



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:40 AM UTC

PDB ID : 3M1I / pdb_00003m1i
Title : Crystal structure of yeast CRM1 (Xpo1p) in complex with yeast RanBP1 (Yrb1p) and yeast RanGTP (Gsp1pGTP)
Authors : Koyama, M.; Matsuura, Y.
Deposited on : 2010-03-05
Resolution : 2.00 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

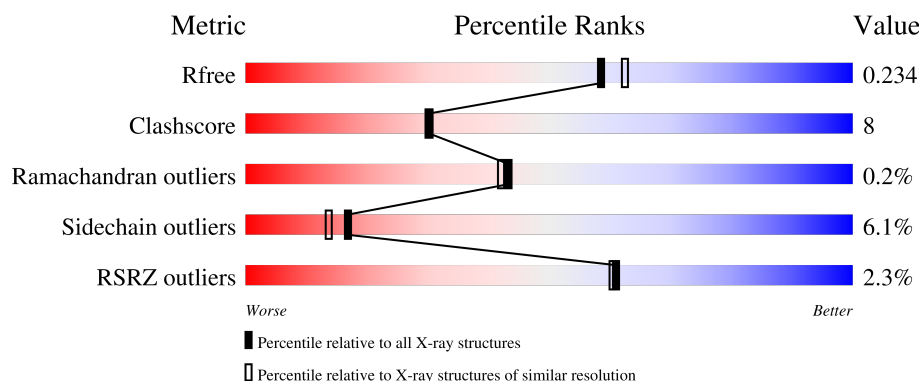
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

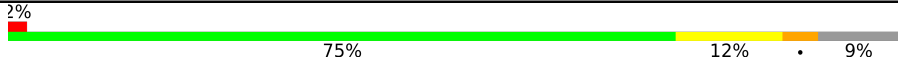
The reported resolution of this entry is 2.00 Å.

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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	219	
2	B	191	
3	C	1049	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11634 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 GSP1/CNR1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1596	1033	268	290	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	71	LEU	GLN	engineered mutation	UNP P32835

- Molecule 2 is a protein called Ran-specific GTPase-activating protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	131	Total	C	N	O	S	0	0	0
			1055	668	184	198	5			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	MET	deletion	UNP P41920
B	?	-	SER	deletion	UNP P41920
B	?	-	SER	deletion	UNP P41920
B	?	-	GLU	deletion	UNP P41920
B	?	-	ASP	deletion	UNP P41920
B	?	-	LYS	deletion	UNP P41920
B	?	-	LYS	deletion	UNP P41920
B	?	-	PRO	deletion	UNP P41920
B	?	-	VAL	deletion	UNP P41920
B	?	-	VAL	deletion	UNP P41920

- Molecule 3 is a protein called Exportin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	1023	Total	C	N	O	S	0	0	0
			8181	5261	1343	1538	39			

There are 39 discrepancies between the modelled and reference sequences:

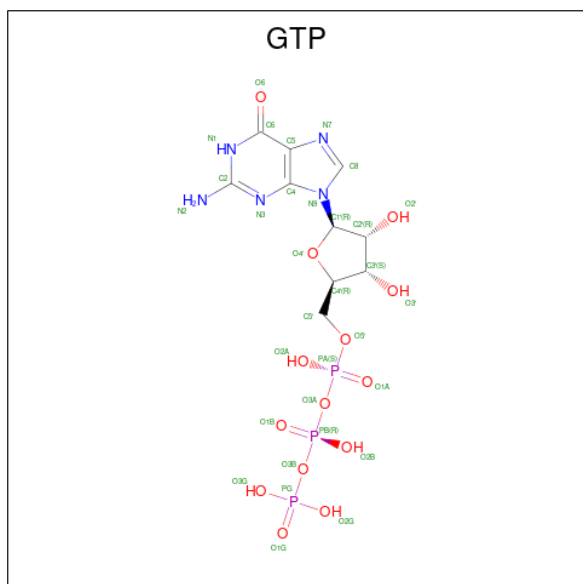
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	expression tag	UNP P30822
C	0	ALA	-	expression tag	UNP P30822
C	?	-	VAL	deletion	UNP P30822
C	?	-	GLN	deletion	UNP P30822
C	?	-	ARG	deletion	UNP P30822
C	?	-	LEU	deletion	UNP P30822
C	?	-	PRO	deletion	UNP P30822
C	?	-	ALA	deletion	UNP P30822
C	?	-	THR	deletion	UNP P30822
C	?	-	GLU	deletion	UNP P30822
C	?	-	MET	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	PRO	deletion	UNP P30822
C	?	-	LEU	deletion	UNP P30822
C	?	-	ILE	deletion	UNP P30822
C	?	-	GLN	deletion	UNP P30822
C	?	-	LEU	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	VAL	deletion	UNP P30822
C	?	-	GLY	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	GLN	deletion	UNP P30822
C	?	-	ALA	deletion	UNP P30822
C	?	-	ILE	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	THR	deletion	UNP P30822
C	?	-	GLY	deletion	UNP P30822
C	?	-	SER	deletion	UNP P30822
C	?	-	GLY	deletion	UNP P30822
C	?	-	ALA	deletion	UNP P30822
C	?	-	LEU	deletion	UNP P30822
C	?	-	ASN	deletion	UNP P30822
C	?	-	PRO	deletion	UNP P30822
C	?	-	GLU	deletion	UNP P30822
C	?	-	TYR	deletion	UNP P30822
C	?	-	MET	deletion	UNP P30822
C	?	-	LYS	deletion	UNP P30822

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	ARG	deletion	UNP P30822
C	?	-	PHE	deletion	UNP P30822

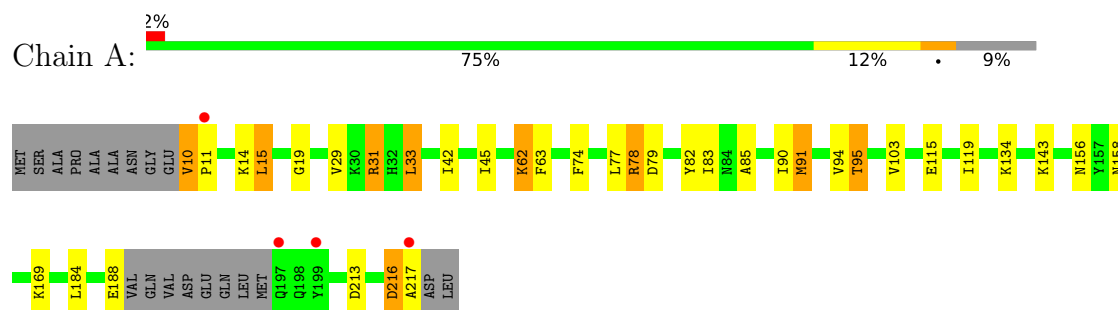
- Molecule 4 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



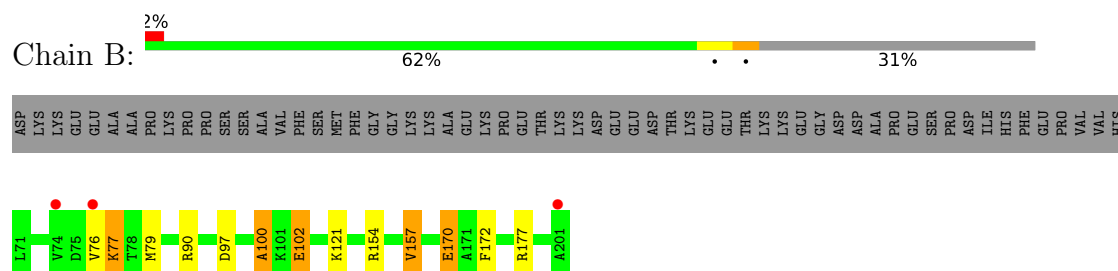
3 Residue-property plots

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 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.

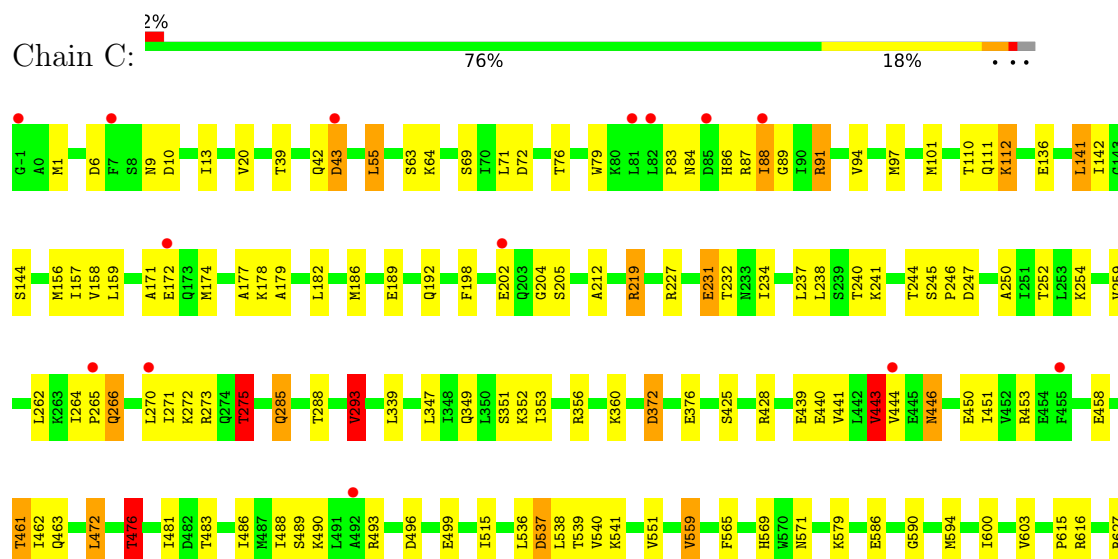
- Molecule 1: GTP-binding nuclear protein GSP1/CNR1

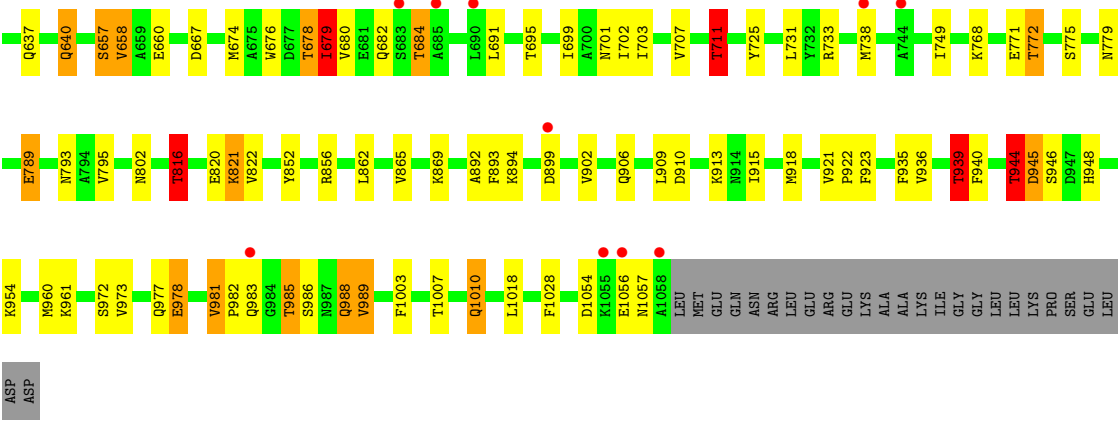


- Molecule 2: Ran-specific GTPase-activating protein 1



- Molecule 3: Exportin-1





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	106.21 Å 106.21 Å 303.60 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.51 – 2.00 43.51 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.51-2.00) 99.9 (43.51-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.00 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.175 , 0.220 (Not available) , 0.234	Depositor DCC
R_{free} test set	5918 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11634	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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: MG, GTP

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	1.51	7/1636 (0.4%)	1.29	9/2219 (0.4%)
2	B	1.18	2/1072 (0.2%)	1.15	4/1433 (0.3%)
3	C	1.40	31/8339 (0.4%)	1.28	43/11317 (0.4%)
All	All	1.40	40/11047 (0.4%)	1.27	56/14969 (0.4%)

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
1	A	1	0
2	B	0	1
3	C	1	1
All	All	2	2

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	103	VAL	CA-CB	8.96	1.59	1.53
3	C	852	TYR	CA-CB	8.60	1.58	1.52
3	C	600	ILE	CA-CB	8.54	1.63	1.54
1	A	119	ILE	C-O	6.84	1.30	1.24
3	C	250	ALA	CA-CB	6.71	1.64	1.53
1	A	94	VAL	CA-CB	6.59	1.62	1.54
1	A	85	ALA	C-O	6.50	1.31	1.23
3	C	795	VAL	CA-CB	6.39	1.62	1.54
3	C	679	ILE	CA-CB	6.27	1.62	1.54
3	C	212	ALA	N-CA	6.18	1.53	1.46
3	C	158	VAL	C-O	-6.17	1.17	1.24
3	C	351	SER	N-CA	6.12	1.54	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	94	VAL	CA-CB	6.08	1.61	1.54
3	C	177	ALA	CA-CB	6.07	1.62	1.53
3	C	69	SER	N-CA	6.04	1.53	1.46
3	C	339	LEU	CA-C	5.94	1.60	1.52
3	C	441	VAL	CA-CB	5.93	1.60	1.53
1	A	31	ARG	CD-NE	-5.92	1.38	1.46
3	C	285	GLN	C-O	-5.76	1.17	1.24
3	C	259	VAL	CA-CB	5.75	1.61	1.54
3	C	481	ILE	CA-CB	5.73	1.61	1.54
3	C	472	LEU	C-O	5.72	1.30	1.24
2	B	90	ARG	CB-CG	-5.72	1.35	1.52
3	C	1010	GLN	CB-CG	-5.72	1.35	1.52
3	C	711	THR	CA-CB	5.67	1.62	1.53
3	C	893	PHE	N-CA	-5.66	1.38	1.46
3	C	640	GLN	CD-OE1	5.65	1.34	1.23
3	C	376	GLU	C-N	5.65	1.41	1.33
3	C	939	THR	CB-CG2	-5.58	1.34	1.52
2	B	157	VAL	CA-CB	5.57	1.62	1.54
3	C	63	SER	C-O	-5.54	1.17	1.24
3	C	667	ASP	CA-C	5.50	1.59	1.52
3	C	353	ILE	C-O	-5.45	1.18	1.24
3	C	20	VAL	CA-C	5.39	1.59	1.52
3	C	936	VAL	CA-CB	5.25	1.60	1.54
3	C	899	ASP	CA-C	5.24	1.59	1.52
1	A	90	ILE	CA-CB	-5.19	1.48	1.54
1	A	19	GLY	N-CA	5.14	1.50	1.45
3	C	89	GLY	N-CA	5.07	1.52	1.45
3	C	288	THR	CA-CB	5.00	1.61	1.53

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	376	GLU	O-C-N	-13.94	108.53	121.35
3	C	945	ASP	CA-C-N	-11.50	105.12	122.83
3	C	945	ASP	C-N-CA	-11.50	105.12	122.83
3	C	376	GLU	CA-C-N	10.13	130.19	119.76
3	C	376	GLU	C-N-CA	10.13	130.19	119.76
3	C	293	VAL	CB-CA-C	-8.85	99.13	112.05
2	B	102	GLU	N-CA-C	7.58	120.89	108.99
1	A	91	MET	CG-SD-CE	7.46	117.31	100.90
3	C	944	THR	N-CA-CB	-7.29	99.94	110.65
3	C	275	THR	OG1-CB-CG2	7.10	123.50	109.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	31	ARG	NE-CZ-NH1	6.97	128.47	121.50
3	C	428	ARG	CG-CD-NE	-6.73	97.20	112.00
3	C	232	THR	CA-C-N	-6.70	111.53	122.65
3	C	232	THR	C-N-CA	-6.70	111.53	122.65
3	C	275	THR	CA-CB-OG1	6.61	119.51	109.60
3	C	789	GLU	CA-C-N	-6.46	112.34	119.19
3	C	789	GLU	C-N-CA	-6.46	112.34	119.19
1	A	95	THR	OG1-CB-CG2	6.45	122.19	109.30
1	A	95	THR	CA-CB-OG1	6.37	119.15	109.60
3	C	141	LEU	N-CA-C	-6.23	104.57	111.36
3	C	559	VAL	CG1-CB-CG2	6.12	124.26	110.80
3	C	219	ARG	CB-CG-CD	5.96	125.00	111.30
3	C	476	THR	OG1-CB-CG2	5.95	121.20	109.30
3	C	893	PHE	CB-CA-C	5.93	121.74	109.95
3	C	893	PHE	N-CA-C	5.91	120.33	113.18
3	C	856	ARG	N-CA-C	-5.81	105.02	111.36
2	B	100	ALA	CA-C-N	-5.81	113.76	122.24
2	B	100	ALA	C-N-CA	-5.81	113.76	122.24
3	C	892	ALA	CA-C-N	-5.77	110.33	121.58
3	C	892	ALA	C-N-CA	-5.77	110.33	121.58
3	C	1010	GLN	CB-CA-C	5.72	120.29	110.79
3	C	682	GLN	N-CA-C	5.68	117.95	111.02
3	C	945	ASP	O-C-N	-5.66	116.27	122.84
3	C	89	GLY	N-CA-C	-5.58	105.63	112.77
3	C	219	ARG	NE-CZ-NH1	5.52	127.02	121.50
3	C	443	VAL	CB-CA-C	5.49	119.16	110.81
2	B	77	LYS	N-CA-C	5.46	118.59	109.96
3	C	91	ARG	CB-CG-CD	-5.46	98.75	111.30
3	C	444	VAL	CB-CA-C	-5.45	101.37	111.18
3	C	701	ASN	N-CA-C	5.40	117.17	111.28
3	C	110	THR	N-CA-C	5.37	119.99	113.38
3	C	725	TYR	N-CA-C	5.37	116.81	111.07
3	C	702	ILE	CB-CA-C	-5.36	105.11	111.97
3	C	944	THR	OG1-CB-CG2	5.35	120.00	109.30
3	C	749	ILE	N-CA-C	5.34	117.72	111.05
1	A	15	LEU	CA-C-N	-5.32	116.59	123.19
1	A	15	LEU	C-N-CA	-5.32	116.59	123.19
3	C	594	MET	CG-SD-CE	-5.15	89.57	100.90
3	C	603	VAL	CG1-CB-CG2	5.11	122.05	110.80
3	C	946	SER	CB-CA-C	5.11	118.98	111.06
3	C	816	THR	N-CA-CB	5.07	117.57	110.12
1	A	103	VAL	CA-C-N	-5.06	114.40	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	VAL	C-N-CA	-5.06	114.40	119.56
3	C	1056	GLU	N-CA-C	-5.05	105.86	112.23
3	C	559	VAL	N-CA-CB	-5.04	103.02	110.58
1	A	158	ASN	CB-CA-C	-5.04	103.47	111.48

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	95	THR	CB
3	C	275	THR	CB

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	76	VAL	Peptide
3	C	204	GLY	Peptide

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	1596	0	1573	18	0
2	B	1055	0	1050	11	0
3	C	8181	0	8201	156	0
4	A	32	0	12	0	0
5	A	1	0	0	0	0
6	A	151	0	0	3	0
6	B	45	0	0	0	0
6	C	573	0	0	29	0
All	All	11634	0	10836	182	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 8.

All (182) 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:43:ASP:HB2	6:C:1551:HOH:O	1.53	1.08
2:B:170:GLU:HG3	2:B:172:PHE:CE2	1.89	1.07
3:C:76:THR:HG22	6:C:1456:HOH:O	1.56	1.05
3:C:174:MET:HE3	3:C:178:LYS:HG2	1.40	1.02
3:C:97:MET:HE2	3:C:101:MET:HE1	1.43	1.00
3:C:156:MET:HE2	3:C:156:MET:HA	1.48	0.95
3:C:97:MET:HE2	3:C:101:MET:CE	1.98	0.93
1:A:217:ALA:HB1	2:B:154:ARG:HH22	1.34	0.92
3:C:171:ALA:O	3:C:172:GLU:HB2	1.75	0.87
3:C:657:SER:HB2	3:C:660:GLU:OE1	1.76	0.86
3:C:101:MET:HE2	6:C:1565:HOH:O	1.76	0.85
3:C:902:VAL:HG13	3:C:954:LYS:HE2	1.59	0.83
3:C:472:LEU:O	3:C:476:THR:HG23	1.80	0.82
3:C:921:VAL:HG13	3:C:922:PRO:HD2	1.62	0.81
3:C:988:GLN:HE21	3:C:988:GLN:H	1.28	0.80
3:C:935:PHE:O	3:C:939:THR:HG23	1.83	0.78
3:C:174:MET:CE	3:C:178:LYS:HG2	2.13	0.78
3:C:1054:ASP:HA	3:C:1057:ASN:HB2	1.68	0.76
2:B:170:GLU:HG3	2:B:172:PHE:CZ	2.21	0.75
3:C:244:THR:HG22	3:C:285:GLN:OE1	1.85	0.74
3:C:1007:THR:H	3:C:1010:GLN:HE21	1.32	0.74
3:C:271:ILE:O	3:C:275:THR:HG23	1.87	0.74
3:C:768:LYS:O	3:C:772:THR:HG23	1.88	0.74
3:C:174:MET:HE2	3:C:179:ALA:N	2.02	0.73
1:A:10:VAL:N	6:A:681:HOH:O	2.20	0.72
1:A:213:ASP:HB3	1:A:216:ASP:HB2	1.72	0.72
1:A:29:VAL:HG12	1:A:33:LEU:HD22	1.73	0.70
3:C:779:ASN:C	3:C:779:ASN:HD22	2.00	0.70
3:C:657:SER:CB	3:C:660:GLU:OE1	2.40	0.69
3:C:349:GLN:HE22	3:C:352:LYS:NZ	1.90	0.69
3:C:174:MET:HE3	3:C:178:LYS:CG	2.21	0.68
3:C:707:VAL:O	3:C:711:THR:HG23	1.94	0.67
3:C:174:MET:CE	3:C:182:LEU:HD12	2.25	0.67
3:C:691:LEU:HD21	3:C:738:MET:HE2	1.75	0.67
3:C:539:THR:HG22	3:C:551:VAL:HG12	1.78	0.66
3:C:816:THR:HG21	6:C:1417:HOH:O	1.95	0.66
3:C:539:THR:HG22	3:C:551:VAL:CG1	2.26	0.66
3:C:72:ASP:O	3:C:76:THR:HG23	1.94	0.65
1:A:31:ARG:HD3	1:A:156:ASN:OD1	1.97	0.64
3:C:101:MET:CE	6:C:1565:HOH:O	2.41	0.63
2:B:170:GLU:HG3	2:B:172:PHE:HE2	1.55	0.63
3:C:238:LEU:O	3:C:252:THR:HG21	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:174:MET:HE2	3:C:178:LYS:C	2.25	0.62
3:C:458:GLU:OE2	3:C:461:THR:HG21	2.00	0.62
3:C:816:THR:CG2	6:C:1417:HOH:O	2.47	0.61
3:C:674:MET:O	3:C:678:THR:HG23	2.01	0.61
3:C:97:MET:HE2	3:C:101:MET:HE3	1.83	0.61
3:C:443:VAL:HG13	3:C:451:ILE:CG2	2.30	0.61
2:B:154:ARG:NH2	2:B:177:ARG:HD3	2.15	0.61
3:C:960:MET:HG3	3:C:1028:PHE:CE2	2.35	0.60
3:C:676:TRP:O	3:C:680:VAL:HG23	2.00	0.60
3:C:918:MET:HE1	3:C:923:PHE:HD2	1.66	0.60
3:C:174:MET:HE1	3:C:182:LEU:HD12	1.83	0.59
3:C:189:GLU:O	3:C:192:GLN:HG2	2.02	0.59
3:C:6:ASP:OD2	3:C:9:ASN:ND2	2.32	0.59
2:B:170:GLU:CG	2:B:172:PHE:CE2	2.78	0.59
3:C:771:GLU:O	3:C:775:SER:HB2	2.03	0.59
3:C:156:MET:HA	3:C:156:MET:CE	2.28	0.58
3:C:453:ARG:HG2	3:C:453:ARG:HH11	1.67	0.58
3:C:940:PHE:O	3:C:944:THR:HB	2.03	0.58
3:C:439:GLU:HB3	3:C:462:ILE:HG12	1.86	0.58
3:C:156:MET:HE2	3:C:156:MET:CA	2.29	0.57
3:C:458:GLU:OE2	3:C:461:THR:CG2	2.52	0.57
3:C:820:GLU:OE1	3:C:821:LYS:HE3	2.03	0.57
1:A:10:VAL:HB	1:A:11:PRO:HD2	1.86	0.57
3:C:171:ALA:O	3:C:172:GLU:CB	2.41	0.57
3:C:982:PRO:O	3:C:985:THR:HG23	2.05	0.56
3:C:446:ASN:ND2	3:C:450:GLU:H	2.02	0.56
3:C:944:THR:HG22	6:C:1215:HOH:O	2.05	0.56
1:A:216:ASP:O	1:A:217:ALA:C	2.49	0.56
3:C:463:GLN:HG2	6:C:1613:HOH:O	2.05	0.56
3:C:156:MET:CE	3:C:159:LEU:HD12	2.36	0.56
3:C:83:PRO:HG2	3:C:86:HIS:CD2	2.41	0.56
3:C:793:ASN:HB2	6:C:1207:HOH:O	2.05	0.55
3:C:1:MET:HB2	6:C:1562:HOH:O	2.07	0.55
3:C:865:VAL:HG23	6:C:1177:HOH:O	2.06	0.55
3:C:266:GLN:O	3:C:272:LYS:HE3	2.07	0.55
3:C:802:ASN:ND2	6:C:1277:HOH:O	2.37	0.55
3:C:43:ASP:OD1	3:C:43:ASP:C	2.51	0.54
3:C:935:PHE:O	3:C:939:THR:CG2	2.53	0.54
3:C:816:THR:HG22	3:C:862:LEU:HB2	1.90	0.54
3:C:493:ARG:HH11	3:C:499:GLU:CD	2.16	0.54
3:C:986:SER:OG	3:C:989:VAL:HG13	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:733:ARG:NH2	3:C:793:ASN:OD1	2.40	0.54
3:C:237:LEU:HD22	3:C:241:LYS:HD3	1.89	0.54
3:C:84:ASN:O	3:C:88:ILE:HG13	2.08	0.54
3:C:349:GLN:HE22	3:C:352:LYS:HZ1	1.56	0.54
3:C:472:LEU:O	3:C:476:THR:CG2	2.52	0.54
3:C:174:MET:CE	3:C:178:LYS:C	2.81	0.53
3:C:738:MET:HE3	3:C:738:MET:HA	1.91	0.53
3:C:453:ARG:HG2	3:C:453:ARG:NH1	2.23	0.53
3:C:1018:LEU:HD22	3:C:1028:PHE:CE1	2.44	0.53
3:C:909:LEU:HD22	3:C:913:LYS:HE3	1.89	0.53
3:C:446:ASN:C	3:C:446:ASN:HD22	2.17	0.53
3:C:174:MET:CE	3:C:179:ALA:HA	2.39	0.52
3:C:227:ARG:O	3:C:231:GLU:HB2	2.10	0.52
3:C:91:ARG:HD3	6:C:1602:HOH:O	2.09	0.52
3:C:768:LYS:O	3:C:772:THR:CG2	2.57	0.52
3:C:174:MET:HE1	3:C:179:ALA:HA	1.91	0.51
3:C:894:LYS:O	3:C:948:HIS:HE1	1.95	0.50
3:C:496:ASP:OD1	3:C:496:ASP:C	2.55	0.50
3:C:678:THR:HG22	6:C:1508:HOH:O	2.11	0.50
3:C:198:PHE:O	3:C:202:GLU:HG2	2.10	0.49
3:C:443:VAL:HG13	3:C:451:ILE:HG23	1.93	0.49
3:C:538:LEU:O	3:C:541:LYS:O	2.30	0.49
1:A:134:LYS:HE2	6:A:761:HOH:O	2.13	0.49
3:C:571:ASN:HB2	6:C:1213:HOH:O	2.11	0.49
3:C:972:SER:HA	6:C:1376:HOH:O	2.11	0.49
2:B:170:GLU:CG	2:B:172:PHE:HE2	2.23	0.49
3:C:349:GLN:HE22	3:C:352:LYS:HZ2	1.58	0.48
3:C:178:LYS:HD3	6:C:1224:HOH:O	2.14	0.48
1:A:217:ALA:CB	2:B:154:ARG:HH22	2.14	0.48
3:C:142:ILE:CD1	3:C:192:GLN:HB2	2.44	0.48
3:C:921:VAL:CG1	3:C:922:PRO:HD2	2.40	0.48
3:C:79:TRP:CE2	3:C:87:ARG:HD3	2.49	0.47
3:C:537:ASP:C	3:C:537:ASP:OD2	2.57	0.47
3:C:446:ASN:HD21	3:C:450:GLU:H	1.62	0.47
3:C:944:THR:CG2	6:C:1215:HOH:O	2.63	0.47
3:C:779:ASN:C	3:C:779:ASN:ND2	2.72	0.46
3:C:270:LEU:HD23	6:C:1618:HOH:O	2.14	0.46
3:C:793:ASN:CB	6:C:1207:HOH:O	2.60	0.46
3:C:678:THR:CG2	6:C:1508:HOH:O	2.64	0.46
2:B:100:ALA:HB3	2:B:102:GLU:HG2	1.98	0.45
3:C:486:ILE:O	3:C:490:LYS:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:156:MET:HE2	3:C:159:LEU:HD12	1.97	0.45
3:C:157:ILE:HD12	6:C:1339:HOH:O	2.17	0.45
3:C:538:LEU:C	3:C:540:VAL:H	2.25	0.45
3:C:906:GLN:NE2	3:C:910:ASP:OD1	2.49	0.45
3:C:91:ARG:CD	6:C:1602:HOH:O	2.62	0.45
3:C:174:MET:CE	3:C:178:LYS:CG	2.89	0.45
3:C:372:ASP:OD1	6:C:1528:HOH:O	2.21	0.45
3:C:699:ILE:HG21	3:C:731:LEU:HD21	1.99	0.45
3:C:679:ILE:HD13	3:C:679:ILE:HA	1.78	0.44
3:C:476:THR:HG21	3:C:515:ILE:HG21	2.00	0.44
3:C:55:LEU:O	3:C:64:LYS:HE2	2.18	0.44
1:A:78:ARG:HD2	6:A:704:HOH:O	2.17	0.44
3:C:43:ASP:OD1	6:C:1359:HOH:O	2.21	0.43
3:C:615:PRO:O	3:C:616:ARG:CB	2.64	0.43
1:A:45:ILE:HA	1:A:74:PHE:HB3	2.01	0.43
3:C:476:THR:HB	3:C:483:THR:HG21	2.00	0.43
3:C:565:PHE:O	3:C:569:HIS:HD2	2.01	0.43
3:C:39:THR:O	3:C:43:ASP:HB3	2.19	0.43
3:C:112:LYS:HA	3:C:112:LYS:HD2	1.80	0.43
3:C:43:ASP:CB	6:C:1551:HOH:O	2.35	0.43
3:C:680:VAL:O	3:C:684:THR:HB	2.18	0.42
3:C:349:GLN:NE2	3:C:352:LYS:NZ	2.64	0.42
3:C:657:SER:OG	3:C:660:GLU:OE1	2.38	0.42
1:A:169:LYS:HD3	1:A:169:LYS:HA	1.84	0.42
2:B:97:ASP:O	2:B:100:ALA:O	2.37	0.42
3:C:174:MET:CE	3:C:179:ALA:N	2.79	0.42
3:C:356:ARG:HG2	3:C:360:LYS:HE3	2.02	0.42
3:C:977:GLN:O	3:C:978:GLU:C	2.62	0.42
3:C:240:THR:HG22	3:C:241:LYS:N	2.35	0.42
3:C:707:VAL:O	3:C:711:THR:CG2	2.63	0.42
3:C:627:ARG:HH11	3:C:627:ARG:HD3	1.75	0.42
1:A:82:TYR:N	1:A:82:TYR:CD1	2.86	0.41
3:C:264:ILE:HA	3:C:265:PRO:HD2	1.78	0.41
3:C:679:ILE:HD11	3:C:695:THR:HG23	2.01	0.41
3:C:156:MET:HE1	3:C:159:LEU:HD12	2.00	0.41
1:A:143:LYS:O	3:C:254:LYS:NZ	2.43	0.41
3:C:981:VAL:HG13	3:C:985:THR:HG23	2.02	0.41
3:C:944:THR:HG23	3:C:1003:PHE:HZ	1.85	0.41
1:A:14:LYS:NZ	1:A:83:ILE:O	2.41	0.41
1:A:62:LYS:HD2	1:A:63:PHE:N	2.36	0.41
3:C:174:MET:CE	3:C:179:ALA:CA	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:347:LEU:HD23	3:C:347:LEU:HA	1.88	0.41
3:C:539:THR:HG22	3:C:551:VAL:HG11	1.99	0.41
3:C:869:LYS:HE3	6:C:1577:HOH:O	2.19	0.41
3:C:988:GLN:HE21	3:C:988:GLN:N	2.07	0.41
3:C:13:ILE:HD12	3:C:13:ILE:HG23	1.75	0.41
3:C:536:LEU:HD23	3:C:536:LEU:HA	1.83	0.41
3:C:182:LEU:HD23	3:C:182:LEU:HA	1.92	0.41
3:C:293:VAL:HG23	6:C:1609:HOH:O	2.21	0.41
2:B:79:MET:HE2	2:B:79:MET:HB3	2.01	0.40
3:C:960:MET:HG3	3:C:1028:PHE:CD2	2.56	0.40
1:A:15:LEU:C	1:A:15:LEU:HD23	2.46	0.40
3:C:245:SER:HA	3:C:246:PRO:HD3	1.99	0.40
1:A:79:ASP:N	1:A:79:ASP:OD2	2.55	0.40
3:C:453:ARG:HG3	3:C:590:GLY:HA2	2.04	0.40
3:C:658:VAL:HG12	6:C:1604:HOH:O	2.21	0.40
3:C:945:ASP:OD1	3:C:945:ASP:C	2.65	0.40
3:C:493:ARG:HD2	3:C:499:GLU:CD	2.46	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	A	196/219 (90%)	191 (97%)	4 (2%)	1 (0%)	24	21
2	B	129/191 (68%)	121 (94%)	8 (6%)	0	100	100
3	C	1021/1049 (97%)	992 (97%)	27 (3%)	2 (0%)	43	42
All	All	1346/1459 (92%)	1304 (97%)	39 (3%)	3 (0%)	43	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	216	ASP
3	C	978	GLU
3	C	983	GLN

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	169/188 (90%)	158 (94%)	11 (6%)	15	12
2	B	108/164 (66%)	104 (96%)	4 (4%)	30	30
3	C	914/955 (96%)	856 (94%)	58 (6%)	16	13
All	All	1191/1307 (91%)	1118 (94%)	73 (6%)	17	14

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

Mol	Chain	Res	Type
1	A	10	VAL
1	A	33	LEU
1	A	42	ILE
1	A	62	LYS
1	A	77	LEU
1	A	78	ARG
1	A	91	MET
1	A	95	THR
1	A	115	GLU
1	A	184	LEU
1	A	188	GLU
2	B	77	LYS
2	B	121	LYS
2	B	157	VAL
2	B	170	GLU
3	C	10	ASP
3	C	42	GLN
3	C	43	ASP
3	C	55	LEU
3	C	71	LEU

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Mol	Chain	Res	Type
3	C	88	ILE
3	C	111	GLN
3	C	112	LYS
3	C	136	GLU
3	C	141	LEU
3	C	144	SER
3	C	186	MET
3	C	205	SER
3	C	219	ARG
3	C	231	GLU
3	C	234	ILE
3	C	247	ASP
3	C	262	LEU
3	C	266	GLN
3	C	273	ARG
3	C	275	THR
3	C	293	VAL
3	C	372	ASP
3	C	425	SER
3	C	440	GLU
3	C	443	VAL
3	C	446	ASN
3	C	461	THR
3	C	476	THR
3	C	488	ILE
3	C	489	SER
3	C	537	ASP
3	C	559	VAL
3	C	579	LYS
3	C	586	GLU
3	C	637	GLN
3	C	640	GLN
3	C	657	SER
3	C	658	VAL
3	C	678	THR
3	C	679	ILE
3	C	684	THR
3	C	703	ILE
3	C	711	THR
3	C	772	THR
3	C	789	GLU
3	C	816	THR

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Mol	Chain	Res	Type
3	C	821	LYS
3	C	822	VAL
3	C	915	ILE
3	C	939	THR
3	C	944	THR
3	C	961	LYS
3	C	973	VAL
3	C	981	VAL
3	C	985	THR
3	C	988	GLN
3	C	989	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	ASN
1	A	147	GLN
1	A	158	ASN
1	A	198	GLN
3	C	31	GLN
3	C	35	GLN
3	C	56	GLN
3	C	86	HIS
3	C	111	GLN
3	C	154	ASN
3	C	173	GLN
3	C	305	ASN
3	C	349	GLN
3	C	446	ASN
3	C	477	HIS
3	C	569	HIS
3	C	779	ASN
3	C	813	ASN
3	C	824	HIS
3	C	906	GLN
3	C	928	HIS
3	C	948	HIS
3	C	988	GLN
3	C	997	ASN
3	C	1005	HIS
3	C	1010	GLN
3	C	1038	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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
4	GTP	A	1177	5	33,34,34	1.88	4 (12%)	50,54,54	1.52	10 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GTP	A	1177	5	-	5/22/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1177	GTP	PB-O3A	-7.56	1.51	1.59
4	A	1177	GTP	PB-O3B	5.04	1.64	1.59
4	A	1177	GTP	C2-N1	2.15	1.42	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1177	GTP	O5'-C5'	-2.01	1.37	1.44

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1177	GTP	C2-N3-C4	3.41	118.17	112.30
4	A	1177	GTP	C5-C4-N3	-3.40	122.98	128.39
4	A	1177	GTP	O3'-C3'-C2'	-2.72	103.11	111.82
4	A	1177	GTP	N9-C4-N3	2.67	131.29	125.95
4	A	1177	GTP	O3B-PG-O1G	-2.49	97.96	111.04
4	A	1177	GTP	O3G-PG-O3B	2.37	112.60	104.64
4	A	1177	GTP	O6-C6-N1	-2.28	115.82	120.11
4	A	1177	GTP	C5-C6-N1	2.28	119.06	113.25
4	A	1177	GTP	N1-C2-N3	-2.25	119.20	123.32
4	A	1177	GTP	O2B-PB-O3A	2.21	113.25	107.27

There are no chirality outliers.

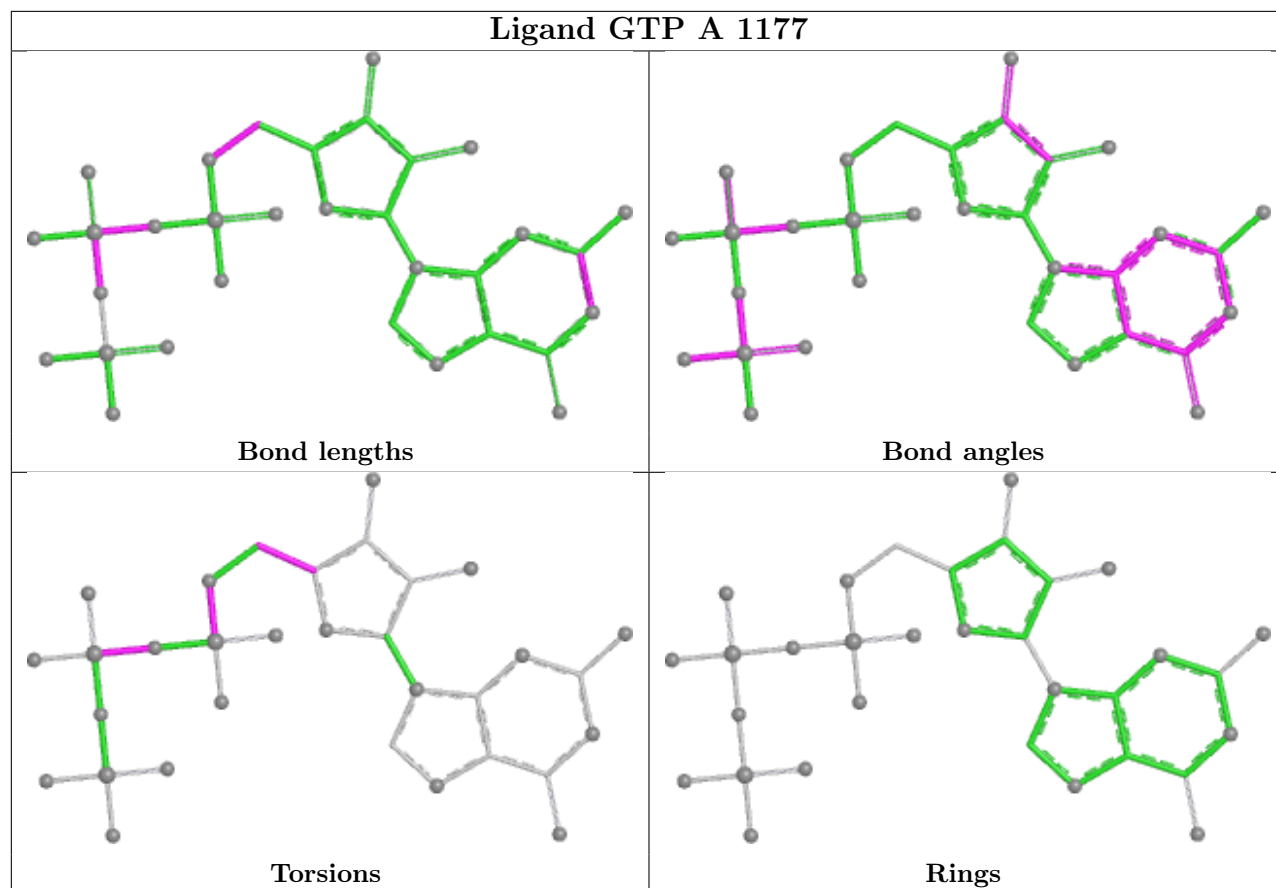
All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1177	GTP	C5'-O5'-PA-O3A
4	A	1177	GTP	C5'-O5'-PA-O1A
4	A	1177	GTP	O4'-C4'-C5'-O5'
4	A	1177	GTP	C3'-C4'-C5'-O5'
4	A	1177	GTP	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

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	200/219 (91%)	-0.18	4 (2%) 65 64	20, 33, 68, 95	0
2	B	131/191 (68%)	0.19	3 (2%) 61 60	30, 47, 80, 110	0
3	C	1023/1049 (97%)	0.00	24 (2%) 61 60	20, 39, 64, 102	0
All	All	1354/1459 (92%)	-0.01	31 (2%) 61 60	20, 39, 67, 110	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	76	VAL	6.2
1	A	217	ALA	3.9
1	A	197	GLN	3.5
3	C	43	ASP	3.3
1	A	199	TYR	3.2
3	C	81	LEU	3.0
3	C	88	ILE	3.0
3	C	683	SER	3.0
3	C	265	PRO	3.0
3	C	1058	ALA	2.9
3	C	899	ASP	2.8
3	C	270	LEU	2.8
3	C	455	PHE	2.7
3	C	85	ASP	2.7
3	C	738	MET	2.7
3	C	172	GLU	2.5
3	C	82	LEU	2.5
3	C	7	PHE	2.5
2	B	201	ALA	2.5
3	C	492	ALA	2.4
3	C	444	VAL	2.4
3	C	-1	GLY	2.4
3	C	685	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
3	C	202	GLU	2.3
2	B	74	VAL	2.3
1	A	11	PRO	2.2
3	C	744	ALA	2.2
3	C	1055	LYS	2.2
3	C	983	GLN	2.1
3	C	1056	GLU	2.1
3	C	690	LEU	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](#)

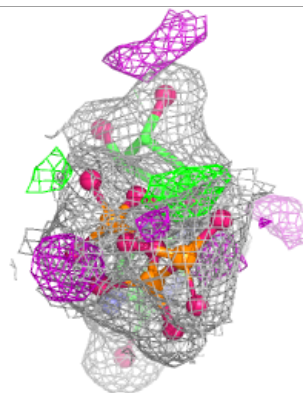
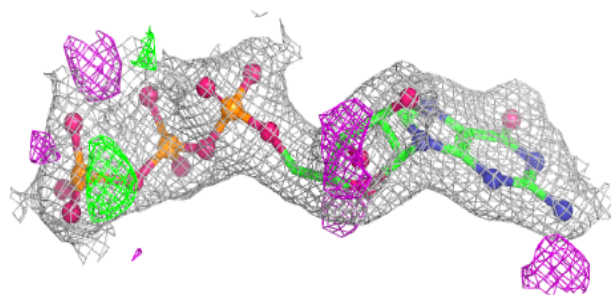
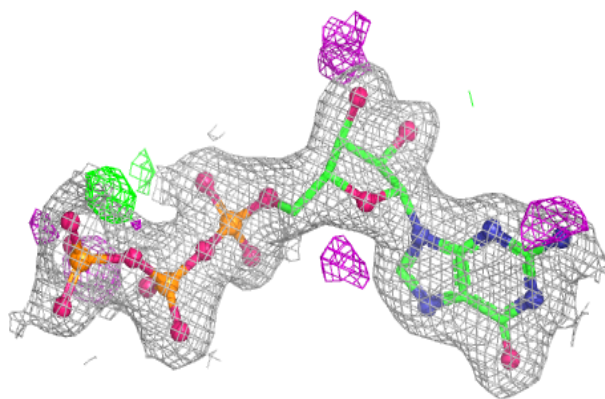
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GTP	A	1177	32/32	0.98	0.05	21,26,31,34	0
5	MG	A	1178	1/1	0.99	0.11	18,18,18,18	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 1177:

$2mF_o-DF_c$ (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.