



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:53 AM UTC

PDB ID : 5TDE / pdb_00005tde
Title : TEV Cleaved Human ATP Citrate Lyase Bound to Citrate
Authors : Hu, J.; Fraser, M.E.
Deposited on : 2016-09-19
Resolution : 1.70 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
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

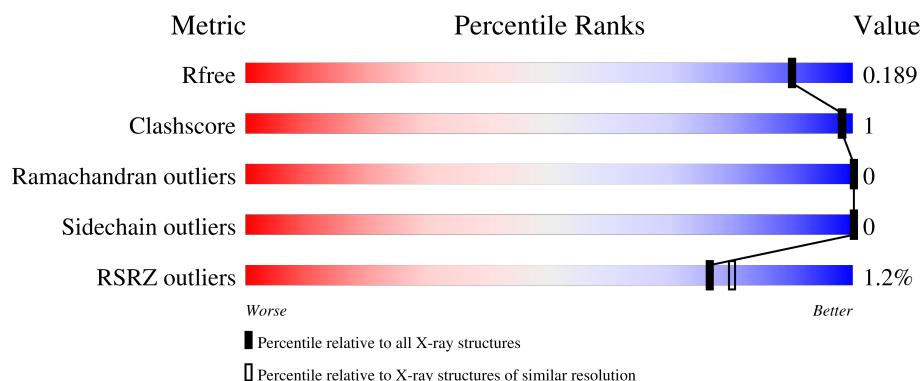
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

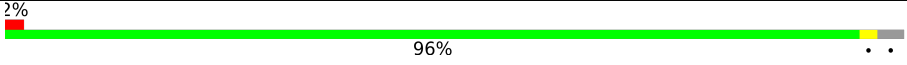
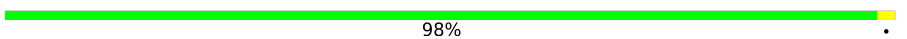
The reported resolution of this entry is 1.70 Å.

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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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 7 unique types of molecules in this entry. The entry contains 12540 atoms, of which 5928 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	420	Total	C	H	N	O	S	0	4	0
			6664	2141	3345	557	612	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	324	Total	C	H	N	O	S	0	9	0
			5048	1582	2546	432	468	20			

There is a discrepancy between the modelled and reference sequences:

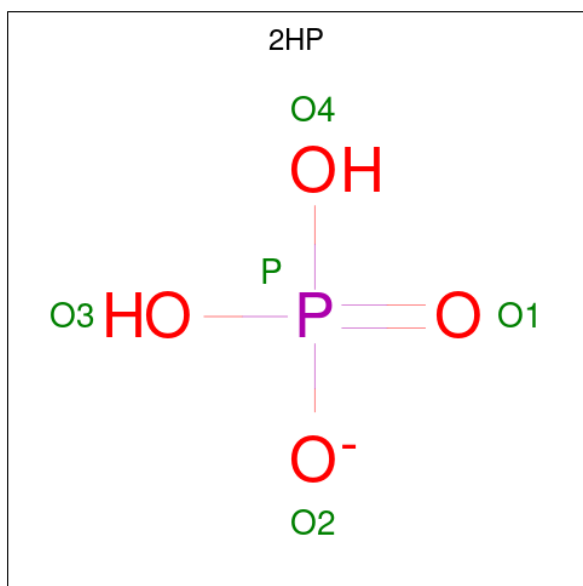
Chain	Residue	Modelled	Actual	Comment	Reference
B	487	SER	-	SEE REMARK 999	UNP P53396

- Molecule 3 is CITRATE ANION (CCD ID: FLC) (formula: C₆H₅O₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			18	6	5	7		

- Molecule 4 is DIHYDROGENPHOSPHATE ION (CCD ID: 2HP) (formula: $\text{H}_2\text{O}_4\text{P}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Na	0	0
			2	2		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).

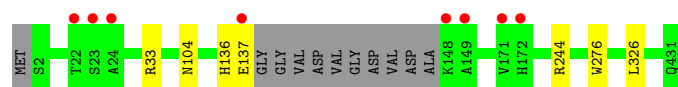


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			14	3	8	3		
6	A	1	Total	C	H	O	0	0
			14	3	8	3		
6	B	1	Total	C	H	O	0	0
			14	3	8	3		
6	B	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	401	Total	O	0	0
			401	401		
7	B	336	Total	O	0	0
			336	336		

- Molecule 1: ATP-citrate synthase



- Chain B: 98%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.83Å 84.08Å 194.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.39 – 1.70 51.39 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.1 (51.39-1.70) 94.4 (51.39-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 1.70Å)	Xtriage
Refinement program	PHENIX (1.10_2152: ???)	Depositor
R, R_{free}	0.155 , 0.184 0.158 , 0.189	Depositor DCC
R_{free} test set	5021 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	16.3	Xtriage
Anisotropy	0.683	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	12540	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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: NA, 2HP, FLC, GOL

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	0.40	0/3417	0.54	0/4632
2	B	0.42	0/2595	0.59	0/3502
All	All	0.41	0/6012	0.56	0/8134

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	3319	3345	3336	4	0
2	B	2502	2546	2528	4	0
3	A	13	5	5	0	0
4	A	5	0	0	0	0
4	B	10	0	0	0	0
5	A	2	0	0	0	0
6	A	12	16	16	0	0
6	B	12	16	16	0	0
7	A	401	0	0	0	4
7	B	336	0	0	1	5
All	All	6612	5928	5901	7	5

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 1.

All (7) 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	2.36	0.52
1:A:33:ARG:NE	1:A:104:ASN:OD1	2.48	0.46
2:B:511:GLY:HA3	7:B:1067:HOH:O	2.15	0.46
2:B:505:GLN:HG3	2:B:572:PHE:CG	2.54	0.43
1:A:136:HIS:O	1:A:137:GLU:HB2	2.20	0.42
1:A:276:TRP:CH2	1:A:326:LEU:HB3	2.55	0.42
2:B:743:TRP:CH2	2:B:789:SER:HA	2.56	0.40

All (5) 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)
7:A:856:HOH:O	7:B:1036:HOH:O[3_655]	1.75	0.45
7:A:905:HOH:O	7:B:1104:HOH:O[3_645]	1.88	0.32
7:A:883:HOH:O	7:B:1284:HOH:O[3_655]	2.07	0.13
7:B:1294:HOH:O	7:B:1320:HOH:O[3_655]	2.09	0.11
7:A:856:HOH:O	7:B:1015:HOH:O[3_655]	2.19	0.01

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	420/431 (97%)	414 (99%)	6 (1%)	0	100	100
2	B	331/324 (102%)	327 (99%)	4 (1%)	0	100	100
All	All	751/755 (100%)	741 (99%)	10 (1%)	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	355/358 (99%)	355 (100%)	0	100	100
2	B	273/264 (103%)	273 (100%)	0	100	100
All	All	628/622 (101%)	628 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	638	ASN

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 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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	FLC	A	501	5	12,12,12	0.97	0	17,17,17	1.24	1 (5%)
6	GOL	A	505	-	5,5,5	0.49	0	5,5,5	0.51	0
4	2HP	B	902	-	4,4,4	3.70	1 (25%)	6,6,6	2.20	2 (33%)
6	GOL	B	903	-	5,5,5	0.47	0	5,5,5	0.48	0
6	GOL	B	904	-	5,5,5	0.36	0	5,5,5	0.49	0
4	2HP	B	901	5	4,4,4	2.49	2 (50%)	6,6,6	1.15	0
4	2HP	A	502	-	4,4,4	2.76	1 (25%)	6,6,6	1.40	1 (16%)
6	GOL	A	506	-	5,5,5	0.56	0	5,5,5	0.73	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	GOL	A	505	-	-	0/4/4/4	-
3	FLC	A	501	5	-	2/16/16/16	-
6	GOL	B	903	-	-	3/4/4/4	-
6	GOL	B	904	-	-	0/4/4/4	-
6	GOL	A	506	-	-	4/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	902	2HP	P-O4	7.23	1.75	1.54
4	A	502	2HP	P-O4	5.19	1.69	1.54
4	B	901	2HP	P-O4	3.98	1.66	1.54
4	B	901	2HP	P-O3	-2.41	1.47	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	902	2HP	O4-P-O2	-4.13	95.06	107.91
3	A	501	FLC	OB1-CBC-CB	-2.68	116.91	122.09
4	A	502	2HP	O4-P-O2	-2.29	100.78	107.91
4	B	902	2HP	O3-P-O2	2.21	114.80	107.91

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	506	GOL	O1-C1-C2-C3
6	A	506	GOL	C1-C2-C3-O3
6	B	903	GOL	O1-C1-C2-O2
6	B	903	GOL	O1-C1-C2-C3
6	A	506	GOL	O1-C1-C2-O2
6	A	506	GOL	O2-C2-C3-O3
6	B	903	GOL	C1-C2-C3-O3
3	A	501	FLC	CB-CA-CAC-OA1
3	A	501	FLC	CB-CA-CAC-OA2

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	420/431 (97%)	-0.22	8 (1%) 66 71	11, 24, 42, 77	4 (0%)
2	B	324/324 (100%)	-0.53	1 (0%) 90 91	9, 19, 32, 59	10 (3%)
All	All	744/755 (98%)	-0.36	9 (1%) 76 80	9, 22, 39, 77	14 (1%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	LYS	3.1
1	A	24	ALA	3.0
1	A	137	GLU	3.0
1	A	171	VAL	2.6
1	A	149	ALA	2.5
1	A	23	SER	2.4
1	A	172	HIS	2.3
2	B	807	ASN	2.3
1	A	22	THR	2.1

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	A	506	6/6	0.80	0.16	35,45,56,56	0
6	GOL	B	903	6/6	0.89	0.13	23,41,47,54	0
6	GOL	A	505	6/6	0.90	0.11	23,37,48,54	0
6	GOL	B	904	6/6	0.90	0.10	31,40,48,51	0
4	2HP	B	902	5/5	0.94	0.09	32,35,42,45	0
5	NA	A	503	1/1	0.98	0.08	23,23,23,23	0
5	NA	A	504	1/1	0.98	0.04	31,31,31,31	0
3	FLC	A	501	13/13	0.98	0.04	13,16,24,25	0
4	2HP	B	901	5/5	0.99	0.06	14,14,18,20	0
4	2HP	A	502	5/5	0.99	0.04	18,19,21,23	0

6.5 Other polymers ⓘ

There are no such residues in this entry.