



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 09:57 pm BST

PDB ID : 5TDM  
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](mailto: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.

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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

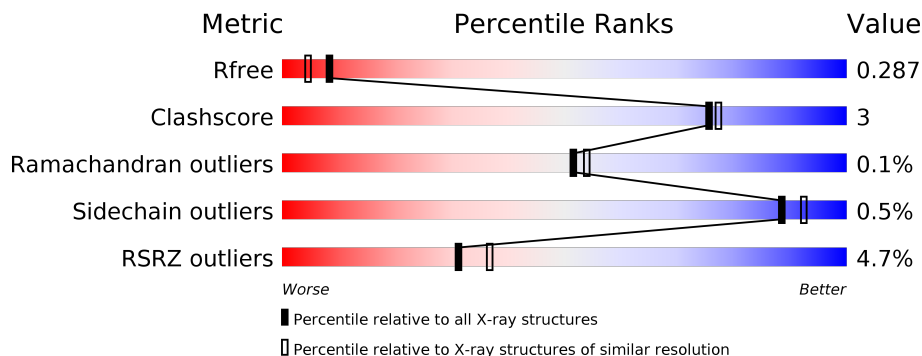
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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 [i](#)

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
			Total	C	H	N	O	S				
1	A	427	6696	2156	3353	560	618	9	0	0	0	

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

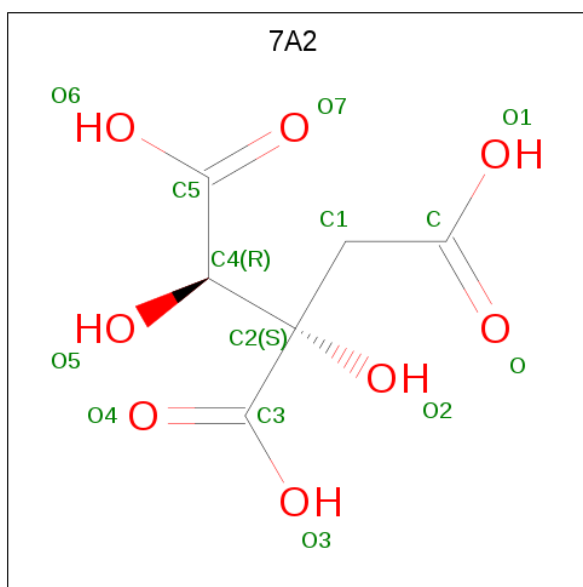
- Molecule 2 is a protein called ATP-citrate synthase.

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

There is a discrepancy between the modelled and reference sequences:

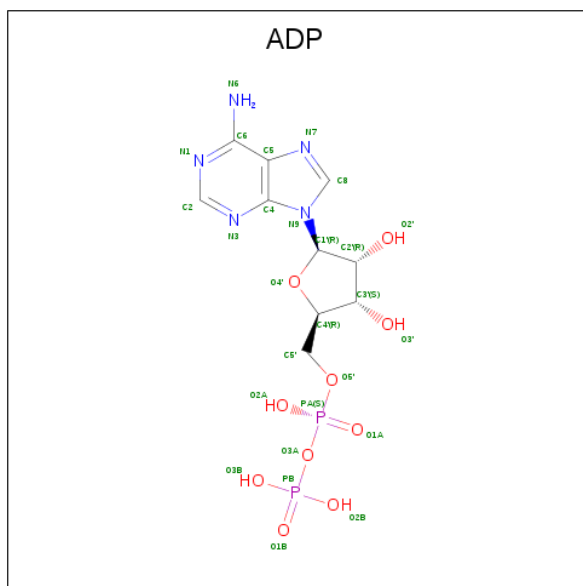
Chain	Residue	Modelled	Actual	Comment	Reference
B	487	SER	-	SEE REMARK 999	UNP P53396

- Molecule 3 is 3-C-carboxy-2-deoxy-L-threo-pentonic acid (three-letter code: 7A2) (formula: C<sub>6</sub>H<sub>8</sub>O<sub>8</sub>).



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

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

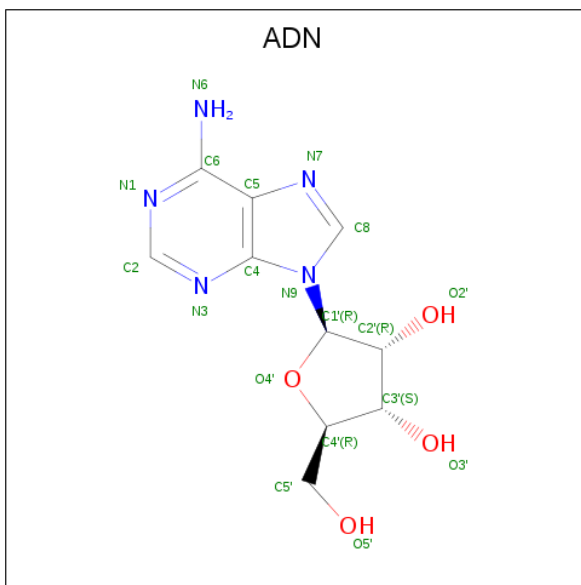


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
4	A	1	42	10	15	5	10	2	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

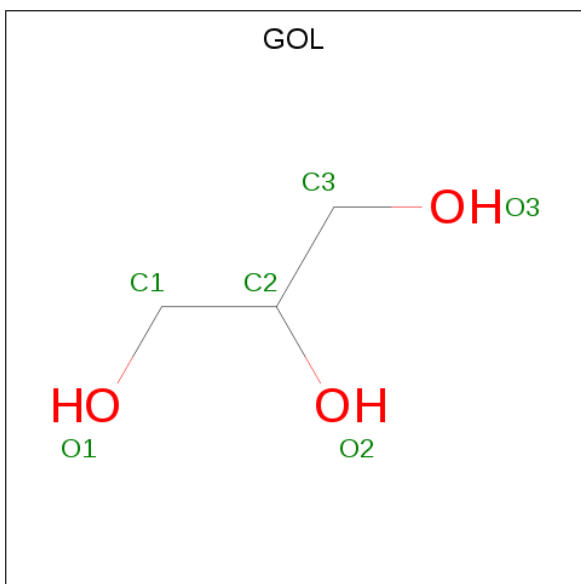
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Mg 1 1	0	0
5	A	1	Total Mg 1 1	0	0

- Molecule 6 is ADENOSINE (three-letter code: ADN) (formula: C<sub>10</sub>H<sub>13</sub>N<sub>5</sub>O<sub>4</sub>).



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

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

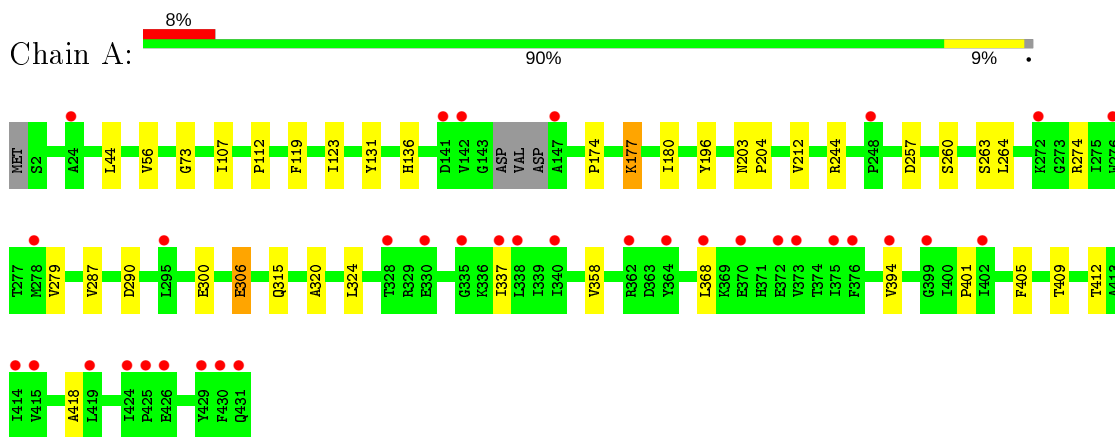
- Molecule 8 is water.

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

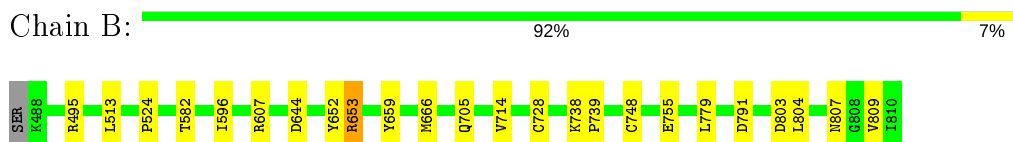
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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, $\alpha$ , $\beta$ , $\gamma$	55.38Å 84.81Å 193.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.77 – 2.10 63.83 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.8 (63.77-2.10) 75.4 (63.83-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.25 (at 2.10Å)	Xtrriage
Refinement program	PHENIX (1.10_2152: ???)	Depositor
R, $R_{free}$	0.216 , 0.286 0.217 , 0.287	Depositor DCC
$R_{free}$ test set	2678 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.0	Xtrriage
Anisotropy	0.809	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 46.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	11959	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	65.0	wwPDB-VP

Xtrriage'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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, MG, ADP, ADN, 7A2, 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	0.66	1/3421 (0.0%)	0.68	1/4639 (0.0%)
2	B	0.63	0/2511	0.71	1/3388 (0.0%)
All	All	0.65	1/5932 (0.0%)	0.70	2/8027 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	196	TYR	CD1-CE1	7.86	1.51	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	607	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	A	44	LEU	CB-CG-CD2	-5.43	101.77	111.00

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	3343	3353	3352	23	0
2	B	2474	2480	2476	15	0
3	A	14	5	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	36	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 3.

All (36) 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
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: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:OD1	2:B:644:ASP:N	2.49	0.46
1:A:180:ILE:HD11	1:A:212:VAL:HG23	1.99	0.45
1:A:56:VAL:HA	1:A:73:GLY:O	2.15	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
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:405:PHE:HD2	1:A:409:THR:HG1	1.66	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:279:VAL:O	1:A:306:GLU:HA	2.20	0.42
3:A:901:7A2:C1	3:A:901:7A2:O7	2.67	0.42
1:A:257:ASP:HB2	1:A:264:LEU:O	2.19	0.42
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

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%)	47	49
2	B	321/324 (99%)	315 (98%)	6 (2%)	0	100	100
All	All	744/755 (98%)	722 (97%)	21 (3%)	1 (0%)	51	54

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%)	352 (99%)	2 (1%)	86	90
2	B	263/263 (100%)	262 (100%)	1 (0%)	91	94
All	All	617/621 (99%)	614 (100%)	3 (0%)	88	92

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	177	LYS
1	A	306	GLU
2	B	653	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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	2,5	10,14,15	4.81	4 (40%)	5,20,22	1.60	1 (20%)

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	2,5	-	1/5/12/14	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	760	NEP	P-O3P	13.23	1.58	1.47
2	B	760	NEP	P-O1P	5.71	1.66	1.54
2	B	760	NEP	P-O2P	-4.13	1.46	1.54
2	B	760	NEP	CE1-ND1	-2.02	1.31	1.35

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	760	NEP	O2P-P-O3P	-2.69	107.62	113.44

There are no chirality outliers.

All (1) torsion outliers are listed below:

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

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates 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
6	ADN	A	904	7	18,21,21	0.62	0	18,31,31	0.78	1 (5%)
7	GOL	B	902	-	5,5,5	0.44	0	5,5,5	0.27	0
4	ADP	A	902	5	24,29,29	0.61	0	29,45,45	0.87	1 (3%)
3	7A2	A	901	-	2,13,13	0.55	0	2,19,19	0.87	0
7	GOL	A	905	6	5,5,5	0.41	0	5,5,5	0.79	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
6	ADN	A	904	7	-	0/2/22/22	0/3/3/3
7	GOL	B	902	-	-	0/4/4/4	-
4	ADP	A	902	5	-	1/12/32/32	0/3/3/3
3	7A2	A	901	-	-	1/9/21/21	-
7	GOL	A	905	6	-	2/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	902	ADP	C5-C6-N6	2.68	124.43	120.35
6	A	904	ADN	C5-C6-N6	2.21	123.72	120.35

There are no chirality outliers.

All (4) torsion outliers are listed below:

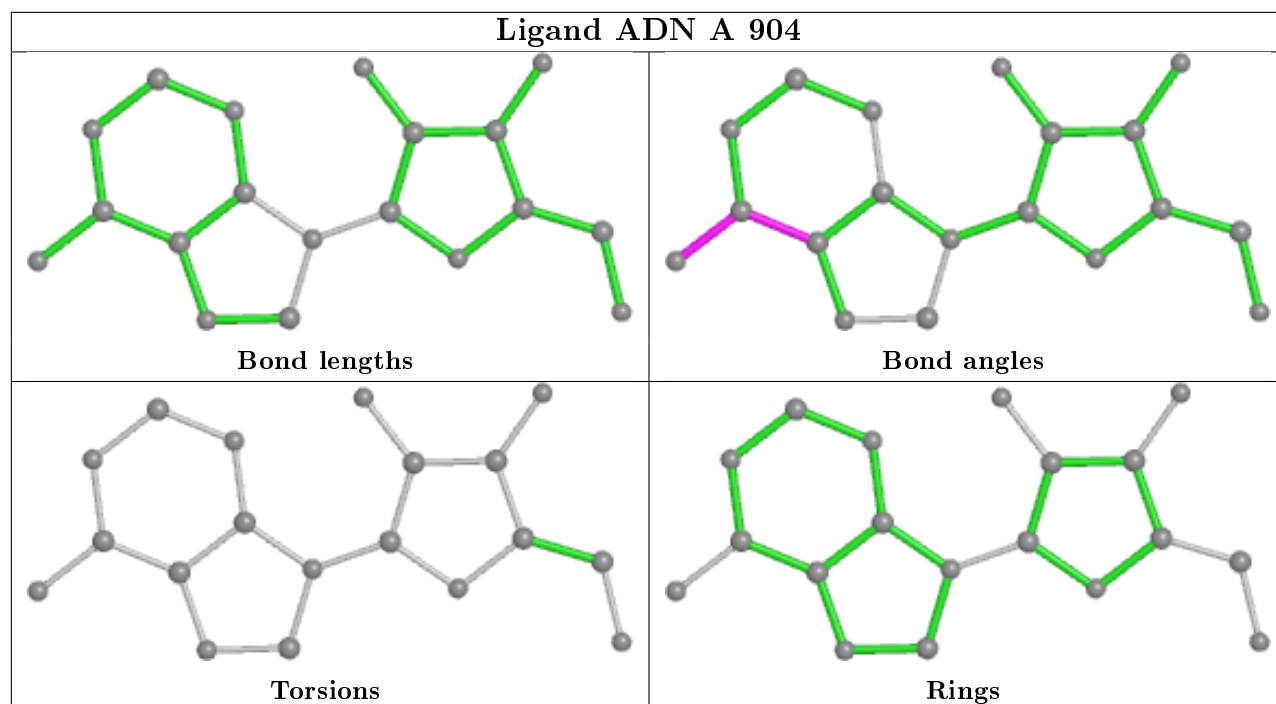
Mol	Chain	Res	Type	Atoms
3	A	901	7A2	C3-C2-C4-O5
7	A	905	GOL	O1-C1-C2-O2
7	A	905	GOL	O1-C1-C2-C3
4	A	902	ADP	C5'-O5'-PA-O1A

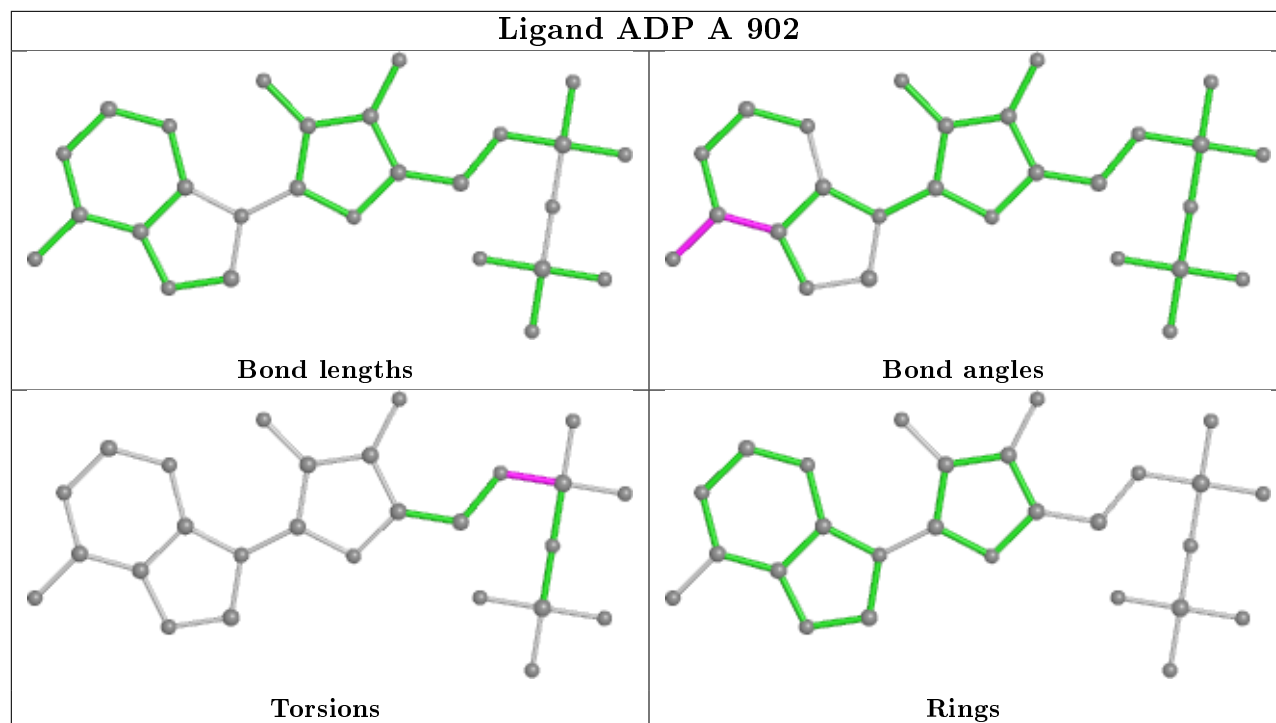
There are no ring outliers.

3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	904	ADN	1	0
3	A	901	7A2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	427/431 (99%)	0.58	35 (8%) <span style="border: 1px solid red; padding: 2px;">11</span> <span style="border: 1px solid red; padding: 2px;">15</span>	28, 59, 96, 123	0
2	B	322/324 (99%)	0.32	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	36, 54, 76, 93	0
All	All	749/755 (99%)	0.47	35 (4%) <span style="border: 1px solid red; padding: 2px;">31</span> <span style="border: 1px solid red; padding: 2px;">37</span>	28, 55, 89, 123	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	430	PHE	7.2
1	A	429	TYR	4.8
1	A	147	ALA	4.7
1	A	368	LEU	4.6
1	A	373	VAL	4.5
1	A	335	GLY	4.4
1	A	364	TYR	4.4
1	A	276	TRP	4.3
1	A	375	ILE	4.1
1	A	340	ILE	4.1
1	A	425	PRO	3.8
1	A	141	ASP	3.7
1	A	419	LEU	3.6
1	A	338	LEU	3.6
1	A	337	ILE	3.4
1	A	402	ILE	3.3
1	A	142	VAL	3.1
1	A	376	PHE	2.9
1	A	295	LEU	2.8
1	A	330	GLU	2.6
1	A	272	LYS	2.6
1	A	431	GLN	2.6
1	A	328	THR	2.5
1	A	370	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	24	ALA	2.4
1	A	399	GLY	2.3
1	A	394	VAL	2.3
1	A	278	MET	2.3
1	A	415	VAL	2.3
1	A	426	GLU	2.2
1	A	424	ILE	2.1
1	A	248	PRO	2.0
1	A	372	GLU	2.0
1	A	414	ILE	2.0
1	A	362	ARG	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	NEP	B	760	14/15	0.95	0.15	41,65,89,94	0

## 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, 95<sup>th</sup> 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.

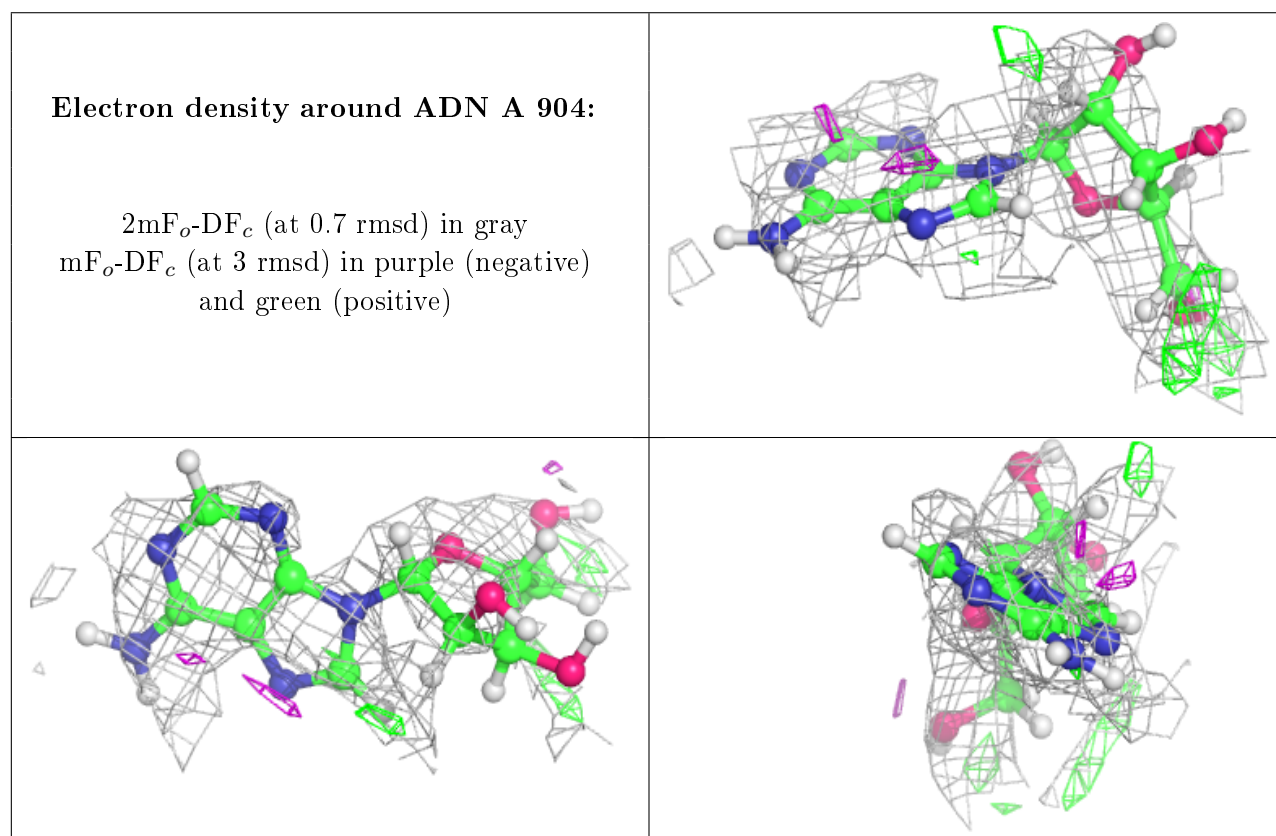
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	MG	A	903	1/1	0.50	0.12	90,90,90,90	0
7	GOL	A	905	6/6	0.65	0.37	57,100,163,165	0
6	ADN	A	904	19/19	0.71	0.36	72,144,204,234	0
5	MG	B	901	1/1	0.80	0.09	63,63,63,63	0
7	GOL	B	902	6/6	0.86	0.20	58,102,128,133	0
3	7A2	A	901	14/14	0.90	0.14	44,84,115,129	0

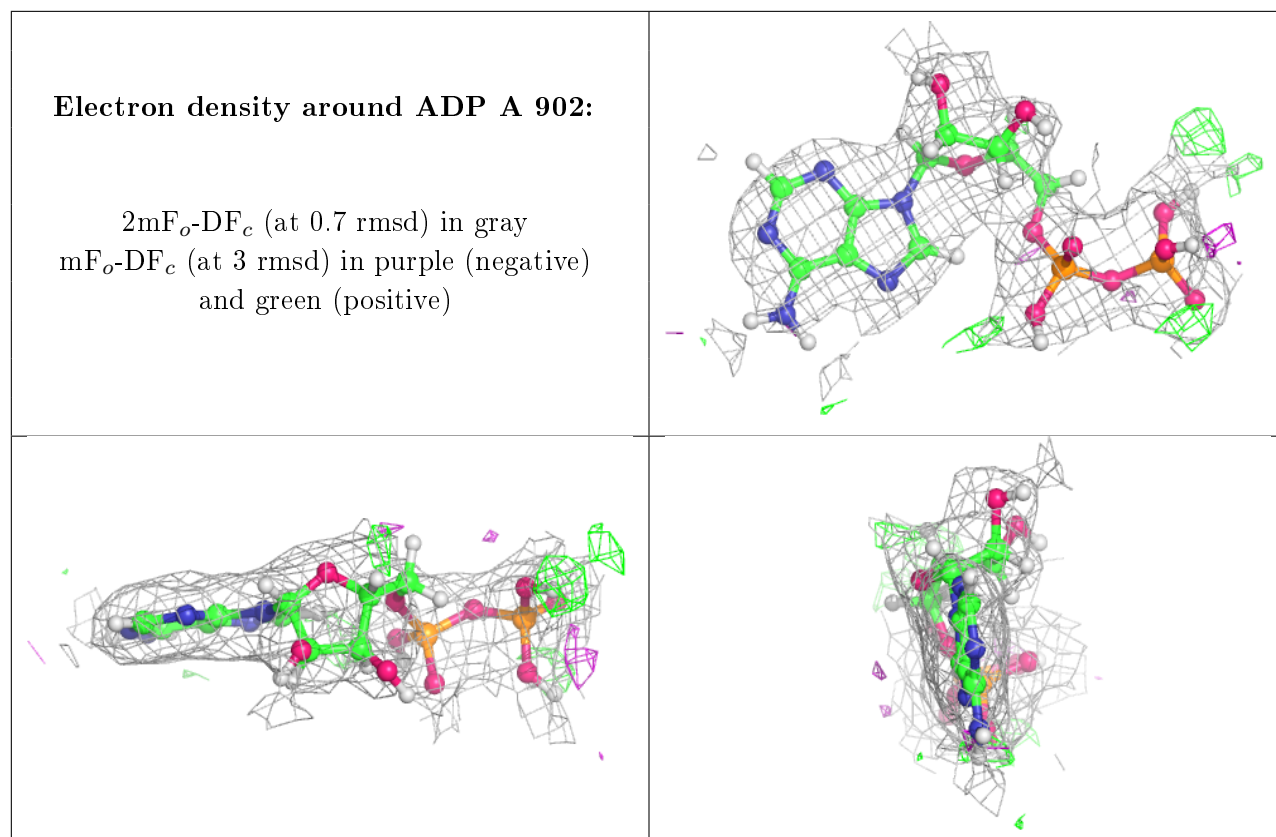
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ADP	A	902	27/27	0.94	0.11	27,72,113,123	0

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.