



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:09 PM UTC

PDB ID : 9ZOK / pdb_00009zok
Title : Crystal structure of Phosphoribosylaminoimidazole carboxylase from Burkholderia xenovorans (ADP complex)
Authors : Seattle Structural Genomics Center for Infectious Disease; Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2025-12-15
Resolution : 2.09 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

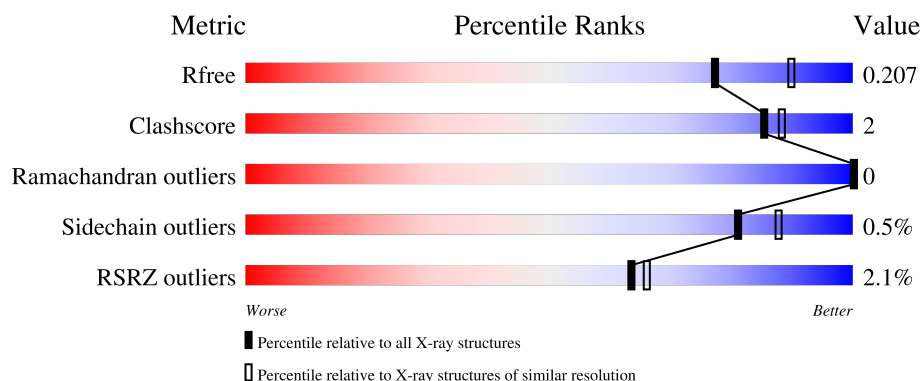
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	397	94%
1	B	397	92%
1	C	397	92%
1	D	397	92%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	GOL	B	403	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 12425 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 N5-carboxyaminoimidazole ribonucleotide synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	7	0
			2943	1848	533	544	18			
1	B	385	Total	C	N	O	S	0	7	0
			2907	1829	527	534	17			
1	C	388	Total	C	N	O	S	0	5	0
			2883	1812	515	536	20			
1	D	386	Total	C	N	O	S	0	5	0
			2844	1794	505	525	20			

There are 32 discrepancies between the modelled and reference sequences:

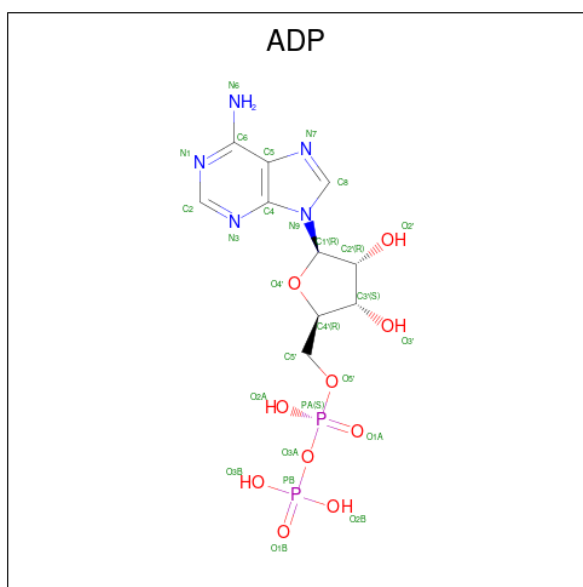
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q13UJ9
A	2	ALA	-	expression tag	UNP Q13UJ9
A	3	HIS	-	expression tag	UNP Q13UJ9
A	4	HIS	-	expression tag	UNP Q13UJ9
A	5	HIS	-	expression tag	UNP Q13UJ9
A	6	HIS	-	expression tag	UNP Q13UJ9
A	7	HIS	-	expression tag	UNP Q13UJ9
A	8	HIS	-	expression tag	UNP Q13UJ9
B	1	MET	-	expression tag	UNP Q13UJ9
B	2	ALA	-	expression tag	UNP Q13UJ9
B	3	HIS	-	expression tag	UNP Q13UJ9
B	4	HIS	-	expression tag	UNP Q13UJ9
B	5	HIS	-	expression tag	UNP Q13UJ9
B	6	HIS	-	expression tag	UNP Q13UJ9
B	7	HIS	-	expression tag	UNP Q13UJ9
B	8	HIS	-	expression tag	UNP Q13UJ9
C	1	MET	-	expression tag	UNP Q13UJ9
C	2	ALA	-	expression tag	UNP Q13UJ9
C	3	HIS	-	expression tag	UNP Q13UJ9
C	4	HIS	-	expression tag	UNP Q13UJ9
C	5	HIS	-	expression tag	UNP Q13UJ9

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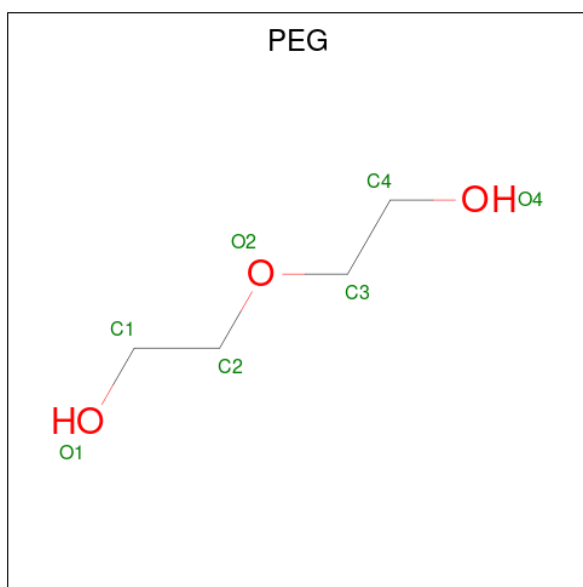
Chain	Residue	Modelled	Actual	Comment	Reference
C	6	HIS	-	expression tag	UNP Q13UJ9
C	7	HIS	-	expression tag	UNP Q13UJ9
C	8	HIS	-	expression tag	UNP Q13UJ9
D	1	MET	-	expression tag	UNP Q13UJ9
D	2	ALA	-	expression tag	UNP Q13UJ9
D	3	HIS	-	expression tag	UNP Q13UJ9
D	4	HIS	-	expression tag	UNP Q13UJ9
D	5	HIS	-	expression tag	UNP Q13UJ9
D	6	HIS	-	expression tag	UNP Q13UJ9
D	7	HIS	-	expression tag	UNP Q13UJ9
D	8	HIS	-	expression tag	UNP Q13UJ9

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).

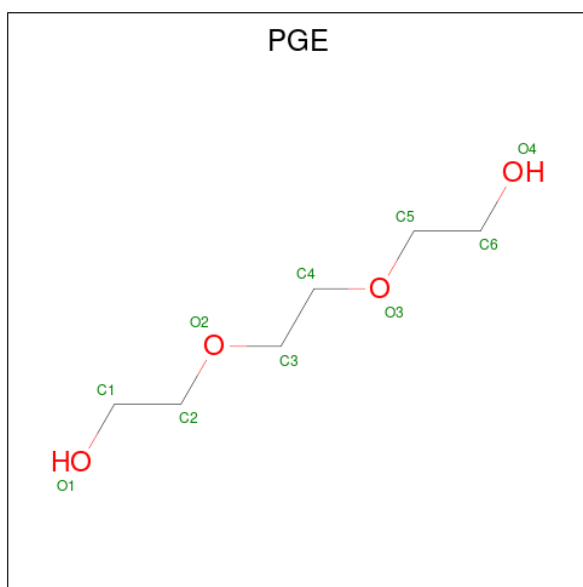


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		

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

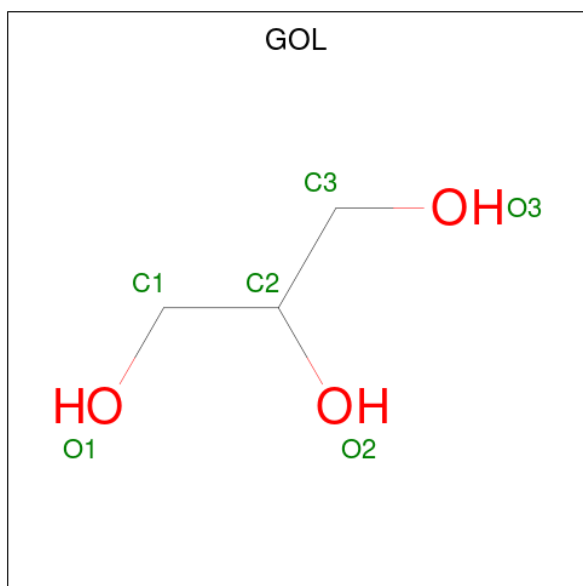
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		
4	B	1	Total	Mg	0	0
			1	1		
4	C	2	Total	Mg	0	0
			2	2		
4	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



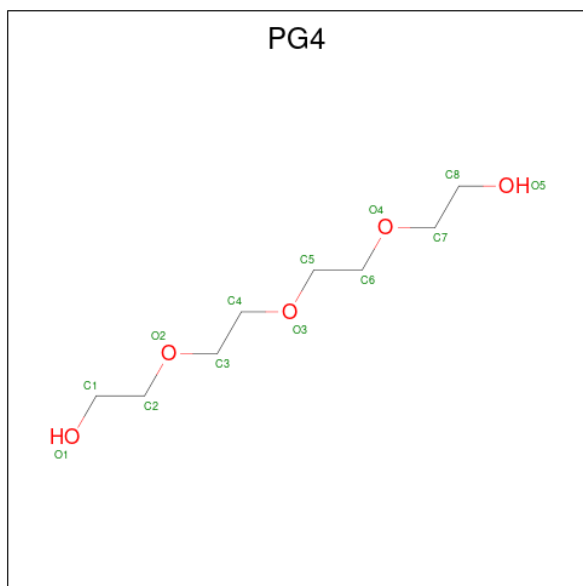
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	221	Total	O	0	0
			221	221		
8	B	207	Total	O	0	0
			207	207		
8	C	157	Total	O	0	0
			157	157		
8	D	101	Total	O	0	0
			101	101		

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: N5-carboxyaminoimidazole ribonucleotide synthase



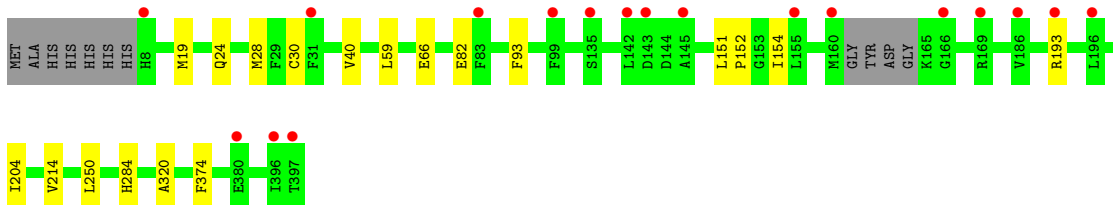
- Molecule 1: N5-carboxyaminoimidazole ribonucleotide synthase



- Molecule 1: N5-carboxyaminoimidazole ribonucleotide synthase



- Molecule 1: N5-carboxyaminoimidazole ribonucleotide synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.91Å 169.04Å 170.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.09 – 2.09 47.09 – 2.09	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.09-2.09) 100.0 (47.09-2.09)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.08Å)	Xtriage
Refinement program	PHENIX (2.0_5765: ???)	Depositor
R, R_{free}	0.168 , 0.204 0.173 , 0.207	Depositor DCC
R_{free} test set	5605 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	34.5	Xtriage
Anisotropy	0.673	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.013 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12425	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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: ADP, MG, PEG, GOL, PGE, PG4

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	0.45	0/3024	0.57	0/4114
1	B	0.44	0/2987	0.59	0/4066
1	C	0.41	0/2955	0.57	0/4023
1	D	0.32	0/2918	0.50	0/3981
All	All	0.41	0/11884	0.56	0/16184

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	2943	0	2945	10	0
1	B	2907	0	2915	16	0
1	C	2883	0	2850	14	0
1	D	2844	0	2797	12	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
2	C	27	0	12	0	0
2	D	27	0	12	0	0
3	A	7	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	2	0	0	0	0
4	B	1	0	0	0	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
5	B	10	0	14	1	0
6	B	6	0	8	5	0
6	C	6	0	8	2	0
6	D	6	0	8	1	0
7	B	13	0	18	0	0
8	A	221	0	0	1	0
8	B	207	0	0	1	0
8	C	157	0	0	0	0
8	D	101	0	0	0	0
All	All	12425	0	11621	52	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 2.

All (52) 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:195:PRO:HB2	1:B:270:GLU:HG3	1.75	0.68
1:B:285:ASN:H	6:B:403:GOL:H32	1.63	0.62
1:D:284:HIS:CD2	6:D:402:GOL:H31	2.34	0.62
1:B:19:MET:SD	1:B:82[A]:GLU:HG2	2.39	0.62
1:B:117:LYS:HE3	1:B:191:GLU:OE2	2.00	0.62
1:D:66:GLU:HG3	1:D:93:PHE:CZ	2.36	0.60
1:C:47:GLU:HB3	1:C:60:ARG:CZ	2.34	0.57
1:D:204:ILE:HG22	1:D:214:VAL:HG12	1.87	0.56
1:A:160:MET:HA	1:A:160:MET:HE2	1.88	0.54
1:B:118[A]:ARG:CZ	1:B:132:VAL:HG21	2.39	0.52
1:C:350:MET:HE1	1:C:385:ALA:O	2.10	0.52
1:C:47:GLU:HB3	1:C:60:ARG:NH1	2.25	0.52
1:B:117:LYS:HD3	8:B:505:HOH:O	2.10	0.51
1:B:284:HIS:CD2	6:B:403:GOL:H31	2.46	0.51
1:B:59:LEU:HD12	1:B:59:LEU:N	2.25	0.50
1:B:285:ASN:H	6:B:403:GOL:C3	2.24	0.50
1:D:30[A]:CYS:SG	1:D:40:VAL:HG11	2.52	0.50
1:D:151:LEU:HA	1:D:152:PRO:C	2.38	0.49
1:B:282:ARG:HH12	6:B:403:GOL:H31	1.77	0.49
1:A:350:MET:HE1	1:A:385:ALA:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:HIS:CE1	1:A:185:GLY:HA2	2.49	0.48
1:B:28:MET:HE1	1:B:288:HIS:CE1	2.49	0.47
1:A:198:PHE:CE2	1:A:268:ILE:HG13	2.49	0.47
1:C:19:MET:SD	1:C:82[A]:GLU:HG2	2.55	0.47
1:C:151:LEU:HA	1:C:152:PRO:C	2.40	0.46
1:D:19:MET:SD	1:D:82[A]:GLU:HG2	2.56	0.45
1:D:204:ILE:C	1:D:204:ILE:HD12	2.41	0.45
1:A:204:ILE:C	1:A:204:ILE:HD12	2.41	0.45
1:D:214:VAL:CG1	1:D:250:LEU:HD21	2.47	0.45
1:D:154:ILE:HD11	1:D:193:ARG:CZ	2.46	0.45
1:B:11:ILE:HD13	1:B:78:ALA:HB2	1.99	0.44
1:B:117:LYS:NZ	1:B:277:ASN:O	2.48	0.44
1:C:204:ILE:HD12	1:C:204:ILE:C	2.42	0.44
1:B:284:HIS:CD2	6:B:403:GOL:C3	3.02	0.42
1:C:47:GLU:HG3	1:C:48:ASN:CG	2.44	0.42
1:D:320:ALA:HA	1:D:374:PHE:O	2.20	0.42
1:C:30[A]:CYS:SG	1:C:40:VAL:HG11	2.59	0.42
1:C:200:VAL:HA	1:C:218:ALA:O	2.19	0.42
1:D:59:LEU:HD12	1:D:59:LEU:N	2.34	0.42
1:A:147:LEU:O	1:A:151:LEU:HG	2.20	0.42
1:C:284:HIS:CD2	6:C:402:GOL:H12	2.53	0.42
1:B:19:MET:HE1	1:B:80:SER:HB3	2.02	0.42
1:C:57:ARG:HG2	1:C:59:LEU:HD12	2.01	0.42
1:C:284:HIS:CD2	6:C:402:GOL:C1	3.03	0.42
1:C:360:LYS:O	1:C:362:GLU:N	2.53	0.41
1:D:24:GLN:C	1:D:28:MET:HE3	2.45	0.41
1:A:71[A]:GLU:HG3	1:A:74[A]:ARG:HH22	1.86	0.40
1:B:383:ARG:HE	5:B:402:PGE:H6	1.86	0.40
1:A:25:LEU:CD2	1:A:288:HIS:HE1	2.34	0.40
1:A:146:LYS:HD2	8:A:653:HOH:O	2.20	0.40
1:A:200:VAL:HA	1:A:218:ALA:O	2.21	0.40
1:C:320:ALA:HA	1:C:374:PHE:O	2.21	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	395/397 (100%)	392 (99%)	3 (1%)	0	100	100
1	B	388/397 (98%)	384 (99%)	4 (1%)	0	100	100
1	C	389/397 (98%)	384 (99%)	5 (1%)	0	100	100
1	D	387/397 (98%)	383 (99%)	4 (1%)	0	100	100
All	All	1559/1588 (98%)	1543 (99%)	16 (1%)	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	302/305 (99%)	300 (99%)	2 (1%)	76	83
1	B	299/305 (98%)	298 (100%)	1 (0%)	86	91
1	C	292/305 (96%)	289 (99%)	3 (1%)	68	76
1	D	285/305 (93%)	285 (100%)	0	100	100
All	All	1178/1220 (97%)	1172 (100%)	6 (0%)	81	88

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	GLU
1	A	160	MET
1	B	270	GLU
1	C	105	ARG
1	C	146	LYS
1	C	360	LYS

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

Mol	Chain	Res	Type
1	B	180	HIS
1	B	243	GLN
1	B	247	GLN
1	B	251	GLN
1	B	302	GLN
1	C	8	HIS
1	C	180	HIS
1	C	302	GLN
1	D	111	GLN
1	D	180	HIS
1	D	243	GLN
1	D	247	GLN
1	D	251	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 6 are monoatomic - leaving 10 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	$\# Z > 2$
6	GOL	D	402	-	5,5,5	0.32	0	5,5,5	0.35	0
6	GOL	C	402	-	5,5,5	0.31	0	5,5,5	0.92	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PEG	A	402	-	6,6,6	0.27	0	5,5,5	0.77	0
7	PG4	B	404	-	12,12,12	0.28	0	11,11,11	0.33	0
5	PGE	B	402	-	9,9,9	0.33	0	8,8,8	0.72	0
2	ADP	B	401	-	28,29,29	1.30	5 (17%)	43,45,45	1.85	11 (25%)
6	GOL	B	403	-	5,5,5	0.18	0	5,5,5	0.73	0
2	ADP	C	401	4	28,29,29	1.45	5 (17%)	43,45,45	1.88	10 (23%)
2	ADP	D	401	-	28,29,29	1.41	5 (17%)	43,45,45	1.78	9 (20%)
2	ADP	A	401	4	28,29,29	1.43	5 (17%)	43,45,45	1.78	10 (23%)

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
6	GOL	D	402	-	-	2/4/4/4	-
6	GOL	C	402	-	-	4/4/4/4	-
3	PEG	A	402	-	-	2/4/4/4	-
7	PG4	B	404	-	-	4/10/10/10	-
5	PGE	B	402	-	-	5/7/7/7	-
2	ADP	B	401	-	-	2/16/32/32	0/3/3/3
6	GOL	B	403	-	-	2/4/4/4	-
2	ADP	C	401	4	-	0/16/32/32	0/3/3/3
2	ADP	D	401	-	-	1/16/32/32	0/3/3/3
2	ADP	A	401	4	-	0/16/32/32	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	ADP	C5-C4	4.78	1.47	1.39
2	C	401	ADP	C5-C4	4.68	1.47	1.39
2	B	401	ADP	C5-C4	4.53	1.47	1.39
2	A	401	ADP	C5-C4	4.16	1.46	1.39
2	A	401	ADP	PA-O3A	3.28	1.63	1.59
2	C	401	ADP	C5-N7	-3.12	1.33	1.39
2	D	401	ADP	C5-N7	-2.62	1.34	1.39
2	A	401	ADP	C5-N7	-2.57	1.34	1.39
2	D	401	ADP	C5-C6	2.34	1.47	1.41
2	B	401	ADP	C5-C6	2.33	1.47	1.41
2	B	401	ADP	C4-N9	-2.32	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ADP	C5-N7	-2.26	1.35	1.39
2	D	401	ADP	C8-N7	2.22	1.35	1.31
2	C	401	ADP	PA-O3A	2.20	1.61	1.59
2	C	401	ADP	C5-C6	2.17	1.47	1.41
2	A	401	ADP	C5-C6	2.15	1.47	1.41
2	A	401	ADP	C4-N9	-2.10	1.33	1.37
2	B	401	ADP	C8-N7	2.06	1.35	1.31
2	C	401	ADP	C4-N9	-2.05	1.33	1.37
2	D	401	ADP	C4-N9	-2.03	1.33	1.37

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	ADP	C5-C4-N3	-5.51	119.13	126.72
2	D	401	ADP	C5-C4-N3	-5.26	119.47	126.72
2	A	401	ADP	C5-C4-N3	-5.18	119.58	126.72
2	B	401	ADP	C5-C4-N3	-5.10	119.69	126.72
2	C	401	ADP	N3-C4-N9	4.89	135.49	127.17
2	D	401	ADP	N3-C4-N9	4.64	135.06	127.17
2	B	401	ADP	N3-C4-N9	4.55	134.90	127.17
2	A	401	ADP	N3-C4-N9	4.39	134.64	127.17
2	C	401	ADP	N3-C2-N1	-3.88	122.70	128.58
2	B	401	ADP	C2-N3-C4	3.74	120.98	111.83
2	B	401	ADP	N3-C2-N1	-3.71	122.97	128.58
2	D	401	ADP	N3-C2-N1	-3.70	122.98	128.58
2	A	401	ADP	N3-C2-N1	-3.68	123.01	128.58
2	C	401	ADP	C2-N3-C4	3.58	120.58	111.83
2	B	401	ADP	C4-N9-C8	3.58	109.50	105.74
2	D	401	ADP	C2-N3-C4	3.35	120.02	111.83
2	A	401	ADP	C2-N3-C4	3.26	119.80	111.83
2	D	401	ADP	C4-N9-C8	3.08	108.98	105.74
2	A	401	ADP	C4-N9-C8	2.91	108.80	105.74
2	C	401	ADP	C4-N9-C8	2.85	108.73	105.74
2	D	401	ADP	C2-N1-C6	2.83	123.38	118.73
2	A	401	ADP	C4-C5-N7	-2.80	107.38	110.58
2	C	401	ADP	C2'-C1'-N9	-2.79	106.38	113.30
2	D	401	ADP	C4-C5-N7	-2.75	107.43	110.58
2	A	401	ADP	C2-N1-C6	2.68	123.13	118.73
2	B	401	ADP	C4-C5-N7	-2.64	107.56	110.58
2	B	401	ADP	N9-C8-N7	-2.63	110.21	113.94
2	C	401	ADP	C4-C5-N7	-2.58	107.63	110.58
2	B	401	ADP	C5-N7-C8	2.45	107.30	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ADP	C5-N7-C8	2.41	107.24	103.45
2	C	401	ADP	C5-N7-C8	2.35	107.14	103.45
2	D	401	ADP	C5-N7-C8	2.27	107.01	103.45
2	A	401	ADP	N9-C8-N7	-2.21	110.80	113.94
2	B	401	ADP	C2-N1-C6	2.18	122.31	118.73
2	C	401	ADP	C2-N1-C6	2.18	122.31	118.73
2	A	401	ADP	C2'-C1'-N9	-2.17	107.91	113.30
2	B	401	ADP	C6-C5-N7	2.13	136.20	132.09
2	C	401	ADP	O3B-PB-O2B	2.11	115.70	107.80
2	B	401	ADP	C2'-C1'-N9	-2.07	108.16	113.30
2	D	401	ADP	O2A-PA-O1A	2.02	121.83	112.44

There are no chirality outliers.

All (22) torsion outliers are listed below:

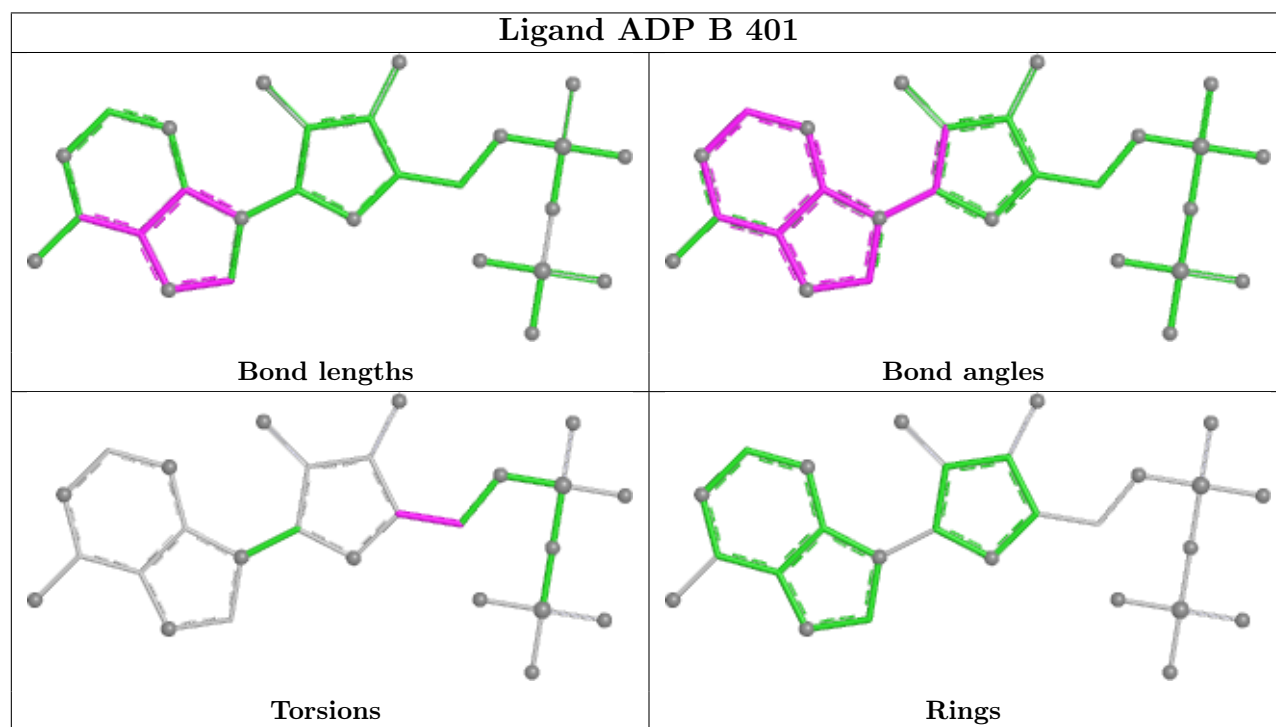
Mol	Chain	Res	Type	Atoms
6	B	403	GOL	O1-C1-C2-C3
5	B	402	PGE	O2-C3-C4-O3
3	A	402	PEG	O2-C3-C4-O4
6	C	402	GOL	O1-C1-C2-C3
6	C	402	GOL	C1-C2-C3-O3
3	A	402	PEG	O1-C1-C2-O2
6	B	403	GOL	O1-C1-C2-O2
6	C	402	GOL	O2-C2-C3-O3
5	B	402	PGE	O1-C1-C2-O2
7	B	404	PG4	O1-C1-C2-O2
6	D	402	GOL	O1-C1-C2-C3
7	B	404	PG4	O2-C3-C4-O3
7	B	404	PG4	O4-C7-C8-O5
7	B	404	PG4	C6-C5-O3-C4
5	B	402	PGE	C4-C3-O2-C2
6	C	402	GOL	O1-C1-C2-O2
6	D	402	GOL	O1-C1-C2-O2
2	B	401	ADP	O4'-C4'-C5'-O5'
5	B	402	PGE	C6-C5-O3-C4
2	B	401	ADP	C3'-C4'-C5'-O5'
5	B	402	PGE	O3-C5-C6-O4
2	D	401	ADP	PB-O3A-PA-O1A

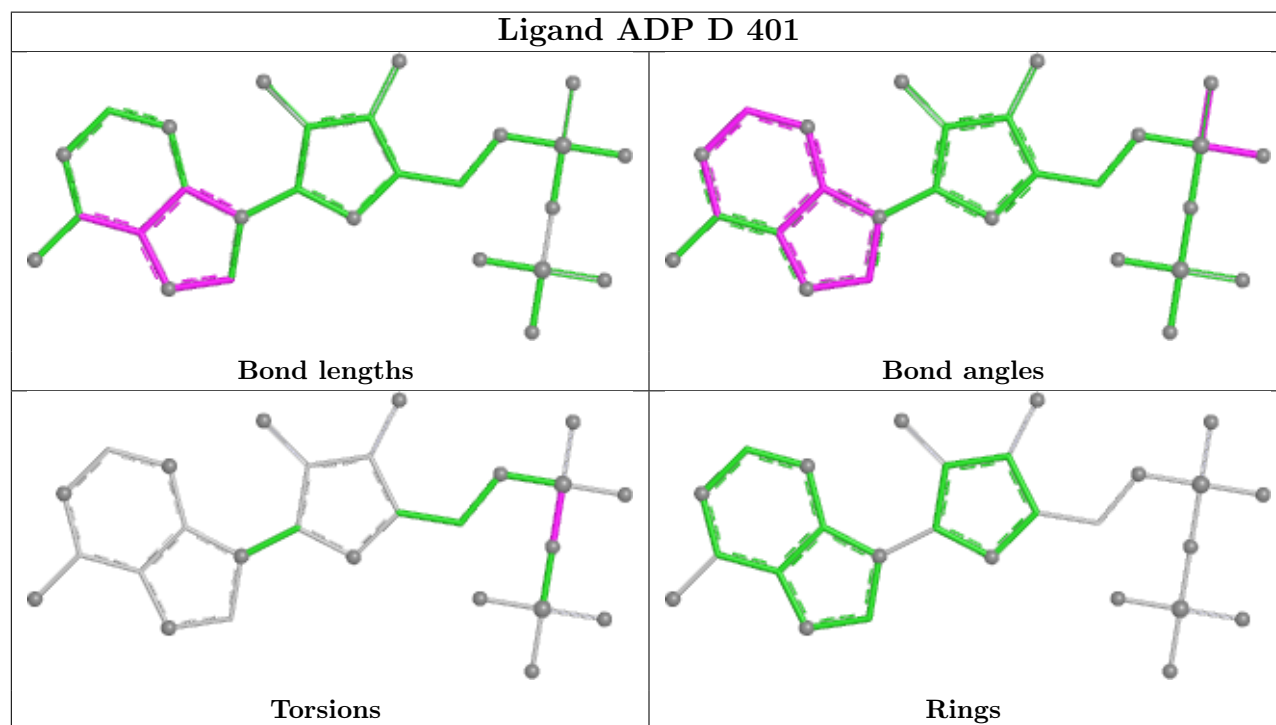
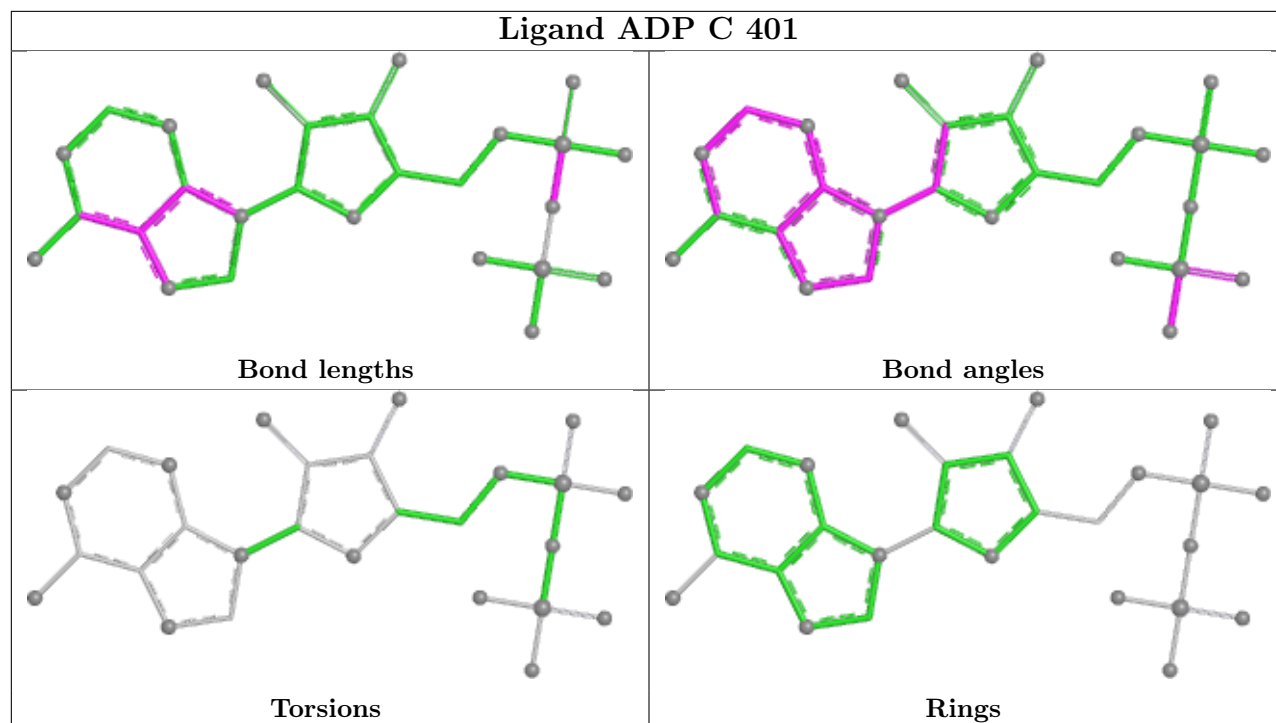
There are no ring outliers.

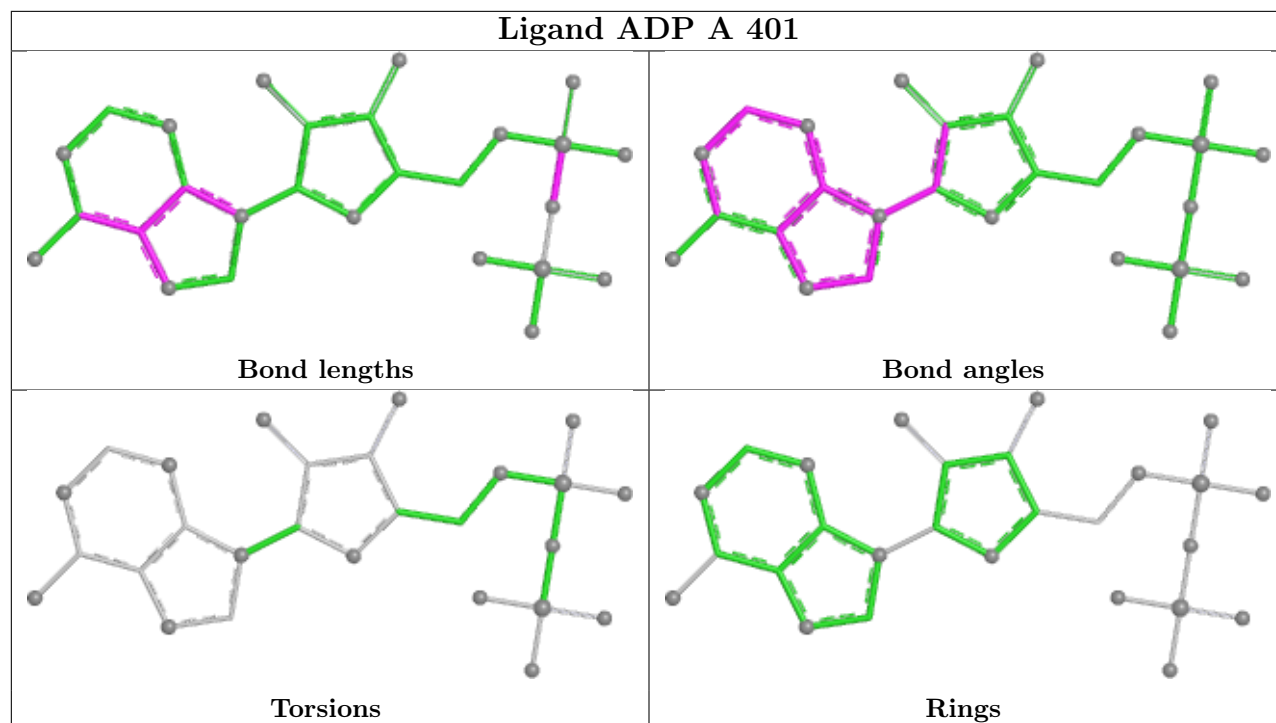
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	402	GOL	1	0
6	C	402	GOL	2	0
5	B	402	PGE	1	0
6	B	403	GOL	5	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	390/397 (98%)	-0.09	5 (1%) 75 77	21, 36, 54, 90	7 (1%)
1	B	385/397 (96%)	-0.12	4 (1%) 79 81	19, 36, 57, 92	7 (1%)
1	C	388/397 (97%)	0.08	6 (1%) 72 74	21, 41, 75, 121	5 (1%)
1	D	386/397 (97%)	0.50	18 (4%) 36 38	22, 53, 108, 139	5 (1%)
All	All	1549/1588 (97%)	0.09	33 (2%) 63 66	19, 40, 85, 139	24 (1%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	99[A]	PHE	5.0
1	D	8	HIS	3.7
1	B	160	MET	3.5
1	C	396	ILE	3.3
1	B	8	HIS	3.3
1	C	338	PRO	3.2
1	D	397	THR	3.2
1	B	99[A]	PHE	3.2
1	C	333	ASP	3.1
1	D	186	VAL	3.1
1	C	339	ALA	3.0
1	C	362	GLU	2.9
1	A	397	THR	2.7
1	B	397	THR	2.6
1	D	142	LEU	2.6
1	D	143	ASP	2.5
1	D	193	ARG	2.5
1	D	380	GLU	2.4
1	A	99	PHE	2.4
1	D	169	ARG	2.4
1	D	160	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	83	PHE	2.3
1	A	8	HIS	2.3
1	D	135	SER	2.3
1	D	31	PHE	2.2
1	C	359	GLY	2.2
1	D	145	ALA	2.2
1	D	155	LEU	2.2
1	D	196	LEU	2.1
1	D	396	ILE	2.1
1	D	166	GLY	2.1
1	A	46	ASP	2.0
1	A	243	GLN	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 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.

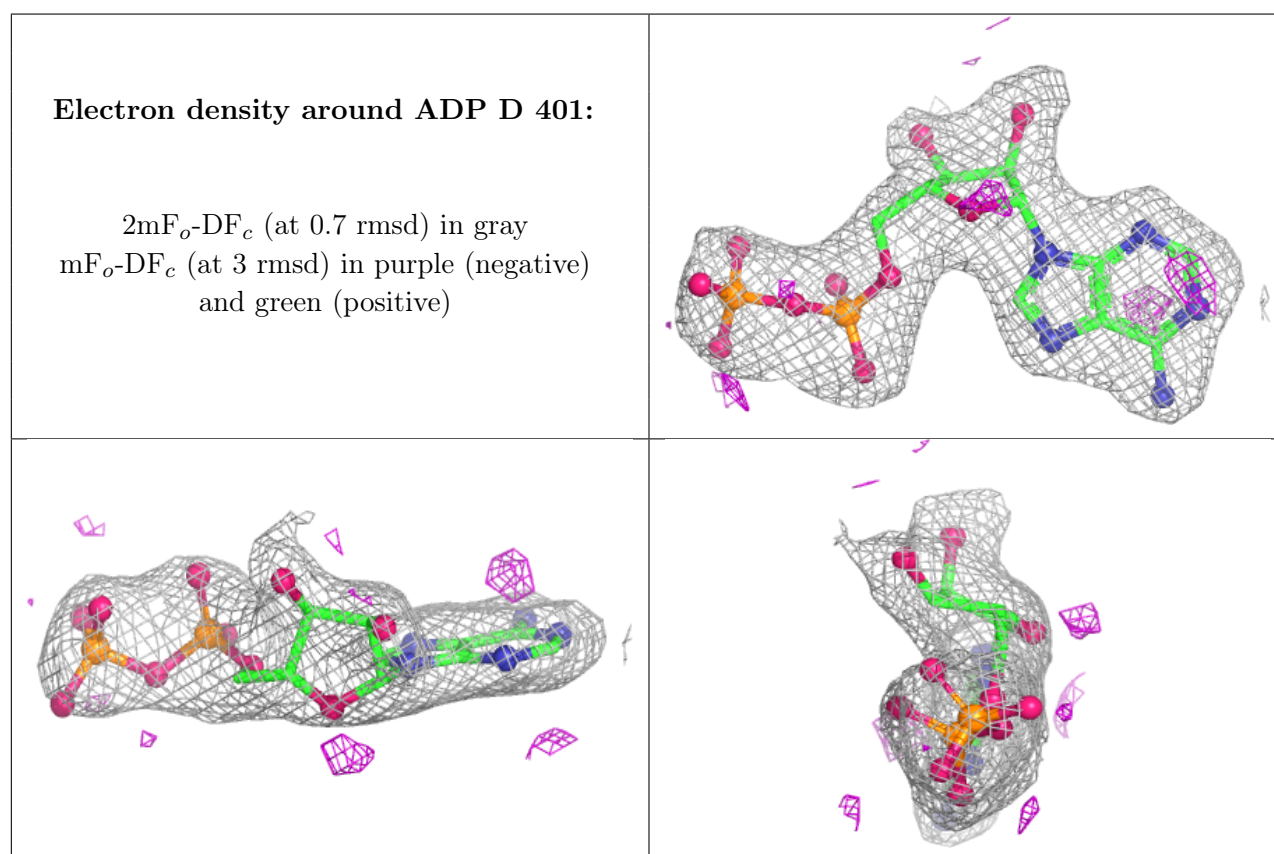
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	GOL	C	402	6/6	0.75	0.17	45,50,63,66	0
5	PGE	B	402	10/10	0.85	0.14	50,60,67,69	0
6	GOL	D	402	6/6	0.85	0.12	52,61,65,68	0
2	ADP	D	401	27/27	0.86	0.10	66,72,101,110	0
6	GOL	B	403	6/6	0.86	0.12	41,48,53,57	0
2	ADP	B	401	27/27	0.90	0.09	41,49,84,100	0
3	PEG	A	402	7/7	0.90	0.13	46,49,56,61	0
7	PG4	B	404	13/13	0.90	0.12	48,53,61,63	0
4	MG	D	403	1/1	0.93	0.07	62,62,62,62	0
4	MG	A	403	1/1	0.93	0.07	45,45,45,45	0
2	ADP	A	401	27/27	0.95	0.07	30,38,44,46	0

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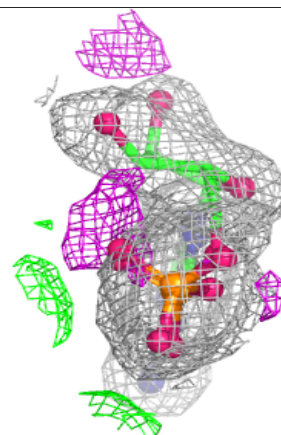
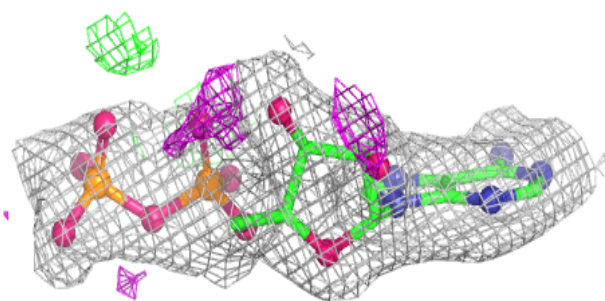
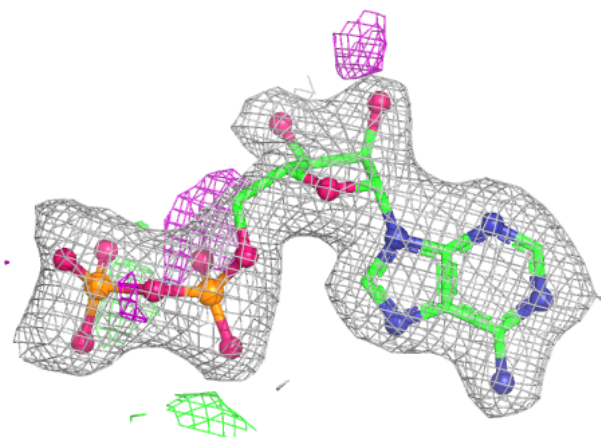
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	B	405	1/1	0.96	0.05	38,38,38,38	0
4	MG	C	403	1/1	0.97	0.12	40,40,40,40	0
4	MG	C	404	1/1	0.97	0.06	44,44,44,44	0
2	ADP	C	401	27/27	0.98	0.05	33,38,44,46	0
4	MG	A	404	1/1	0.99	0.09	36,36,36,36	0

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.

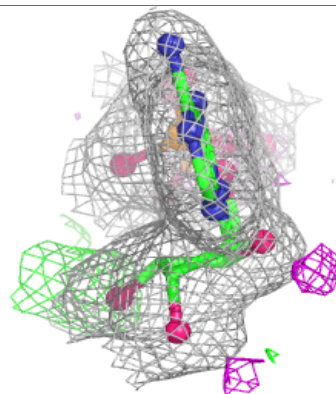
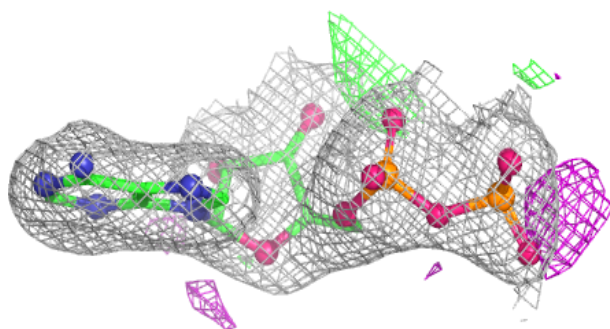
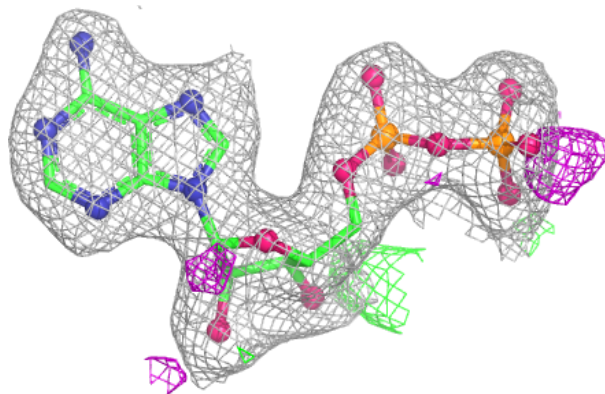


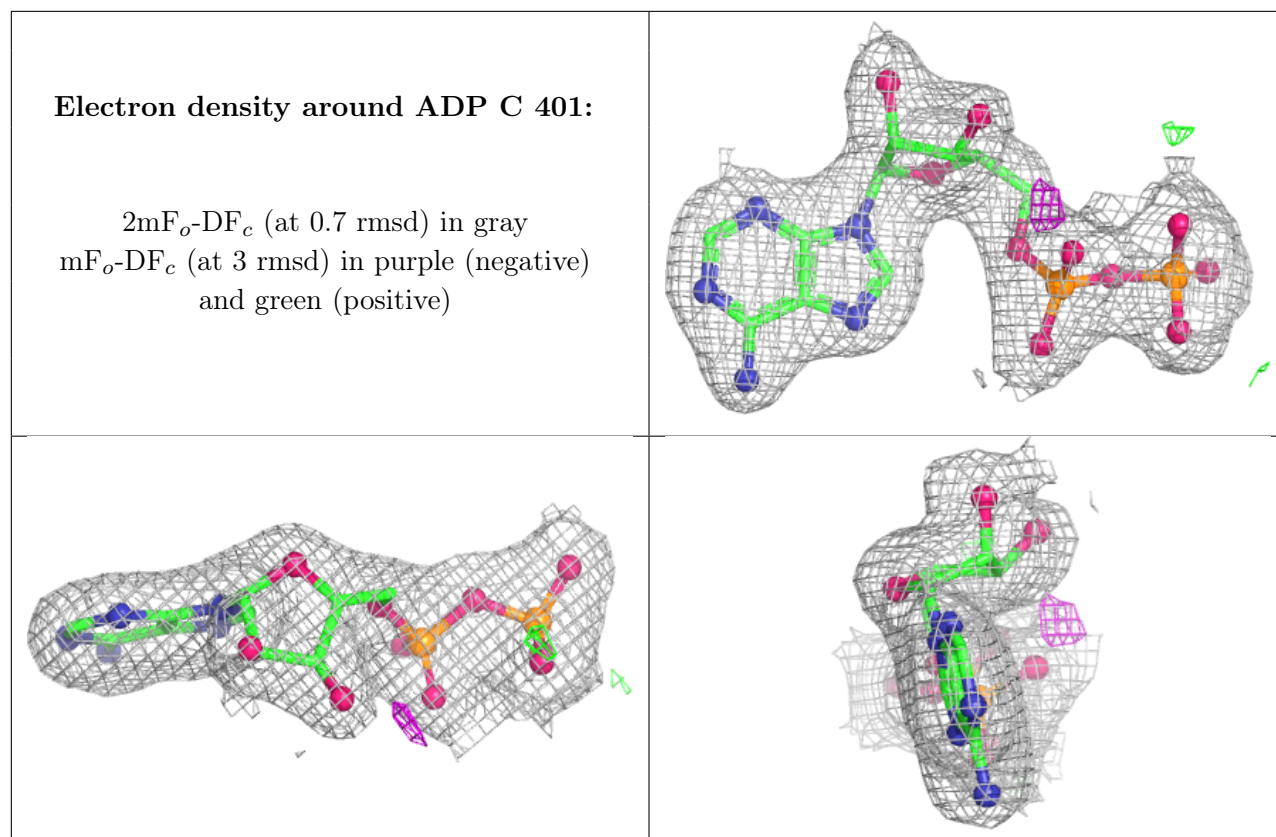
Electron density around ADP B 401:

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

**Electron density around ADP A 401:**

$2mF_o-DF_c$ (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.