



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 09:53 pm BST

PDB ID : 4CDG
Title : Crystal structure of the Bloom's syndrome helicase BLM in complex with Nanobody
Authors : Newman, J.A.; Savitsky, P.; Allerston, C.K.; Pike, A.C.W.; Pardon, E.; Steyaert, J.; Arrowsmith, C.H.; von Delft, F.; Bountra, C.; Edwards, A.; Gileadi, O.
Deposited on : 2013-10-31
Resolution : 2.79 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

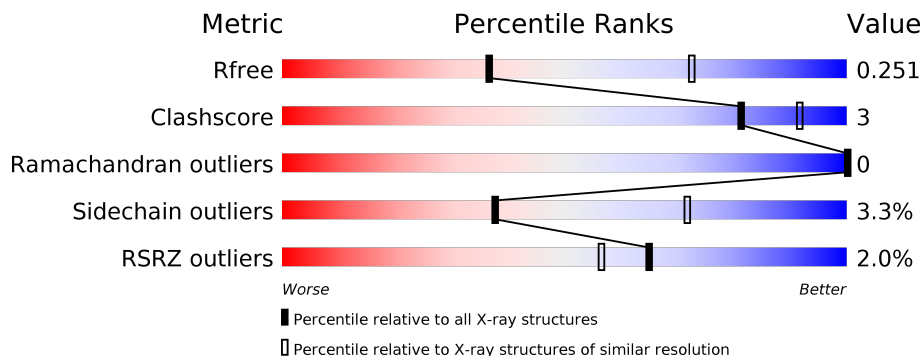
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	673	<p>2% 84% 8% 7%</p>
1	B	673	<p>0% 84% 8% 8%</p>
2	C	147	<p>2% 74% 9% 17%</p>
2	D	147	<p>2% 71% 12% 17%</p>

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 11711 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 BLOOM SYNDROME PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	625	Total	C	N	O	S	0	0	0
			4910	3127	834	916	33			
1	B	616	Total	C	N	O	S	0	0	0
			4858	3091	823	910	34			

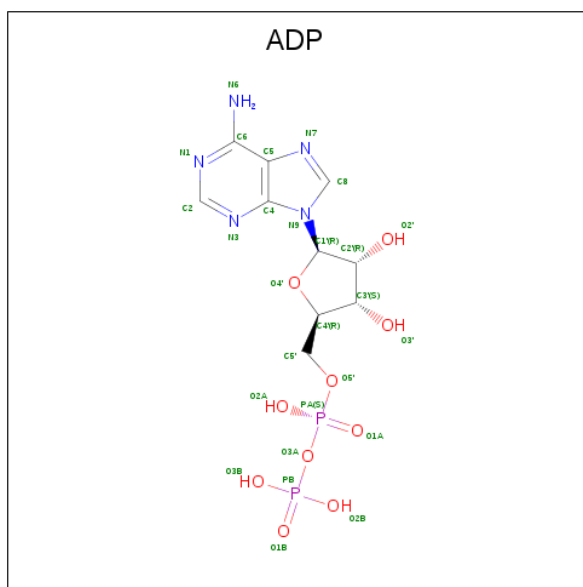
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	634	SER	-	expression tag	UNP P54132
A	635	MET	-	expression tag	UNP P54132
A	1299	MET	-	expression tag	UNP P54132
A	1300	SER	-	expression tag	UNP P54132
A	1301	GLU	-	expression tag	UNP P54132
A	1302	ASN	-	expression tag	UNP P54132
A	1303	LEU	-	expression tag	UNP P54132
A	1304	TYR	-	expression tag	UNP P54132
A	1305	PHE	-	expression tag	UNP P54132
A	1306	GLN	-	expression tag	UNP P54132
B	634	SER	-	expression tag	UNP P54132
B	635	MET	-	expression tag	UNP P54132
B	1299	MET	-	expression tag	UNP P54132
B	1300	SER	-	expression tag	UNP P54132
B	1301	GLU	-	expression tag	UNP P54132
B	1302	ASN	-	expression tag	UNP P54132
B	1303	LEU	-	expression tag	UNP P54132
B	1304	TYR	-	expression tag	UNP P54132
B	1305	PHE	-	expression tag	UNP P54132
B	1306	GLN	-	expression tag	UNP P54132

- Molecule 2 is a protein called NANOBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	122	Total	C	N	O	S	0	0	0
			942	587	168	181	6			
2	D	122	Total	C	N	O	S	0	0	0
			945	589	168	182	6			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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



- Molecule 2: NANOBODY



HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.67Å 88.42Å 95.12Å 69.39° 89.84° 77.07°	Depositor
Resolution (Å)	42.88 – 2.79 57.95 – 2.79	Depositor EDS
% Data completeness (in resolution range)	98.5 (42.88-2.79) 98.5 (57.95-2.79)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.81Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.206 , 0.245 0.217 , 0.251	Depositor DCC
R_{free} test set	2174 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	67.5	Xtrriage
Anisotropy	0.333	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11711	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.07% 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: ZN, ADP

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	0.21	0/5007	0.39	0/6771
1	B	0.22	0/4953	0.39	0/6692
2	C	0.22	0/962	0.37	0/1304
2	D	0.22	0/964	0.38	0/1305
All	All	0.22	0/11886	0.38	0/16072

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 [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	4910	0	4821	24	0
1	B	4858	0	4761	23	0
2	C	942	0	891	6	0
2	D	945	0	903	7	0
3	A	27	0	12	0	0
3	B	27	0	12	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	11711	0	11400	59	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (59) 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:676:ILE:HD13	1:B:701:LEU:HD23	1.69	0.74
1:A:1211:ARG:NH2	1:A:1251:GLU:O	2.27	0.67
1:B:859:ARG:NH1	1:B:978:GLY:O	2.27	0.67
1:A:859:ARG:NH1	1:A:978:GLY:O	2.29	0.66
2:D:59:TYR:OH	2:D:69:ARG:NH1	2.32	0.63
2:D:58:TRP:NE1	2:D:100:MET:SD	2.72	0.63
1:A:643:ARG:NH1	1:B:823:SER:O	2.36	0.58
1:B:1211:ARG:NH2	1:B:1251:GLU:O	2.36	0.58
2:C:59:TYR:OH	2:C:69:ARG:NH1	2.37	0.58
2:C:84:SER:O	2:C:88:ARG:NH1	2.36	0.57
2:C:58:TRP:NE1	2:C:100:MET:SD	2.78	0.56
1:A:1116:ASP:OD1	1:A:1139:ARG:NH2	2.35	0.56
1:A:733:PRO:HB2	1:A:760:ILE:HA	1.89	0.55
1:B:1221:GLU:OE2	1:B:1285:TYR:OH	2.25	0.53
1:B:966:LEU:HD22	1:B:967:PRO:HD2	1.91	0.53
1:B:1046:GLY:O	1:B:1048:ASN:ND2	2.42	0.53
2:D:28:GLU:OE2	2:D:117:CYS:N	2.37	0.52
1:A:671:ASN:ND2	1:A:854:SER:O	2.42	0.52
1:A:1221:GLU:OE2	1:A:1285:TYR:OH	2.26	0.52
1:B:1116:ASP:OD1	1:B:1139:ARG:NH2	2.33	0.52
2:D:84:SER:O	2:D:88:ARG:NH1	2.43	0.51
1:B:1194:ASN:OD1	1:B:1195:SER:N	2.42	0.51
1:A:655:MET:HG3	1:A:701:LEU:HD11	1.93	0.51
1:A:735:THR:HG21	1:A:749:ILE:HD12	1.94	0.50
1:B:671:ASN:ND2	1:B:854:SER:O	2.45	0.50
1:B:662:LYS:NZ	1:B:731:ASP:O	2.45	0.49
2:C:58:TRP:CD1	2:C:102:LEU:HB2	2.49	0.48
2:C:88:ARG:NH2	2:C:111:ASP:OD2	2.46	0.48
1:A:701:LEU:HB3	1:A:702:PRO:HD3	1.96	0.48
2:D:58:TRP:CD1	2:D:102:LEU:HB2	2.49	0.47
1:A:998:VAL:HG12	1:A:1002:LYS:HE3	1.97	0.47
1:A:893:ILE:HD13	1:A:904:MET:HE2	1.97	0.46
1:B:1175:ASN:OD1	1:B:1176:LYS:N	2.49	0.45
1:A:735:THR:OG1	1:A:736:TYR:N	2.48	0.45
1:A:954:ASP:HB2	1:A:979:ARG:NH2	2.32	0.45
1:A:909:GLN:NE2	1:A:915:ALA:O	2.50	0.44
1:B:994:THR:HG22	1:B:996:HIS:H	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1064:ASP:OD1	1:B:1065:ASN:N	2.51	0.43
1:B:665:LEU:HD13	3:B:1634:ADP:C6	2.54	0.43
1:B:735:THR:HG22	1:B:760:ILE:HD12	2.00	0.43
1:A:699:TYR:HB3	1:A:827:MET:HE1	2.00	0.43
1:B:735:THR:OG1	1:B:736:TYR:N	2.51	0.43
1:A:1064:ASP:OD1	1:A:1065:ASN:N	2.51	0.43
2:D:104:MET:HE2	2:D:107:LEU:HD21	2.00	0.42
1:A:863:LYS:HG2	1:A:1061:VAL:HA	2.01	0.42
1:B:898:ARG:HG3	1:B:919:HIS:CG	2.54	0.42
1:A:710:THR:HB	1:A:762:LEU:HD13	2.02	0.42
1:B:859:ARG:NH2	1:B:861:ASN:OD1	2.53	0.42
2:D:82:VAL:HG12	2:D:85:VAL:HG22	2.01	0.41
1:B:701:LEU:HB3	1:B:702:PRO:HD3	2.03	0.41
1:A:735:THR:HG22	1:A:760:ILE:HD12	2.02	0.41
1:B:1122:LYS:HE3	1:B:1122:LYS:HB3	1.90	0.41
1:B:893:ILE:HD13	1:B:904:MET:HE2	2.03	0.41
2:C:74:THR:HG23	2:C:76:ALA:H	1.86	0.41
1:A:1277:GLU:N	1:A:1277:GLU:OE1	2.45	0.40
1:A:714:SER:O	1:A:767:PRO:HD3	2.21	0.40
1:A:1226:CYS:SG	1:A:1246:LEU:HD21	2.61	0.40
1:A:834:ASN:HA	1:A:835:PRO:HD3	1.91	0.40
1:B:1277:GLU:N	1:B:1277:GLU:OE1	2.47	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	619/673 (92%)	594 (96%)	25 (4%)	0	100	100
1	B	606/673 (90%)	584 (96%)	22 (4%)	0	100	100
2	C	120/147 (82%)	116 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	120/147 (82%)	116 (97%)	4 (3%)	0	100	100
All	All	1465/1640 (89%)	1410 (96%)	55 (4%)	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	531/601 (88%)	512 (96%)	19 (4%)	35	69
1	B	528/601 (88%)	512 (97%)	16 (3%)	41	75
2	C	99/118 (84%)	96 (97%)	3 (3%)	41	75
2	D	100/118 (85%)	96 (96%)	4 (4%)	31	65
All	All	1258/1438 (88%)	1216 (97%)	42 (3%)	38	72

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	662	LYS
1	A	688	LEU
1	A	709	VAL
1	A	714	SER
1	A	722	ASP
1	A	756	LYS
1	A	789	LEU
1	A	874	VAL
1	A	899	ARG
1	A	910	ARG
1	A	949	PHE
1	A	955	LYS
1	A	994	THR
1	A	1108	ARG
1	A	1121	SER
1	A	1139	ARG

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Mol	Chain	Res	Type
1	A	1156	ASP
1	A	1164	ASN
1	A	1205	VAL
1	B	709	VAL
1	B	714	SER
1	B	722	ASP
1	B	757	ASP
1	B	789	LEU
1	B	816	MET
1	B	899	ARG
1	B	910	ARG
1	B	949	PHE
1	B	955	LYS
1	B	1033	ILE
1	B	1121	SER
1	B	1122	LYS
1	B	1139	ARG
1	B	1156	ASP
1	B	1199	LYS
2	C	41	ARG
2	C	62	THR
2	C	142	HIS
2	D	41	ARG
2	D	62	THR
2	D	75	SER
2	D	99	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	802	GLN
1	A	1021	ASN
1	A	1058	HIS
1	A	1089	GLN
1	B	802	GLN
1	B	975	GLN
1	B	1048	ASN
1	B	1162	ASN
2	C	118	ASN
2	C	136	GLN
2	D	23	GLN
2	D	118	ASN

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Mol	Chain	Res	Type
2	D	136	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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
3	ADP	A	1634	-	24,29,29	0.95	1 (4%)	29,45,45	1.47	4 (13%)
3	ADP	B	1634	-	24,29,29	0.96	1 (4%)	29,45,45	1.39	3 (10%)

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	ADP	A	1634	-	-	3/12/32/32	0/3/3/3
3	ADP	B	1634	-	-	2/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1634	ADP	C5-C4	2.45	1.47	1.40
3	A	1634	ADP	C5-C4	2.43	1.47	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1634	ADP	PA-O3A-PB	-3.64	120.35	132.83
3	B	1634	ADP	PA-O3A-PB	-3.55	120.64	132.83
3	A	1634	ADP	N3-C2-N1	-3.38	123.39	128.68
3	B	1634	ADP	N3-C2-N1	-3.28	123.56	128.68
3	B	1634	ADP	C4-C5-N7	-2.81	106.47	109.40
3	A	1634	ADP	C4-C5-N7	-2.68	106.61	109.40
3	A	1634	ADP	C3'-C2'-C1'	2.67	105.00	100.98

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1634	ADP	C5'-O5'-PA-O1A
3	A	1634	ADP	C5'-O5'-PA-O3A
3	A	1634	ADP	C5'-O5'-PA-O2A
3	B	1634	ADP	PA-O3A-PB-O2B
3	B	1634	ADP	C5'-O5'-PA-O1A

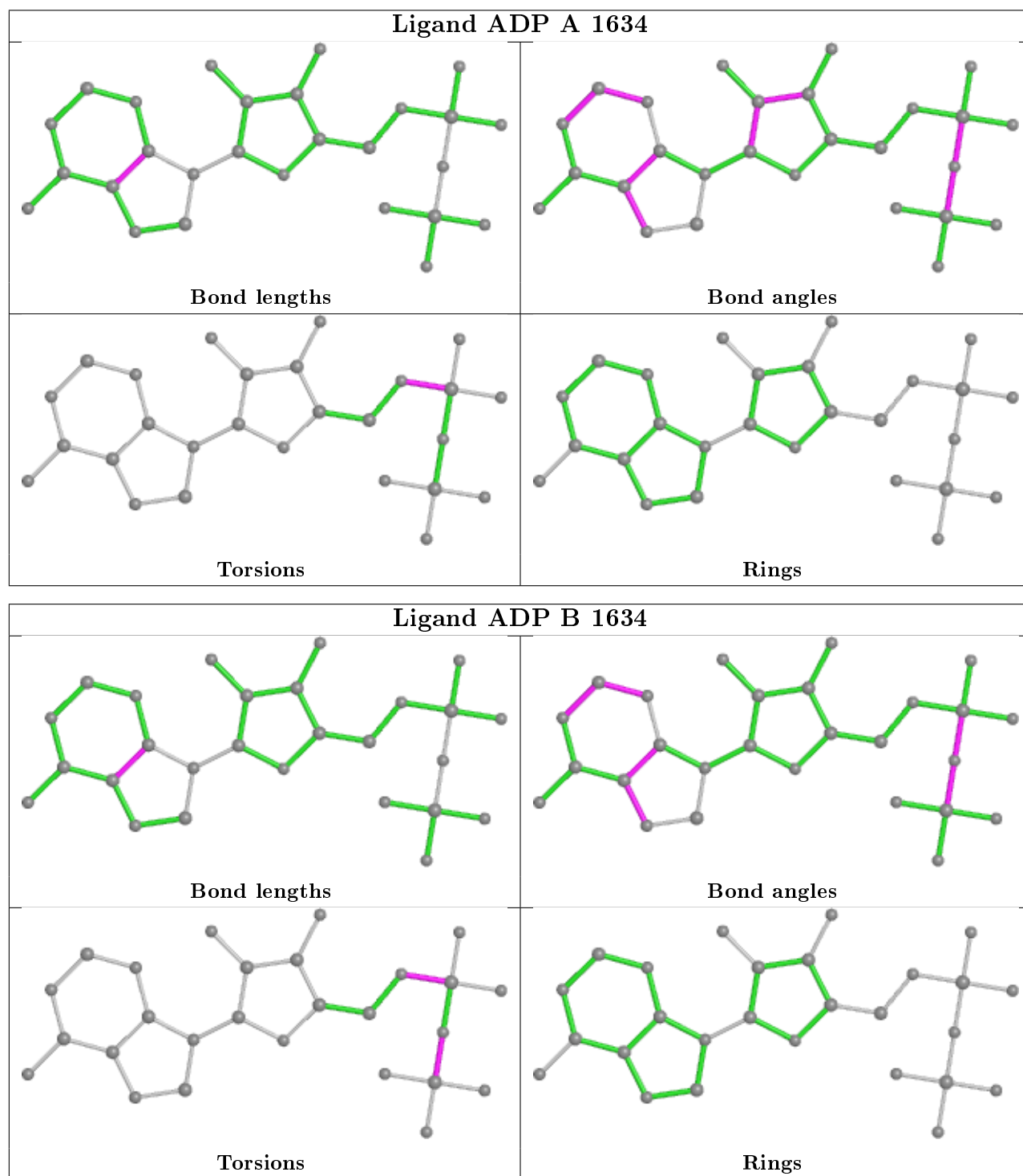
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1634	ADP	1	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

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	625/673 (92%)	0.19	15 (2%) 59 49	37, 67, 111, 134	0
1	B	616/673 (91%)	0.14	8 (1%) 77 72	33, 61, 98, 115	0
2	C	122/147 (82%)	0.24	3 (2%) 57 47	41, 81, 109, 123	0
2	D	122/147 (82%)	0.19	3 (2%) 57 47	43, 70, 93, 119	0
All	All	1485/1640 (90%)	0.18	29 (1%) 65 56	33, 66, 107, 134	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1171	VAL	3.8
1	A	1017	GLU	3.7
1	B	772	ALA	3.5
1	B	1160	TYR	3.4
1	A	1128	SER	3.3
1	A	1109	PHE	3.1
1	A	1170	TYR	3.1
1	B	1128	SER	3.0
1	A	746	ALA	2.9
2	D	66	GLN	2.8
2	C	144	HIS	2.8
1	A	1002	LYS	2.6
1	A	811	TYR	2.5
1	B	848	LEU	2.4
2	C	143	HIS	2.4
2	D	144	HIS	2.4
1	A	1172	MET	2.4
1	A	1160	TYR	2.3
1	A	1126	ILE	2.3
1	A	792	PHE	2.3
1	A	1088	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	1171	VAL	2.2
1	A	644	PHE	2.2
1	A	744	SER	2.2
2	C	34	VAL	2.2
1	B	814	MET	2.1
2	D	65	LYS	2.1
1	B	643	ARG	2.0
1	B	1109	PHE	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](#)

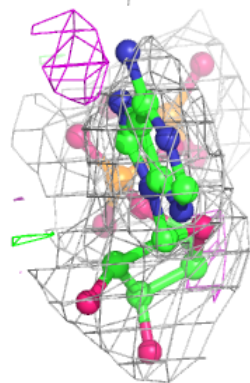
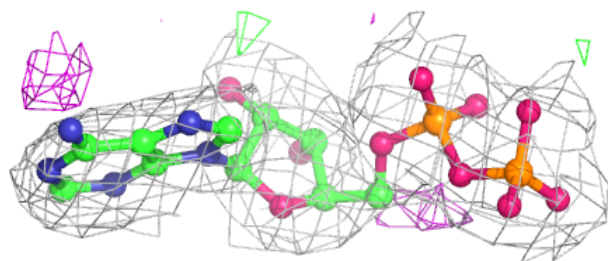
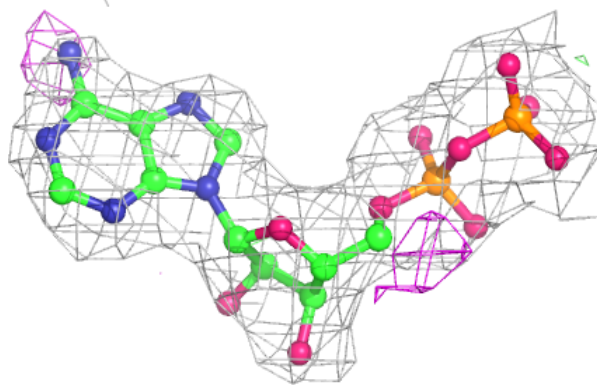
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ADP	B	1634	27/27	0.96	0.19	29,47,55,62	0
3	ADP	A	1634	27/27	0.97	0.20	34,52,64,84	0
4	ZN	B	1643	1/1	0.99	0.21	55,55,55,55	0
4	ZN	A	1643	1/1	1.00	0.19	54,54,54,54	0

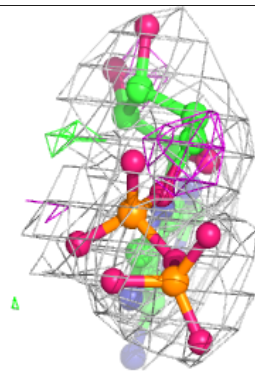
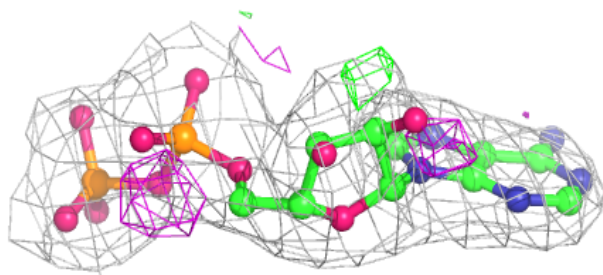
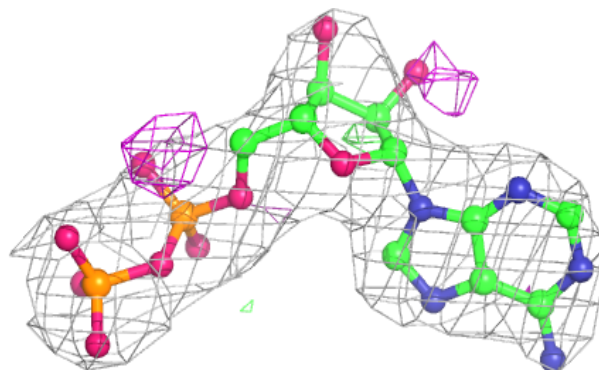
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP B 1634:

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

**Electron density around ADP A 1634:**

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



6.5 Other polymers

There are no such residues in this entry.