



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 7A8R / pdb_00007a8r
Title : Structure of RecQL from Bos taurus
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Deposited on : 2020-08-30
Resolution : 2.31 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

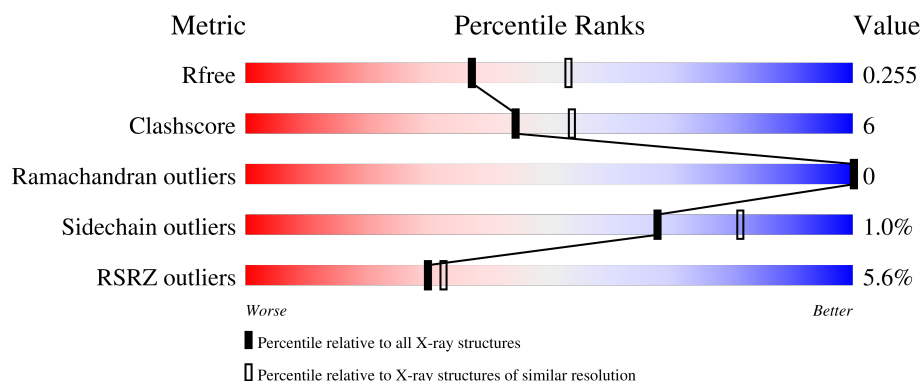
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.31 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	531	6% 84% 15%
1	B	531	5% 82% 17%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8618 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 ATP-dependent DNA helicase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	531	Total	C	N	O	S	0	0	0
			4237	2704	727	770	36			
1	B	530	Total	C	N	O	S	0	0	0
			4229	2699	726	769	35			

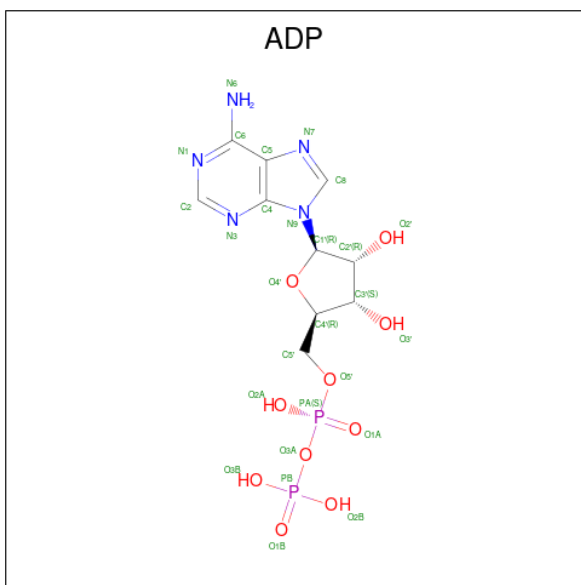
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	62	MET	-	initiating methionine	UNP A0JN36
B	62	MET	-	initiating methionine	UNP A0JN36

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

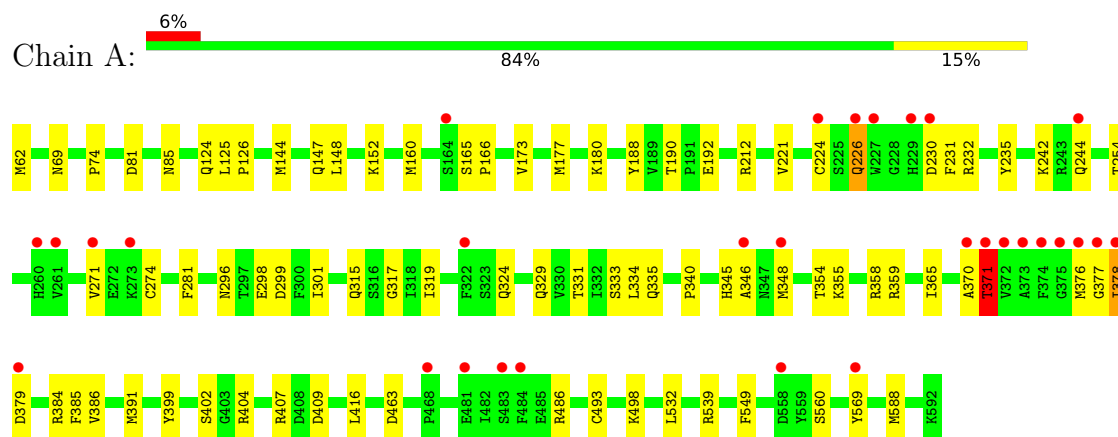
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	57	Total O 57 57	0	0
5	B	37	Total O 37 37	0	0

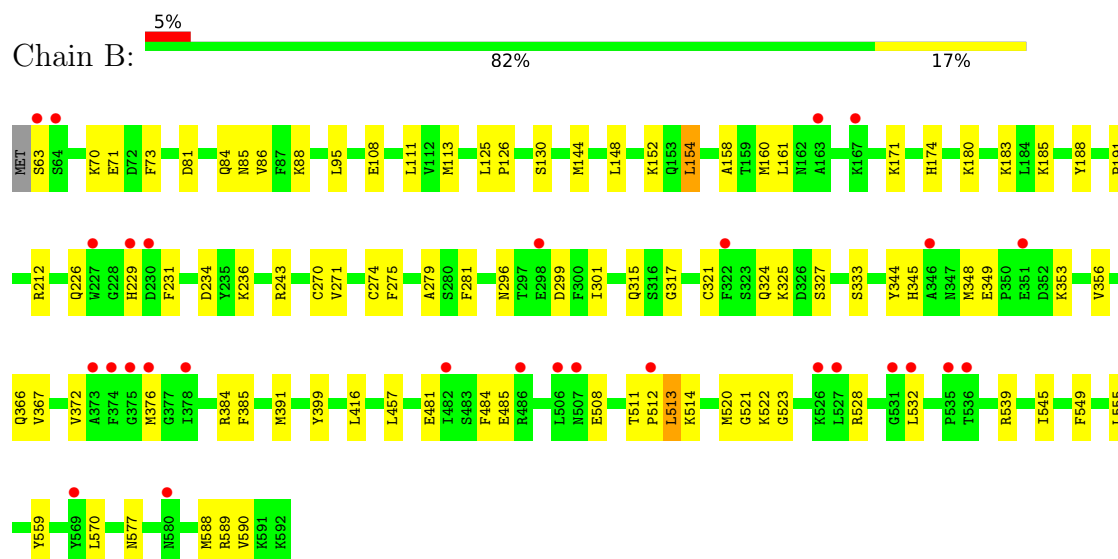
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: ATP-dependent DNA helicase



• Molecule 1: ATP-dependent DNA helicase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.58Å 117.22Å 203.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.16 – 2.31 23.16 – 2.31	Depositor EDS
% Data completeness (in resolution range)	75.2 (23.16-2.31) 75.2 (23.16-2.31)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.31Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.201 , 0.253 0.202 , 0.255	Depositor DCC
R_{free} test set	2299 reflections (3.86%)	wwPDB-VP
Wilson B-factor (Å ²)	49.3	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8618	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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: ADP, ZN, MG

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	0.20	0/4325	0.50	0/5830
1	B	0.19	0/4317	0.48	0/5820
All	All	0.19	0/8642	0.49	0/11650

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	3
All	All	0	7

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	370	ALA	Peptide
1	A	371	THR	Peptide
1	A	377	GLY	Peptide
1	A	378	ILE	Peptide
1	B	229	HIS	Peptide
1	B	481	GLU	Peptide
1	B	508	GLU	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4237	0	4295	47	2
1	B	4229	0	4286	55	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	27	0	12	0	0
3	B	27	0	12	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	57	0	0	6	0
5	B	37	0	0	6	0
All	All	8618	0	8605	102	2

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 (102) 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:463:ASP:OD1	5:A:1101:HOH:O	1.90	0.88
1:A:69:ASN:OD1	5:A:1102:HOH:O	1.95	0.85
1:B:243:ARG:NE	1:B:270:CYS:SG	2.57	0.77
1:A:180:LYS:HA	1:A:212:ARG:HD2	1.67	0.76
1:B:348:MET:HB2	1:B:353:LYS:HD3	1.69	0.74
1:B:174:HIS:NE2	5:B:1102:HOH:O	2.20	0.73
1:A:62:MET:N	5:A:1104:HOH:O	2.28	0.67
1:A:301:ILE:HD13	1:A:333:SER:HB2	1.76	0.67
1:B:301:ILE:HD13	1:B:333:SER:HB2	1.76	0.67
1:B:144:MET:HG2	1:B:188:TYR:HB3	1.76	0.66
1:A:144:MET:HG2	1:A:188:TYR:HB3	1.78	0.66
1:A:329:GLN:NE2	5:A:1103:HOH:O	2.30	0.65
1:B:81:ASP:OD1	1:B:85:ASN:ND2	2.30	0.64
1:A:81:ASP:OD1	1:A:85:ASN:ND2	2.31	0.62
1:B:522:LYS:HD3	1:B:523:GLY:H	1.63	0.61
1:B:148:LEU:HD11	1:B:160:MET:HB3	1.82	0.61
1:B:366:GLN:HG3	1:B:367:VAL:HG13	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:GLU:OE2	5:A:1103:HOH:O	2.16	0.60
1:A:296:ASN:HB3	1:A:299:ASP:HB2	1.85	0.59
1:B:511:THR:HG22	1:B:514:LYS:HB2	1.84	0.59
1:A:190:THR:HB	1:A:192:GLU:HG2	1.86	0.58
1:B:457:LEU:O	5:B:1101:HOH:O	2.17	0.58
1:A:539:ARG:NH2	5:A:1106:HOH:O	2.37	0.57
1:B:512:PRO:O	1:B:513:LEU:HB3	2.04	0.57
1:B:521:GLY:HA2	1:B:528:ARG:HD2	1.88	0.56
1:B:296:ASN:HB3	1:B:299:ASP:HB2	1.86	0.56
1:B:108:GLU:HG2	1:B:271:VAL:HG13	1.88	0.55
1:B:520:MET:SD	5:B:1117:HOH:O	2.58	0.55
1:A:124:GLN:NE2	1:A:147:GLN:OE1	2.40	0.54
1:B:324:GLN:HE21	1:B:345:HIS:CD2	2.25	0.54
1:A:271:VAL:O	1:A:274:CYS:HB2	2.08	0.54
1:B:549:PHE:HB3	1:B:555:LEU:HD13	1.89	0.53
1:A:498:LYS:HE3	1:A:532:LEU:HD21	1.91	0.52
1:B:324:GLN:HA	1:B:327:SER:HB3	1.90	0.52
1:A:144:MET:HE2	1:A:160:MET:HB2	1.91	0.52
1:B:231:PHE:HZ	1:B:236:LYS:HD3	1.75	0.52
1:A:345:HIS:HB2	1:A:348:MET:HG3	1.90	0.52
1:B:152:LYS:NZ	1:B:158:ALA:H	2.08	0.52
1:B:317:GLY:HA3	1:B:385:PHE:CZ	2.45	0.52
1:A:315:GLN:HB3	1:A:384:ARG:HG3	1.92	0.51
1:B:174:HIS:CE1	5:B:1102:HOH:O	2.61	0.51
1:A:331:THR:O	1:A:335:GLN:HG3	2.10	0.51
1:A:148:LEU:HD11	1:A:160:MET:HB3	1.92	0.51
1:B:376:MET:N	1:B:376:MET:SD	2.84	0.51
1:B:344:TYR:CE2	1:B:353:LYS:HD2	2.46	0.51
1:B:391:MET:HA	1:B:416:LEU:HD11	1.94	0.50
1:B:111:LEU:HD11	1:B:113:MET:HE3	1.92	0.50
1:B:315:GLN:HB3	1:B:384:ARG:HG3	1.93	0.50
1:A:242:LYS:HG2	1:A:271:VAL:HG12	1.94	0.50
1:A:386:VAL:HG11	1:A:402:SER:HA	1.94	0.50
1:B:325:LYS:HD2	1:B:325:LYS:H	1.78	0.49
1:B:70:LYS:HE2	1:B:73:PHE:CZ	2.47	0.49
1:B:63:SER:HA	1:B:275:PHE:HD1	1.79	0.47
1:B:324:GLN:HE21	1:B:345:HIS:CG	2.33	0.47
1:B:539:ARG:NH2	5:B:1107:HOH:O	2.47	0.47
1:A:224:CYS:HA	1:A:235:TYR:HB3	1.97	0.46
1:A:493:CYS:HB2	1:A:549:PHE:CZ	2.50	0.46
1:A:391:MET:HA	1:A:416:LEU:HD11	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:GLN:O	1:B:88:LYS:HA	2.16	0.46
1:A:144:MET:CE	1:A:160:MET:HB2	2.46	0.45
1:B:545:ILE:HA	1:B:588:MET:HE1	1.97	0.45
1:B:183:LYS:HA	1:B:183:LYS:HD3	1.85	0.45
1:B:191:PRO:HB2	1:B:234:ASP:HB3	1.99	0.45
1:B:231:PHE:CZ	1:B:236:LYS:HD3	2.51	0.45
1:A:340:PRO:HB2	1:A:365:ILE:HA	1.99	0.45
1:A:354:THR:HG22	1:A:358:ARG:HD2	1.98	0.45
1:B:171:LYS:HD3	1:B:171:LYS:HA	1.63	0.45
1:A:486:ARG:HA	1:A:588:MET:O	2.17	0.44
1:A:324:GLN:HG2	1:A:371:THR:CG2	2.48	0.44
1:A:317:GLY:HA3	1:A:385:PHE:CZ	2.52	0.44
1:B:344:TYR:HD2	1:B:356:VAL:HG11	1.82	0.44
1:B:144:MET:HE2	1:B:160:MET:HB2	1.99	0.44
1:B:321:CYS:O	1:B:372:VAL:HG23	2.18	0.44
1:A:226:GLN:H	1:A:226:GLN:HG2	1.58	0.44
1:A:226:GLN:HA	1:A:231:PHE:CG	2.53	0.44
1:A:177:MET:O	1:A:212:ARG:NH1	2.49	0.43
1:A:407:ARG:HD3	1:A:407:ARG:HA	1.38	0.43
1:B:485:GLU:HG2	1:B:590:VAL:HG23	2.01	0.43
1:B:130:SER:O	1:B:185:LYS:NZ	2.51	0.43
1:B:271:VAL:HG12	1:B:274:CYS:HB2	2.00	0.43
1:A:125:LEU:HB3	1:A:126:PRO:HD3	2.00	0.43
1:B:71:GLU:OE1	5:B:1103:HOH:O	2.21	0.43
1:A:221:VAL:HG12	1:A:254:THR:HB	2.01	0.43
1:B:180:LYS:HA	1:B:212:ARG:HD2	2.00	0.42
1:A:346:ALA:HB2	1:A:376:MET:HE2	2.00	0.42
1:B:86:VAL:O	1:B:154:LEU:HD12	2.19	0.42
1:B:484:PHE:CD1	1:B:589:ARG:HD3	2.55	0.42
1:A:319:ILE:HD12	1:A:334:LEU:HD11	2.02	0.42
1:B:281:PHE:O	1:B:399:TYR:OH	2.30	0.41
1:B:95:LEU:HD12	1:B:279:ALA:HB2	2.01	0.41
1:A:407:ARG:C	1:A:409:ASP:H	2.29	0.41
1:B:125:LEU:HB3	1:B:126:PRO:HD3	2.03	0.41
1:A:560:SER:HB2	1:A:569:TYR:HE2	1.86	0.41
1:B:512:PRO:HG3	1:B:570:LEU:HG	2.01	0.41
1:A:165:SER:HA	1:A:166:PRO:HD3	1.93	0.41
1:A:230:ASP:HB3	1:A:232:ARG:NH1	2.36	0.41
1:B:144:MET:CE	1:B:160:MET:HB2	2.51	0.41
1:A:355:LYS:O	1:A:359:ARG:HG3	2.20	0.41
1:A:173:VAL:O	1:A:177:MET:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:PHE:O	1:A:399:TYR:OH	2.26	0.40
1:A:379:ASP:CG	1:A:404:ARG:HE	2.29	0.40
1:B:161:LEU:HD23	1:B:161:LEU:HA	1.94	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:PRO:O	1:A:152:LYS:NZ[4_556]	2.01	0.19
1:A:244:GLN:NE2	1:B:559:TYR:O[3_655]	2.12	0.08

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	529/531 (100%)	528 (100%)	1 (0%)	0	100	100
1	B	528/531 (99%)	528 (100%)	0	0	100	100
All	All	1057/1062 (100%)	1056 (100%)	1 (0%)	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	469/469 (100%)	466 (99%)	3 (1%)	78	88
1	B	468/469 (100%)	462 (99%)	6 (1%)	61	76
All	All	937/938 (100%)	928 (99%)	9 (1%)	68	81

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	226	GLN
1	A	371	THR
1	A	378	ILE
1	B	154	LEU
1	B	226	GLN
1	B	349	GLU
1	B	513	LEU
1	B	532	LEU
1	B	577	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	A	389	HIS
1	A	553	GLN
1	A	580	ASN
1	B	181	ASN
1	B	324	GLN
1	B	335	GLN
1	B	389	HIS
1	B	548	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 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	B	1002	4	28,29,29	1.42	5 (17%)	43,45,45	1.87	10 (23%)
3	ADP	A	1002	4	28,29,29	1.42	4 (14%)	43,45,45	1.81	11 (25%)

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	B	1002	4	-	2/16/32/32	0/3/3/3
3	ADP	A	1002	4	-	2/16/32/32	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	ADP	C5-C4	4.67	1.47	1.39
3	B	1002	ADP	C5-C4	4.62	1.47	1.39
3	B	1002	ADP	C5-C6	2.80	1.48	1.41
3	A	1002	ADP	C5-C6	2.72	1.48	1.41
3	B	1002	ADP	C8-N7	2.44	1.36	1.31
3	A	1002	ADP	C8-N7	2.41	1.36	1.31
3	A	1002	ADP	C5-N7	-2.23	1.35	1.39
3	B	1002	ADP	PA-O3A	2.19	1.61	1.59
3	B	1002	ADP	C5-N7	-2.04	1.35	1.39

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1002	ADP	C5-C4-N3	-5.82	118.71	126.72
3	A	1002	ADP	C5-C4-N3	-5.69	118.89	126.72
3	B	1002	ADP	N3-C4-N9	4.51	134.84	127.17
3	A	1002	ADP	N3-C4-N9	4.26	134.41	127.17
3	A	1002	ADP	C4-C5-N7	-3.82	106.21	110.58
3	B	1002	ADP	C2-N3-C4	3.82	121.15	111.83
3	B	1002	ADP	C4-C5-N7	-3.71	106.34	110.58
3	A	1002	ADP	C2-N3-C4	3.49	120.36	111.83
3	B	1002	ADP	N3-C2-N1	-3.41	123.41	128.58
3	A	1002	ADP	N3-C2-N1	-3.01	124.03	128.58
3	B	1002	ADP	C4-N9-C8	2.77	108.65	105.74
3	A	1002	ADP	C5-N7-C8	2.70	107.69	103.45
3	B	1002	ADP	C5-N7-C8	2.69	107.69	103.45
3	A	1002	ADP	C4-N9-C8	2.48	108.34	105.74
3	B	1002	ADP	C6-C5-N7	2.31	136.54	132.09
3	B	1002	ADP	N9-C8-N7	-2.16	110.86	113.94
3	A	1002	ADP	C6-C5-N7	2.13	136.19	132.09
3	B	1002	ADP	O4'-C1'-N9	2.08	112.09	108.09
3	A	1002	ADP	N9-C8-N7	-2.03	111.06	113.94
3	A	1002	ADP	O4'-C1'-N9	2.02	111.97	108.09
3	A	1002	ADP	O3B-PB-O2B	2.00	115.31	107.80

There are no chirality outliers.

All (4) torsion outliers are listed below:

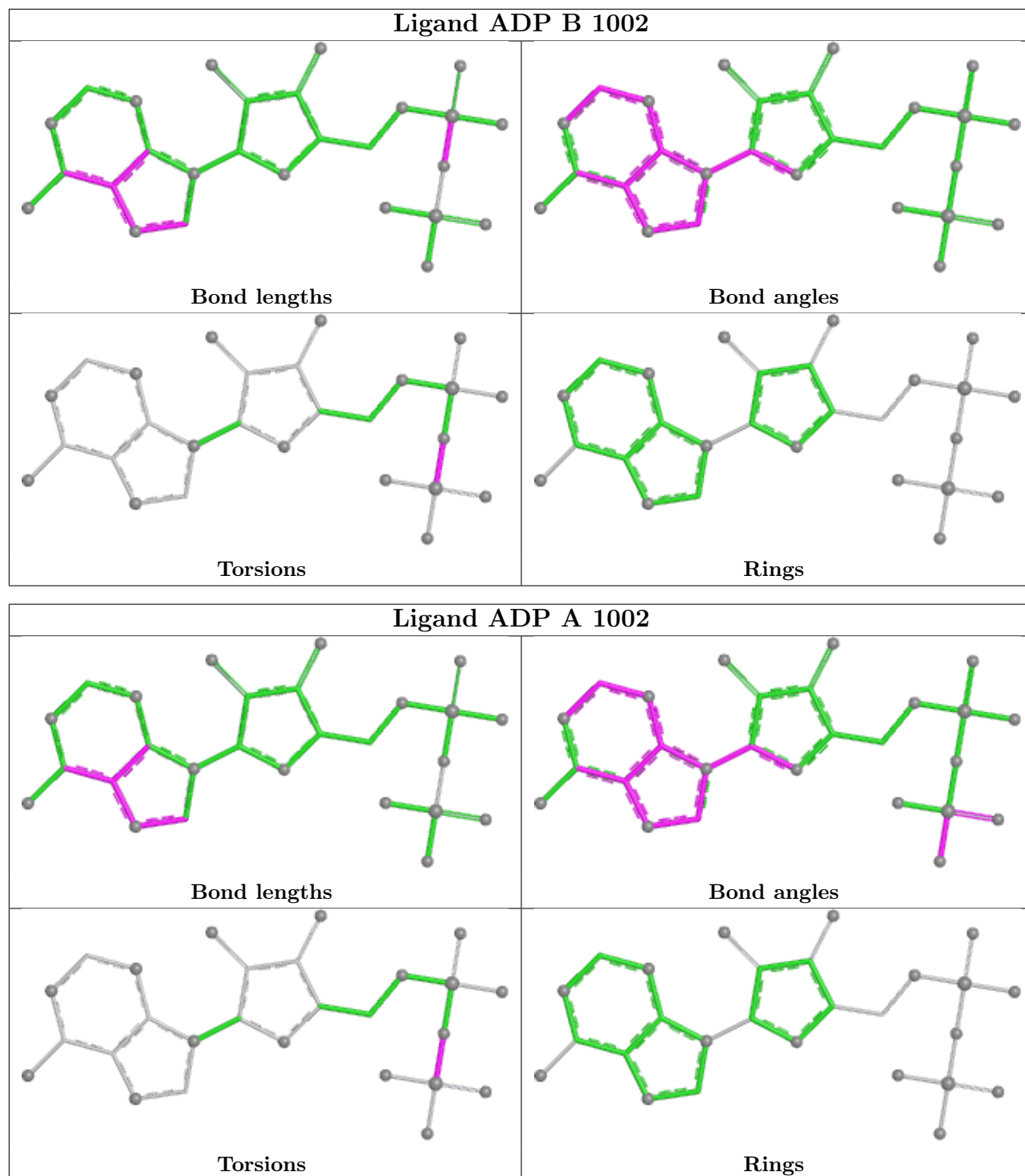
Mol	Chain	Res	Type	Atoms
3	A	1002	ADP	PA-O3A-PB-O2B
3	A	1002	ADP	PA-O3A-PB-O3B
3	B	1002	ADP	PA-O3A-PB-O2B
3	B	1002	ADP	PA-O3A-PB-O3B

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

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

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	531/531 (100%)	0.28	30 (5%) 30 32	32, 56, 85, 114	0
1	B	530/531 (99%)	0.41	29 (5%) 30 33	35, 63, 102, 117	0
All	All	1061/1062 (99%)	0.35	59 (5%) 30 32	32, 59, 96, 117	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	374	PHE	5.4
1	A	260	HIS	4.9
1	A	375	GLY	4.6
1	A	370	ALA	4.3
1	B	63	SER	4.1
1	A	371	THR	4.0
1	B	373	ALA	4.0
1	B	375	GLY	3.9
1	B	227	TRP	3.9
1	B	229	HIS	3.9
1	A	373	ALA	3.8
1	B	374	PHE	3.7
1	A	229	HIS	3.6
1	A	378	ILE	3.5
1	A	484	PHE	3.5
1	A	227	TRP	3.4
1	B	526	LYS	3.4
1	B	351	GLU	3.2
1	B	527	LEU	3.2
1	B	163	ALA	3.1
1	B	580	ASN	3.1
1	B	482	ILE	3.0
1	B	536	THR	3.0
1	B	298	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	322	PHE	2.9
1	A	377	GLY	2.8
1	B	531	GLY	2.8
1	A	322	PHE	2.8
1	A	376	MET	2.8
1	B	346	ALA	2.7
1	A	379	ASP	2.7
1	A	261	VAL	2.7
1	A	483	SER	2.6
1	A	372	VAL	2.6
1	B	167	LYS	2.6
1	A	226	GLN	2.5
1	A	346	ALA	2.5
1	A	271	VAL	2.5
1	B	378	ILE	2.4
1	A	569	TYR	2.4
1	A	273	LYS	2.3
1	A	558	ASP	2.3
1	B	64	SER	2.3
1	A	230	ASP	2.3
1	A	164	SER	2.3
1	B	535	PRO	2.2
1	B	486	ARG	2.2
1	B	507	ASN	2.2
1	A	348	MET	2.1
1	B	230	ASP	2.1
1	A	244	GLN	2.1
1	A	224	CYS	2.1
1	B	376	MET	2.1
1	B	506	LEU	2.1
1	B	532	LEU	2.1
1	B	512	PRO	2.1
1	B	569	TYR	2.1
1	A	468	PRO	2.0
1	A	481	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

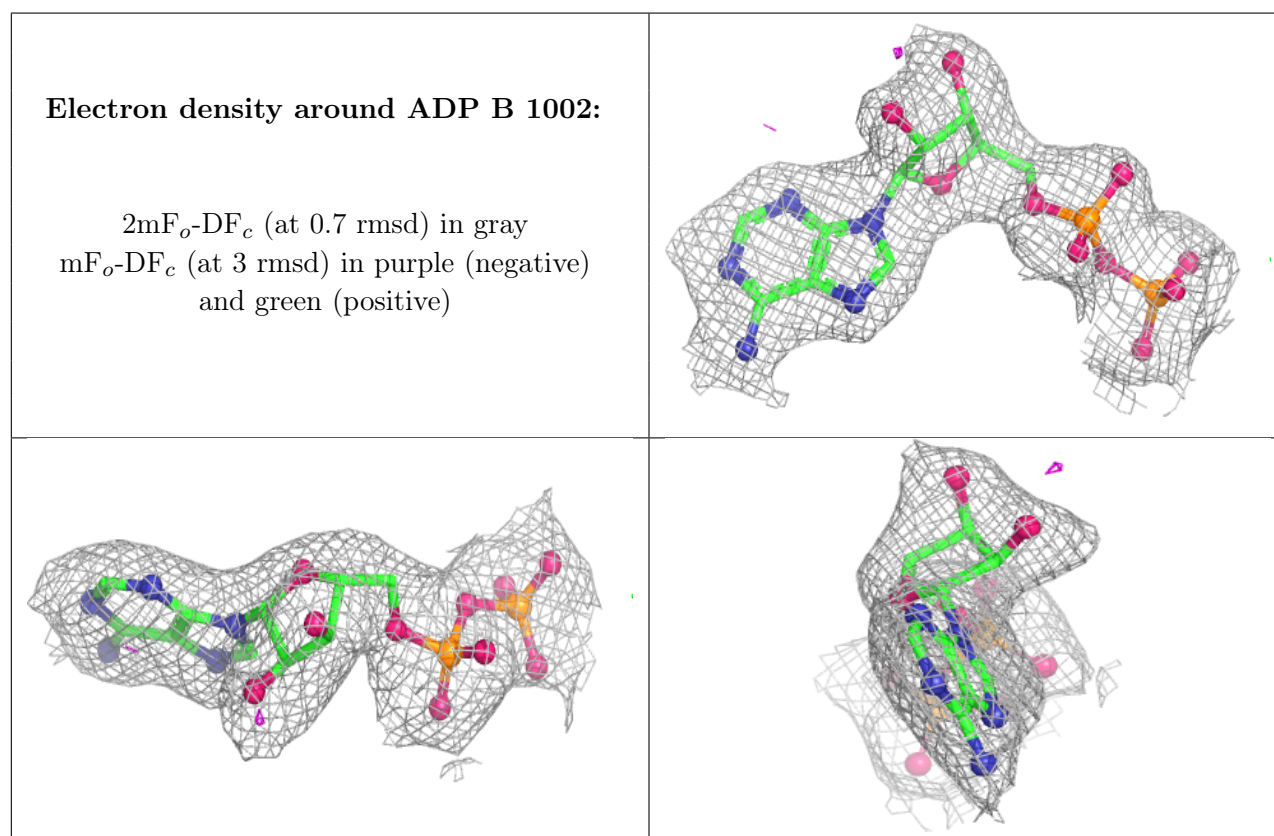
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

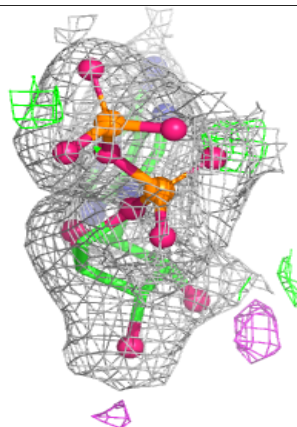
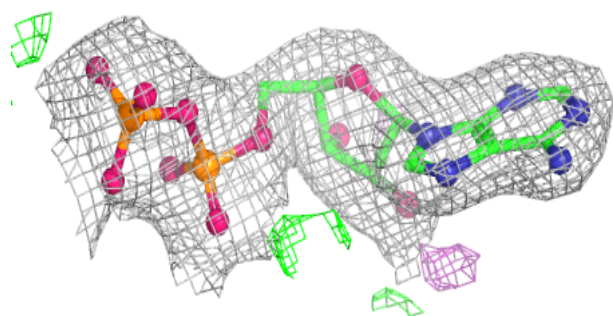
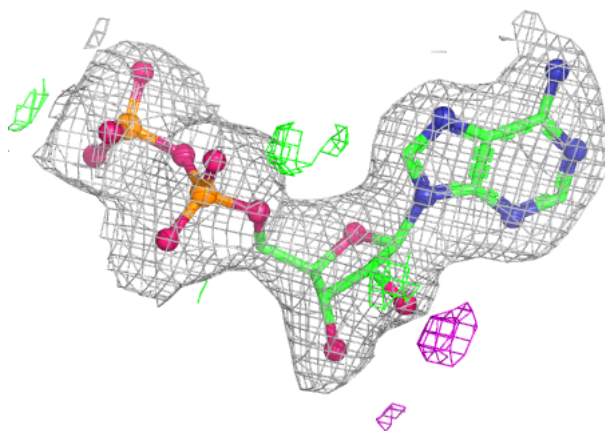
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	B	1003	1/1	0.95	0.10	60,60,60,60	0
3	ADP	B	1002	27/27	0.97	0.06	38,46,53,65	0
4	MG	A	1003	1/1	0.97	0.13	44,44,44,44	0
3	ADP	A	1002	27/27	0.97	0.07	37,46,54,59	0
2	ZN	B	1001	1/1	0.98	0.05	41,41,41,41	0
2	ZN	A	1001	1/1	1.00	0.05	44,44,44,44	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.



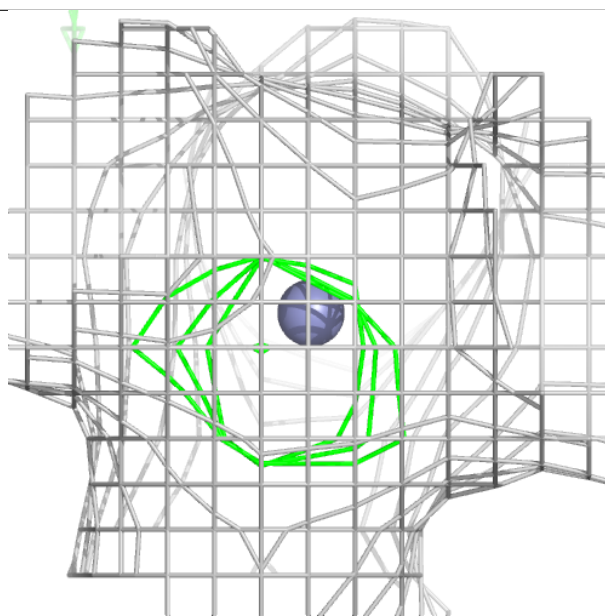
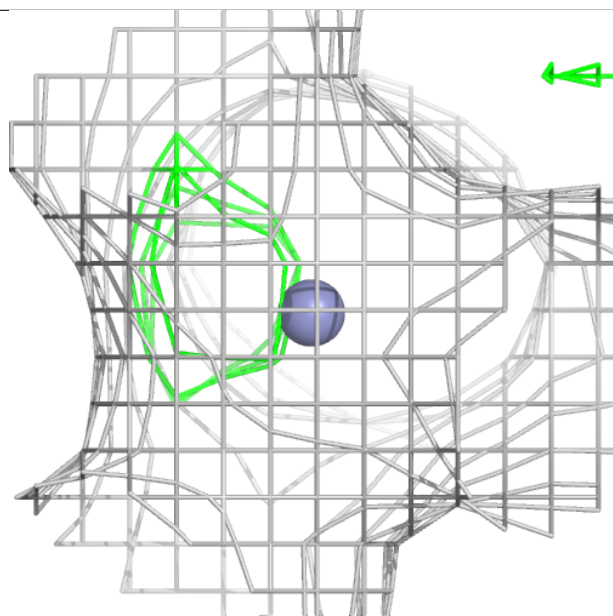
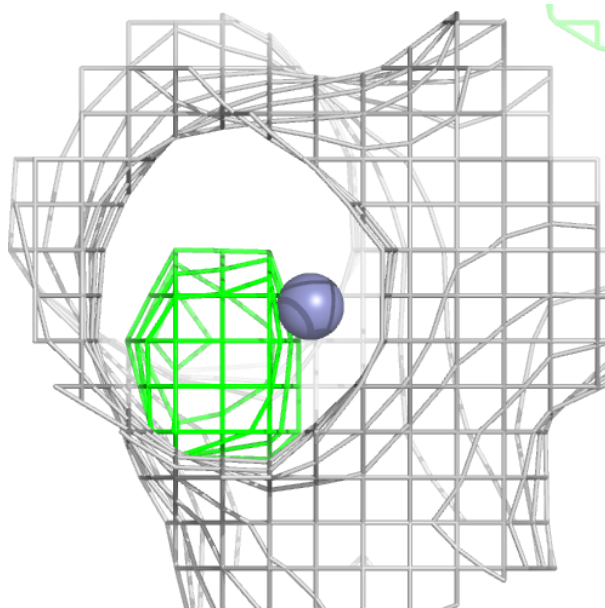
Electron density around ADP A 1002:

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



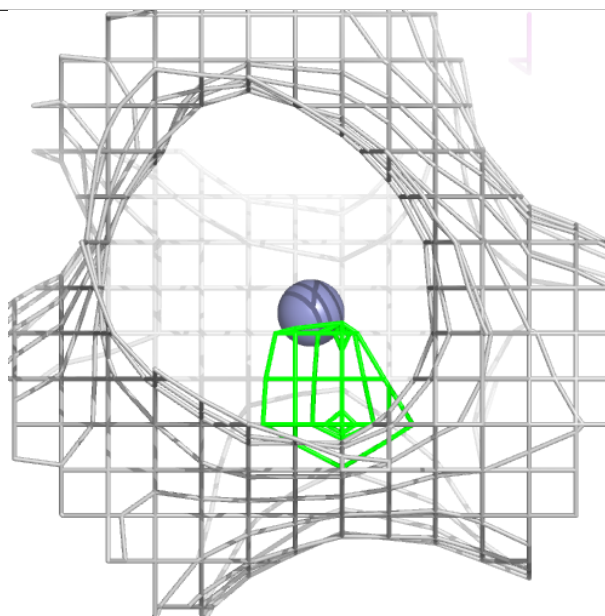
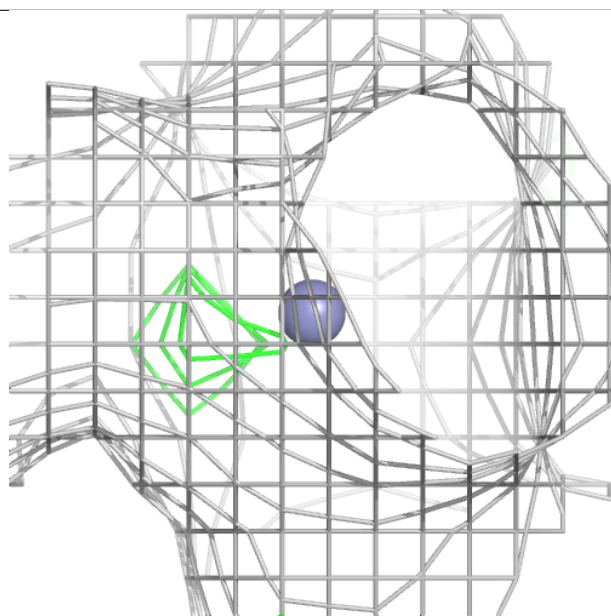
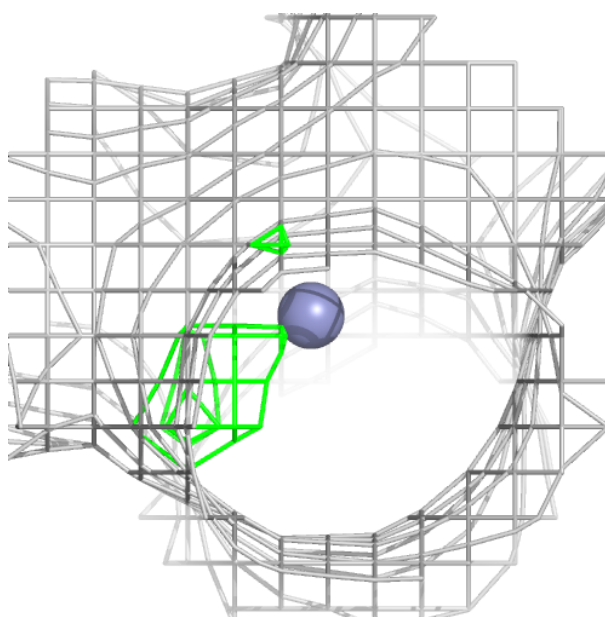
Electron density around ZN B 1001:

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



Electron density around ZN A 1001:

$2mF_o-DF_c$ (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.