



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:43 AM UTC

PDB ID : 1R6B / pdb_00001r6b
Title : High resolution crystal structure of ClpA
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Deposited on : 2003-10-15
Resolution : 2.25 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

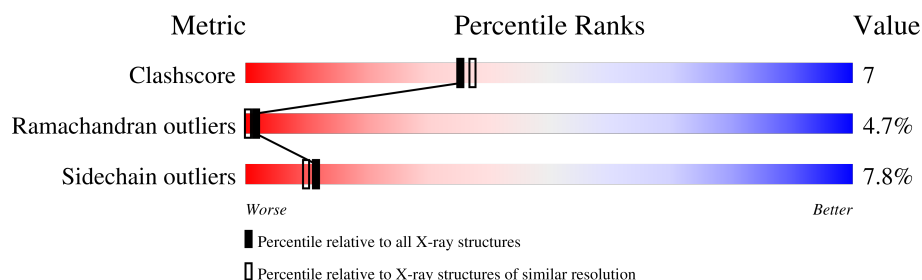
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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))
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	X	758	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5623 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 ClpA protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	704	Total	C	N	O	S	0	0	0
			5506	3466	986	1037	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	169	LEU	MET	engineered mutation	UNP P0ABH9
X	367	ALA	GLY	SEE REMARK 999	UNP P0ABH9
X	411	LEU	VAL	SEE REMARK 999	UNP P0ABH9
X	533	LEU	VAL	SEE REMARK 999	UNP P0ABH9

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	X	3	Total	Mg	0	0
			3	3		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

- Molecule 4 is water.

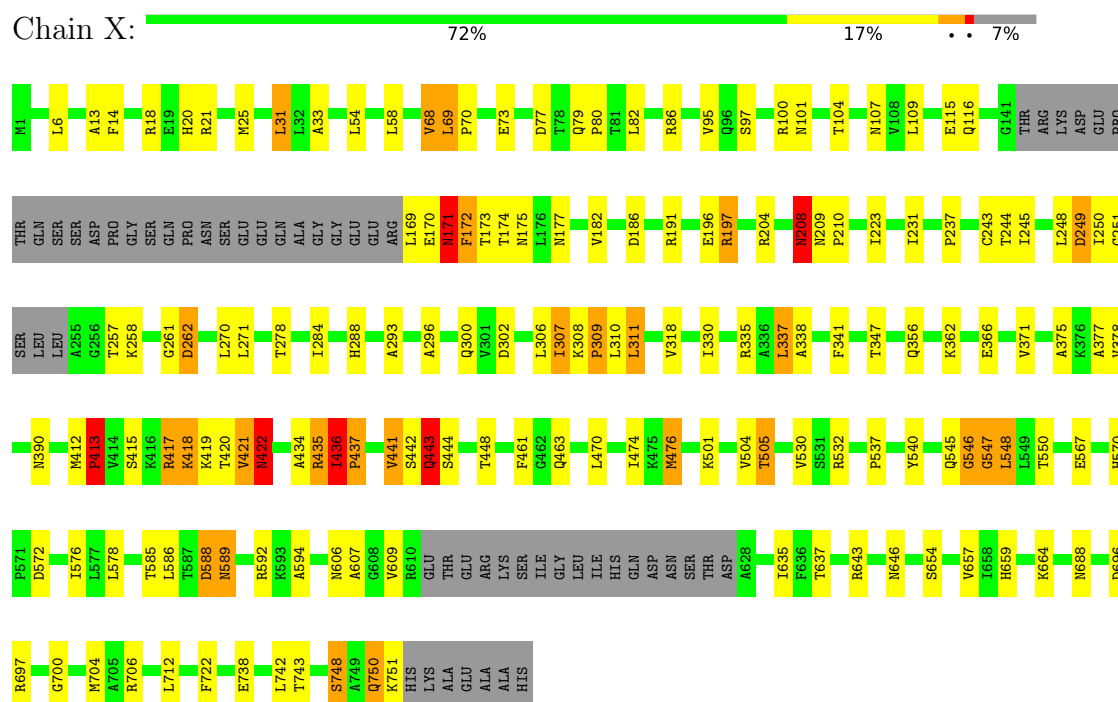
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	X	60	Total	O	0	0
			60	60		

3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

- Molecule 1: ClpA protein



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	124.11Å 124.11Å 97.05Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.25	Depositor
% Data completeness (in resolution range)	93.8 (20.00-2.25)	Depositor
R_{merge}	0.03	Depositor
R_{sym}	0.03	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.234 , 0.278	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5623	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	X	1.06	11/5586 (0.2%)	0.93	13/7540 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	436	ILE	CA-C	9.21	1.62	1.52
1	X	284	ILE	CA-CB	8.07	1.64	1.53
1	X	704	MET	SD-CE	7.75	1.99	1.79
1	X	537	PRO	CA-C	7.04	1.58	1.52
1	X	422	ASN	CB-CG	-5.98	1.37	1.52
1	X	318	VAL	CA-CB	5.94	1.62	1.54
1	X	378	VAL	CA-CB	5.73	1.61	1.54
1	X	97	SER	C-O	-5.55	1.16	1.24
1	X	637	THR	CA-CB	5.27	1.61	1.53
1	X	208	ASN	CB-CG	-5.17	1.39	1.52
1	X	375	ALA	CA-CB	5.06	1.61	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	443	GLN	OE1-CD-NE2	-9.81	112.79	122.60
1	X	20	HIS	N-CA-C	-7.51	104.11	113.28
1	X	443	GLN	CG-CD-NE2	6.67	126.40	116.40
1	X	537	PRO	N-CA-C	6.48	118.61	110.70
1	X	307	ILE	N-CA-C	5.91	118.48	111.09
1	X	171	ASN	N-CA-C	5.90	123.36	110.80
1	X	750	GLN	N-CA-C	-5.79	103.94	111.71
1	X	548	LEU	N-CA-C	5.61	122.76	110.80
1	X	390	ASN	N-CA-C	5.28	119.43	112.89
1	X	116	GLN	N-CA-C	5.26	117.70	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	69	LEU	N-CA-C	5.20	121.31	109.81
1	X	101	ASN	N-CA-C	5.19	117.02	111.36
1	X	73	GLU	N-CA-C	5.03	118.58	109.32

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	X	5506	0	5598	83	0
2	X	3	0	0	0	0
3	X	54	0	23	0	0
4	X	60	0	0	3	0
All	All	5623	0	5621	83	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 7.

All (83) 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:X:442:SER:O	1:X:443:GLN:CG	1.77	1.32
1:X:420:THR:O	1:X:421:VAL:HG23	1.08	1.24
1:X:420:THR:O	1:X:421:VAL:CG2	1.94	1.16
1:X:442:SER:O	1:X:443:GLN:HG3	0.85	1.02
1:X:550:THR:HG23	1:X:594:ALA:HB2	1.40	0.99
1:X:436:ILE:H	1:X:437:PRO:CD	1.82	0.93
1:X:442:SER:C	1:X:443:GLN:HG3	1.96	0.90
1:X:250:ILE:HG22	1:X:251:GLY:H	1.41	0.86
1:X:420:THR:C	1:X:421:VAL:HG23	2.02	0.84
1:X:104:THR:H	1:X:107:ASN:HD22	1.27	0.81
1:X:436:ILE:H	1:X:437:PRO:HD2	1.43	0.81
1:X:377:ALA:HB2	1:X:422:ASN:HB3	1.66	0.78
1:X:434:ALA:O	1:X:435:ARG:HB2	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:250:ILE:HG22	1:X:251:GLY:N	2.04	0.72
1:X:607:ALA:HB1	1:X:635:ILE:HD13	1.72	0.70
1:X:418:LYS:HD3	1:X:422:ASN:OD1	1.94	0.67
1:X:418:LYS:HE2	1:X:422:ASN:HD21	1.63	0.64
1:X:436:ILE:N	1:X:437:PRO:CD	2.59	0.63
1:X:177:ASN:ND2	1:X:245:ILE:H	1.95	0.63
1:X:13:ALA:HB2	1:X:33:ALA:HB2	1.80	0.62
1:X:191:ARG:HD3	1:X:223:ILE:HD11	1.82	0.62
1:X:169:LEU:HD12	1:X:170:GLU:H	1.65	0.62
1:X:261:GLY:O	1:X:262:ASP:HB2	1.99	0.61
1:X:69:LEU:N	1:X:70:PRO:HD3	2.14	0.61
1:X:271:LEU:HD13	1:X:307:ILE:HG13	1.83	0.61
1:X:442:SER:O	1:X:443:GLN:CB	2.49	0.60
1:X:589:ASN:HB3	1:X:592:ARG:HD2	1.83	0.60
1:X:250:ILE:CG2	1:X:251:GLY:H	2.11	0.60
1:X:171:ASN:O	1:X:172:PHE:HB2	2.01	0.60
1:X:436:ILE:N	1:X:437:PRO:HD2	2.15	0.60
1:X:550:THR:HG21	1:X:588:ASP:OD1	2.02	0.60
1:X:546:GLY:O	1:X:589:ASN:ND2	2.35	0.59
1:X:14:PHE:O	1:X:18:ARG:HG3	2.03	0.59
1:X:337:LEU:HD12	1:X:341:PHE:HE1	1.69	0.57
1:X:417:ARG:O	1:X:418:LYS:HB2	2.04	0.57
1:X:570:HIS:HD2	1:X:572:ASP:HB2	1.71	0.55
1:X:204:ARG:O	1:X:208:ASN:ND2	2.38	0.55
1:X:659:HIS:CE1	1:X:688:ASN:HD22	2.25	0.54
1:X:249:ASP:N	1:X:249:ASP:OD1	2.41	0.54
1:X:31:LEU:HD13	1:X:58:LEU:HD11	1.90	0.53
1:X:588:ASP:CG	1:X:589:ASN:N	2.66	0.52
1:X:68:VAL:HG12	1:X:70:PRO:HD3	1.92	0.51
1:X:308:LYS:HB2	1:X:309:PRO:HD3	1.93	0.51
1:X:191:ARG:NH1	4:X:901:HOH:O	2.43	0.51
1:X:306:LEU:HD23	1:X:311:LEU:HD13	1.92	0.50
1:X:545:GLN:O	1:X:547:GLY:N	2.45	0.50
1:X:748:SER:C	1:X:750:GLN:H	2.20	0.50
1:X:578:LEU:HD22	1:X:643:ARG:HD2	1.94	0.48
1:X:371:VAL:HG22	1:X:418:LYS:HD2	1.95	0.48
1:X:436:ILE:H	1:X:437:PRO:HD3	1.71	0.48
1:X:588:ASP:OD1	1:X:588:ASP:C	2.58	0.47
1:X:659:HIS:CE1	1:X:688:ASN:ND2	2.82	0.47
1:X:659:HIS:HE1	1:X:688:ASN:ND2	2.13	0.47
1:X:412:MET:O	1:X:413:PRO:C	2.57	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:13:ALA:HB1	1:X:25:MET:HE1	1.98	0.46
1:X:309:PRO:HB2	1:X:310:LEU:H	1.58	0.46
1:X:25:MET:HB3	1:X:80:PRO:HA	1.97	0.46
1:X:288:HIS:HB2	1:X:330:ILE:HD11	1.98	0.45
1:X:308:LYS:HE3	1:X:337:LEU:HD13	1.97	0.45
1:X:86:ARG:NH1	1:X:115:GLU:OE2	2.50	0.45
1:X:337:LEU:HD12	1:X:341:PHE:CE1	2.49	0.45
1:X:570:HIS:CD2	1:X:572:ASP:H	2.35	0.44
1:X:191:ARG:NH2	1:X:347:THR:O	2.50	0.44
1:X:197:ARG:HD3	4:X:941:HOH:O	2.18	0.44
1:X:169:LEU:HD23	1:X:174:THR:HG23	2.01	0.43
1:X:95:VAL:HG13	1:X:100:ARG:HB2	2.00	0.42
1:X:362:LYS:HD3	1:X:366:GLU:OE2	2.19	0.42
1:X:540:TYR:HD2	1:X:540:TYR:HA	1.71	0.42
1:X:589:ASN:CB	1:X:592:ARG:HD2	2.49	0.42
1:X:209:ASN:HA	1:X:210:PRO:HD3	1.90	0.42
1:X:434:ALA:O	1:X:435:ARG:CB	2.64	0.42
1:X:79:GLN:HA	1:X:80:PRO:HD3	1.92	0.42
1:X:169:LEU:HB3	1:X:174:THR:HG21	2.01	0.42
1:X:476:MET:HB3	1:X:476:MET:HE3	1.74	0.42
1:X:501:LYS:O	1:X:505:THR:HG23	2.20	0.41
1:X:177:ASN:HD21	1:X:245:ILE:H	1.68	0.41
1:X:576:ILE:HG23	1:X:586:LEU:HD21	2.02	0.41
1:X:250:ILE:CG2	1:X:251:GLY:N	2.72	0.40
1:X:470:LEU:O	1:X:474:ILE:HG12	2.21	0.40
1:X:664:LYS:NZ	4:X:982:HOH:O	2.52	0.40
1:X:231:ILE:HG21	1:X:243:CYS:O	2.21	0.40
1:X:461:PHE:HB2	1:X:657:VAL:HG13	2.03	0.40
1:X:696:ASP:O	1:X:700:GLY:N	2.51	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	X	696/758 (92%)	604 (87%)	59 (8%)	33 (5%)	2 0

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	171	ASN
1	X	172	PHE
1	X	262	ASP
1	X	413	PRO
1	X	415	SER
1	X	417	ARG
1	X	421	VAL
1	X	435	ARG
1	X	436	ILE
1	X	443	GLN
1	X	175	ASN
1	X	296	ALA
1	X	300	GLN
1	X	309	PRO
1	X	418	LYS
1	X	419	LYS
1	X	422	ASN
1	X	437	PRO
1	X	546	GLY
1	X	548	LEU
1	X	293	ALA
1	X	335	ARG
1	X	441	VAL
1	X	258	LYS
1	X	444	SER
1	X	530	VAL
1	X	532	ARG
1	X	173	THR
1	X	237	PRO
1	X	338	ALA
1	X	609	VAL
1	X	182	VAL
1	X	547	GLY

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	X	592/639 (93%)	546 (92%)	46 (8%)	11	10

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	6	LEU
1	X	21	ARG
1	X	31	LEU
1	X	54	LEU
1	X	68	VAL
1	X	77	ASP
1	X	82	LEU
1	X	109	LEU
1	X	186	ASP
1	X	196	GLU
1	X	197	ARG
1	X	208	ASN
1	X	244	THR
1	X	248	LEU
1	X	249	ASP
1	X	257	THR
1	X	270	LEU
1	X	278	THR
1	X	302	ASP
1	X	311	LEU
1	X	337	LEU
1	X	356	GLN
1	X	413	PRO
1	X	422	ASN
1	X	441	VAL
1	X	448	THR
1	X	463	GLN
1	X	476	MET
1	X	504	VAL
1	X	505	THR
1	X	567	GLU
1	X	585	THR
1	X	588	ASP
1	X	589	ASN

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Mol	Chain	Res	Type
1	X	606	ASN
1	X	646	ASN
1	X	654	SER
1	X	697	ARG
1	X	706	ARG
1	X	712	LEU
1	X	722	PHE
1	X	738	GLU
1	X	742	LEU
1	X	743	THR
1	X	748	SER
1	X	751	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	4	GLN
1	X	22	HIS
1	X	64	GLN
1	X	107	ASN
1	X	116	GLN
1	X	177	ASN
1	X	200	GLN
1	X	233	GLN
1	X	273	GLN
1	X	276	GLN
1	X	356	GLN
1	X	422	ASN
1	X	528	HIS
1	X	570	HIS
1	X	575	ASN
1	X	606	ASN
1	X	659	HIS
1	X	660	GLN
1	X	688	ASN

5.3.3 RNA ⓘ

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 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	ADP	X	781	2	28,29,29	1.21	4 (14%)	43,45,45	1.80	7 (16%)
3	ADP	X	780	2	28,29,29	1.39	5 (17%)	43,45,45	1.73	11 (25%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ADP	X	781	2	-	0/16/32/32	0/3/3/3
3	ADP	X	780	2	-	1/16/32/32	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	X	780	ADP	O3'-C3'	-3.40	1.34	1.43
3	X	781	ADP	C5-N7	-3.20	1.33	1.39
3	X	781	ADP	C2-N1	2.62	1.38	1.33
3	X	780	ADP	C5-N7	-2.58	1.34	1.39
3	X	780	ADP	C8-N7	2.54	1.36	1.31
3	X	780	ADP	C4-N9	-2.30	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	X	780	ADP	C2'-C3'	-2.25	1.47	1.53
3	X	781	ADP	PA-O2A	-2.01	1.46	1.55
3	X	781	ADP	C5-C6	-2.01	1.35	1.41

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	X	781	ADP	N3-C2-N1	-5.43	120.36	128.58
3	X	780	ADP	N3-C2-N1	-5.07	120.91	128.58
3	X	781	ADP	C5-C4-N3	-4.94	119.92	126.72
3	X	780	ADP	N9-C8-N7	-4.32	107.81	113.94
3	X	780	ADP	C5-C4-N3	-4.31	120.78	126.72
3	X	781	ADP	N3-C4-N9	3.63	133.34	127.17
3	X	780	ADP	C2-N3-C4	3.42	120.17	111.83
3	X	781	ADP	N9-C8-N7	-3.38	109.15	113.94
3	X	781	ADP	C2-N3-C4	3.36	120.05	111.83
3	X	781	ADP	C5-N7-C8	3.33	108.68	103.45
3	X	780	ADP	C5-N7-C8	2.94	108.08	103.45
3	X	781	ADP	C6-C5-C4	2.70	120.87	117.18
3	X	780	ADP	O3'-C3'-C4'	-2.37	104.28	111.08
3	X	780	ADP	O3'-C3'-C2'	-2.35	104.27	111.82
3	X	780	ADP	N3-C4-N9	2.28	131.05	127.17
3	X	780	ADP	C4-C5-N7	-2.27	107.98	110.58
3	X	780	ADP	C5-C4-N9	2.17	108.17	105.81
3	X	780	ADP	C4-N9-C8	2.10	107.95	105.74

There are no chirality outliers.

All (1) torsion outliers are listed below:

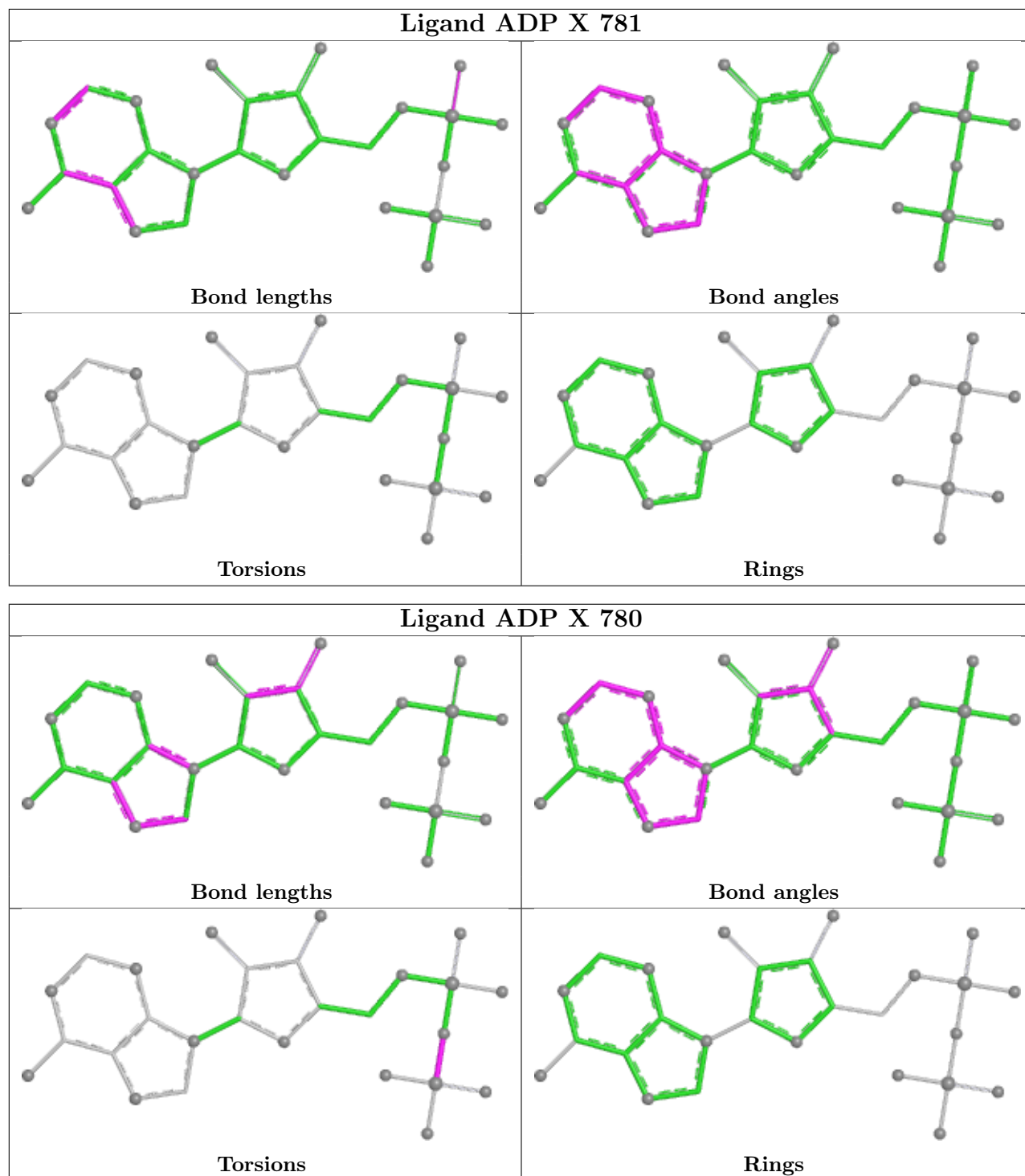
Mol	Chain	Res	Type	Atoms
3	X	780	ADP	PA-O3A-PB-O2B

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.