



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 09:25 PM EDT

PDB ID : 4HNA / pdb_00004hna
Title : Kinesin motor domain in the ADP-MG-ALFX state in complex with tubulin and a DARPIN
Authors : Gigant, B.; Knossow, M.
Deposited on : 2012-10-19
Resolution : 3.19 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

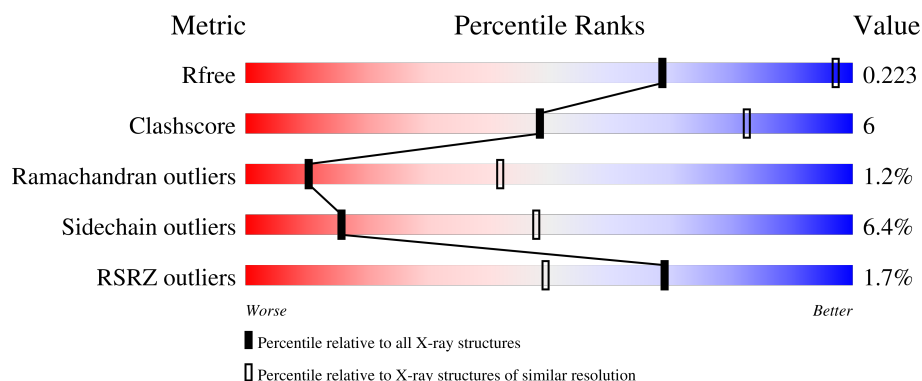
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	451	 2% 75% 18% • 5%
2	B	445	 2% 82% 12% • •
3	D	169	 2% 79% 14% • 6%
4	K	349	 % 67% 25% • 5%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 10607 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 Tubulin alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	0	0
			3349	2123	571	633	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	232	SER	GLY	SEE REMARK 999	UNP D0VWZ0
A	340	SER	THR	SEE REMARK 999	UNP D0VWZ0

- Molecule 2 is a protein called Tubulin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	431	Total	C	N	O	S	0	0	0
			3378	2118	579	655	26			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	318	ILE	VAL	SEE REMARK 999	UNP D0VWY9

- Molecule 3 is a protein called DESIGNED ANKYRIN REPEAT PROTEIN (DARPIN) D2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	159	Total	C	N	O	S	0	0	0
			1185	743	202	238	2			

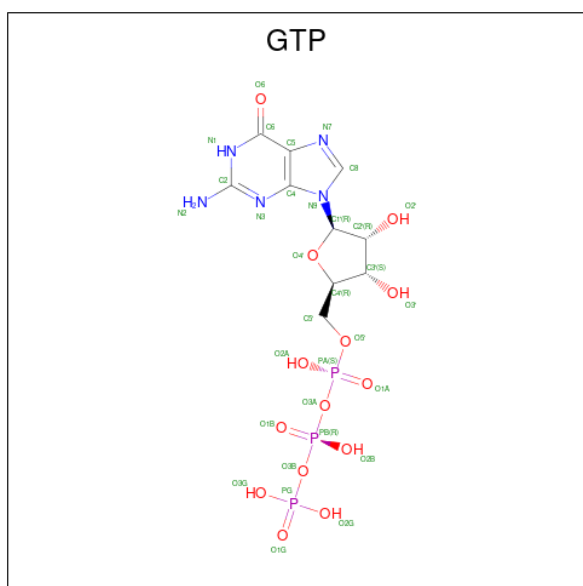
- Molecule 4 is a protein called Kinesin-1 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	K	333	Total	C	N	O	S	0	0	0
			2593	1616	447	520	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	7	SER	CYS	SEE REMARK 999	UNP P33176
K	65	ALA	CYS	SEE REMARK 999	UNP P33176
K	168	ALA	CYS	SEE REMARK 999	UNP P33176
K	174	SER	CYS	SEE REMARK 999	UNP P33176
K	294	ALA	CYS	SEE REMARK 999	UNP P33176
K	330	SER	CYS	SEE REMARK 999	UNP P33176

- Molecule 5 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).

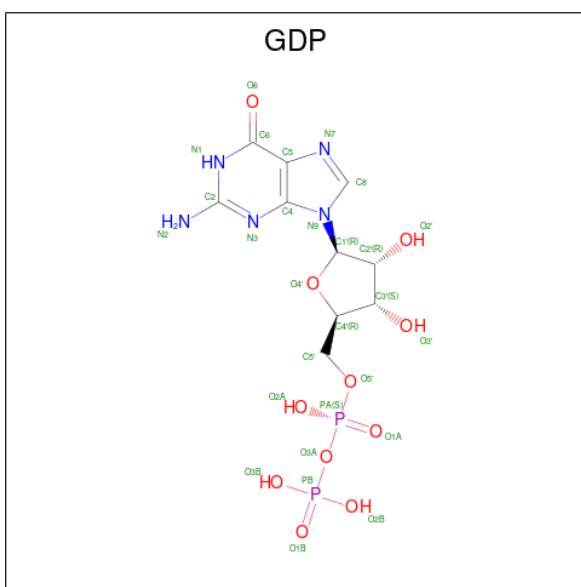


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

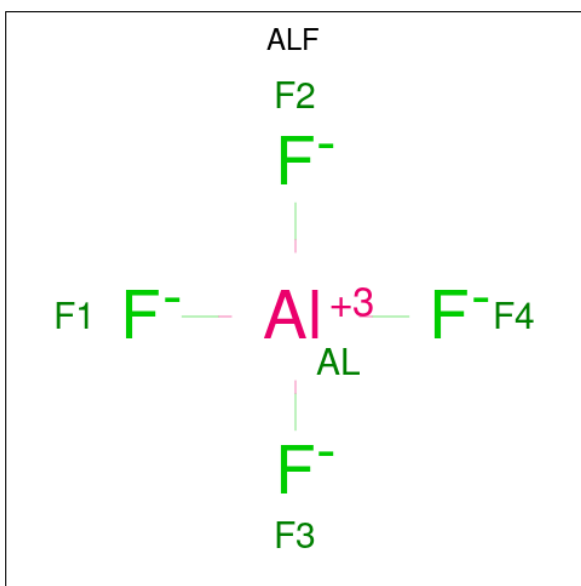
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		
6	B	1	Total	Mg	0	0
			1	1		
6	K	1	Total	Mg	0	0
			1	1		

- Molecule 7 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



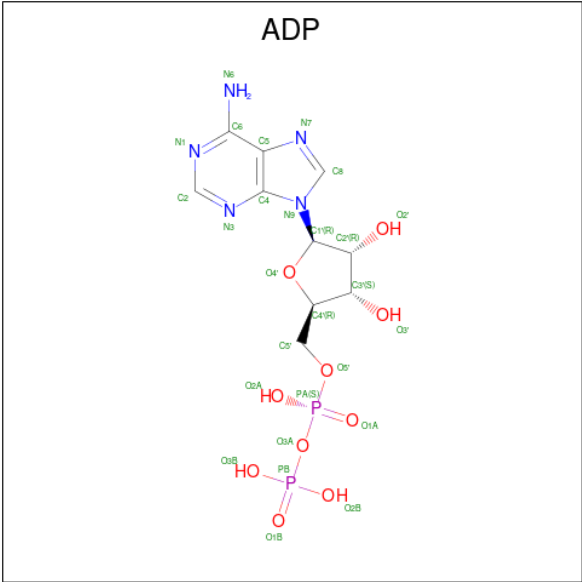
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 8 is TETRAFLUOROALUMINATE ION (CCD ID: ALF) (formula: AlF_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total 5	Al 1	F 4	0	0
8	K	1	Total 5	Al 1	F 4	0	0

- Molecule 9 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	K	1	Total	C	N	O	P	
			27	10	5	10	2	

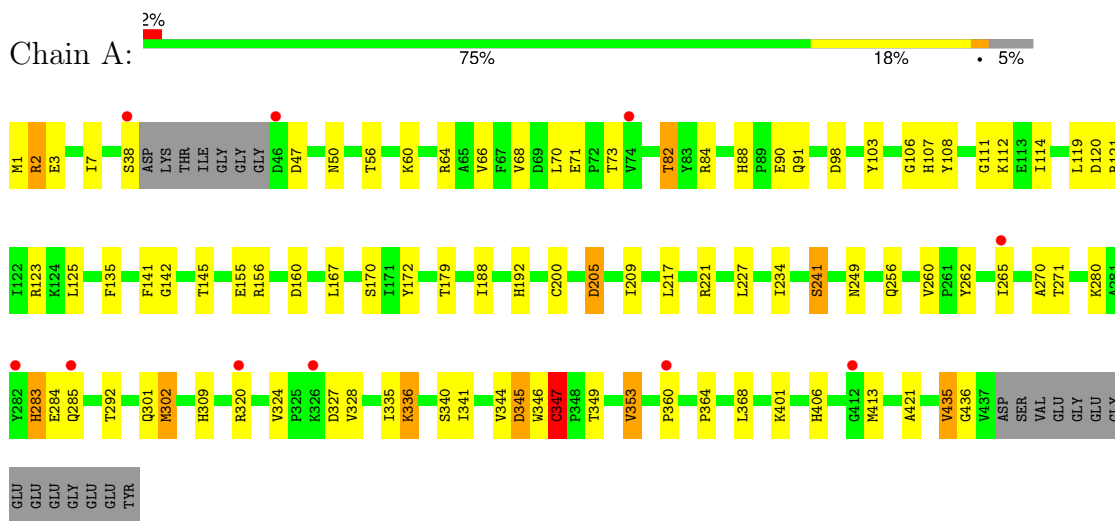
- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	K	2	Total O		
			2 2	0	0

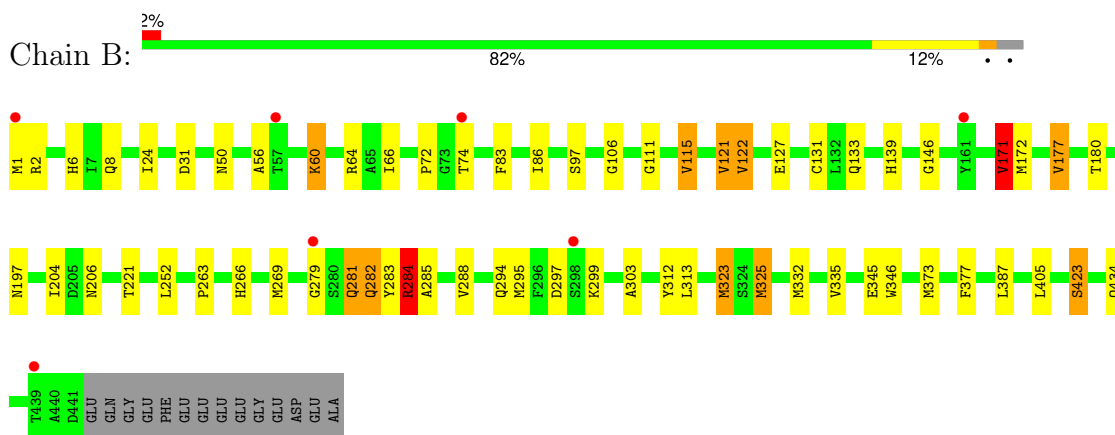
3 Residue-property plots [i](#)

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.

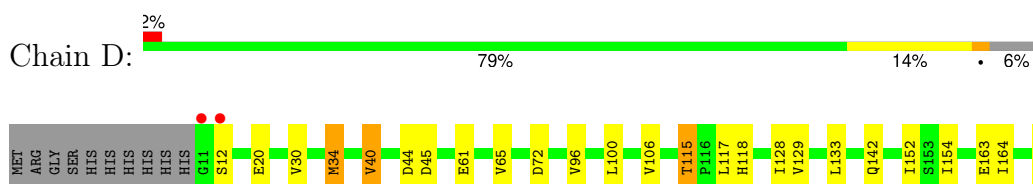
- Molecule 1: Tubulin alpha chain



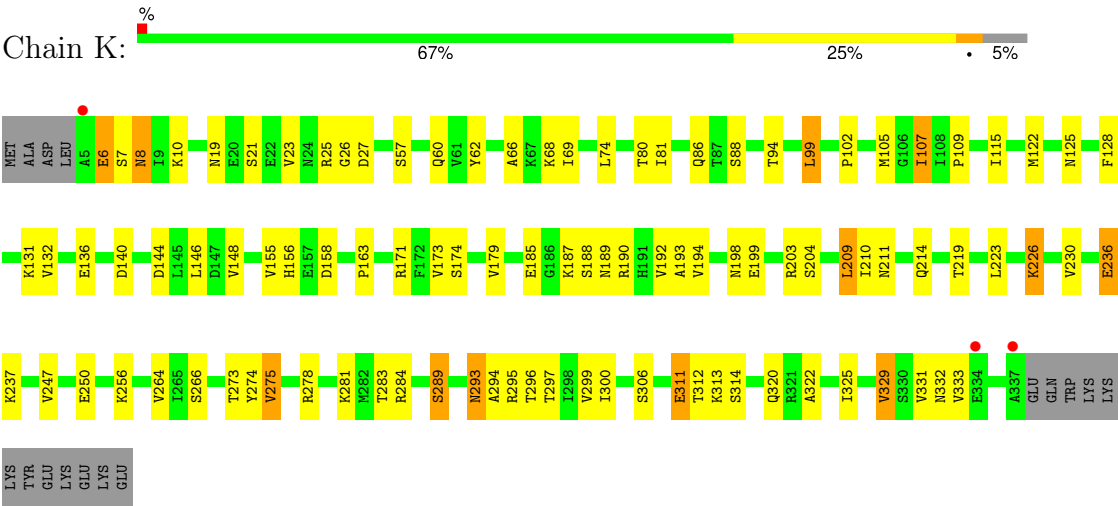
- Molecule 2: Tubulin beta chain



- Molecule 3: DESIGNED ANKYRIN REPEAT PROTEIN (DARPIN) D2



● Molecule 4: Kinesin-1 heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.05Å 160.54Å 174.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.14 – 3.19 59.14 – 3.19	Depositor EDS
% Data completeness (in resolution range)	80.2 (59.14-3.19) 80.3 (59.14-3.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.08 (at 3.19Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.177 , 0.211 0.200 , 0.223	Depositor DCC
R_{free} test set	1519 reflections (4.09%)	wwPDB-VP
Wilson B-factor (Å ²)	62.5	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 63.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	10607	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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: GTP, ADP, MG, ALF, GDP

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	0.84	1/3424 (0.0%)	1.37	16/4649 (0.3%)
2	B	0.80	0/3452	1.34	23/4676 (0.5%)
3	D	0.85	0/1200	1.50	12/1626 (0.7%)
4	K	0.89	0/2632	1.49	26/3551 (0.7%)
All	All	0.84	1/10708 (0.0%)	1.41	77/14502 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	347	CYS	CA-C	-5.23	1.48	1.52

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	K	333	VAL	N-CA-C	10.12	119.93	110.42
1	A	82	THR	N-CA-C	-10.10	95.33	110.28
1	A	345	ASP	CA-CB-CG	8.97	121.57	112.60
2	B	297	ASP	CA-CB-CG	7.97	120.57	112.60
2	B	97	SER	N-CA-C	7.13	118.85	111.14
4	K	158	ASP	CA-CB-CG	7.07	119.67	112.60
1	A	73	THR	N-CA-C	6.95	119.74	111.33
2	B	171	VAL	N-CA-CB	6.89	118.86	111.00
2	B	180	THR	CA-C-N	6.88	129.48	120.60
2	B	180	THR	C-N-CA	6.88	129.48	120.60
3	D	44	ASP	CA-CB-CG	6.87	119.47	112.60
2	B	281	GLN	CA-C-N	6.87	134.06	121.70
2	B	281	GLN	C-N-CA	6.87	134.06	121.70
4	K	329	VAL	N-CA-C	6.84	117.92	108.89
1	A	285	GLN	N-CA-C	6.78	120.72	109.46
4	K	275	VAL	N-CA-CB	6.75	120.66	111.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	122	VAL	N-CA-CB	6.73	120.68	110.58
4	K	230	VAL	N-CA-C	6.64	118.40	108.23
4	K	306	SER	CA-C-N	6.55	129.06	120.28
4	K	306	SER	C-N-CA	6.55	129.06	120.28
2	B	171	VAL	CB-CA-C	6.46	118.78	110.96
2	B	115	VAL	N-CA-C	6.39	117.07	110.36
3	D	115	THR	CB-CA-C	6.37	118.65	108.91
1	A	340	SER	CA-C-N	6.20	129.41	122.59
1	A	340	SER	C-N-CA	6.20	129.41	122.59
4	K	144	ASP	CA-CB-CG	6.17	118.77	112.60
4	K	125	ASN	CA-CB-CG	6.07	118.67	112.60
4	K	332	ASN	CA-C-N	6.07	128.12	120.72
4	K	332	ASN	C-N-CA	6.07	128.12	120.72
4	K	8	ASN	CA-CB-CG	6.01	118.61	112.60
1	A	347	CYS	N-CA-C	5.98	119.94	109.48
1	A	205	ASP	CA-CB-CG	5.98	118.58	112.60
2	B	177	VAL	CB-CA-C	5.92	116.27	111.06
1	A	142	GLY	N-CA-C	5.85	122.63	113.86
2	B	60	LYS	N-CA-C	5.85	119.44	110.20
1	A	327	ASP	CA-CB-CG	5.85	118.45	112.60
4	K	25	ARG	N-CA-C	-5.81	100.72	108.74
2	B	197	ASN	N-CA-C	5.79	121.24	113.72
3	D	12	SER	CA-C-N	5.76	128.27	120.38
3	D	12	SER	C-N-CA	5.76	128.27	120.38
2	B	31	ASP	CA-CB-CG	5.74	118.33	112.60
4	K	189	ASN	N-CA-C	-5.73	108.09	114.62
4	K	236	GLU	N-CA-C	5.63	118.61	108.48
4	K	281	LYS	CA-C-N	5.59	128.09	120.54
4	K	281	LYS	C-N-CA	5.59	128.09	120.54
4	K	189	ASN	CA-C-N	5.51	128.58	120.82
4	K	189	ASN	C-N-CA	5.51	128.58	120.82
1	A	364	PRO	N-CA-C	5.40	123.60	112.47
4	K	300	ILE	N-CA-CB	5.38	117.14	111.00
2	B	74	THR	CA-C-N	5.33	127.43	120.28
2	B	74	THR	C-N-CA	5.33	127.43	120.28
1	A	309	HIS	CA-CB-CG	5.33	119.12	113.80
4	K	320	GLN	CA-C-N	5.24	128.07	120.79
4	K	320	GLN	C-N-CA	5.24	128.07	120.79
3	D	154	ILE	CA-C-N	5.19	127.23	120.28
3	D	154	ILE	C-N-CA	5.19	127.23	120.28
2	B	2	ARG	N-CA-C	5.16	119.03	111.34
1	A	353	VAL	N-CA-C	5.15	115.38	108.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	PHE	N-CA-C	5.14	117.79	109.40
2	B	127	GLU	CB-CG-CD	5.13	121.32	112.60
3	D	163	GLU	CA-C-N	5.12	127.48	120.46
3	D	163	GLU	C-N-CA	5.12	127.48	120.46
3	D	72	ASP	CA-CB-CG	5.12	117.72	112.60
4	K	266	SER	CA-C-N	5.12	127.44	120.54
4	K	266	SER	C-N-CA	5.12	127.44	120.54
2	B	313	LEU	N-CA-C	-5.08	105.20	111.40
4	K	185	GLU	CA-C-N	5.07	125.56	119.94
4	K	185	GLU	C-N-CA	5.07	125.56	119.94
2	B	284	ARG	CA-C-N	5.06	131.20	121.54
2	B	284	ARG	C-N-CA	5.06	131.20	121.54
2	B	72	PRO	CA-C-N	5.05	125.69	120.03
2	B	72	PRO	C-N-CA	5.05	125.69	120.03
3	D	133	LEU	CA-C-N	5.05	127.01	120.44
3	D	133	LEU	C-N-CA	5.05	127.01	120.44
1	A	283	HIS	CA-C-N	5.04	131.17	121.54
1	A	283	HIS	C-N-CA	5.04	131.17	121.54
3	D	30	VAL	N-CA-C	-5.01	105.82	110.53

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	3349	0	3264	44	0
2	B	3378	0	3253	36	0
3	D	1185	0	1171	7	0
4	K	2593	0	2563	35	0
5	A	32	0	12	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	K	1	0	0	0	0
7	B	28	0	12	1	0
8	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	K	5	0	0	0	0
9	K	27	0	12	0	0
10	K	2	0	0	0	0
All	All	10607	0	10287	115	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 6.

All (115) 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)
4:K:88:SER:HB3	4:K:198:ASN:HB2	1.61	0.83
1:A:221:ARG:HG2	2:B:325:MET:HB3	1.64	0.80
1:A:71:GLU:HB3	1:A:98:ASP:HB3	1.70	0.74
4:K:194:VAL:HG13	4:K:199:GLU:HA	1.71	0.72
2:B:281:GLN:HA	2:B:282:GLN:CB	2.20	0.71
1:A:88:HIS:HB3	1:A:91:GLN:HE21	1.56	0.71
1:A:262:TYR:HB2	1:A:265:ILE:HD12	1.72	0.71
2:B:1:MET:HB3	2:B:131:CYS:SG	2.31	0.70
1:A:265:ILE:HD13	1:A:435:VAL:HG21	1.73	0.70
2:B:66:ILE:HG12	2:B:121:VAL:HG13	1.75	0.69
1:A:70:LEU:HD13	1:A:145:THR:HG23	1.74	0.68
1:A:209:ILE:HD11	1:A:302:MET:HE2	1.77	0.67
2:B:6:HIS:HE1	2:B:8:GLN:HG3	1.62	0.64
4:K:122:MET:HE3	4:K:128:PHE:HZ	1.63	0.63
1:A:108:TYR:HD2	1:A:112:LYS:HZ1	1.45	0.62
2:B:263:PRO:O	2:B:266:HIS:HD2	1.83	0.61
2:B:1:MET:HG2	2:B:133:GLN:HG2	1.81	0.61
2:B:206:ASN:HD21	7:B:501:GDP:HN22	1.49	0.61
2:B:50:ASN:O	2:B:64:ARG:NH2	2.33	0.61
4:K:94:THR:HG22	4:K:107:ILE:HD11	1.82	0.61
1:A:344:VAL:HG23	1:A:347:CYS:HB2	1.82	0.60
1:A:70:LEU:HD13	1:A:145:THR:CG2	2.32	0.60
4:K:105:MET:HB2	4:K:109:PRO:HG2	1.84	0.59
1:A:50:ASN:O	1:A:64:ARG:NH1	2.36	0.59
2:B:1:MET:CB	2:B:131:CYS:SG	2.90	0.59
2:B:288:VAL:HG22	2:B:323:MET:HE3	1.85	0.58
2:B:6:HIS:CE1	2:B:8:GLN:HG3	2.39	0.58
1:A:103:TYR:CD2	1:A:413:MET:HE1	2.38	0.57
2:B:133:GLN:NE2	2:B:252:LEU:H	2.03	0.57
1:A:292:THR:HG22	1:A:335:ILE:HD12	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:ILE:HG21	1:A:302:MET:SD	2.45	0.57
4:K:322:ALA:HA	4:K:325:ILE:HD12	1.87	0.57
2:B:6:HIS:HE1	2:B:8:GLN:HE21	1.52	0.56
1:A:7:ILE:HG23	1:A:66:VAL:HG13	1.87	0.56
1:A:103:TYR:HD2	1:A:413:MET:HE1	1.69	0.56
2:B:312:TYR:CE2	2:B:377:PHE:HZ	2.24	0.56
4:K:284:ARG:HH11	4:K:284:ARG:HG2	1.71	0.56
3:D:34:MET:HA	3:D:34:MET:HE2	1.87	0.55
4:K:19:ASN:O	4:K:23:VAL:HG23	2.06	0.55
4:K:289:SER:HA	4:K:294:ALA:HB3	1.87	0.55
2:B:172:MET:HG3	2:B:387:LEU:HD11	1.88	0.55
1:A:256:GLN:O	1:A:260:VAL:HG22	2.08	0.53
1:A:292:THR:HG22	1:A:335:ILE:CD1	2.39	0.53
1:A:270:ALA:HB3	1:A:302:MET:HG3	1.91	0.52
2:B:139:HIS:HD2	2:B:146:GLY:O	1.91	0.52
1:A:265:ILE:CD1	1:A:435:VAL:HG21	2.39	0.52
1:A:167:LEU:HG	1:A:200:CYS:HB3	1.92	0.52
2:B:106:GLY:O	2:B:111:GLY:HA3	2.10	0.52
2:B:295:MET:CG	2:B:377:PHE:HB2	2.40	0.51
2:B:434:GLN:NE2	4:K:274:TYR:HB3	2.26	0.51
1:A:88:HIS:HD2	1:A:90:GLU:H	1.57	0.51
4:K:295:ARG:HG2	4:K:329:VAL:HG21	1.93	0.51
4:K:131:LYS:HG3	4:K:211:ASN:HB3	1.93	0.50
1:A:271:THR:CG2	1:A:301:GLN:HG2	2.41	0.50
4:K:156:HIS:O	4:K:163:PRO:HA	2.11	0.49
4:K:193:ALA:HB3	4:K:203:ARG:HD2	1.93	0.49
1:A:241:SER:HB2	1:A:249:ASN:O	2.13	0.49
1:A:401:LYS:HG3	2:B:346:TRP:CE3	2.48	0.49
2:B:323:MET:HE3	2:B:373:MET:HE3	1.95	0.49
1:A:119:LEU:HD11	1:A:156:ARG:HB3	1.94	0.49
2:B:6:HIS:CE1	2:B:8:GLN:HE21	2.29	0.48
4:K:66:ALA:C	4:K:68:LYS:H	2.21	0.48
3:D:118:HIS:HD2	3:D:152:ILE:HD11	1.78	0.48
4:K:247:VAL:HA	4:K:250:GLU:OE1	2.14	0.47
1:A:217:LEU:HD21	1:A:368:LEU:HD23	1.96	0.47
4:K:311:GLU:H	4:K:311:GLU:HG2	1.36	0.47
1:A:346:TRP:HZ2	1:A:435:VAL:HG13	1.79	0.47
1:A:2:ARG:HH11	1:A:2:ARG:HA	1.80	0.47
2:B:434:GLN:NE2	4:K:274:TYR:CB	2.78	0.47
2:B:269:MET:HG3	2:B:303:ALA:HB3	1.97	0.47
4:K:136:GLU:HB2	4:K:204:SER:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:HIS:HE1	1:A:155:GLU:OE2	1.98	0.46
2:B:221:THR:HG21	3:D:20:GLU:CD	2.40	0.46
3:D:61:GLU:O	3:D:65:VAL:HG23	2.16	0.46
1:A:344:VAL:HG12	1:A:436:GLY:HA3	1.98	0.46
4:K:194:VAL:HG22	4:K:199:GLU:O	2.15	0.46
4:K:264:VAL:HG21	4:K:283:THR:HB	1.97	0.46
2:B:171:VAL:HA	2:B:204:ILE:O	2.15	0.46
4:K:80:THR:HB	4:K:296:THR:HG23	1.98	0.46
1:A:172:TYR:HB3	1:A:205:ASP:HA	1.98	0.45
3:D:117:LEU:HD11	3:D:129:VAL:HG13	1.97	0.45
4:K:284:ARG:HG2	4:K:284:ARG:NH1	2.32	0.45
2:B:56:ALA:HB3	2:B:60:LYS:HG3	1.99	0.44
1:A:336:LYS:HA	1:A:341:ILE:HD11	1.99	0.44
1:A:141:PHE:CE2	1:A:170:SER:HB3	2.53	0.44
2:B:133:GLN:HE22	2:B:252:LEU:H	1.64	0.44
2:B:332:MET:O	2:B:335:VAL:HG12	2.18	0.44
1:A:280:LYS:HE3	1:A:283:HIS:CB	2.48	0.44
4:K:62:TYR:HD1	4:K:107:ILE:HG23	1.82	0.44
1:A:106:GLY:O	1:A:111:GLY:HA3	2.18	0.43
1:A:406:HIS:CG	2:B:263:PRO:HD3	2.53	0.43
2:B:83:PHE:O	2:B:86:ILE:HG22	2.18	0.43
1:A:123:ARG:HH21	1:A:160:ASP:HB3	1.83	0.43
1:A:328:VAL:HG11	1:A:353:VAL:HG11	1.99	0.43
1:A:1:MET:N	1:A:3:GLU:OE2	2.49	0.43
4:K:173:VAL:HG21	4:K:179:VAL:HG22	1.99	0.43
3:D:34:MET:HE1	3:D:40:VAL:HG13	2.01	0.42
2:B:423:SER:HB3	4:K:278:ARG:HH22	1.83	0.42
2:B:295:MET:HG3	2:B:377:PHE:HB2	2.00	0.42
2:B:295:MET:HE2	2:B:295:MET:HB3	1.96	0.42
4:K:102:PRO:HA	4:K:105:MET:HE3	2.02	0.42
4:K:99:LEU:HD11	4:K:187:LYS:HD2	2.01	0.42
4:K:214:GLN:HG3	4:K:223:LEU:HB2	2.01	0.42
4:K:203:ARG:HH22	4:K:236:GLU:CD	2.27	0.42
2:B:345:GLU:H	2:B:345:GLU:HG2	1.61	0.42
1:A:192:HIS:CG	1:A:421:ALA:HA	2.55	0.41
1:A:209:ILE:HG22	1:A:227:LEU:HD22	2.01	0.41
1:A:320:ARG:HG3	1:A:360:PRO:HD3	2.03	0.41
3:D:96:VAL:HG11	3:D:128:ILE:HG23	2.02	0.41
4:K:209:LEU:HD11	4:K:226:LYS:HG2	2.02	0.41
1:A:2:ARG:HA	1:A:2:ARG:NH1	2.36	0.41
4:K:74:LEU:HD13	4:K:122:MET:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:10:LYS:O	4:K:297:THR:HA	2.21	0.40
4:K:69:ILE:HG13	4:K:81:ILE:HD11	2.02	0.40
4:K:132:VAL:HG12	4:K:210:ILE:HA	2.03	0.40

There are no symmetry-related clashes.

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	426/451 (94%)	406 (95%)	18 (4%)	2 (0%)	24	59
2	B	429/445 (96%)	416 (97%)	8 (2%)	5 (1%)	10	42
3	D	157/169 (93%)	155 (99%)	2 (1%)	0	100	100
4	K	331/349 (95%)	299 (90%)	23 (7%)	9 (3%)	4	25
All	All	1343/1414 (95%)	1276 (95%)	51 (4%)	16 (1%)	10	42

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	284	GLU
2	B	282	GLN
2	B	285	ALA
4	K	7	SER
4	K	27	ASP
1	A	47	ASP
2	B	283	TYR
2	B	284	ARG
4	K	6	GLU
4	K	99	LEU
4	K	273	THR
4	K	26	GLY

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Mol	Chain	Res	Type
4	K	293	ASN
2	B	279	GLY
4	K	188	SER
4	K	275	VAL

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	359/379 (95%)	338 (94%)	21 (6%)	18	51
2	B	370/383 (97%)	357 (96%)	13 (4%)	32	64
3	D	121/131 (92%)	111 (92%)	10 (8%)	10	38
4	K	291/309 (94%)	262 (90%)	29 (10%)	7	30
All	All	1141/1202 (95%)	1068 (94%)	73 (6%)	16	48

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	38	SER
1	A	56	THR
1	A	60	LYS
1	A	68	VAL
1	A	82	THR
1	A	84	ARG
1	A	114	ILE
1	A	120	ASP
1	A	121	ARG
1	A	125	LEU
1	A	179	THR
1	A	188	ILE
1	A	241	SER
1	A	302	MET
1	A	324	VAL
1	A	336	LYS

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Mol	Chain	Res	Type
1	A	345	ASP
1	A	347	CYS
1	A	349	THR
1	A	435	VAL
2	B	24	ILE
2	B	115	VAL
2	B	121	VAL
2	B	122	VAL
2	B	171	VAL
2	B	177	VAL
2	B	284	ARG
2	B	294	GLN
2	B	299	LYS
2	B	323	MET
2	B	325	MET
2	B	405	LEU
2	B	423	SER
3	D	34	MET
3	D	40	VAL
3	D	45	ASP
3	D	100	LEU
3	D	106	VAL
3	D	115	THR
3	D	142	GLN
3	D	164	ILE
3	D	167	LYS
3	D	168	LEU
4	K	6	GLU
4	K	8	ASN
4	K	21	SER
4	K	57	SER
4	K	60	GLN
4	K	86	GLN
4	K	107	ILE
4	K	115	ILE
4	K	140	ASP
4	K	146	LEU
4	K	148	VAL
4	K	155	VAL
4	K	171	ARG
4	K	174	SER
4	K	190	ARG

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Mol	Chain	Res	Type
4	K	192	VAL
4	K	209	LEU
4	K	219	THR
4	K	226	LYS
4	K	237	LYS
4	K	256	LYS
4	K	289	SER
4	K	293	ASN
4	K	299	VAL
4	K	311	GLU
4	K	312	THR
4	K	313	LYS
4	K	314	SER
4	K	331	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	88	HIS
1	A	91	GLN
1	A	139	HIS
1	A	197	HIS
1	A	216	ASN
1	A	256	GLN
1	A	301	GLN
1	A	329	ASN
2	B	6	HIS
2	B	8	GLN
2	B	14	ASN
2	B	50	ASN
2	B	133	GLN
2	B	139	HIS
2	B	206	ASN
2	B	266	HIS
2	B	294	GLN
2	B	433	GLN
2	B	434	GLN
4	K	34	GLN
4	K	63	ASN
4	K	78	ASN
4	K	93	HIS

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Mol	Chain	Res	Type
4	K	125	ASN
4	K	191	HIS
4	K	287	GLN
4	K	332	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 8 ligands modelled in this entry, 3 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	$\# Z > 2$
7	GDP	B	501	6,8	29,30,30	0.46	0	45,47,47	0.75	1 (2%)
8	ALF	B	503	7	4,4,4	1.41	0	-		
8	ALF	K	403	9	4,4,4	1.32	0	-		
5	GTP	A	600	6	33,34,34	2.17	2 (6%)	50,54,54	0.47	0
9	ADP	K	401	6,8	28,29,29	0.44	0	43,45,45	0.63	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
7	GDP	B	501	6,8	-	3/16/32/32	0/3/3/3
9	ADP	K	401	6,8	-	0/16/32/32	0/3/3/3
5	GTP	A	600	6	-	7/22/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	600	GTP	PA-O3A	-8.72	1.50	1.59
5	A	600	GTP	PB-O3B	-8.45	1.50	1.59

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	501	GDP	O3A-PB-O1B	-2.02	100.43	111.04

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	600	GTP	C5'-O5'-PA-O3A
5	A	600	GTP	C5'-O5'-PA-O1A
5	A	600	GTP	C5'-O5'-PA-O2A
7	B	501	GDP	C5'-O5'-PA-O3A
7	B	501	GDP	C5'-O5'-PA-O1A
7	B	501	GDP	C5'-O5'-PA-O2A
5	A	600	GTP	PB-O3B-PG-O3G
5	A	600	GTP	PB-O3A-PA-O2A
5	A	600	GTP	PB-O3B-PG-O1G
5	A	600	GTP	PB-O3A-PA-O1A

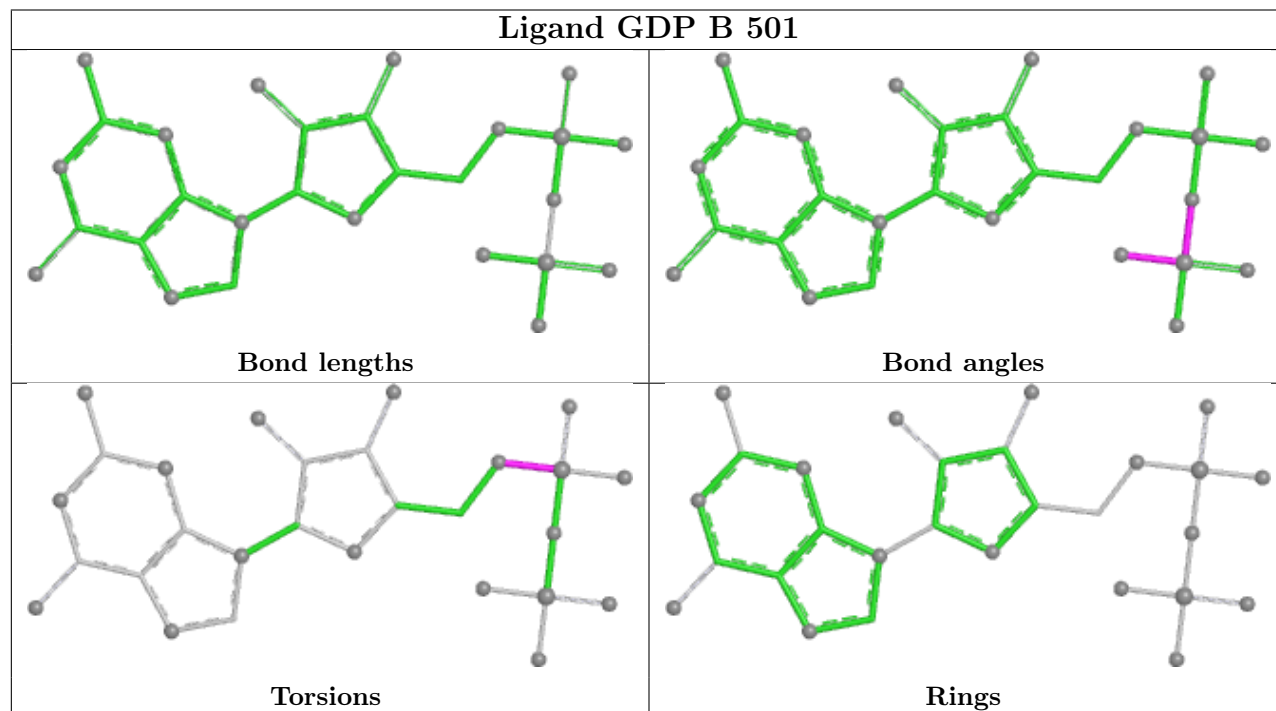
There are no ring outliers.

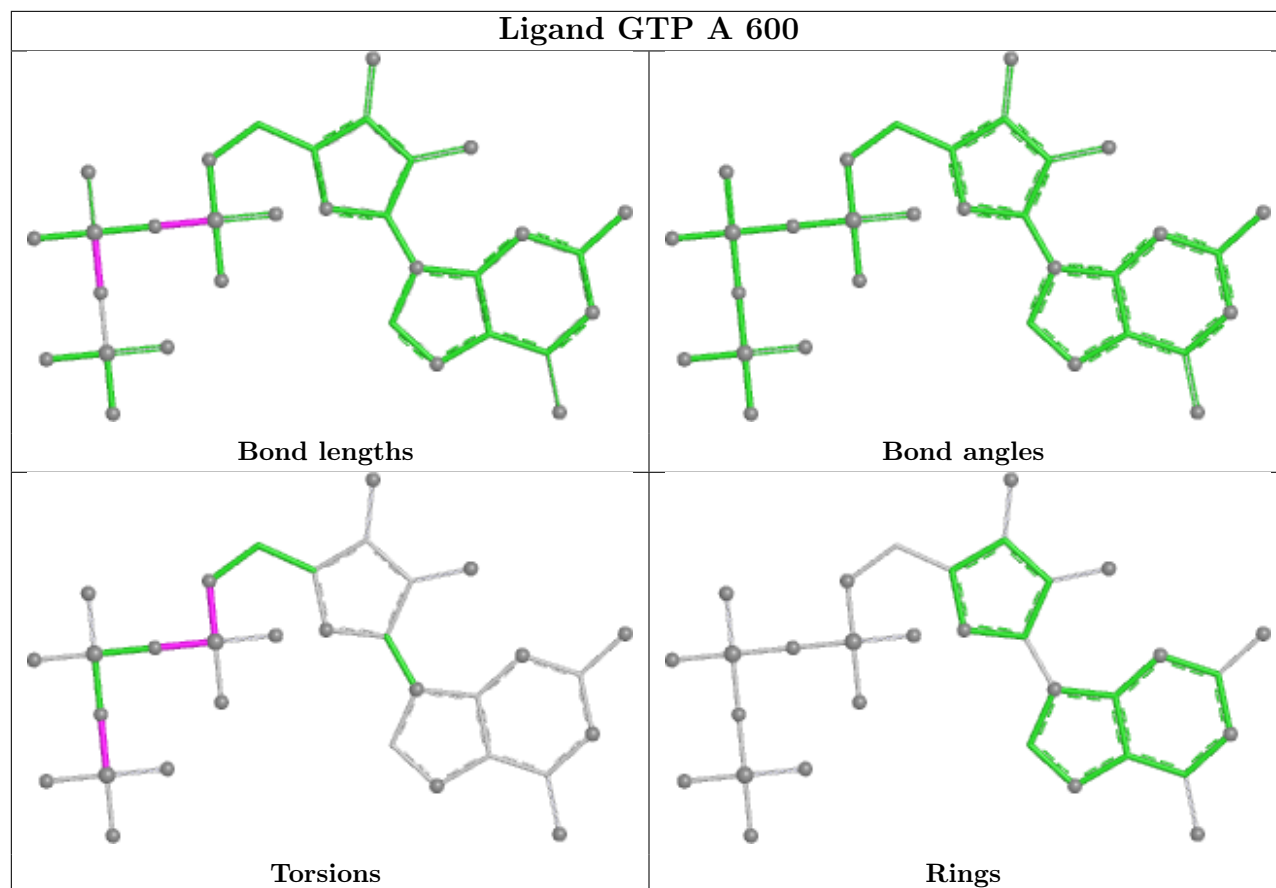
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	501	GDP	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	430/451 (95%)	0.30	10 (2%) 61 41	45, 87, 139, 161	0
2	B	431/445 (96%)	0.19	7 (1%) 70 51	42, 71, 117, 154	0
3	D	159/169 (94%)	0.18	3 (1%) 66 46	45, 75, 114, 158	0
4	K	333/349 (95%)	-0.05	3 (0%) 81 64	44, 70, 118, 144	0
All	All	1353/1414 (95%)	0.17	23 (1%) 69 49	42, 75, 127, 161	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	K	337	ALA	5.3
3	D	12	SER	3.1
3	D	169	ASN	2.9
4	K	5	ALA	2.9
3	D	11	GLY	2.7
1	A	282	TYR	2.7
1	A	326	LYS	2.6
2	B	279	GLY	2.6
4	K	334	GLU	2.4
2	B	57	THR	2.4
2	B	439	THR	2.4
1	A	320	ARG	2.4
2	B	74	THR	2.3
2	B	1	MET	2.2
1	A	74	VAL	2.2
1	A	38	SER	2.2
1	A	285	GLN	2.2
2	B	298	SER	2.1
1	A	360	PRO	2.1
1	A	412	GLY	2.1
1	A	265	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	46	ASP	2.0
2	B	161	TYR	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](#)

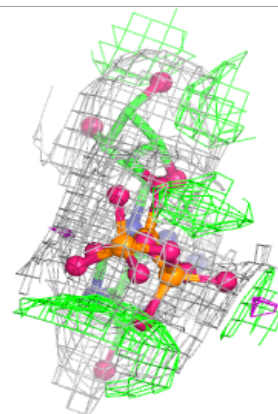
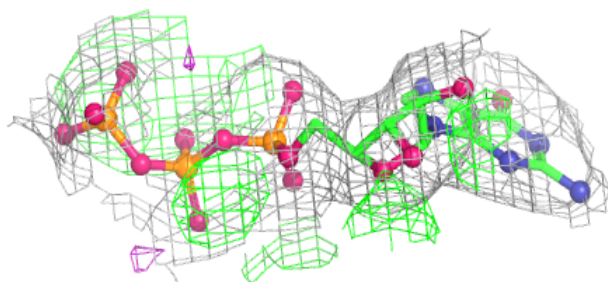
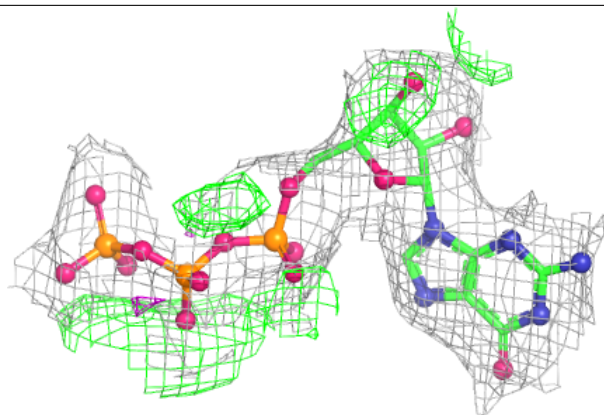
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	ALF	B	503	5/5	0.85	0.15	86,93,98,99	0
6	MG	A	601	1/1	0.87	0.29	28,28,28,28	0
6	MG	B	502	1/1	0.92	0.20	55,55,55,55	0
5	GTP	A	600	32/32	0.92	0.10	51,63,70,74	0
7	GDP	B	501	28/28	0.94	0.10	64,71,79,85	0
8	ALF	K	403	5/5	0.97	0.08	50,51,54,55	0
9	ADP	K	401	27/27	0.97	0.08	55,60,65,69	0
6	MG	K	402	1/1	1.00	0.04	27,27,27,27	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.

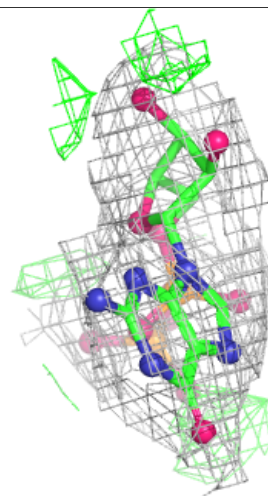
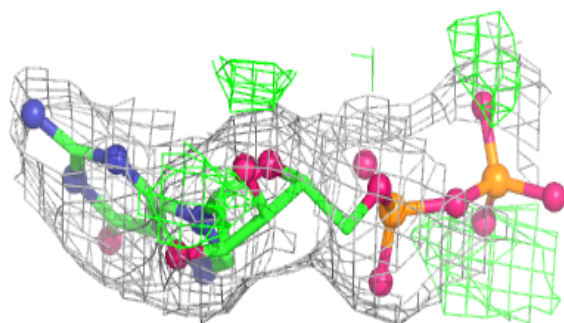
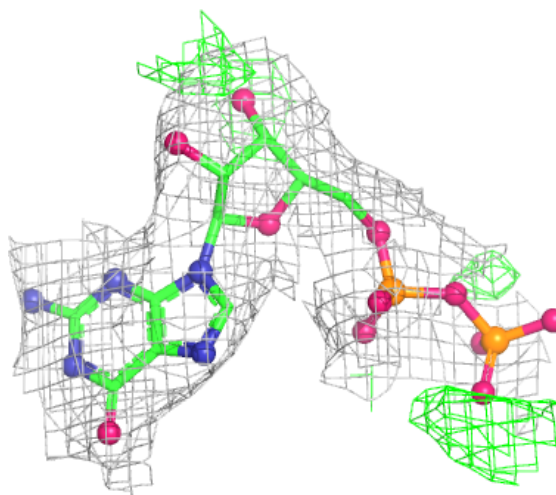
Electron density around GTP A 600:

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



Electron density around GDP B 501:

$2mF_o-DF_c$ (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.