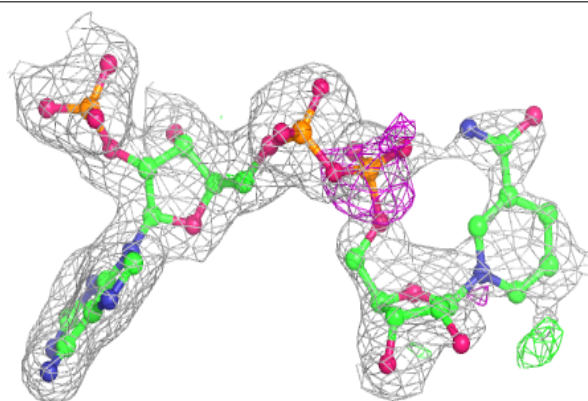
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NDP A 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 NDP B 502:**

$2mF_o-DF_c$ (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.