Full wwPDB X-ray Structure Validation Report (i)

Oct 17, 2023 – 03:22 AM EDT

PDB ID	:	2IKG
Title	:	Aldose reductase complexed with nitrophenyl-oxadiazol type inhibitor at 1.43 A
Authors	:	Steuber, H.; Koch, C.; Heine, A.; Klebe, G.
Deposited on	:	2006-10-02
Resolution	:	1.43 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 (i) 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 (i)) were used in the production of this report:

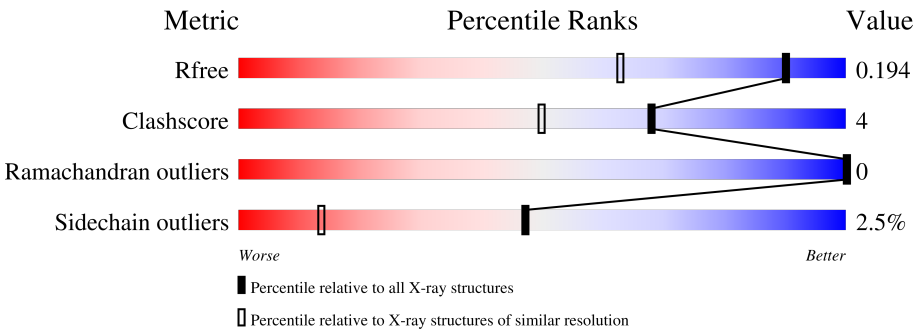
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.43 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	2021 (1.46-1.42)
Clashscore	141614	2086 (1.46-1.42)
Ramachandran outliers	138981	2047 (1.46-1.42)
Sidechain outliers	138945	2047 (1.46-1.42)

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	316	87% 12% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2936 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.

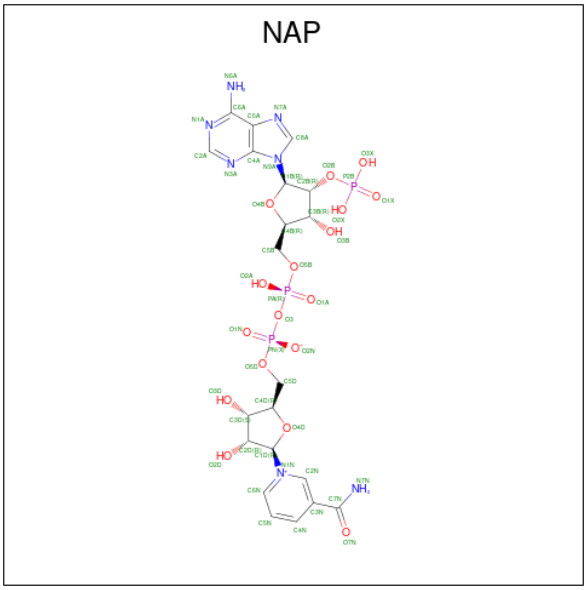
- Molecule 1 is a protein called Aldose reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	316	2545	1633	427	472	13	0	6	0

There is a discrepancy between the modelled and reference sequences:

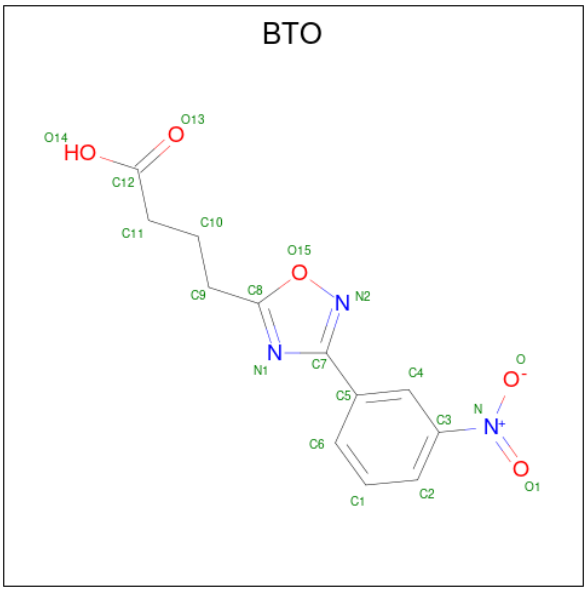
Chain	Residue	Modelled	Actual	Comment	Reference
A	4	ILE	LEU	SEE REMARK 999	UNP P15121

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



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

- Molecule 3 is 4-[3-(3-NITROPHENYL)-1,2,4-OXADIAZOL-5-YL]BUTANOIC ACID (three-letter code: BTO) (formula: $C_{12}H_{11}N_3O_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	20	12	3	5	0	0

- Molecule 4 is water.

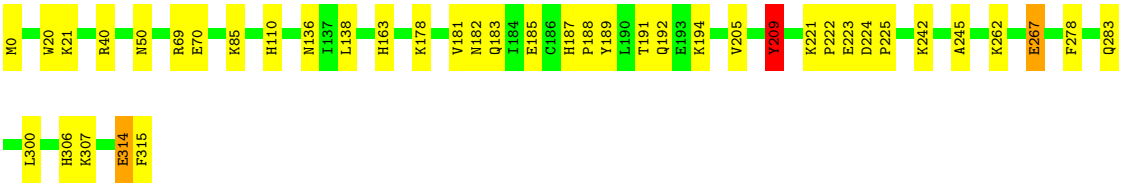
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	323	323	323	0	0

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. 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. 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: Aldose reductase

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	49.28Å 66.75Å 47.15Å 90.00° 92.59° 90.00°	Depositor
Resolution (Å)	25.00 – 1.43 27.62 – 1.43	Depositor EDS
% Data completeness (in resolution range)	98.6 (25.00-1.43) 89.5 (27.62-1.43)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) >$ [1]	3.12 (at 1.43Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.130 , 0.200 0.178 , 0.194	Depositor DCC
R_{free} test set	2546 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	14.1	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 41.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.002 for -l,k,h 0.031 for -h,-k,l 0.024 for l,-k,h	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	2936	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.67% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: BTO, NAP

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	0/2631	1.20	11/3572 (0.3%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	283	GLN	CG-CD-OE1	9.73	141.06	121.60
1	A	209	TYR	CB-CG-CD1	9.38	126.63	121.00
1	A	40	ARG	NE-CZ-NH2	-8.33	116.14	120.30
1	A	209	TYR	CA-CB-CG	7.26	127.20	113.40
1	A	306	HIS	CG-ND1-CE1	6.55	117.37	108.20
1	A	307	LYS	C-N-CA	6.43	137.77	121.70
1	A	209	TYR	CG-CD1-CE1	5.93	126.04	121.30
1	A	69	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	A	314	GLU	N-CA-CB	5.60	120.69	110.60
1	A	188	PRO	O-C-N	-5.24	114.31	122.70
1	A	245	ALA	O-C-N	-5.00	114.70	122.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2545	0	2551	20	0
2	A	48	0	25	2	0
3	A	20	0	10	2	0
4	A	323	0	0	3	0
All	All	2936	0	2586	20	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 4.

All (20) 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:222:PRO:HG2	1:A:223:GLU:OE1	1.88	0.74
1:A:189:TYR:CE2	1:A:225:PRO:HB2	2.33	0.63
1:A:267:GLU:HG2	4:A:469:HOH:O	1.98	0.62
1:A:185:GLU:OE2	1:A:187:HIS:HE1	1.89	0.55
1:A:194:LYS:HD3	1:A:315:PHE:CE2	2.43	0.54
1:A:183:GLN:OE1	2:A:316:NAP:H2N	2.10	0.50
1:A:300:LEU:HG	3:A:317:BTO:N2	2.28	0.49
1:A:70:GLU:O	1:A:70:GLU:HG3	2.13	0.47
1:A:136:ASN:ND2	1:A:138:LEU:H	2.14	0.46
1:A:221:LYS:O	1:A:224:ASP:HB2	2.16	0.46
1:A:262:LYS:O	2:A:316:NAP:H8A	2.16	0.46
1:A:110:HIS:NE2	3:A:317:BTO:O14	2.50	0.45
1:A:20:TRP:CE3	1:A:21:LYS:HE3	2.53	0.44
1:A:178:LYS:NZ	4:A:584:HOH:O	2.50	0.43
1:A:181:VAL:HA	1:A:205:VAL:O	2.18	0.43
1:A:242:LYS:HE2	1:A:278:PHE:CD1	2.54	0.43
1:A:314:GLU:HG2	1:A:315:PHE:N	2.32	0.42
1:A:185:GLU:HB2	1:A:209:TYR:CE1	2.54	0.42
1:A:191:THR:O	1:A:192:GLN:HB2	2.20	0.42
1:A:136:ASN:HB2	4:A:448:HOH:O	2.21	0.41

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	320/316 (101%)	316 (99%)	4 (1%)	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	287/281 (102%)	280 (98%)	7 (2%)	49	15

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	MET
1	A	50	ASN
1	A	85	LYS
1	A	163	HIS
1	A	182	ASN
1	A	209	TYR
1	A	267	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	93	GLN
1	A	136	ASN
1	A	182	ASN
1	A	187	HIS
1	A	283	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	BTO	A	317	-	16,21,21	2.44	7 (43%)	14,28,28	2.03	4 (28%)
2	NAP	A	316	-	45,52,52	1.84	11 (24%)	56,80,80	2.05	11 (19%)

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	BTO	A	317	-	-	3/7/14/14	0/2/2/2
2	NAP	A	316	-	-	4/31/67/67	0/5/5/5

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	317	BTO	C6-C5	5.47	1.55	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	316	NAP	C2N-N1N	5.39	1.41	1.35
2	A	316	NAP	C2A-N3A	4.18	1.38	1.32
3	A	317	BTO	C7-N1	4.14	1.42	1.34
2	A	316	NAP	C2A-N1A	4.10	1.41	1.33
2	A	316	NAP	C4A-N3A	3.47	1.40	1.35
3	A	317	BTO	C7-N2	-3.35	1.28	1.34
2	A	316	NAP	C3N-C7N	3.26	1.55	1.50
2	A	316	NAP	C6N-N1N	2.88	1.42	1.35
3	A	317	BTO	O14-C12	-2.83	1.21	1.30
2	A	316	NAP	O4D-C1D	2.63	1.44	1.41
2	A	316	NAP	P2B-O2B	2.63	1.64	1.59
3	A	317	BTO	C2-C3	2.60	1.43	1.38
2	A	316	NAP	C4N-C3N	2.49	1.43	1.39
3	A	317	BTO	C4-C3	2.48	1.42	1.36
2	A	316	NAP	C5N-C4N	2.26	1.43	1.38
3	A	317	BTO	C3-N	-2.22	1.39	1.45
2	A	316	NAP	C3D-C4D	2.08	1.58	1.53

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	316	NAP	C3N-C2N-N1N	-6.93	113.66	120.43
2	A	316	NAP	N3A-C2A-N1A	-4.99	120.88	128.68
2	A	316	NAP	C3N-C7N-N7N	-4.68	112.13	117.75
2	A	316	NAP	C5A-C6A-N6A	4.67	127.45	120.35
2	A	316	NAP	C2N-C3N-C4N	4.32	123.15	118.26
3	A	317	BTO	C2-C3-N	4.30	122.61	119.38
3	A	317	BTO	C4-C3-N	-3.86	115.34	118.75
2	A	316	NAP	C6N-N1N-C2N	3.57	125.23	121.97
2	A	316	NAP	C6N-C5N-C4N	-3.25	114.72	119.44
2	A	316	NAP	O4D-C4D-C3D	-3.03	99.12	105.11
3	A	317	BTO	C9-C10-C11	-2.92	104.72	114.34
2	A	316	NAP	PN-O3-PA	2.78	142.37	132.83
2	A	316	NAP	C2A-N1A-C6A	2.59	123.18	118.75
2	A	316	NAP	C5A-C6A-N1A	-2.46	114.77	120.35
3	A	317	BTO	O14-C12-C11	2.42	121.79	114.03

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	316	NAP	C4D-C5D-O5D-PN

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Mol	Chain	Res	Type	Atoms
2	A	316	NAP	PA-O3-PN-O5D
3	A	317	BTO	C10-C11-C12-O13
2	A	316	NAP	C2B-O2B-P2B-O3X
3	A	317	BTO	C10-C11-C12-O14
2	A	316	NAP	PA-O3-PN-O2N
3	A	317	BTO	C11-C10-C9-C8

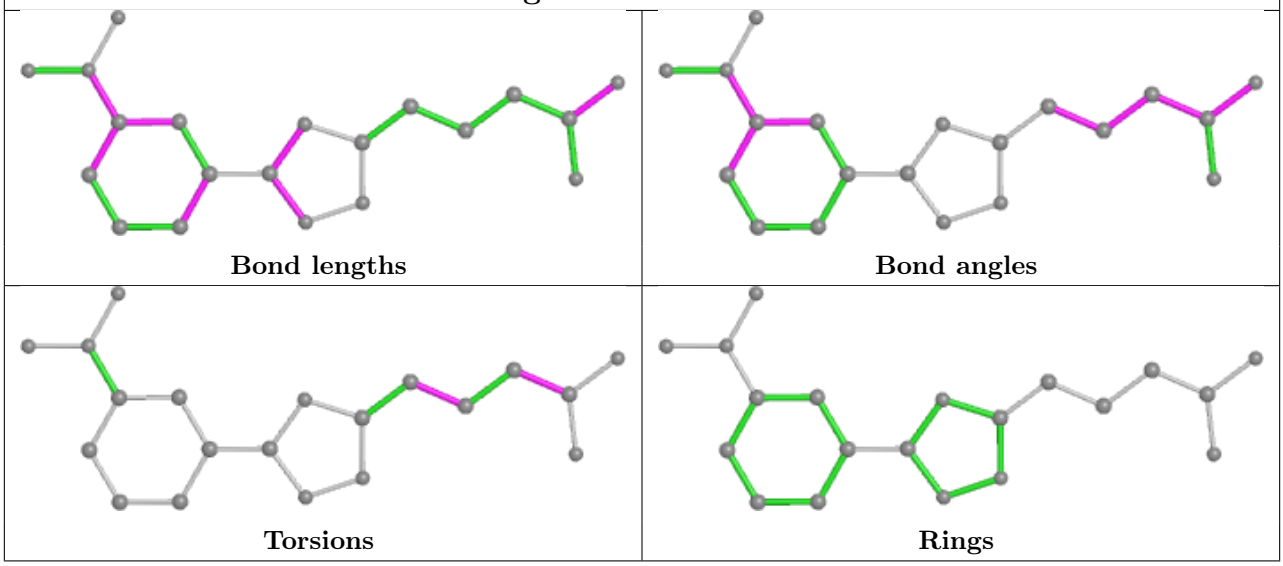
There are no ring outliers.

2 monomers are involved in 4 short contacts:

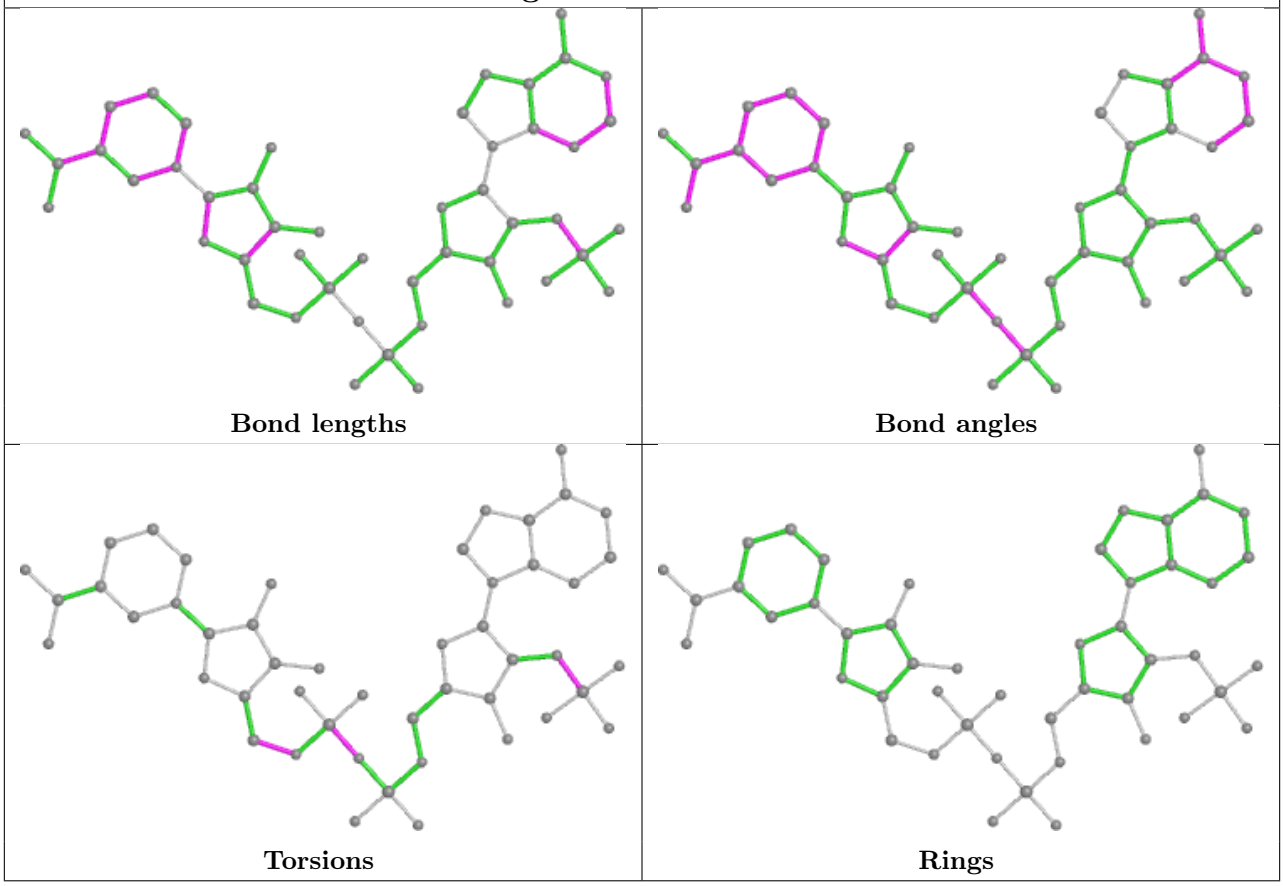
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	317	BTO	2	0
2	A	316	NAP	2	0

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 then 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.

Ligand BTO A 317



Ligand NAP A 316



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 (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

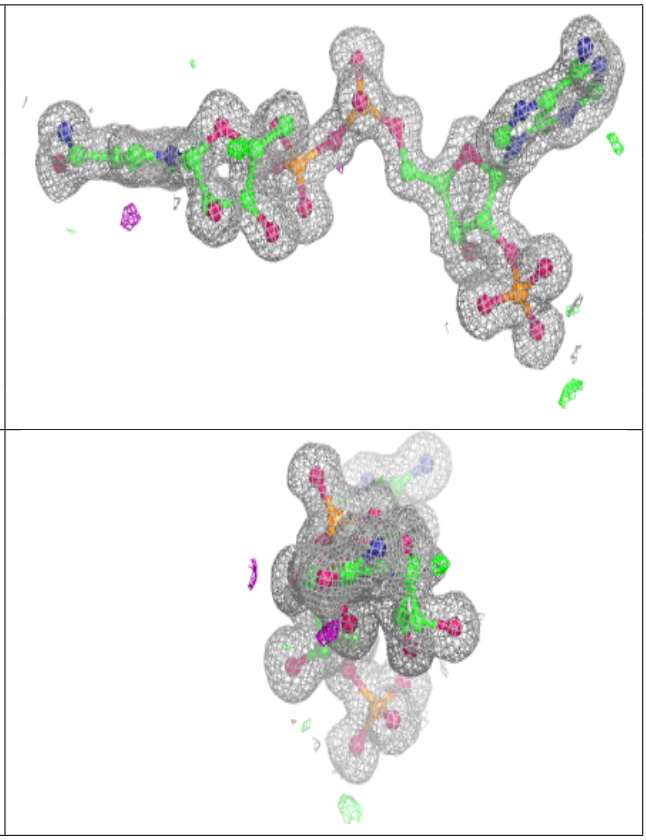
6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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.

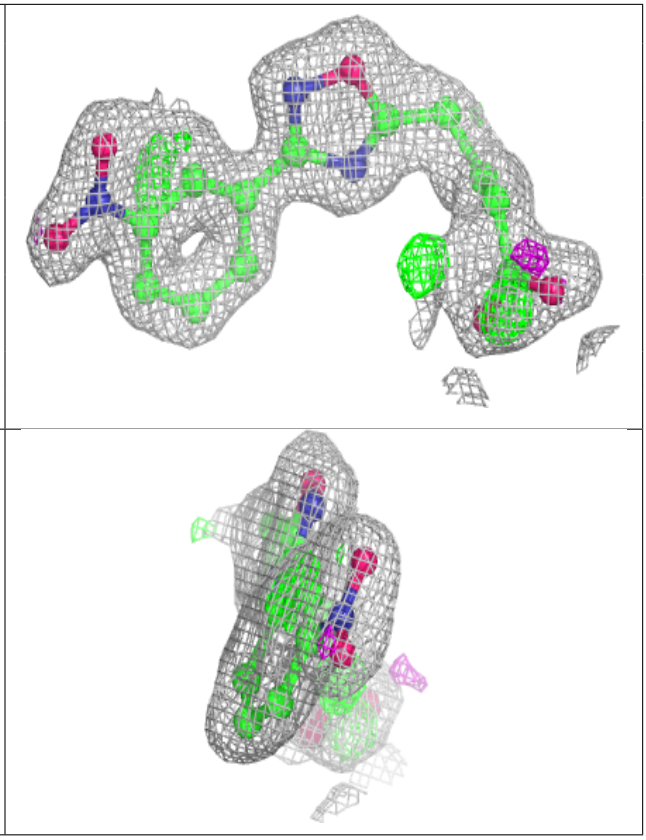
Electron density around NAP A 316:

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



Electron density around BTO A 317:

$2mF_o$-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)

Unable to reproduce the depositors R factor - this section is therefore empty.