



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 01:33 AM UTC

PDB ID : 8HQ6 / pdb_00008hq6
Title : KL2 in complex with CRM1-Ran-RanBP1
Authors : Sun, Q.; Jian, L.
Deposited on : 2022-12-13
Resolution : 2.03 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

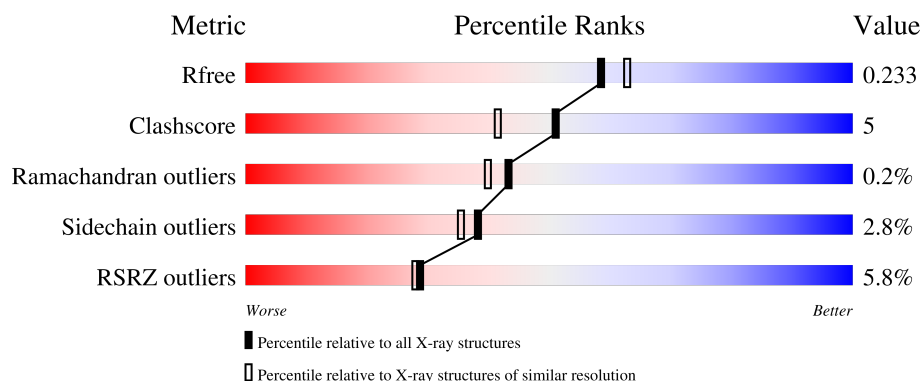
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	216	9% 86% 10% .
2	B	140	8% 83% 9% 8%
3	C	1003	5% 86% 12% ..

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 11668 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 GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	1	0
			1671	1078	286	301	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	69	LEU	GLN	engineered mutation	UNP P62826
A	182	ALA	LEU	engineered mutation	UNP P62826

- Molecule 2 is a protein called YRB1 isoform 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	129	Total	C	N	O	S	0	0	0
			1054	669	182	198	5			

- Molecule 3 is a protein called CRM1 isoform 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	995	Total	C	N	O	S	0	4	0
			8072	5188	1328	1512	44			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP A0A6A5PZI8
C	-1	GLY	-	expression tag	UNP A0A6A5PZI8
C	0	SER	-	expression tag	UNP A0A6A5PZI8
C	27	GLU	SER	engineered mutation	UNP A0A6A5PZI8
C	49	GLU	GLN	engineered mutation	UNP A0A6A5PZI8
C	51	VAL	ALA	engineered mutation	UNP A0A6A5PZI8
C	?	-	VAL	deletion	UNP A0A6A5PZI8
C	?	-	GLN	deletion	UNP A0A6A5PZI8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	ARG	deletion	UNP A0A6A5PZI8
C	?	-	LEU	deletion	UNP A0A6A5PZI8
C	?	-	PRO	deletion	UNP A0A6A5PZI8
C	?	-	ALA	deletion	UNP A0A6A5PZI8
C	?	-	THR	deletion	UNP A0A6A5PZI8
C	?	-	GLU	deletion	UNP A0A6A5PZI8
C	?	-	MET	deletion	UNP A0A6A5PZI8
C	?	-	SER	deletion	UNP A0A6A5PZI8
C	?	-	PRO	deletion	UNP A0A6A5PZI8
C	?	-	LEU	deletion	UNP A0A6A5PZI8
C	?	-	ILE	deletion	UNP A0A6A5PZI8
C	?	-	GLN	deletion	UNP A0A6A5PZI8
C	?	-	LEU	deletion	UNP A0A6A5PZI8
C	?	-	SER	deletion	UNP A0A6A5PZI8
C	?	-	VAL	deletion	UNP A0A6A5PZI8
C	?	-	GLY	deletion	UNP A0A6A5PZI8
C	?	-	SER	deletion	UNP A0A6A5PZI8
C	?	-	GLN	deletion	UNP A0A6A5PZI8
C	?	-	ALA	deletion	UNP A0A6A5PZI8
C	?	-	ILE	deletion	UNP A0A6A5PZI8
C	?	-	SER	deletion	UNP A0A6A5PZI8
C	?	-	THR	deletion	UNP A0A6A5PZI8
C	?	-	GLY	deletion	UNP A0A6A5PZI8
C	?	-	SER	deletion	UNP A0A6A5PZI8
C	?	-	GLY	deletion	UNP A0A6A5PZI8
C	?	-	ALA	deletion	UNP A0A6A5PZI8
C	?	-	LEU	deletion	UNP A0A6A5PZI8
C	?	-	ASN	deletion	UNP A0A6A5PZI8
C	?	-	PRO	deletion	UNP A0A6A5PZI8
C	?	-	GLU	deletion	UNP A0A6A5PZI8
C	?	-	TYR	deletion	UNP A0A6A5PZI8
C	?	-	MET	deletion	UNP A0A6A5PZI8
C	?	-	LYS	deletion	UNP A0A6A5PZI8
C	?	-	ARG	deletion	UNP A0A6A5PZI8
C	?	-	PHE	deletion	UNP A0A6A5PZI8
C	?	-	VAL	deletion	UNP A0A6A5PZI8
C	?	-	LEU	deletion	UNP A0A6A5PZI8
C	?	-	VAL	deletion	UNP A0A6A5PZI8
C	?	-	VAL	deletion	UNP A0A6A5PZI8
C	?	-	GLU	deletion	UNP A0A6A5PZI8
C	?	-	ASN	deletion	UNP A0A6A5PZI8
C	?	-	ASP	deletion	UNP A0A6A5PZI8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	GLU	deletion	UNP A0A6A5PZI8
C	?	-	GLY	deletion	UNP A0A6A5PZI8
C	?	-	GLU	deletion	UNP A0A6A5PZI8
C	?	-	ILE	deletion	UNP A0A6A5PZI8
C	?	-	VAL	deletion	UNP A0A6A5PZI8
C	?	-	ARG	deletion	UNP A0A6A5PZI8
C	?	-	GLU	deletion	UNP A0A6A5PZI8
C	?	-	PHE	deletion	UNP A0A6A5PZI8
C	?	-	VAL	deletion	UNP A0A6A5PZI8
C	?	-	LYS	deletion	UNP A0A6A5PZI8
C	?	-	GLU	deletion	UNP A0A6A5PZI8
C	?	-	SER	deletion	UNP A0A6A5PZI8
C	?	-	ASP	deletion	UNP A0A6A5PZI8
C	?	-	THR	deletion	UNP A0A6A5PZI8
C	537	GLY	ASP	engineered mutation	UNP A0A6A5PZI8
C	539	CYS	THR	engineered mutation	UNP A0A6A5PZI8
C	540	GLU	VAL	engineered mutation	UNP A0A6A5PZI8
C	541	GLN	LYS	engineered mutation	UNP A0A6A5PZI8
C	553	ARG	SER	engineered mutation	UNP A0A6A5PZI8
C	561	GLU	GLN	engineered mutation	UNP A0A6A5PZI8
C	741	THR	ALA	engineered mutation	UNP A0A6A5PZI8
C	1022	CYS	TYR	engineered mutation	UNP A0A6A5PZI8

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0

- Molecule 5 is NITRATE ION (CCD ID: NO3) (formula: NO₃).



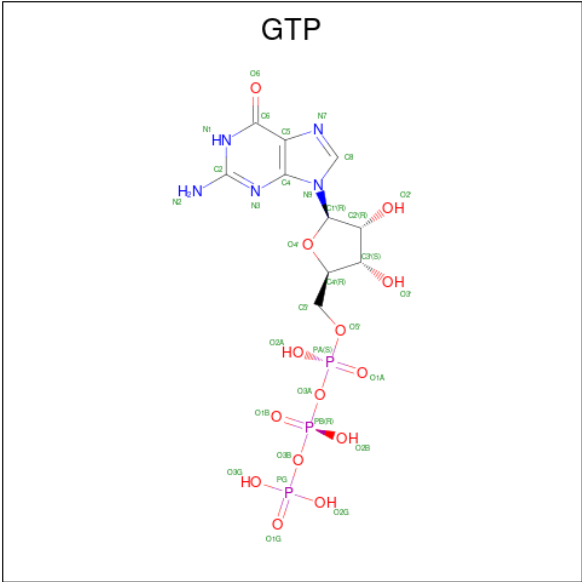
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	N	O	0	0
			4	1	3		
5	A	1	Total	N	O	0	0
			4	1	3		
5	A	1	Total	N	O	0	0
			4	1	3		
5	C	1	Total	N	O	0	0
			4	1	3		
5	C	1	Total	N	O	0	0
			4	1	3		
5	C	1	Total	N	O	0	0
			4	1	3		
5	C	1	Total	N	O	0	0
			4	1	3		
5	C	1	Total	N	O	0	0
			4	1	3		
5	C	1	Total	N	O	0	0
			4	1	3		
5	C	1	Total	N	O	0	0
			4	1	3		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



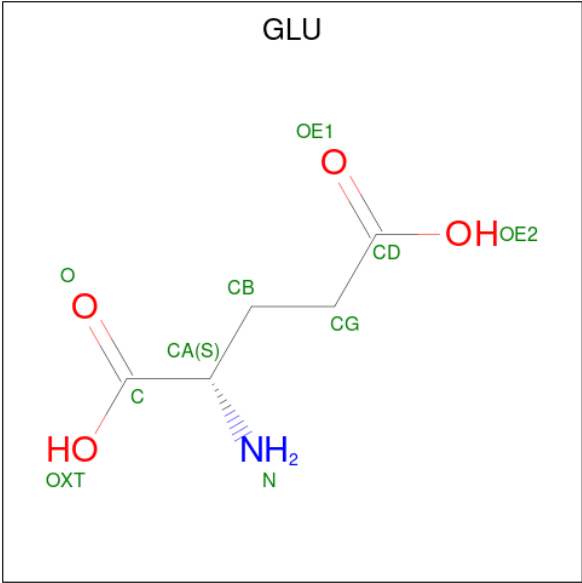
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



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

- Molecule 8 is GLUTAMIC ACID (CCD ID: GLU) (formula: C₅H₉NO₄).

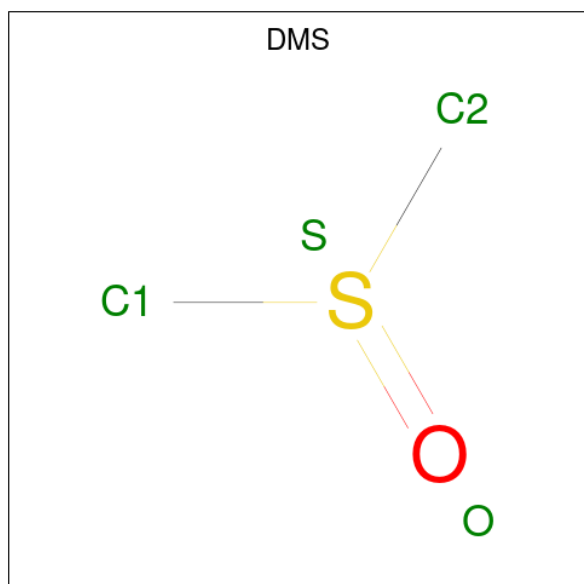


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	C	1	Total	C	N	O	0	0
			9	5	1	3		

- Molecule 9 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

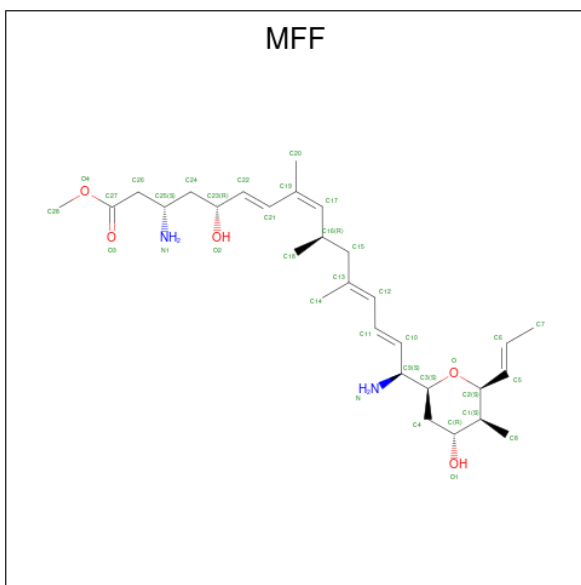
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	C	2	Total Cl 2 2	0	0

- Molecule 10 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	C	1	Total C O S 4 2 1 1	0	0
10	C	1	Total C O S 4 2 1 1	0	0
10	C	1	Total C O S 4 2 1 1	0	0
10	C	1	Total C O S 4 2 1 1	0	0

- Molecule 11 is methyl (3 {S},5 {R},6 {E},8 {Z},10 {R},12 {E},14 {E},16 {S})-3,16-bis(azanyl)-8,10,12-trimethyl-16-[(2 {S},4 {R},5 {S},6 {S})-5-methyl-4-oxidanyl-6-[({E})-prop-1-enyl]oxan-2-yl]-5-oxidanyl-hexadeca-6,8,12,14-tetraenoate (CCD ID: MFF) (formula: $C_{29}H_{48}N_2O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	C	1	Total	C	N	O	0	0
			36	29	2	5		

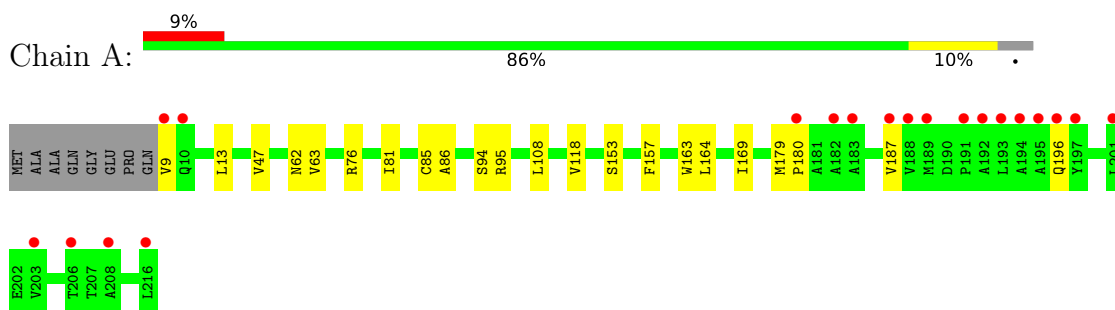
- Molecule 12 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	A	134	Total O 134 134	0	0
12	B	42	Total O 42 42	0	0
12	C	513	Total O 513 513	0	0

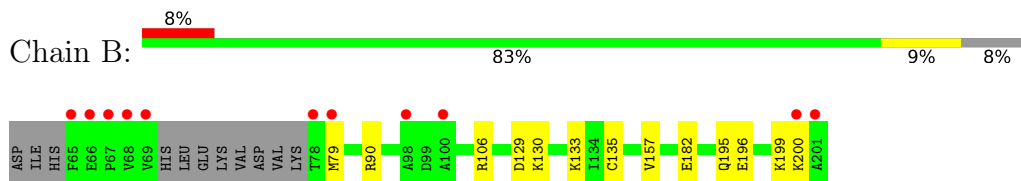
3 Residue-property plots [i](#)

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.

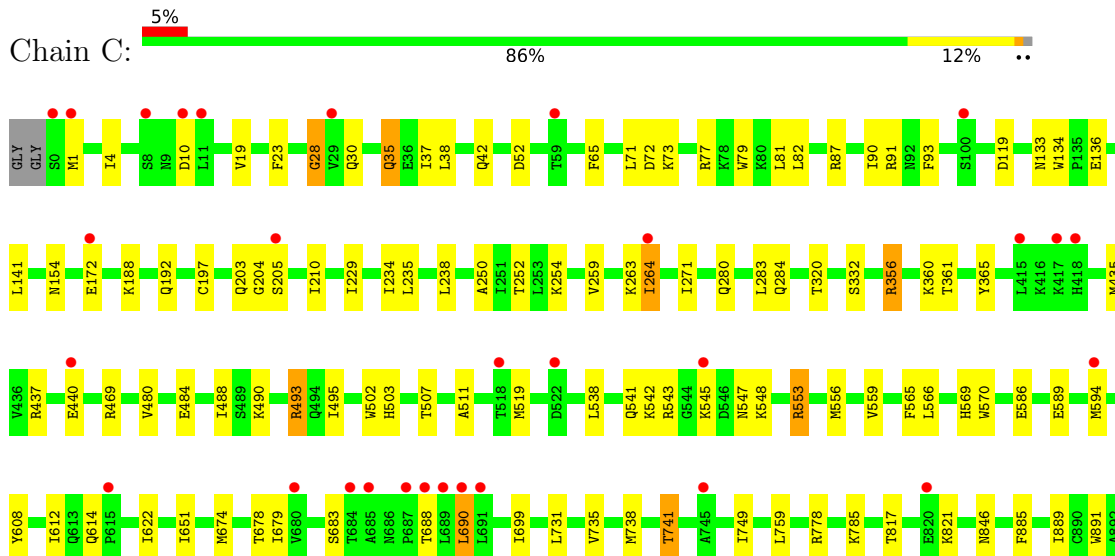
- Molecule 1: GTP-binding nuclear protein Ran

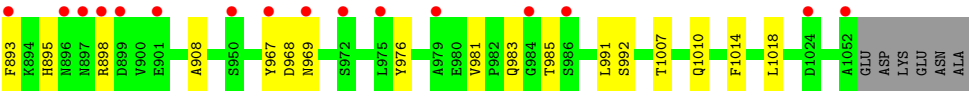


- Molecule 2: YRB1 isoform 1



- Molecule 3: CRM1 isoform 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	104.75Å 104.75Å 302.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.19 – 2.03 99.00 – 2.03	Depositor EDS
% Data completeness (in resolution range)	99.8 (99.19-2.03) 99.8 (99.00-2.03)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.03Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.191 , 0.224 0.206 , 0.233	Depositor DCC
R_{free} test set	5503 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.316	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11668	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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, MG, GOL, MFF, DMS, CL, NO3

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	0.96	0/1712	1.28	0/2321
2	B	0.92	0/1072	1.26	0/1431
3	C	0.98	0/8228	1.41	6/11148 (0.1%)
All	All	0.97	0/11012	1.37	6/14900 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	C	0	1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	263	LYS	CB-CA-C	5.80	118.96	110.26
3	C	28	GLY	CA-C-N	5.35	127.30	120.56
3	C	28	GLY	C-N-CA	5.35	127.30	120.56
3	C	741	THR	CA-CB-OG1	-5.11	101.94	109.60
3	C	361	THR	CA-CB-OG1	-5.06	102.01	109.60
3	C	817	THR	CA-CB-OG1	-5.03	102.05	109.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	469	ARG	Sidechain

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	1671	0	1671	17	1
2	B	1054	0	1052	10	0
3	C	8072	0	8158	96	1
4	A	1	0	0	0	0
5	A	12	0	0	0	0
5	C	44	0	0	2	0
6	A	6	0	8	1	0
6	C	24	0	32	2	0
7	A	32	0	12	0	0
8	C	9	0	5	0	0
9	C	2	0	0	0	0
10	C	16	0	24	3	0
11	C	36	0	0	1	0
12	A	134	0	0	3	0
12	B	42	0	0	1	0
12	C	513	0	0	13	0
All	All	11668	0	10962	115	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 5.

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)
3:C:556[A]:MET:HE3	3:C:594[A]:MET:HG3	1.30	1.09
3:C:556[A]:MET:HE3	3:C:594[A]:MET:CG	1.83	1.07
3:C:503:HIS:O	3:C:507:THR:HG23	1.71	0.90
1:A:94:SER:HB2	3:C:898:ARG:HH21	1.39	0.88
3:C:556[A]:MET:CE	3:C:594[A]:MET:HG3	2.06	0.84
3:C:556[A]:MET:CE	3:C:594[A]:MET:CG	2.55	0.83
3:C:356:ARG:HA	6:C:1106:GOL:H12	1.63	0.80
3:C:495:ILE:HG23	3:C:541:GLN:HE22	1.48	0.79
3:C:52[A]:ASP:OD1	3:C:93:PHE:CE2	2.42	0.72
3:C:556[A]:MET:HE3	3:C:594[A]:MET:HG2	1.71	0.71
3:C:556[A]:MET:HE1	3:C:594[A]:MET:HB3	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:197[A]:CYS:SG	12:C:1294:HOH:O	2.54	0.66
3:C:502:TRP:HE1	3:C:547:ASN:HD22	1.43	0.66
2:B:90:ARG:NH2	12:B:301:HOH:O	2.21	0.66
3:C:23:PHE:CE1	3:C:35:GLN:HG3	2.33	0.64
3:C:250:ALA:HA	12:C:1214:HOH:O	1.99	0.63
3:C:90:ILE:HA	10:C:1115:DMS:H23	1.82	0.62
1:A:179:MET:HE3	2:B:133:LYS:HG2	1.82	0.61
3:C:235:LEU:HD21	3:C:259:VAL:HG11	1.83	0.61
3:C:437:ARG:NH1	5:C:1114:NO3:O2	2.34	0.60
3:C:502:TRP:HE1	3:C:547:ASN:ND2	2.00	0.59
3:C:679:ILE:O	3:C:683:SER:HB3	2.03	0.59
3:C:495:ILE:HG23	3:C:541:GLN:NE2	2.17	0.58
3:C:493:ARG:HD3	12:C:1467:HOH:O	2.02	0.58
3:C:23:PHE:HE1	3:C:35:GLN:HG3	1.70	0.57
3:C:1:MET:HE3	3:C:37:ILE:HD12	1.85	0.57
3:C:79:TRP:NE1	3:C:87:ARG:HD2	2.21	0.56
3:C:846:ASN:HB2	10:C:1109:DMS:H21	1.87	0.56
3:C:556[B]:MET:HE2	11:C:1123:MFF:C20	2.36	0.55
3:C:119:ASP:OD2	3:C:154:ASN:ND2	2.38	0.54
3:C:976:TYR:OH	3:C:985:THR:O	2.24	0.54
3:C:93:PHE:CD1	10:C:1115:DMS:H11	2.43	0.53
3:C:356:ARG:NH2	3:C:440:GLU:OE1	2.42	0.53
1:A:94:SER:HB2	3:C:898:ARG:NH2	2.17	0.53
3:C:967:TYR:HE1	3:C:991:LEU:HD23	1.74	0.52
1:A:153:SER:OG	6:A:304:GOL:H11	2.09	0.52
3:C:893:PHE:HZ	3:C:908:ALA:HB2	1.75	0.52
3:C:542:LYS:O	3:C:548:LYS:HE2	2.10	0.51
3:C:4:ILE:HD11	3:C:37:ILE:HG22	1.91	0.51
3:C:73:LYS:O	3:C:77:ARG:HG3	2.10	0.51
3:C:570:TRP:CD2	3:C:614:GLN:HG3	2.46	0.51
3:C:699:ILE:HG21	3:C:731:LEU:HD21	1.92	0.51
1:A:81:ILE:HD11	3:C:65:PHE:CG	2.46	0.50
3:C:556[A]:MET:HE1	3:C:594[A]:MET:CB	2.41	0.50
3:C:91:ARG:HG3	3:C:134:TRP:CZ3	2.47	0.50
2:B:129:ASP:C	2:B:130:LYS:HG2	2.38	0.49
3:C:891:TRP:CD1	3:C:895:HIS:HE1	2.30	0.49
3:C:735:VAL:HG13	3:C:759:LEU:HB3	1.95	0.49
1:A:157:PHE:CE1	2:B:130:LYS:HG3	2.47	0.49
3:C:204:GLY:HA2	3:C:210:ILE:HD11	1.93	0.49
1:A:9:VAL:HG21	1:A:169:ILE:HD11	1.95	0.48
1:A:76:ARG:HD3	12:A:423:HOH:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:507:THR:HG22	12:C:1215:HOH:O	2.12	0.48
2:B:195:GLN:O	2:B:199:LYS:HG3	2.13	0.48
3:C:1007:THR:H	3:C:1010:GLN:HE21	1.60	0.48
2:B:196:GLU:O	2:B:200:LYS:HD3	2.14	0.48
3:C:553:ARG:HD2	3:C:594[B]:MET:HB2	1.96	0.47
2:B:196:GLU:HA	2:B:199:LYS:HD3	1.96	0.47
3:C:91:ARG:HG3	3:C:134:TRP:CH2	2.50	0.47
3:C:821:LYS:HD2	12:C:1295:HOH:O	2.14	0.47
3:C:480:VAL:CG2	3:C:519:MET:HG2	2.46	0.46
3:C:674:MET:O	3:C:678:THR:HG23	2.15	0.46
3:C:735:VAL:HA	3:C:738:MET:HE3	1.97	0.46
1:A:118:VAL:HG22	1:A:163:TRP:CE3	2.51	0.46
3:C:556[A]:MET:CE	3:C:594[A]:MET:HG2	2.39	0.46
3:C:690:LEU:HG	3:C:738:MET:HE1	1.98	0.46
3:C:141:LEU:C	3:C:141:LEU:HD23	2.40	0.46
3:C:19:VAL:HG12	3:C:38:LEU:HG	1.99	0.45
3:C:360:LYS:HD2	5:C:1101:NO3:O3	2.16	0.45
3:C:622:ILE:HB	3:C:651:ILE:HG21	1.99	0.45
3:C:254:LYS:NZ	12:C:1235:HOH:O	2.49	0.44
3:C:967:TYR:CE1	3:C:991:LEU:HD23	2.53	0.44
3:C:507:THR:CG2	12:C:1215:HOH:O	2.66	0.44
3:C:885:PHE:O	3:C:889:ILE:HG13	2.18	0.44
3:C:72:ASP:OD2	12:C:1201:HOH:O	2.21	0.44
3:C:264:ILE:HG23	3:C:271:ILE:HG21	1.98	0.43
3:C:570:TRP:CE2	3:C:614:GLN:HG3	2.52	0.43
3:C:1014:PHE:CZ	3:C:1018:LEU:HD22	2.54	0.43
1:A:47:VAL:HA	1:A:63:VAL:O	2.19	0.43
1:A:13:LEU:C	1:A:13:LEU:HD23	2.43	0.43
3:C:559:VAL:HG13	3:C:566:LEU:HD21	1.99	0.43
2:B:106:ARG:HD3	2:B:135:CYS:O	2.19	0.43
3:C:356:ARG:HD3	6:C:1106:GOL:O2	2.18	0.43
3:C:968:ASP:O	3:C:969:ASN:HB2	2.18	0.43
3:C:238:LEU:O	3:C:252:THR:HG21	2.19	0.43
12:A:453:HOH:O	3:C:254:LYS:HE3	2.18	0.42
3:C:493:ARG:CD	12:C:1467:HOH:O	2.64	0.42
3:C:192:GLN:HG2	12:C:1581:HOH:O	2.18	0.42
3:C:690:LEU:CG	3:C:738:MET:HE1	2.50	0.42
3:C:82:LEU:O	3:C:87:ARG:NH1	2.52	0.42
3:C:548:LYS:NZ	12:C:1256:HOH:O	2.52	0.42
1:A:86:ALA:CB	1:A:108:LEU:HD21	2.49	0.42
3:C:235:LEU:HD23	3:C:235:LEU:HA	1.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:280:GLN:O	3:C:284:GLN:HB2	2.20	0.42
3:C:679:ILE:O	3:C:683:SER:CB	2.66	0.42
3:C:893:PHE:CZ	3:C:908:ALA:HB2	2.52	0.42
3:C:229:ILE:HA	3:C:234:ILE:CG2	2.50	0.41
3:C:608:TYR:CZ	3:C:612:ILE:HG21	2.55	0.41
3:C:77:ARG:HD2	12:C:1687:HOH:O	2.20	0.41
3:C:480:VAL:HG21	3:C:519:MET:HG2	2.02	0.41
12:A:453:HOH:O	3:C:254:LYS:CE	2.68	0.41
1:A:47:VAL:CG1	1:A:62[A]:ASN:HD21	2.33	0.41
3:C:320:THR:HB	3:C:365:TYR:HB2	2.03	0.41
3:C:565:PHE:O	3:C:569:HIS:HD2	2.03	0.41
3:C:821:LYS:NZ	12:C:1262:HOH:O	2.53	0.41
1:A:85:CYS:HB2	1:A:164:LEU:HD22	2.03	0.41
3:C:435:MET:HG3	3:C:511:ALA:HA	2.02	0.41
1:A:157:PHE:HE1	2:B:130:LYS:HG3	1.85	0.41
3:C:30:GLN:HA	3:C:30:GLN:NE2	2.36	0.41
3:C:1018:LEU:HD12	3:C:1018:LEU:HA	1.96	0.41
3:C:484:GLU:O	3:C:488:ILE:HG12	2.21	0.40
3:C:749:ILE:HD12	3:C:749:ILE:HA	1.94	0.40
3:C:821:LYS:HA	3:C:821:LYS:HD3	1.93	0.40
1:A:81:ILE:HD11	3:C:65:PHE:CD1	2.57	0.40
1:A:180:PRO:HB3	2:B:79:MET:HG3	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)
1:A:95:ARG:NH2	3:C:10:ASP:OD2[5_444]	2.14	0.06

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	207/216 (96%)	200 (97%)	7 (3%)	0	100	100
2	B	125/140 (89%)	118 (94%)	7 (6%)	0	100	100
3	C	997/1003 (99%)	976 (98%)	19 (2%)	2 (0%)	43	40
All	All	1329/1359 (98%)	1294 (97%)	33 (2%)	2 (0%)	43	40

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	205	SER
3	C	28	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	A	180/184 (98%)	178 (99%)	2 (1%)	65	68
2	B	110/121 (91%)	108 (98%)	2 (2%)	51	52
3	C	914/915 (100%)	885 (97%)	29 (3%)	34	30
All	All	1204/1220 (99%)	1171 (97%)	33 (3%)	38	37

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	187	VAL
1	A	196	GLN
2	B	157	VAL
2	B	182	GLU
3	C	35	GLN
3	C	42	GLN
3	C	71	LEU
3	C	81	LEU
3	C	133	ASN
3	C	136	GLU
3	C	172	GLU

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Mol	Chain	Res	Type
3	C	188	LYS
3	C	203	GLN
3	C	264	ILE
3	C	283	LEU
3	C	332	SER
3	C	356	ARG
3	C	490	LYS
3	C	493	ARG
3	C	538	LEU
3	C	543	ARG
3	C	545	LYS
3	C	553	ARG
3	C	586	GLU
3	C	589	GLU
3	C	688	THR
3	C	690	LEU
3	C	741	THR
3	C	778	ARG
3	C	785	LYS
3	C	981	VAL
3	C	983	GLN
3	C	992	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	154	ASN
3	C	9	ASN
3	C	30	GLN
3	C	31	GLN
3	C	35	GLN
3	C	42	GLN
3	C	111	GLN
3	C	203	GLN
3	C	274	GLN
3	C	345	GLN
3	C	426	GLN
3	C	503	HIS
3	C	541	GLN
3	C	547	ASN
3	C	569	HIS
3	C	928	HIS

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Mol	Chain	Res	Type
3	C	1010	GLN

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 29 ligands modelled in this entry, 3 are monoatomic - leaving 26 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	$\# Z > 2$
5	NO3	A	302	-	1,3,3	0.08	0	0,3,3	-	-
11	MFF	C	1123	-	34,36,36	1.17	2 (5%)	37,47,47	2.78	16 (43%)
10	DMS	C	1120	-	3,3,3	0.30	0	3,3,3	0.10	0
5	NO3	C	1122	-	1,3,3	0.48	0	0,3,3	-	-
5	NO3	C	1107	-	1,3,3	0.10	0	0,3,3	-	-
5	NO3	A	303	-	1,3,3	0.57	0	0,3,3	-	-
5	NO3	C	1111	-	1,3,3	0.37	0	0,3,3	-	-
6	GOL	A	304	-	5,5,5	0.10	0	5,5,5	0.29	0
8	GLU	C	1102	-	7,8,9	0.97	0	4,9,11	1.11	0
10	DMS	C	1118	-	3,3,3	0.33	0	3,3,3	0.11	0
6	GOL	C	1117	-	5,5,5	0.12	0	5,5,5	0.37	0
6	GOL	C	1106	-	5,5,5	0.18	0	5,5,5	0.51	0
5	NO3	C	1101	-	1,3,3	1.01	0	0,3,3	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NO3	C	1119	-	1,3,3	0.12	0	0,3,3	-	-
5	NO3	C	1114	-	1,3,3	0.01	0	0,3,3	-	-
7	GTP	A	306	4	33,34,34	0.56	0	50,54,54	0.50	0
10	DMS	C	1115	-	3,3,3	0.25	0	3,3,3	0.15	0
5	NO3	C	1121	-	1,3,3	0.25	0	0,3,3	-	-
5	NO3	C	1105	-	1,3,3	0.76	0	0,3,3	-	-
5	NO3	A	305	-	1,3,3	0.05	0	0,3,3	-	-
5	NO3	C	1113	-	1,3,3	0.02	0	0,3,3	-	-
10	DMS	C	1109	-	3,3,3	0.36	0	3,3,3	0.26	0
5	NO3	C	1112	-	1,3,3	0.01	0	0,3,3	-	-
5	NO3	C	1110	-	1,3,3	0.30	0	0,3,3	-	-
6	GOL	C	1108	-	5,5,5	0.18	0	5,5,5	0.45	0
6	GOL	C	1116	-	5,5,5	0.21	0	5,5,5	0.43	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	C	1108	-	-	2/4/4/4	-
7	GTP	A	306	4	-	3/22/38/38	0/3/3/3
6	GOL	A	304	-	-	2/4/4/4	-
11	MFF	C	1123	-	-	10/33/52/52	0/1/1/1
8	GLU	C	1102	-	-	0/6/7/9	-
6	GOL	C	1117	-	-	4/4/4/4	-
6	GOL	C	1116	-	-	3/4/4/4	-
6	GOL	C	1106	-	-	4/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	C	1123	MFF	O4-C27	4.96	1.48	1.33
11	C	1123	MFF	C21-C19	2.26	1.50	1.46

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	1123	MFF	O4-C27-C26	6.34	125.15	111.47
11	C	1123	MFF	C22-C21-C19	-5.40	114.72	126.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	1123	MFF	C23-C24-C25	5.30	123.10	114.27
11	C	1123	MFF	C24-C23-C22	-5.05	102.06	112.64
11	C	1123	MFF	C21-C19-C17	4.14	129.91	118.82
11	C	1123	MFF	C20-C19-C17	-4.05	111.74	123.21
11	C	1123	MFF	C25-C26-C27	4.03	118.69	112.44
11	C	1123	MFF	C23-C22-C21	3.66	132.35	124.73
11	C	1123	MFF	C11-C12-C13	-3.58	122.66	127.69
11	C	1123	MFF	C12-C11-C10	-3.47	116.39	124.43
11	C	1123	MFF	C28-O4-C27	2.75	124.86	116.07
11	C	1123	MFF	O4-C27-O3	-2.57	115.52	123.20
11	C	1123	MFF	C18-C16-C17	-2.53	103.26	110.04
11	C	1123	MFF	O-C2-C5	-2.44	101.23	108.68
11	C	1123	MFF	C18-C16-C15	2.40	112.59	109.83
11	C	1123	MFF	O3-C27-C26	-2.30	119.30	124.65

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	1106	GOL	O1-C1-C2-C3
6	C	1106	GOL	C1-C2-C3-O3
6	C	1116	GOL	O1-C1-C2-C3
6	C	1117	GOL	O1-C1-C2-C3
6	C	1117	GOL	C1-C2-C3-O3
11	C	1123	MFF	N1-C25-C26-C27
11	C	1123	MFF	C24-C25-C26-C27
11	C	1123	MFF	C23-C24-C25-N1
11	C	1123	MFF	O-C3-C9-C10
11	C	1123	MFF	C4-C3-C9-C10
11	C	1123	MFF	O3-C27-O4-C28
11	C	1123	MFF	C26-C27-O4-C28
6	C	1106	GOL	O1-C1-C2-O2
6	A	304	GOL	C1-C2-C3-O3
6	C	1108	GOL	O1-C1-C2-C3
6	C	1106	GOL	O2-C2-C3-O3
6	C	1108	GOL	O1-C1-C2-O2
6	C	1116	GOL	O1-C1-C2-O2
6	C	1117	GOL	O1-C1-C2-O2
6	C	1117	GOL	O2-C2-C3-O3
7	A	306	GTP	PA-O3A-PB-O1B
11	C	1123	MFF	C13-C15-C16-C18
11	C	1123	MFF	C11-C10-C9-C3

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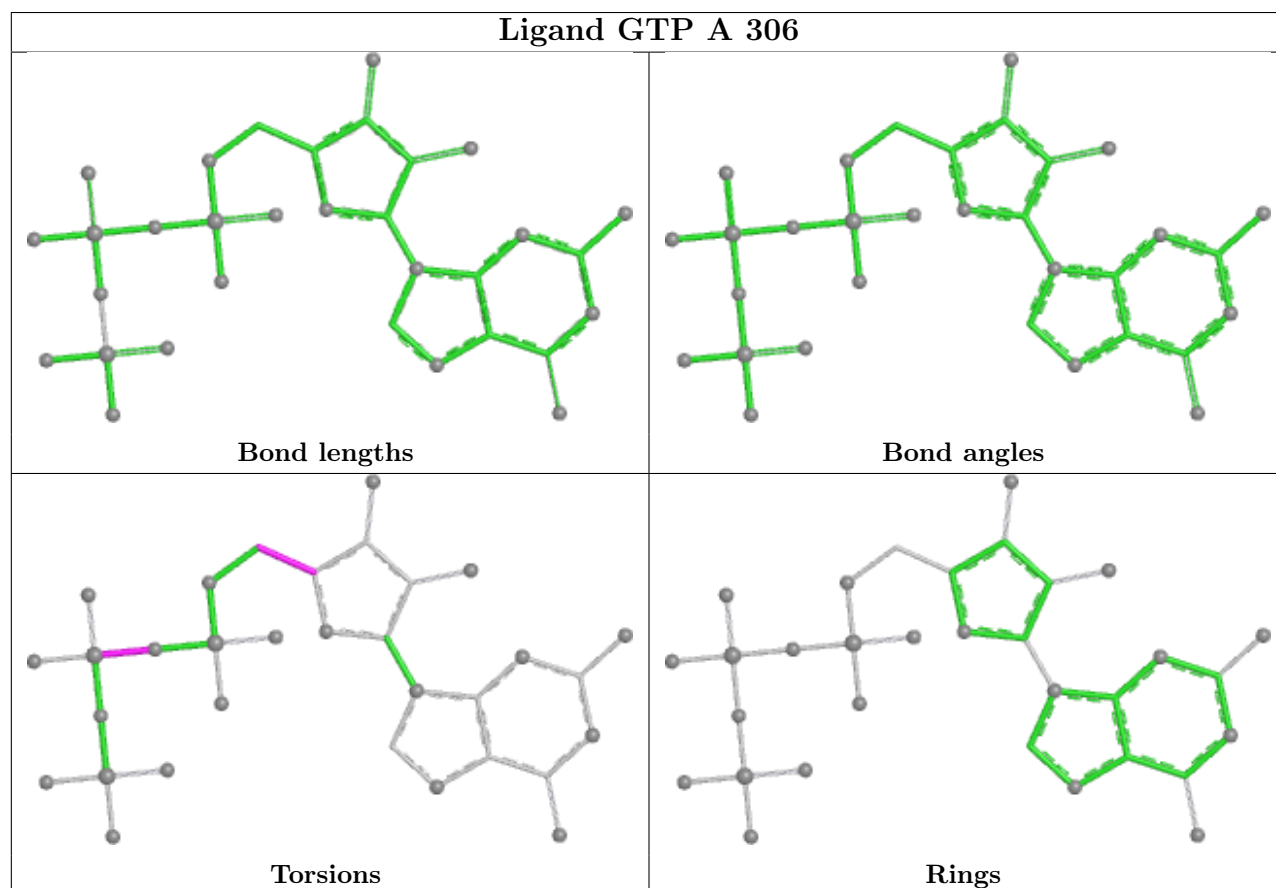
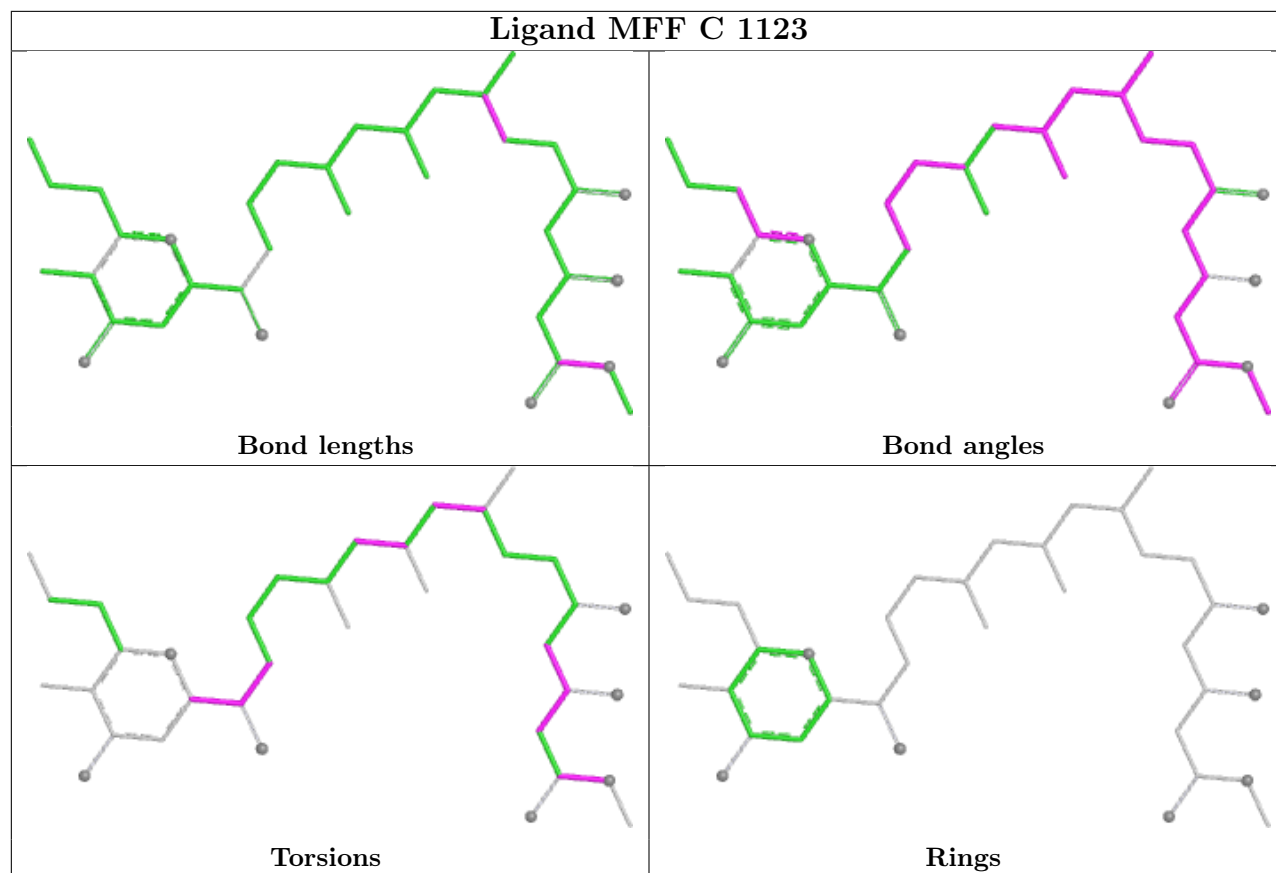
Mol	Chain	Res	Type	Atoms
6	A	304	GOL	O2-C2-C3-O3
7	A	306	GTP	PA-O3A-PB-O2B
11	C	1123	MFF	C16-C17-C19-C20
6	C	1116	GOL	C1-C2-C3-O3
7	A	306	GTP	O4'-C4'-C5'-O5'

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	C	1123	MFF	1	0
6	A	304	GOL	1	0
6	C	1106	GOL	2	0
5	C	1101	NO3	1	0
5	C	1114	NO3	1	0
10	C	1115	DMS	2	0
10	C	1109	DMS	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	208/216 (96%)	0.51	20 (9%)	13 12	22, 42, 103, 118	1 (0%)
2	B	129/140 (92%)	0.71	11 (8%)	16 16	30, 54, 90, 145	0
3	C	995/1003 (99%)	0.46	46 (4%)	37 36	20, 46, 76, 107	4 (0%)
All	All	1332/1359 (98%)	0.49	77 (5%)	29 28	20, 47, 80, 145	5 (0%)

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	201	ALA	9.4
3	C	0	SER	6.8
2	B	69	VAL	6.2
2	B	65	PHE	5.3
2	B	68	VAL	5.1
3	C	898	ARG	5.0
3	C	1052	ALA	4.8
3	C	967	TYR	4.7
1	A	191	PRO	4.5
3	C	691	LEU	4.3
1	A	195	ALA	4.2
3	C	893	PHE	4.1
3	C	522	ASP	4.0
3	C	972	SER	3.9
3	C	689	LEU	3.8
2	B	66	GLU	3.7
1	A	9	VAL	3.6
1	A	188	VAL	3.5
3	C	896	ASN	3.5
1	A	189	MET	3.5
1	A	197	TYR	3.4
2	B	67	PRO	3.4
1	A	193	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
3	C	418	HIS	3.4
3	C	29	VAL	3.3
1	A	194	ALA	3.3
3	C	897	ASN	3.2
3	C	415	LEU	3.2
3	C	687	PRO	3.1
1	A	183	ALA	3.0
2	B	200	LYS	3.0
1	A	180	PRO	3.0
3	C	8	SER	2.9
1	A	192	ALA	2.9
3	C	690	LEU	2.8
3	C	417	LYS	2.8
1	A	187	VAL	2.8
3	C	1	MET	2.7
2	B	79	MET	2.7
2	B	78	THR	2.7
1	A	203	VAL	2.7
3	C	518	THR	2.6
3	C	11	LEU	2.6
3	C	59	THR	2.6
1	A	206	THR	2.5
1	A	216	LEU	2.5
3	C	820	GLU	2.5
2	B	98	ALA	2.4
1	A	201	LEU	2.4
3	C	979	ALA	2.4
3	C	984	GLY	2.4
3	C	899	ASP	2.4
3	C	100	SER	2.4
3	C	205	SER	2.3
3	C	172	GLU	2.3
3	C	545	LYS	2.2
2	B	100	ALA	2.2
3	C	975	LEU	2.2
3	C	615	PRO	2.2
3	C	10	ASP	2.2
1	A	10	GLN	2.2
3	C	986	SER	2.2
3	C	688	THR	2.2
1	A	182	ALA	2.1
3	C	950	SER	2.1

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Mol	Chain	Res	Type	RSRZ
3	C	685	ALA	2.1
3	C	594[A]	MET	2.1
3	C	440	GLU	2.1
3	C	1024	ASP	2.1
3	C	745	ALA	2.1
3	C	901	GLU	2.1
3	C	969	ASN	2.1
3	C	684	THR	2.1
3	C	264	ILE	2.1
1	A	208	ALA	2.1
1	A	196	GLN	2.0
3	C	680	VAL	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($\text{\AA}^2$)	Q<0.9
8	GLU	C	1102	9/10	0.56	0.19	83,87,91,97	0
10	DMS	C	1109	4/4	0.78	0.25	89,90,92,93	0
10	DMS	C	1120	4/4	0.78	0.28	93,94,96,98	0
10	DMS	C	1118	4/4	0.79	0.25	93,94,95,96	0
6	GOL	A	304	6/6	0.80	0.18	58,66,69,71	0
6	GOL	C	1116	6/6	0.81	0.26	84,87,90,93	0
6	GOL	C	1117	6/6	0.81	0.17	82,86,86,87	0
5	NO3	C	1112	4/4	0.82	0.12	69,70,74,75	0
5	NO3	C	1119	4/4	0.83	0.13	80,84,85,86	0
5	NO3	A	305	4/4	0.84	0.13	63,70,72,73	0
10	DMS	C	1115	4/4	0.85	0.24	109,111,112,114	0

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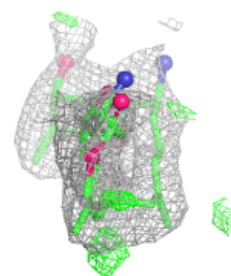
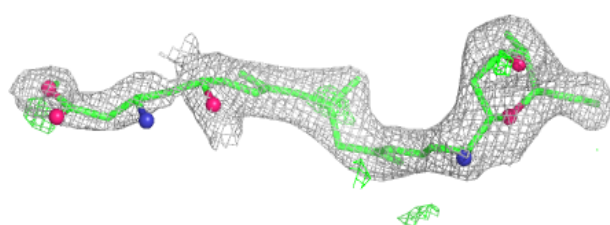
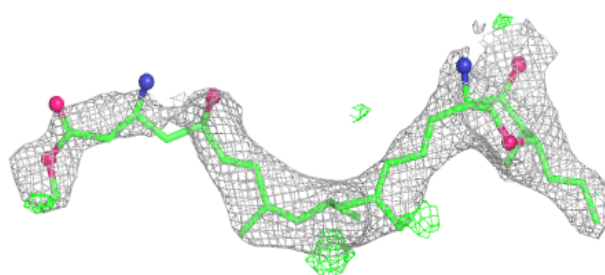
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NO3	C	1114	4/4	0.86	0.12	66,69,70,71	0
5	NO3	C	1122	4/4	0.87	0.10	73,76,77,77	0
5	NO3	C	1113	4/4	0.88	0.14	63,70,73,76	0
6	GOL	C	1108	6/6	0.88	0.15	67,71,72,75	0
5	NO3	C	1105	4/4	0.88	0.24	57,58,58,62	0
11	MFF	C	1123	36/36	0.88	0.16	68,75,89,92	0
5	NO3	C	1101	4/4	0.89	0.30	49,51,55,56	0
5	NO3	A	302	4/4	0.89	0.16	45,47,51,54	0
5	NO3	C	1107	4/4	0.89	0.10	74,83,83,88	0
6	GOL	C	1106	6/6	0.90	0.13	63,65,67,69	0
5	NO3	A	303	4/4	0.90	0.11	62,66,67,73	0
9	CL	C	1103	1/1	0.92	0.09	75,75,75,75	0
5	NO3	C	1110	4/4	0.92	0.10	60,63,67,72	0
5	NO3	C	1111	4/4	0.93	0.10	71,72,74,75	0
7	GTP	A	306	32/32	0.94	0.09	20,20,20,20	0
5	NO3	C	1121	4/4	0.94	0.09	58,62,68,72	0
9	CL	C	1104	1/1	0.96	0.09	67,67,67,67	0
4	MG	A	301	1/1	0.98	0.04	31,31,31,31	0

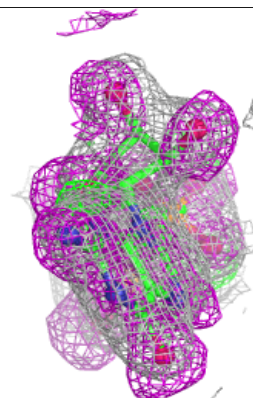
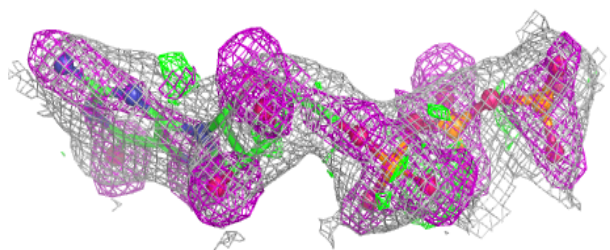
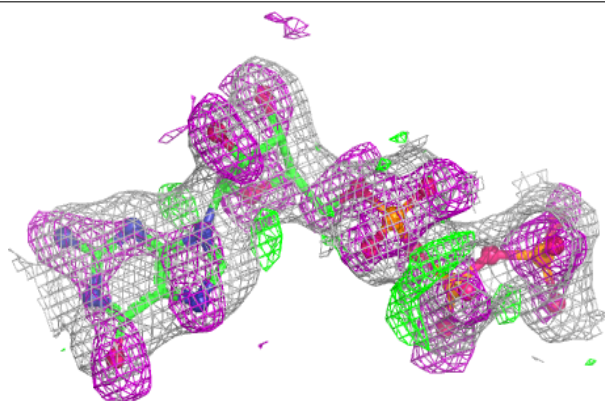
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around MFF C 1123:

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

**Electron density around GTP A 306:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.