



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 04:26 am GMT

PDB ID : 5DLQ
Title : Crystal structure of RanGTP-Exportin 4-eIF5A complex
Authors : Aksu, M.; Trakhanov, S.; Gorlich, D.
Deposited on : 2015-09-07
Resolution : 3.20 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

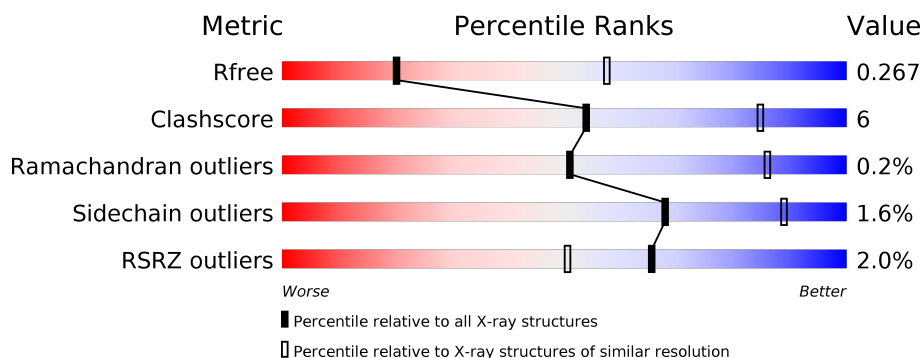
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	1113	<div> <div>3%</div> <div> <div></div> <div>73%</div> <div>16%</div> <div>12%</div> </div> </div>
1	B	1113	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>15%</div> <div>5%</div> </div> </div>
2	C	176	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>16%</div> <div>••</div> </div> </div>
2	D	176	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div>••</div> </div> </div>
3	E	141	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>16%</div> <div>••</div> </div> </div>
3	F	141	<div> <div></div> <div> <div></div> <div>84%</div> <div>13%</div> <div>••</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 20688 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 Exportin-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	1052	Total	C	N	O	S	0	0	0
			8199	5231	1351	1569	48			
1	A	983	Total	C	N	O	S	0	0	0
			7584	4849	1247	1443	45			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	PRO	deletion	UNP Q9ESJ0
B	?	-	LYS	deletion	UNP Q9ESJ0
B	?	-	LEU	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	ARG	deletion	UNP Q9ESJ0
B	?	-	HIS	deletion	UNP Q9ESJ0
B	?	-	TYR	deletion	UNP Q9ESJ0
B	?	-	ILE	deletion	UNP Q9ESJ0
B	?	-	ALA	deletion	UNP Q9ESJ0
B	?	-	MET	deletion	UNP Q9ESJ0
B	?	-	PHE	deletion	UNP Q9ESJ0
B	?	-	GLU	deletion	UNP Q9ESJ0
B	?	-	SER	deletion	UNP Q9ESJ0
B	?	-	SER	deletion	UNP Q9ESJ0
B	?	-	GLN	deletion	UNP Q9ESJ0
B	?	-	ASN	deletion	UNP Q9ESJ0
B	?	-	VAL	deletion	UNP Q9ESJ0
B	?	-	LEU	deletion	UNP Q9ESJ0
B	?	-	LEU	deletion	UNP Q9ESJ0
B	?	-	LYS	deletion	UNP Q9ESJ0
B	?	-	SER	deletion	UNP Q9ESJ0
B	?	-	ASP	deletion	UNP Q9ESJ0
B	?	-	THR	deletion	UNP Q9ESJ0
B	?	-	ASP	deletion	UNP Q9ESJ0
B	?	-	GLU	deletion	UNP Q9ESJ0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	VAL	deletion	UNP Q9ESJ0
B	?	-	PHE	deletion	UNP Q9ESJ0
B	?	-	ARG	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	HIS	deletion	UNP Q9ESJ0
B	?	-	GLU	deletion	UNP Q9ESJ0
B	?	-	PRO	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	GLN	deletion	UNP Q9ESJ0
B	?	-	ALA	deletion	UNP Q9ESJ0
B	?	-	ALA	deletion	UNP Q9ESJ0
B	?	-	GLY	deletion	UNP Q9ESJ0
B	?	-	ARG	deletion	UNP Q9ESJ0
B	1085	TYR	HIS	engineered mutation	UNP Q9ESJ0
A	?	-	PRO	deletion	UNP Q9ESJ0
A	?	-	LYS	deletion	UNP Q9ESJ0
A	?	-	LEU	deletion	UNP Q9ESJ0
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	ARG	deletion	UNP Q9ESJ0
A	?	-	HIS	deletion	UNP Q9ESJ0
A	?	-	TYR	deletion	UNP Q9ESJ0
A	?	-	ILE	deletion	UNP Q9ESJ0
A	?	-	ALA	deletion	UNP Q9ESJ0
A	?	-	MET	deletion	UNP Q9ESJ0
A	?	-	PHE	deletion	UNP Q9ESJ0
A	?	-	GLU	deletion	UNP Q9ESJ0
A	?	-	SER	deletion	UNP Q9ESJ0
A	?	-	SER	deletion	UNP Q9ESJ0
A	?	-	GLN	deletion	UNP Q9ESJ0
A	?	-	ASN	deletion	UNP Q9ESJ0
A	?	-	VAL	deletion	UNP Q9ESJ0
A	?	-	LEU	deletion	UNP Q9ESJ0
A	?	-	LEU	deletion	UNP Q9ESJ0
A	?	-	LYS	deletion	UNP Q9ESJ0
A	?	-	SER	deletion	UNP Q9ESJ0
A	?	-	ASP	deletion	UNP Q9ESJ0
A	?	-	THR	deletion	UNP Q9ESJ0
A	?	-	ASP	deletion	UNP Q9ESJ0
A	?	-	GLU	deletion	UNP Q9ESJ0
A	?	-	VAL	deletion	UNP Q9ESJ0
A	?	-	PHE	deletion	UNP Q9ESJ0
A	?	-	ARG	deletion	UNP Q9ESJ0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	HIS	deletion	UNP Q9ESJ0
A	?	-	GLU	deletion	UNP Q9ESJ0
A	?	-	PRO	deletion	UNP Q9ESJ0
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	GLN	deletion	UNP Q9ESJ0
A	?	-	ALA	deletion	UNP Q9ESJ0
A	?	-	ALA	deletion	UNP Q9ESJ0
A	?	-	GLY	deletion	UNP Q9ESJ0
A	?	-	ARG	deletion	UNP Q9ESJ0
A	1085	TYR	HIS	engineered mutation	UNP Q9ESJ0

- Molecule 2 is a protein called GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	173	Total	C	N	O	S	0	0	0
			1368	890	238	236	4			
2	D	172	Total	C	N	O	S	0	0	0
			1373	894	236	239	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	69	LEU	GLN	engineered mutation	UNP P62826
D	69	LEU	GLN	engineered mutation	UNP P62826

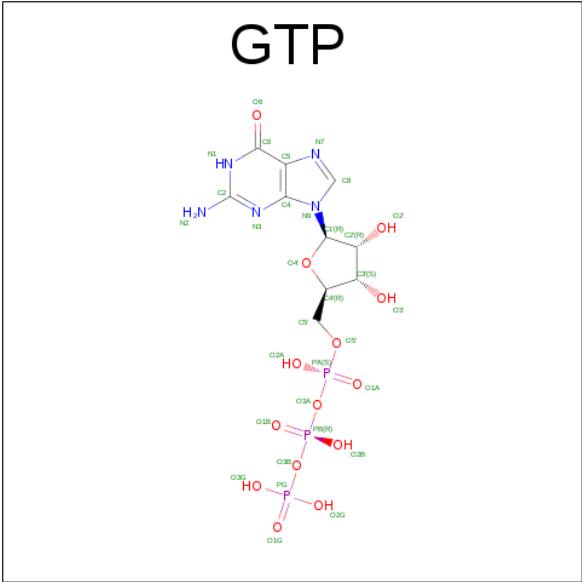
- Molecule 3 is a protein called Eukaryotic translation initiation factor 5A-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	138	Total	C	N	O	S	0	0	0
			1058	663	180	206	9			
3	E	137	Total	C	N	O	S	0	0	0
			1040	651	177	203	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	14	GLY	-	expression tag	UNP P63241
E	14	GLY	-	expression tag	UNP P63241

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



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

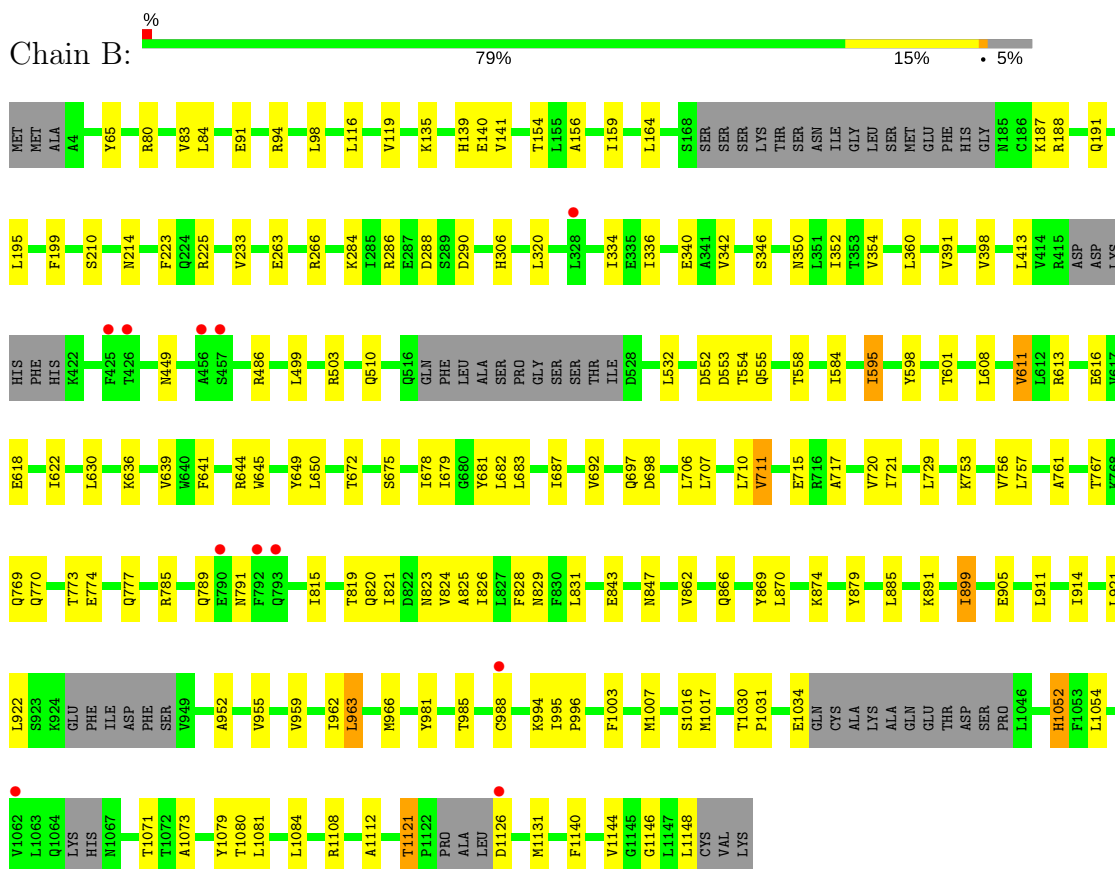
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	1	Total	Mg	0	0
			1	1		
5	C	1	Total	Mg	0	0
			1	1		

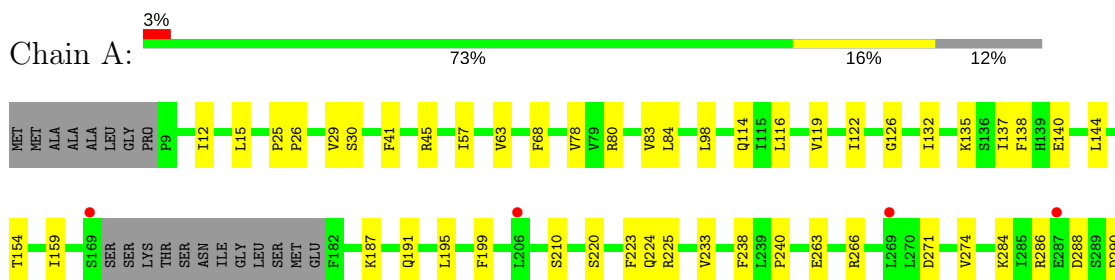
3 Residue-property plots

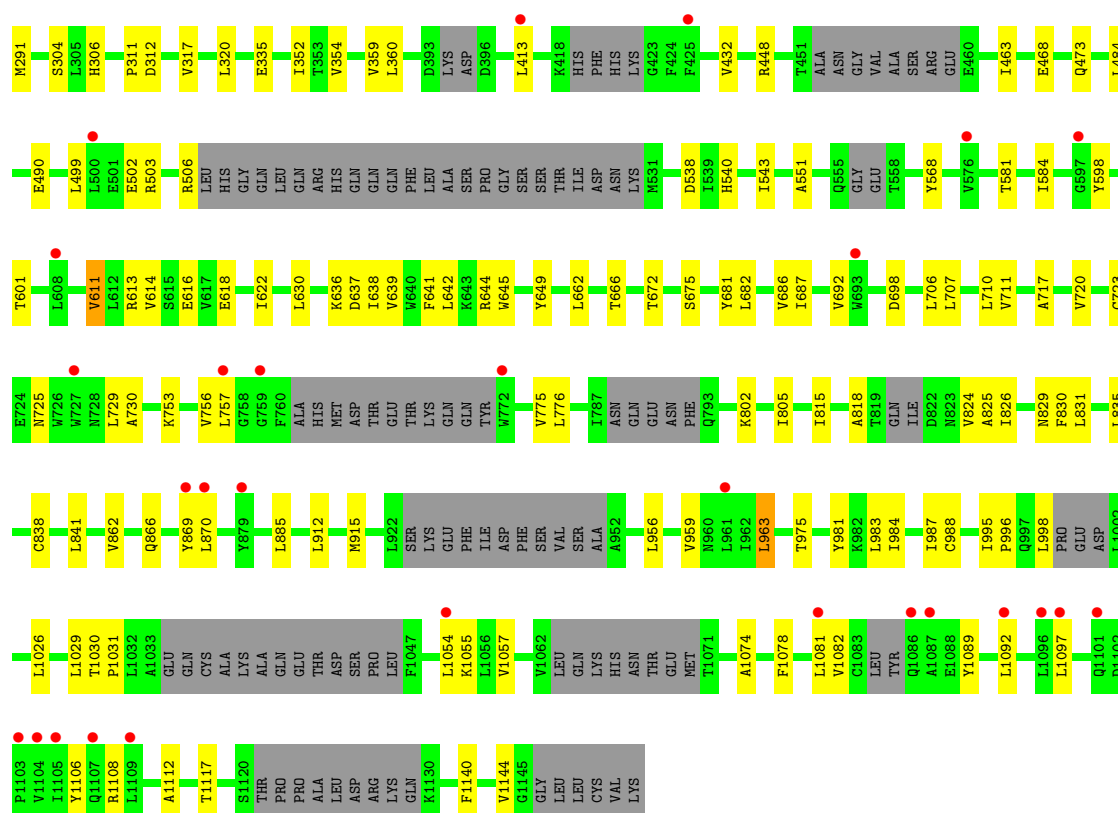
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Exportin-4

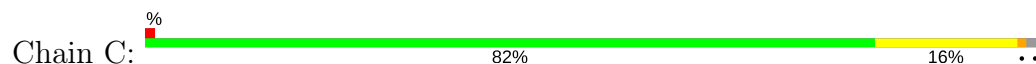


• Molecule 1: Exportin-4

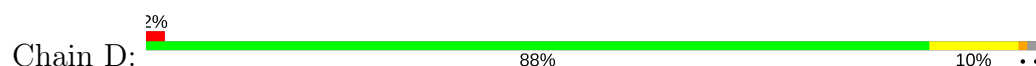




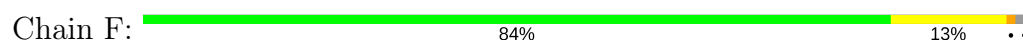
- Molecule 2: GTP-binding nuclear protein Ran



- Molecule 2: GTP-binding nuclear protein Ran

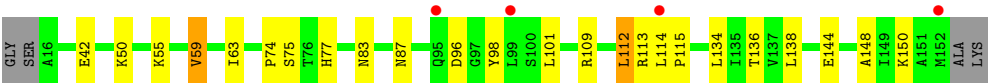


- Molecule 3: Eukaryotic translation initiation factor 5A-1



- Molecule 3: Eukaryotic translation initiation factor 5A-1





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.62Å 98.62Å 726.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.41 – 3.20 49.41 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.41-3.20) 100.0 (49.41-3.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 3.19Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.221 , 0.270 0.216 , 0.267	Depositor DCC
R_{free} test set	3491 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	106.3	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 76.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20688	wwPDB-VP
Average B, all atoms (Å ²)	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% 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, 5CT

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	0.24	0/7708	0.38	0/10473
1	B	0.24	0/8341	0.39	0/11336
2	C	0.25	0/1403	0.44	0/1904
2	D	0.25	0/1408	0.42	0/1908
3	E	0.24	0/1038	0.47	0/1398
3	F	0.25	0/1057	0.48	0/1423
All	All	0.24	0/20955	0.40	0/28442

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	7584	0	7444	105	0
1	B	8199	0	8095	98	0
2	C	1368	0	1353	17	0
2	D	1373	0	1369	13	0
3	E	1040	0	1036	12	0
3	F	1058	0	1057	11	0
4	C	32	0	12	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	32	0	12	1	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
All	All	20688	0	20378	249	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 (249) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1079:TYR:HH	1:B:1126:ASP:N	1.75	0.84
3:E:136:THR:HB	3:E:148:ALA:HB3	1.63	0.79
3:F:136:THR:HB	3:F:148:ALA:HB3	1.65	0.76
1:B:210:SER:O	1:B:284:LYS:NZ	2.22	0.72
1:B:91:GLU:OE2	1:B:94:ARG:NH1	2.22	0.72
1:B:135:LYS:O	1:B:139:HIS:ND1	2.22	0.70
1:B:622:ILE:HD12	1:B:692:VAL:HG21	1.74	0.69
1:A:622:ILE:HD12	1:A:692:VAL:HG21	1.75	0.67
1:B:911:LEU:HD13	1:B:966:MET:HE1	1.77	0.66
1:A:672:THR:HG23	1:A:675:SER:H	1.62	0.64
1:B:645:TRP:CD1	1:B:649:TYR:HB2	2.33	0.63
1:B:959:VAL:O	1:B:963:LEU:HB2	1.98	0.63
1:B:820:GLN:H	1:B:823:ASN:HB2	1.64	0.63
3:E:114:LEU:HD23	3:E:115:PRO:HD2	1.83	0.61
1:A:191:GLN:NE2	1:A:238:PHE:O	2.33	0.61
1:A:210:SER:O	1:A:284:LYS:NZ	2.32	0.60
1:B:553:ASP:O	1:B:555:GLN:N	2.33	0.60
1:B:721:ILE:HD13	1:B:761:ALA:HB3	1.83	0.60
1:B:988:CYS:HA	1:B:995:ILE:HD11	1.82	0.59
1:B:391:VAL:HG11	1:B:449:ASN:HA	1.85	0.59
1:A:645:TRP:CD1	1:A:649:TYR:HB2	2.37	0.58
1:B:352:ILE:HD13	1:B:413:LEU:HD12	1.85	0.58
1:B:636:LYS:HD3	1:B:698:ASP:HB3	1.85	0.58
1:B:825:ALA:O	1:B:829:ASN:ND2	2.37	0.58
1:B:824:VAL:HG11	1:B:869:TYR:HB2	1.86	0.58
1:A:956:LEU:HD13	1:A:998:LEU:HD21	1.86	0.57
1:A:135:LYS:HD2	1:A:138:PHE:CE2	2.38	0.57
1:A:1112:ALA:HB1	1:A:1140:PHE:HD1	1.68	0.57
1:A:581:THR:HA	1:A:601:THR:HG23	1.86	0.56
1:A:1026:LEU:HD13	1:A:1074:ALA:HB2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:VAL:O	1:A:1108:ARG:NH1	2.38	0.56
1:B:306:HIS:HB3	1:B:354:VAL:HG13	1.87	0.56
1:A:835:LEU:HD21	1:A:862:VAL:HG21	1.86	0.56
1:A:956:LEU:HD21	1:A:987:ILE:HG21	1.88	0.56
1:A:618:GLU:HG3	1:A:630:LEU:HD13	1.87	0.56
1:B:320:LEU:HD21	1:B:360:LEU:HD23	1.88	0.55
3:E:55:LYS:HA	3:E:74:PRO:HA	1.87	0.55
1:A:584:ILE:HG23	1:A:613:ARG:HH22	1.71	0.55
1:B:83:VAL:O	1:B:1108:ARG:NH1	2.39	0.55
1:A:45:ARG:O	1:A:80:ARG:NH1	2.40	0.55
1:B:639:VAL:HG13	1:B:706:LEU:HD12	1.89	0.54
1:A:753:LYS:HA	1:A:818:ALA:HB2	1.89	0.54
1:B:116:LEU:HD22	1:B:159:ILE:HG12	1.90	0.54
1:B:707:LEU:O	1:B:711:VAL:HG23	2.07	0.54
1:B:672:THR:HG23	1:B:675:SER:H	1.74	0.53
2:D:151:ALA:HA	2:D:157:PHE:CE2	2.44	0.53
1:A:320:LEU:HD21	1:A:360:LEU:HD23	1.91	0.53
1:A:468:GLU:HG3	1:A:473:GLN:HB2	1.91	0.53
1:A:448:ARG:NH1	1:A:538:ASP:OD1	2.41	0.52
3:F:42:GLU:HB3	3:F:59:VAL:HG13	1.92	0.52
1:B:831:LEU:HD13	1:B:862:VAL:HG22	1.91	0.52
2:D:43:LEU:HA	2:D:72:PHE:HB3	1.92	0.52
2:D:123:LYS:HG2	4:D:217:GTP:C6	2.44	0.52
1:B:1016:SER:OG	1:B:1017:MET:N	2.42	0.52
1:B:899:ILE:HG12	3:F:119:LEU:HD13	1.92	0.52
1:A:191:GLN:HG2	1:A:240:PRO:HD3	1.92	0.51
1:B:981:TYR:O	1:B:985:THR:OG1	2.23	0.51
2:C:39:TYR:CD1	4:C:217:GTP:H5"	2.45	0.51
1:B:611:VAL:HG21	1:B:641:PHE:CZ	2.45	0.51
1:B:819:THR:HG21	1:B:866:GLN:HG2	1.93	0.51
3:E:98:TYR:CE1	3:E:113:ARG:HG2	2.46	0.51
1:A:98:LEU:HD21	1:A:119:VAL:HG11	1.92	0.51
2:D:29:ARG:HG2	2:D:157:PHE:CE1	2.46	0.51
1:A:137:ILE:H	1:A:137:ILE:HD12	1.74	0.51
1:A:838:CYS:HA	1:A:841:LEU:HB2	1.92	0.51
1:A:12:ILE:HG12	1:A:57:ILE:HG12	1.92	0.51
1:A:825:ALA:O	1:A:829:ASN:ND2	2.44	0.50
1:B:346:SER:O	1:B:350:ASN:ND2	2.44	0.50
1:A:187:LYS:O	1:A:191:GLN:HB2	2.11	0.50
1:B:553:ASP:OD1	1:B:644:ARG:NH2	2.42	0.50
1:A:707:LEU:O	1:A:711:VAL:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:725:ASN:O	1:A:729:LEU:HG	2.12	0.50
2:C:15:LEU:HD13	2:C:23:LYS:HB3	1.92	0.50
3:F:55:LYS:HA	3:F:74:PRO:HA	1.93	0.50
1:B:847:ASN:O	3:F:86:ARG:HD2	2.11	0.50
1:A:199:PHE:HB2	1:A:233:VAL:HG11	1.94	0.50
3:F:17:THR:HG22	3:F:82:PRO:HA	1.93	0.50
1:A:502:GLU:OE2	1:A:506:ARG:NE	2.45	0.49
1:B:821:ILE:HG13	1:B:869:TYR:CE1	2.47	0.49
1:A:756:VAL:HG21	1:A:815:ILE:HG12	1.94	0.49
1:A:831:LEU:HD13	1:A:862:VAL:HG12	1.94	0.49
1:A:776:LEU:HB3	1:A:830:PHE:CZ	2.47	0.49
1:A:988:CYS:HA	1:A:995:ILE:HD11	1.93	0.49
1:A:611:VAL:HG21	1:A:641:PHE:CE2	2.48	0.49
2:C:35:PHE:CE2	2:C:37:LYS:HG2	2.48	0.49
1:B:611:VAL:HG21	1:B:641:PHE:CE2	2.47	0.49
1:B:753:LYS:O	1:B:757:LEU:HB2	2.13	0.49
3:E:101:LEU:HD12	3:E:112:LEU:HD23	1.94	0.49
1:A:717:ALA:HA	1:A:720:VAL:HG12	1.95	0.49
1:B:286:ARG:HE	1:B:334:ILE:HG23	1.78	0.49
1:A:639:VAL:HG13	1:A:706:LEU:HD12	1.95	0.49
1:A:912:LEU:HD11	1:A:975:THR:HG22	1.95	0.48
1:B:84:LEU:HA	1:B:1108:ARG:HH12	1.77	0.48
1:A:642:LEU:HB2	1:A:706:LEU:HD11	1.95	0.48
1:B:94:ARG:NH2	1:B:140:GLU:OE1	2.45	0.48
3:F:89:PHE:CE1	3:F:103:GLN:HG2	2.49	0.48
1:B:263:GLU:HA	1:B:266:ARG:HB2	1.96	0.48
2:D:35:PHE:HE2	2:D:37:LYS:HG2	1.78	0.48
1:A:824:VAL:HG11	1:A:869:TYR:HB2	1.95	0.48
2:C:123:LYS:HG2	4:C:217:GTP:C6	2.49	0.48
1:A:271:ASP:O	1:A:274:VAL:HG22	2.13	0.48
2:C:95:ARG:HG2	2:C:99:LYS:HE3	1.95	0.48
1:B:1112:ALA:HB1	1:B:1140:PHE:CD1	2.49	0.48
1:A:68:PHE:CD1	1:A:114:GLN:HG3	2.49	0.47
1:A:753:LYS:O	1:A:757:LEU:HB2	2.13	0.47
2:C:77:ASP:OD1	2:C:77:ASP:N	2.47	0.47
2:C:29:ARG:NH1	3:F:42:GLU:OE1	2.45	0.47
3:F:75:SER:C	3:F:77:HIS:H	2.17	0.47
1:A:802:LYS:HA	1:A:805:ILE:HG22	1.97	0.47
1:B:773:THR:HA	1:B:777:GLN:HB2	1.97	0.47
1:A:866:GLN:O	1:A:870:LEU:HG	2.15	0.47
2:D:77:ASP:N	2:D:77:ASP:OD1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:75:SER:C	3:E:77:HIS:H	2.18	0.47
1:B:584:ILE:HG23	1:B:613:ARG:HH22	1.80	0.47
1:A:263:GLU:HA	1:A:266:ARG:HB2	1.97	0.47
1:A:286:ARG:NH2	1:A:335:GLU:O	2.44	0.47
1:B:682:LEU:HD12	1:B:710:LEU:HD21	1.97	0.47
2:C:151:ALA:HA	2:C:157:PHE:CE2	2.50	0.47
1:A:224:GLN:HG2	1:A:291:MET:HG2	1.97	0.47
1:A:78:VAL:HG11	1:A:122:ILE:HD11	1.97	0.47
1:B:486:ARG:NH2	1:B:552:ASP:OD1	2.42	0.47
1:A:730:ALA:HB1	1:A:775:VAL:HG22	1.97	0.46
1:B:922:LEU:HG	1:B:955:VAL:HG21	1.97	0.46
1:A:686:VAL:HG21	1:A:707:LEU:HD13	1.97	0.46
1:B:1003:PHE:O	1:B:1007:MET:HG2	2.15	0.46
1:A:15:LEU:HD22	1:A:41:PHE:CE1	2.50	0.46
1:A:84:LEU:HD23	1:A:1108:ARG:HH11	1.81	0.46
1:A:306:HIS:HB3	1:A:354:VAL:HG13	1.97	0.46
1:A:614:VAL:HG12	1:A:638:ILE:HD11	1.98	0.46
1:B:952:ALA:HB3	1:B:994:LYS:HZ2	1.79	0.46
3:E:87:ASN:O	3:E:136:THR:HA	2.15	0.46
1:A:682:LEU:HD12	1:A:710:LEU:HD21	1.97	0.46
1:B:1071:THR:HG22	1:B:1073:ALA:H	1.81	0.46
1:B:756:VAL:HG21	1:B:815:ILE:HG12	1.97	0.46
1:A:116:LEU:HD22	1:A:159:ILE:HG12	1.97	0.46
1:B:1112:ALA:HB1	1:B:1140:PHE:HD1	1.81	0.45
2:C:113:GLU:HB3	2:C:114:ASN:H	1.67	0.45
1:A:1030:THR:OG1	1:A:1031:PRO:HD3	2.17	0.45
1:A:352:ILE:HG23	1:A:413:LEU:HD13	1.98	0.45
1:A:68:PHE:CE1	1:A:114:GLN:HG3	2.52	0.45
1:B:717:ALA:HA	1:B:720:VAL:HG12	1.97	0.45
1:A:490:GLU:HG2	1:A:568:TYR:CE1	2.52	0.45
1:A:723:CYS:O	1:A:725:ASN:N	2.42	0.45
1:B:84:LEU:HD23	1:B:1108:ARG:HH11	1.82	0.45
1:A:1097:LEU:HB3	1:A:1106:TYR:HE1	1.82	0.45
1:A:584:ILE:HD12	1:A:601:THR:HG21	1.97	0.45
1:B:828:PHE:CE2	1:B:874:LYS:HB3	2.51	0.45
1:A:805:ILE:HD11	1:A:841:LEU:HG	1.98	0.45
1:A:959:VAL:O	1:A:963:LEU:HB2	2.16	0.45
2:C:35:PHE:HE2	2:C:37:LYS:HG2	1.82	0.45
1:A:1140:PHE:CE1	1:A:1144:VAL:HG21	2.51	0.45
1:B:595:ILE:HG22	1:B:598:TYR:HB3	1.97	0.45
1:B:618:GLU:HG3	1:B:630:LEU:HD13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:LEU:HD21	1:B:119:VAL:HG11	1.97	0.45
1:A:1055:LYS:HE2	1:A:1092:LEU:HD21	1.98	0.45
1:A:618:GLU:O	1:A:622:ILE:HG12	2.17	0.44
2:D:141:LYS:HD2	2:D:144:LEU:HD12	1.99	0.44
1:A:1089:TYR:HE2	1:A:1117:THR:HG22	1.82	0.44
1:B:154:THR:HG23	1:B:225:ARG:HD3	1.98	0.44
2:D:13:LEU:HA	2:D:85:CYS:O	2.17	0.44
1:B:843:GLU:OE2	1:B:891:LYS:NZ	2.41	0.44
1:A:304:SER:HA	1:A:354:VAL:HG21	1.99	0.44
2:C:13:LEU:HA	2:C:85:CYS:O	2.18	0.44
1:A:611:VAL:HG21	1:A:641:PHE:CZ	2.53	0.44
1:B:286:ARG:HG3	1:B:340:GLU:OE2	2.18	0.44
1:A:981:TYR:HA	1:A:984:ILE:HG12	1.99	0.44
1:B:187:LYS:O	1:B:191:GLN:HG3	2.17	0.44
1:A:915:MET:HE2	1:A:915:MET:HB2	1.86	0.44
1:B:342:VAL:HG23	1:B:398:VAL:HG13	1.99	0.44
2:D:35:PHE:CE2	2:D:37:LYS:HG2	2.53	0.44
1:B:608:LEU:HB3	1:B:645:TRP:CH2	2.52	0.44
1:A:317:VAL:HG22	1:A:359:VAL:HG13	2.00	0.43
1:A:824:VAL:C	1:A:826:ILE:H	2.21	0.43
1:B:510:GLN:HG2	1:B:532:LEU:HD13	1.99	0.43
1:B:1054:LEU:HD13	1:B:1081:LEU:HB3	2.00	0.43
1:A:154:THR:HG23	1:A:225:ARG:HD3	2.01	0.43
1:B:1030:THR:HG23	1:B:1080:THR:HG21	2.01	0.43
2:C:68:GLY:N	4:C:217:GTP:O1G	2.46	0.43
1:A:126:GLY:HA3	1:A:132:ILE:HG13	1.99	0.43
1:A:25:PRO:HA	1:A:26:PRO:HD3	1.92	0.43
1:B:199:PHE:HB2	1:B:233:VAL:HG11	1.99	0.43
3:E:42:GLU:HB2	3:E:59:VAL:HG13	1.99	0.43
1:A:432:VAL:HG21	1:A:484:LEU:HD13	2.00	0.43
1:B:1034:GLU:OE1	1:B:1034:GLU:N	2.50	0.43
1:B:164:LEU:HD11	1:B:195:LEU:HD13	2.01	0.43
1:B:905:GLU:OE1	2:C:37:LYS:HD3	2.19	0.43
1:B:914:ILE:HG21	1:B:962:ILE:HG21	1.99	0.43
1:B:824:VAL:C	1:B:826:ILE:H	2.21	0.43
1:A:616:GLU:HG2	1:A:681:TYR:OH	2.19	0.43
1:B:995:ILE:HB	1:B:996:PRO:HD3	2.01	0.43
1:B:584:ILE:HG13	1:B:601:THR:HG21	2.00	0.43
1:A:1029:LEU:HA	1:A:1029:LEU:HD23	1.83	0.42
1:B:687:ILE:HD11	1:B:729:LEU:HD22	2.01	0.42
1:A:540:HIS:ND1	1:A:637:ASP:OD2	2.29	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:650:LEU:HD23	1:B:650:LEU:HA	1.86	0.42
3:E:63:ILE:HG13	3:E:138:LEU:HD21	2.02	0.42
1:B:141:VAL:HG13	1:B:156:ALA:HB1	2.01	0.42
1:B:879:TYR:CE1	1:B:921:LEU:HG	2.54	0.42
3:E:109:ARG:NH2	3:E:144:GLU:OE1	2.53	0.42
1:A:1078:PHE:O	1:A:1082:VAL:HG23	2.19	0.42
1:B:679:ILE:HD11	1:B:720:VAL:HB	2.01	0.42
1:B:1084:LEU:HA	1:B:1084:LEU:HD23	1.93	0.42
1:B:616:GLU:HG2	1:B:681:TYR:OH	2.19	0.42
1:B:770:GLN:NE2	1:B:774:GLU:OE1	2.53	0.42
2:D:151:ALA:HA	2:D:157:PHE:HE2	1.84	0.42
3:E:134:LEU:HB3	3:E:150:LYS:HB2	2.01	0.42
1:B:1030:THR:HB	1:B:1031:PRO:HD3	2.01	0.42
1:B:1140:PHE:O	1:B:1144:VAL:N	2.48	0.42
1:B:263:GLU:HG2	1:B:266:ARG:HE	1.85	0.42
1:B:288:ASP:C	1:B:290:ASP:H	2.23	0.42
1:B:65:TYR:CD1	2:C:81:ILE:HD12	2.54	0.42
1:A:29:VAL:O	1:A:30:SER:OG	2.29	0.42
1:A:995:ILE:HB	1:A:996:PRO:HD3	2.02	0.42
1:B:683:LEU:HD12	1:B:683:LEU:HA	1.88	0.42
1:A:499:LEU:O	1:A:503:ARG:HG2	2.20	0.42
1:B:499:LEU:O	1:B:503:ARG:HG2	2.20	0.42
1:B:678:ILE:O	1:B:682:LEU:HG	2.20	0.42
1:A:288:ASP:C	1:A:290:ASP:H	2.23	0.41
1:A:662:LEU:O	1:A:666:THR:HG22	2.20	0.41
1:B:866:GLN:O	1:B:870:LEU:HG	2.20	0.41
2:C:101:VAL:N	2:C:102:PRO:HD2	2.35	0.41
1:A:543:ILE:HG12	1:A:611:VAL:HG12	2.02	0.41
2:C:11:PHE:CG	2:C:168:LEU:HD13	2.55	0.41
1:B:769:GLN:O	1:B:773:THR:OG1	2.23	0.41
1:B:1146:GLY:O	1:B:1148:LEU:N	2.54	0.41
1:B:80:ARG:HG3	1:B:1146:GLY:O	2.20	0.41
1:A:636:LYS:HD3	1:A:698:ASP:HB3	2.01	0.41
1:A:220:SER:O	1:A:224:GLN:HG3	2.21	0.41
3:F:98:TYR:CE1	3:F:113:ARG:HG2	2.56	0.41
1:B:336:ILE:O	1:B:340:GLU:HB2	2.20	0.41
1:B:785:ARG:O	1:B:789:GLN:HG3	2.21	0.41
2:C:12:LYS:HE3	2:C:64:TRP:CE2	2.56	0.41
1:A:1054:LEU:HD13	1:A:1081:LEU:HB3	2.03	0.41
1:A:195:LEU:HD11	1:A:233:VAL:HG13	2.02	0.41
1:A:584:ILE:HD11	1:A:598:TYR:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:687:ILE:HD11	1:A:729:LEU:CD1	2.51	0.41
1:A:622:ILE:HG23	1:A:692:VAL:HG11	2.02	0.41
1:A:1057:VAL:HG12	1:A:1078:PHE:CZ	2.55	0.41
1:B:1052:HIS:ND1	1:B:1052:HIS:O	2.50	0.40
2:D:29:ARG:HG2	2:D:157:PHE:CZ	2.56	0.40
1:A:551:ALA:HB3	1:A:644:ARG:HD3	2.02	0.40
1:A:915:MET:HG3	1:A:983:LEU:HD12	2.03	0.40
2:D:88:ILE:HD11	2:D:108:LEU:HD12	2.03	0.40
3:F:152:MET:N	3:F:152:MET:SD	2.94	0.40
1:B:188:ARG:CZ	1:A:311:PRO:HB3	2.51	0.40
1:A:776:LEU:HD23	1:A:830:PHE:CE2	2.57	0.40
2:D:101:VAL:N	2:D:102:PRO:HD2	2.36	0.40
1:A:140:GLU:O	1:A:144:LEU:HD13	2.22	0.40
1:A:463:ILE:HD12	3:E:50:5CT:H22	2.02	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	951/1113 (85%)	915 (96%)	35 (4%)	1 (0%)	55	89
1	B	1036/1113 (93%)	985 (95%)	47 (4%)	4 (0%)	38	77
2	C	171/176 (97%)	167 (98%)	4 (2%)	0	100	100
2	D	170/176 (97%)	166 (98%)	4 (2%)	0	100	100
3	E	134/141 (95%)	127 (95%)	6 (4%)	1 (1%)	25	68
3	F	135/141 (96%)	128 (95%)	7 (5%)	0	100	100
All	All	2597/2860 (91%)	2488 (96%)	103 (4%)	6 (0%)	51	86

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	554	THR
1	A	312	ASP
1	B	214	ASN
3	E	96	ASP
1	B	1121	THR
1	B	899	ILE

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	825/994 (83%)	820 (99%)	5 (1%)	89	96
1	B	901/994 (91%)	887 (98%)	14 (2%)	68	89
2	C	143/154 (93%)	137 (96%)	6 (4%)	34	72
2	D	146/154 (95%)	144 (99%)	2 (1%)	71	90
3	E	111/118 (94%)	108 (97%)	3 (3%)	50	82
3	F	114/118 (97%)	109 (96%)	5 (4%)	33	70
All	All	2240/2532 (88%)	2205 (98%)	35 (2%)	68	89

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	223	PHE
1	B	558	THR
1	B	595	ILE
1	B	611	VAL
1	B	697	GLN
1	B	711	VAL
1	B	715	GLU
1	B	767	THR
1	B	791	ASN
1	B	885	LEU
1	B	963	LEU
1	B	1052	HIS
1	B	1121	THR

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Mol	Chain	Res	Type
1	B	1131	MET
2	C	14	VAL
2	C	15	LEU
2	C	30	HIS
2	C	63	VAL
2	C	69	LEU
2	C	109	VAL
3	F	59	VAL
3	F	76	THR
3	F	83	ASN
3	F	114	LEU
3	F	152	MET
1	A	63	VAL
1	A	223	PHE
1	A	611	VAL
1	A	885	LEU
1	A	963	LEU
2	D	30	HIS
2	D	157	PHE
3	E	59	VAL
3	E	83	ASN
3	E	112	LEU

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

Mol	Chain	Res	Type
1	A	191	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	5CT	E	50	3	14,14,15	0.83	0	11,15,17	1.09	2 (18%)
3	5CT	F	50	3	14,14,15	0.82	0	11,15,17	1.13	1 (9%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	5CT	E	50	3	-	0/12/14/16	0/0/0/0
3	5CT	F	50	3	-	0/12/14/16	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	F	50	5CT	C3-C2-C1	-2.63	109.28	112.20
3	E	50	5CT	C3-C2-C1	-2.52	109.41	112.20
3	E	50	5CT	C1-NZ-CE	-2.02	108.99	113.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	50	5CT	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	GTP	C	217	5	27,34,34	1.10	2 (7%)	27,54,54	1.70	6 (22%)
4	GTP	D	217	5	27,34,34	1.35	4 (14%)	27,54,54	1.71	6 (22%)

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	C	217	5	-	0/18/38/38	0/3/3/3
4	GTP	D	217	5	-	0/18/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	217	GTP	C8-N7	-2.26	1.30	1.34
4	D	217	GTP	C8-N7	-2.18	1.30	1.34
4	C	217	GTP	O4'-C1'	2.60	1.44	1.41
4	D	217	GTP	PB-O1B	2.84	1.61	1.50
4	D	217	GTP	PA-O1A	2.85	1.61	1.50
4	D	217	GTP	O4'-C1'	2.96	1.45	1.41

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	217	GTP	N3-C2-N1	-3.80	121.91	127.46
4	C	217	GTP	N3-C2-N1	-3.72	122.03	127.46
4	C	217	GTP	C5-C6-N1	-3.70	118.21	123.48
4	D	217	GTP	C5-C6-N1	-3.65	118.28	123.48
4	C	217	GTP	C4-C5-N7	-2.34	107.15	109.41
4	D	217	GTP	C6-C5-C4	-2.33	118.52	120.84
4	C	217	GTP	C6-C5-C4	-2.32	118.54	120.84
4	D	217	GTP	C4-C5-N7	-2.27	107.22	109.41
4	C	217	GTP	C6-N1-C2	3.00	120.38	116.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	217	GTP	C6-N1-C2	3.03	120.41	116.06
4	C	217	GTP	C2-N3-C4	4.77	120.73	115.16
4	D	217	GTP	C2-N3-C4	4.95	120.94	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	217	GTP	3	0
4	D	217	GTP	1	0

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	983/1113 (88%)	0.21	32 (3%) 47 31	95, 133, 173, 194	0
1	B	1052/1113 (94%)	0.05	11 (1%) 82 72	68, 108, 158, 180	0
2	C	173/176 (98%)	0.15	1 (0%) 89 83	63, 83, 116, 130	0
2	D	172/176 (97%)	0.30	4 (2%) 61 46	84, 102, 136, 149	0
3	E	136/141 (96%)	0.28	4 (2%) 52 37	108, 126, 147, 152	0
3	F	137/141 (97%)	-0.05	0 100 100	71, 94, 133, 149	0
All	All	2653/2860 (92%)	0.14	52 (1%) 65 50	63, 118, 165, 194	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1086	GLN	4.3
1	B	457	SER	3.8
1	B	456	ALA	3.5
1	A	500	LEU	3.2
1	A	169	SER	3.2
1	B	1062	VAL	3.1
1	A	870	LEU	3.1
1	A	1107	GLN	3.1
1	B	425	PHE	3.0
1	B	793	GLN	3.0
1	A	772	TRP	2.9
1	A	425	PHE	2.9
1	B	792	PHE	2.8
1	A	759	GLY	2.8
1	A	1087	ALA	2.8
3	E	152	MET	2.8
1	A	413	LEU	2.8
3	E	99	LEU	2.8
1	A	1081	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	54	THR	2.6
1	A	1103	PRO	2.5
1	A	576	VAL	2.5
1	A	269	LEU	2.5
1	A	727	TRP	2.5
1	B	426	THR	2.5
1	A	1105	ILE	2.4
2	C	179	MET	2.4
1	A	1109	LEU	2.4
1	A	1054	LEU	2.4
1	A	879	TYR	2.3
1	A	1092	LEU	2.3
2	D	59	ILE	2.3
1	A	869	TYR	2.3
2	D	169	ILE	2.3
1	B	790	GLU	2.3
1	A	206	LEU	2.2
1	A	597	GLY	2.2
1	A	1101	GLN	2.2
1	A	693	TRP	2.2
1	A	961	LEU	2.2
1	B	1126	ASP	2.2
3	E	114	LEU	2.1
2	D	52	PHE	2.1
3	E	95	GLN	2.1
1	A	1097	LEU	2.1
1	A	287	GLU	2.1
1	B	328	LEU	2.1
1	A	608	LEU	2.1
1	A	1104	VAL	2.0
1	B	988	CYS	2.0
1	A	757	LEU	2.0
1	A	1096	LEU	2.0

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	5CT	E	50	15/16	0.93	0.21	-	105,111,117,118	0
3	5CT	F	50	15/16	0.94	0.23	-	89,90,92,92	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
5	MG	C	218	1/1	0.98	0.26	1.22	58,58,58,58	0
4	GTP	C	217	32/32	0.97	0.23	-0.37	58,63,74,85	0
4	GTP	D	217	32/32	0.96	0.23	-0.98	81,87,101,104	0
5	MG	D	218	1/1	0.97	0.34	-	82,82,82,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.