



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:42 PM UTC

PDB ID : 5TDM / pdb_00005tdm
Title : TEV Cleaved Human ATP Citrate Lyase Bound to 4R-Hydroxycitrate and ADP
Authors : Hu, J.; Fraser, M.E.
Deposited on : 2016-09-19
Resolution : 2.10 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

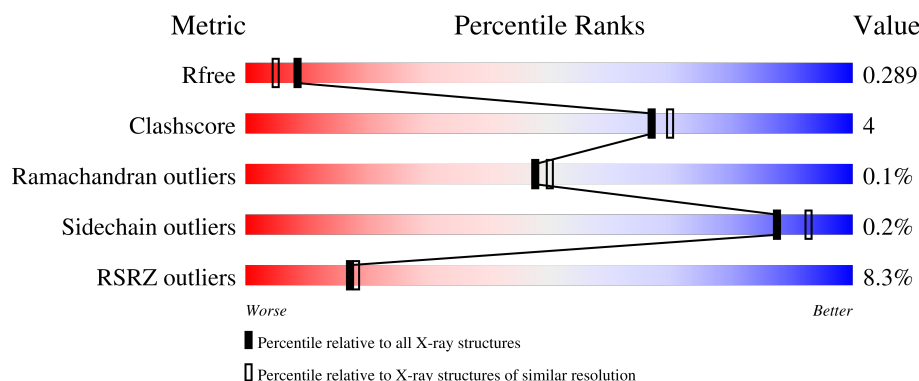
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	431	
2	B	324	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 11959 atoms, of which 5881 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-citrate synthase.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	427	Total	C	H	N	O	S		0	0	0
			6696	2156	3353	560	618	9				

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	426	GLU	-	SEE REMARK 999	UNP P53396
A	427	ASN	-	SEE REMARK 999	UNP P53396
A	428	LEU	-	SEE REMARK 999	UNP P53396
A	429	TYR	-	SEE REMARK 999	UNP P53396
A	430	PHE	-	SEE REMARK 999	UNP P53396
A	431	GLN	-	SEE REMARK 999	UNP P53396

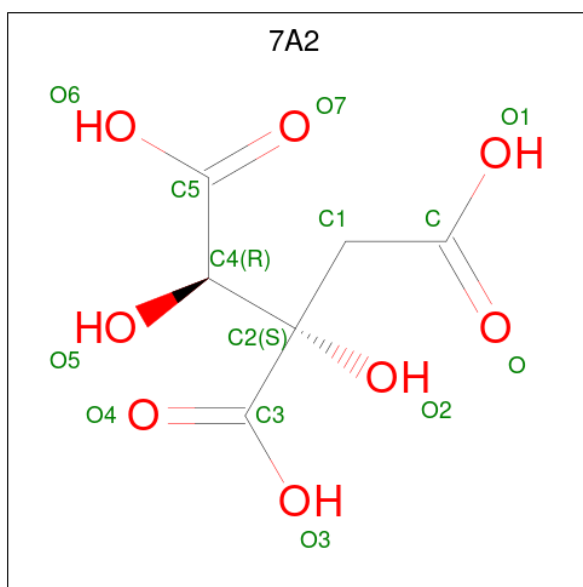
- Molecule 2 is a protein called ATP-citrate synthase.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
2	B	323	Total	C	H	N	O	P	S	0	1	0
			4954	1563	2480	424	466	1	20			

There is a discrepancy between the modelled and reference sequences:

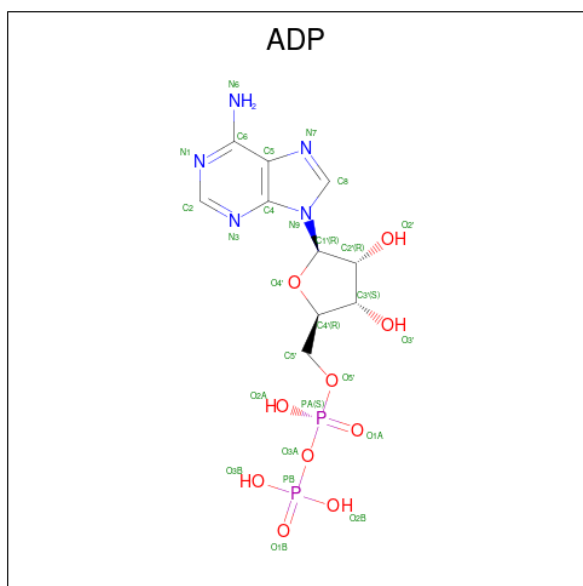
Chain	Residue	Modelled	Actual	Comment	Reference
B	487	SER	-	SEE REMARK 999	UNP P53396

- Molecule 3 is 3-C-carboxy-2-deoxy-L-threo-pentonic acid (CCD ID: 7A2) (formula: C₆H₈O₈).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			19	6	5	8		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

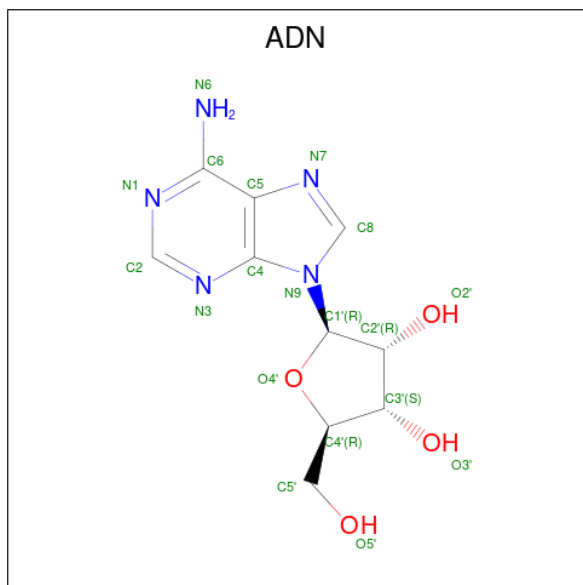


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
4	A	1	Total	C	H	N	O	P	0	0
			42	10	15	5	10	2		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

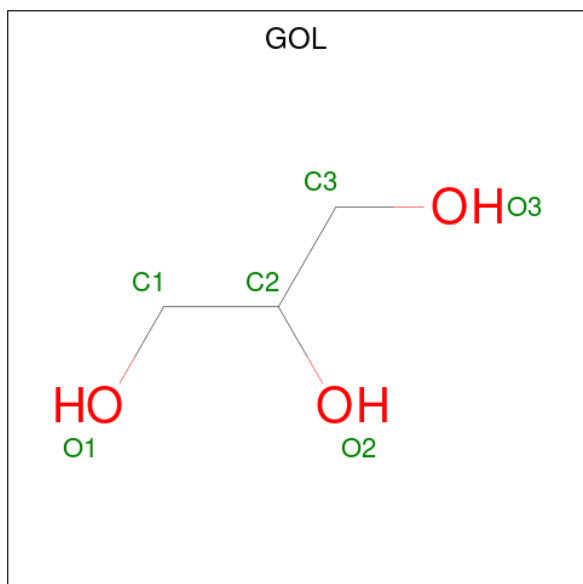
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		

- Molecule 6 is ADENOSINE (CCD ID: ADN) (formula: $C_{10}H_{13}N_5O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
6	A	1	Total	C	H	N	O	0	0
			32	10	13	5	4		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total 13	C 3	H 7	O 3	0	0
7	B	1	Total 14	C 3	H 8	O 3	0	0

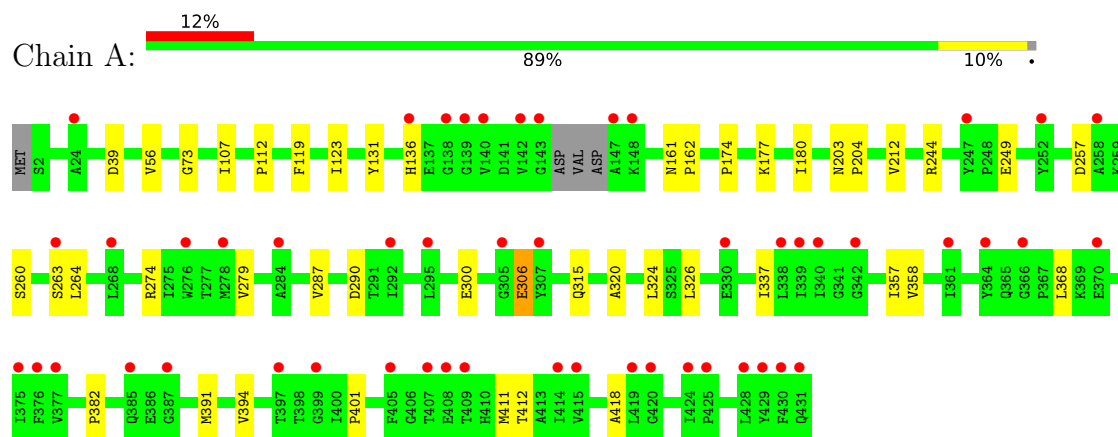
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	113	Total 113	O 113	0	0
8	B	74	Total 74	O 74	0	0

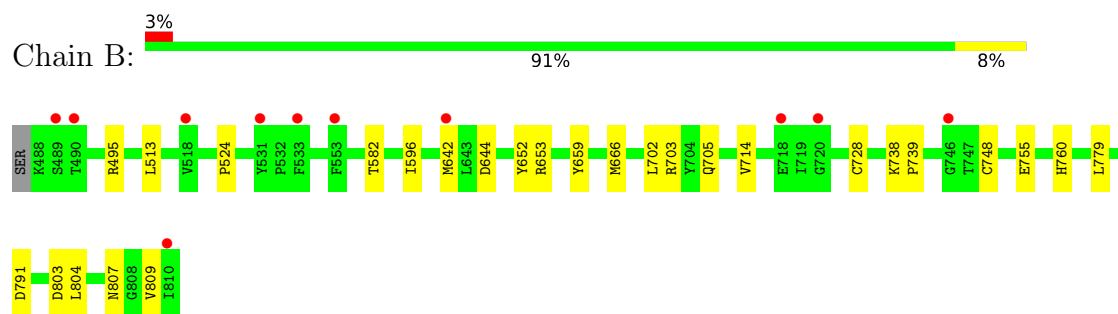
3 Residue-property plots [i](#)

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-citrate synthase



• Molecule 2: ATP-citrate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.38Å 84.81Å 193.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.77 – 2.10 63.77 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.8 (63.77-2.10) 75.4 (63.77-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.10_2152: ???)	Depositor
R, R_{free}	0.216 , 0.286 0.219 , 0.289	Depositor DCC
R_{free} test set	2678 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.809	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11959	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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: GOL, MG, ADN, 7A2, ADP, NEP

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	0.62	0/3421	0.67	0/4639
2	B	0.61	0/2511	0.70	0/3388
All	All	0.62	0/5932	0.69	0/8027

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3343	3353	3352	29	0
2	B	2474	2480	2476	18	0
3	A	14	5	0	1	0
4	A	27	15	12	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	19	13	13	1	0
7	A	6	7	7	1	0
7	B	6	8	8	0	0
8	A	113	0	0	2	0
8	B	74	0	0	1	1
All	All	6078	5881	5868	43	1

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 (43) 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:244:ARG:NH1	2:B:755:GLU:O	2.26	0.69
1:A:290:ASP:HA	2:B:748:CYS:HB3	1.90	0.54
1:A:358:VAL:HG13	1:A:394:VAL:HB	1.91	0.52
2:B:728:CYS:SG	2:B:779:LEU:HD23	2.49	0.52
2:B:804:LEU:HD22	2:B:809:VAL:HG11	1.92	0.51
1:A:382:PRO:HA	2:B:642:MET:HE2	1.92	0.50
2:B:803:ASP:O	2:B:807:ASN:ND2	2.45	0.50
1:A:287:VAL:HG21	2:B:666[B]:MET:SD	2.52	0.49
1:A:260:SER:HB3	1:A:315:GLN:HE22	1.79	0.48
1:A:357:ILE:HG21	1:A:391:MET:HE1	1.95	0.48
1:A:412:THR:OG1	2:B:791:ASP:OD1	2.31	0.47
2:B:513:LEU:HD22	2:B:524:PRO:HB3	1.95	0.47
1:A:123:ILE:HA	1:A:131:TYR:O	2.15	0.47
1:A:257:ASP:OD1	1:A:263:SER:HA	2.15	0.46
2:B:644:ASP:N	2:B:644:ASP:OD1	2.49	0.46
1:A:56:VAL:HA	1:A:73:GLY:O	2.15	0.45
1:A:180:ILE:HD11	1:A:212:VAL:HG23	1.99	0.45
1:A:357:ILE:CG2	1:A:391:MET:HE1	2.46	0.45
1:A:324:LEU:HD13	1:A:368:LEU:HD11	1.98	0.45
2:B:582:THR:HG21	2:B:596:ILE:HD11	1.97	0.45
1:A:320:ALA:O	1:A:324:LEU:HG	2.17	0.45
1:A:112:PRO:HA	8:A:1064:HOH:O	2.15	0.45
2:B:739:PRO:HG2	2:B:804:LEU:HD11	1.98	0.45
1:A:249:GLU:CD	1:A:326:LEU:HD21	2.41	0.44
2:B:705:GLN:HA	2:B:738:LYS:HD2	1.99	0.44
1:A:274:ARG:HG2	1:A:300:GLU:OE2	2.18	0.44
1:A:161:ASN:O	1:A:162:PRO:C	2.61	0.43
6:A:904:ADN:C5	7:A:905:GOL:H2	2.48	0.43
1:A:274:ARG:NH1	8:A:1017:HOH:O	2.50	0.43
2:B:652:TYR:CE1	2:B:653:ARG:HD2	2.53	0.43
2:B:659:TYR:HA	2:B:714:VAL:O	2.19	0.43
1:A:337:ILE:HG21	1:A:418:ALA:HB1	2.00	0.42
2:B:702:LEU:O	2:B:703:ARG:C	2.62	0.42
1:A:279:VAL:O	1:A:306:GLU:HA	2.20	0.42
1:A:39:ASP:OD1	1:A:39:ASP:C	2.63	0.42
3:A:901:7A2:O7	3:A:901:7A2:C1	2.67	0.42
1:A:257:ASP:HB2	1:A:264:LEU:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ASN:HA	1:A:204:PRO:HA	1.81	0.41
2:B:495:ARG:HG2	8:B:1049:HOH:O	2.19	0.41
1:A:119:PHE:CE1	1:A:136:HIS:CD2	3.08	0.41
1:A:174:PRO:HB2	1:A:177:LYS:HD3	2.03	0.40
1:A:56:VAL:O	1:A:107:ILE:HA	2.22	0.40
1:A:411:MET:HE1	2:B:666[A]:MET:HE1	2.04	0.40

All (1) 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)
8:B:1018:HOH:O	8:B:1055:HOH:O[3_655]	2.10	0.10

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	423/431 (98%)	407 (96%)	15 (4%)	1 (0%)	43	44
2	B	321/324 (99%)	315 (98%)	6 (2%)	0	100	100
All	All	744/755 (98%)	722 (97%)	21 (3%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	401	PRO

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	354/358 (99%)	353 (100%)	1 (0%)	86	91
2	B	263/263 (100%)	263 (100%)	0	100	100
All	All	617/621 (99%)	616 (100%)	1 (0%)	87	93

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	306	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	371	HIS
1	A	431	GLN
2	B	563	HIS
2	B	638	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
2	NEP	B	760	5,2	11,14,15	3.05	3 (27%)	8,20,22	1.06	1 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NEP	B	760	5,2	-	2/5/12/14	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	760	NEP	P-O3P	8.37	1.58	1.46
2	B	760	NEP	P-O2P	-3.95	1.46	1.56
2	B	760	NEP	P-O1P	3.70	1.66	1.56

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	760	NEP	O2P-P-O3P	-2.17	107.62	112.95

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	760	NEP	CA-CB-CG-CD2
2	B	760	NEP	CA-CB-CG-ND1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	7A2	A	901	-	12,13,13	0.95	0	13,19,19	2.63	5 (38%)
4	ADP	A	902	5	28,29,29	0.35	0	43,45,45	0.53	0
6	ADN	A	904	7	21,21,21	0.21	0	31,31,31	0.31	0
7	GOL	A	905	6	5,5,5	0.42	0	5,5,5	0.83	0
7	GOL	B	902	-	5,5,5	0.44	0	5,5,5	0.30	0

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	7A2	A	901	-	-	6/21/21/21	-
4	ADP	A	902	5	-	1/16/32/32	0/3/3/3
6	ADN	A	904	7	-	0/6/22/22	0/3/3/3
7	GOL	A	905	6	-	2/4/4/4	-
7	GOL	B	902	-	-	0/4/4/4	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	7A2	C1-C2-C3	-6.18	101.87	110.56
3	A	901	7A2	O3-C3-C2	4.71	121.94	114.06
3	A	901	7A2	O4-C3-C2	-4.20	114.49	122.32
3	A	901	7A2	O6-C5-C4	2.59	123.04	112.80
3	A	901	7A2	O7-C5-C4	-2.23	113.36	122.45

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	901	7A2	O5-C4-C5-O6
7	A	905	GOL	O1-C1-C2-O2
7	A	905	GOL	O1-C1-C2-C3
3	A	901	7A2	O5-C4-C5-O7
4	A	902	ADP	C5'-O5'-PA-O1A

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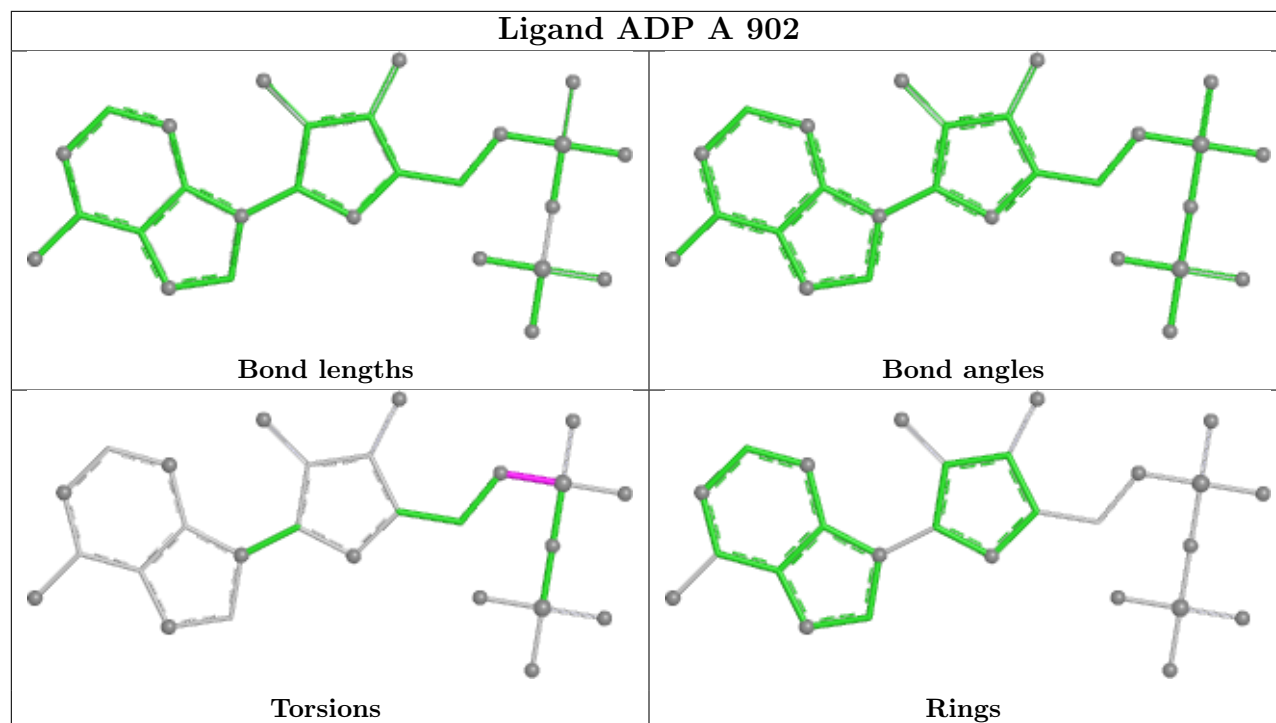
Mol	Chain	Res	Type	Atoms
3	A	901	7A2	O-C-C1-C2
3	A	901	7A2	O1-C-C1-C2
3	A	901	7A2	C2-C4-C5-O6
3	A	901	7A2	C2-C4-C5-O7

There are no ring outliers.

3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	7A2	1	0
6	A	904	ADN	1	0
7	A	905	GOL	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 [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	427/431 (99%)	0.82	51 (11%) 9 9	32, 64, 104, 132	0
2	B	322/324 (99%)	0.61	11 (3%) 48 50	34, 59, 86, 110	1 (0%)
All	All	749/755 (99%)	0.73	62 (8%) 17 18	32, 61, 99, 132	1 (0%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	430	PHE	5.7
1	A	143	GLY	4.4
1	A	142	VAL	4.4
2	B	490	THR	4.2
1	A	305	GLY	3.9
1	A	429	TYR	3.4
1	A	338	LEU	3.4
1	A	385	GLN	3.4
1	A	340	ILE	3.3
1	A	140	VAL	3.1
1	A	295	LEU	3.1
1	A	431	GLN	3.0
1	A	147	ALA	3.0
1	A	428	LEU	3.0
2	B	718	GLU	2.9
1	A	415	VAL	2.9
1	A	330	GLU	2.8
1	A	414	ILE	2.7
1	A	276	TRP	2.7
1	A	407	THR	2.7
1	A	268	LEU	2.6
1	A	24	ALA	2.6
1	A	263	SER	2.5
2	B	533	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	284	ALA	2.5
1	A	364	TYR	2.4
1	A	425	PRO	2.4
1	A	258	ALA	2.4
1	A	375	ILE	2.4
1	A	247	TYR	2.4
1	A	408	GLU	2.4
1	A	138	GLY	2.4
1	A	339	ILE	2.4
2	B	720	GLY	2.3
1	A	377	VAL	2.3
1	A	366	GLY	2.3
2	B	642	MET	2.3
1	A	376	PHE	2.2
2	B	810	ILE	2.2
1	A	342	GLY	2.2
1	A	370	GLU	2.2
1	A	136	HIS	2.2
1	A	405	PHE	2.2
1	A	387	GLY	2.2
2	B	746	GLY	2.2
1	A	307	TYR	2.2
2	B	489	SER	2.2
2	B	553	PHE	2.2
1	A	420	GLY	2.2
1	A	292	ILE	2.1
1	A	252	TYR	2.1
1	A	397	THR	2.1
1	A	424	ILE	2.1
2	B	518	VAL	2.1
1	A	148	LYS	2.1
1	A	278	MET	2.1
1	A	361	ILE	2.1
1	A	419	LEU	2.0
1	A	409	THR	2.0
2	B	531	TYR	2.0
1	A	139	GLY	2.0
1	A	399	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	NEP	B	760	14/15	0.94	0.10	41,65,89,94	0

6.3 Carbohydrates [i](#)

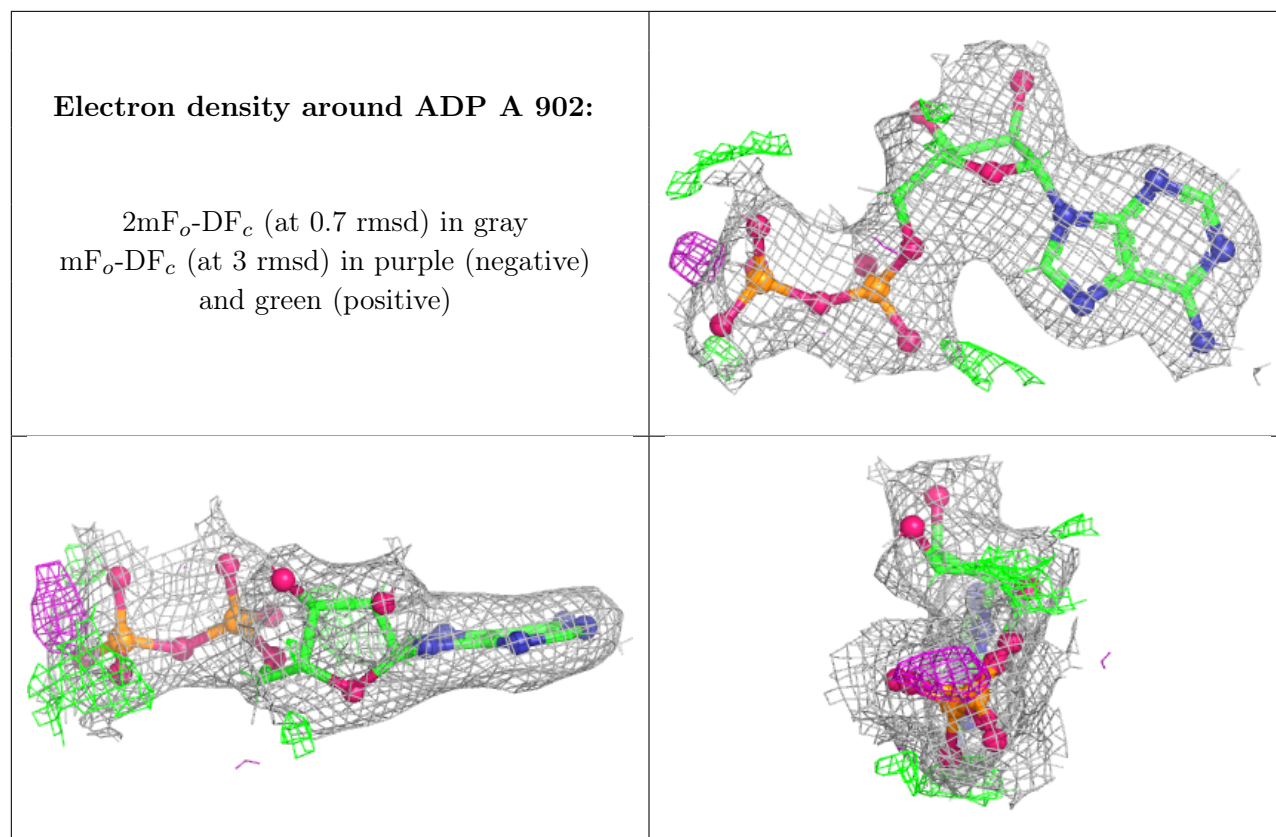
There are no oligosaccharides in this entry.

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	ADN	A	904	19/19	0.63	0.16	72,144,204,234	0
7	GOL	A	905	6/6	0.74	0.23	57,100,163,165	0
5	MG	B	901	1/1	0.81	0.11	63,63,63,63	0
7	GOL	B	902	6/6	0.85	0.18	58,102,128,133	0
3	7A2	A	901	14/14	0.92	0.13	44,84,115,129	0
4	ADP	A	902	27/27	0.93	0.12	27,72,113,123	0
5	MG	A	903	1/1	0.95	0.11	90,90,90,90	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.