



wwPDB X-ray Structure Validation Summary Report ⓘ

May 28, 2020 – 09:56 pm BST

PDB ID : 2FFL
Title : Crystal Structure of Dicer from Giardia intestinalis
Authors : Doudna, J.A.; MacRae, I.J.; Adams, P.D.
Deposited on : 2005-12-19
Resolution : 3.33 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

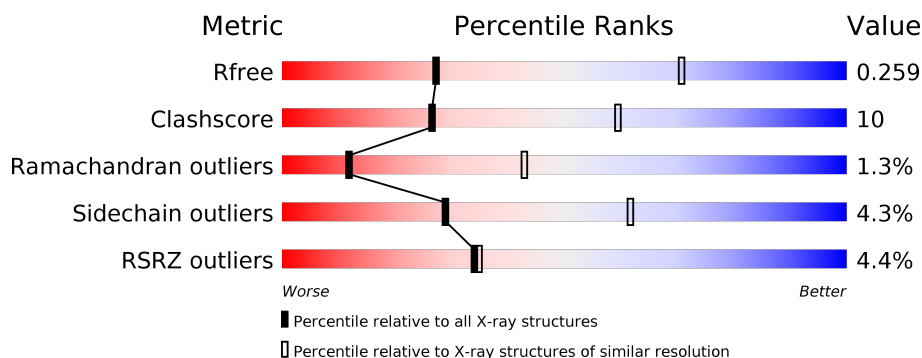
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.33 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}	130704	1060 (3.38-3.30)
Clashscore	141614	1111 (3.38-3.30)
Ramachandran outliers	138981	1090 (3.38-3.30)
Sidechain outliers	138945	1089 (3.38-3.30)
RSRZ outliers	127900	1028 (3.38-3.30)

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 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	756	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 23%, green 71%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 71% 23% • • </div> </div>
1	B	756	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 1%, yellow 22%, green 72%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 72% 22% • • </div> </div>
1	C	756	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 7%, orange 1%, yellow 20%, green 74%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 7% 74% 20% • 5% </div> </div>
1	D	756	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 6%, orange 1%, yellow 23%, green 70%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 6% 70% 23% • 5% </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MN	D	5023	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 22426 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 Dicer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	732	Total	C	N	O	S	0	0	0
			5630	3614	957	1029	30			
1	B	726	Total	C	N	O	S	0	0	0
			5590	3589	950	1021	30			
1	C	718	Total	C	N	O	S	0	0	0
			5539	3560	941	1008	30			
1	D	720	Total	C	N	O	S	0	0	0
			5574	3586	942	1016	30			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	CLONING ARTIFACT	GB 71078474
A	0	ALA	-	CLONING ARTIFACT	GB 71078474
B	-1	GLY	-	CLONING ARTIFACT	GB 71078474
B	0	ALA	-	CLONING ARTIFACT	GB 71078474
C	-1	GLY	-	CLONING ARTIFACT	GB 71078474
C	0	ALA	-	CLONING ARTIFACT	GB 71078474
D	-1	GLY	-	CLONING ARTIFACT	GB 71078474
D	0	ALA	-	CLONING ARTIFACT	GB 71078474

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	9	Total	Mn	0	0
			9	9		
2	A	12	Total	Mn	0	0
			12	12		
2	D	8	Total	Mn	0	0
			8	8		
2	C	10	Total	Mn	0	0
			10	10		

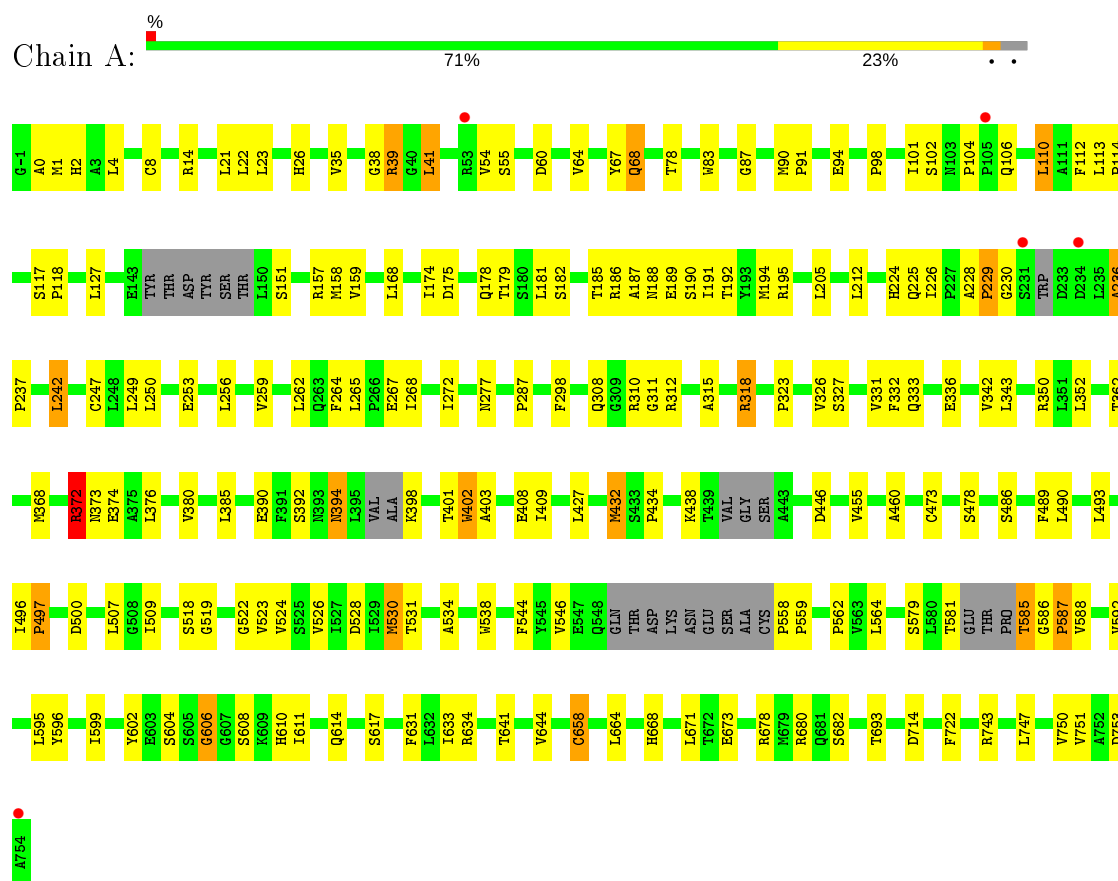
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total 17	O 17	0	0
3	B	21	Total 21	O 21	0	0
3	C	7	Total 7	O 7	0	0
3	D	9	Total 9	O 9	0	0

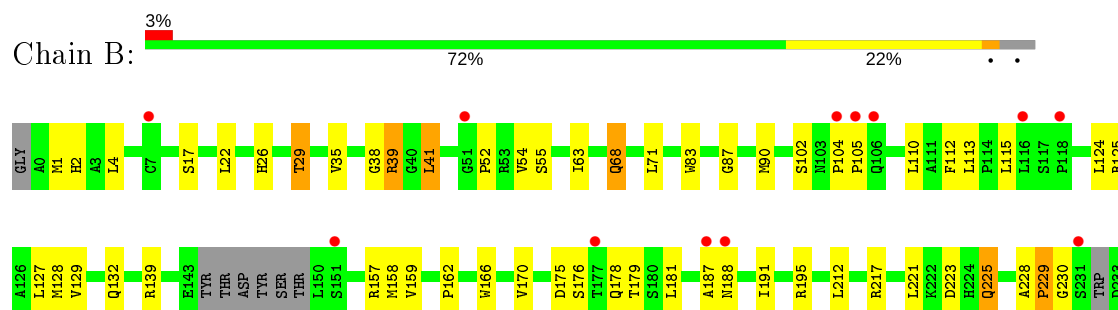
3 Residue-property plots

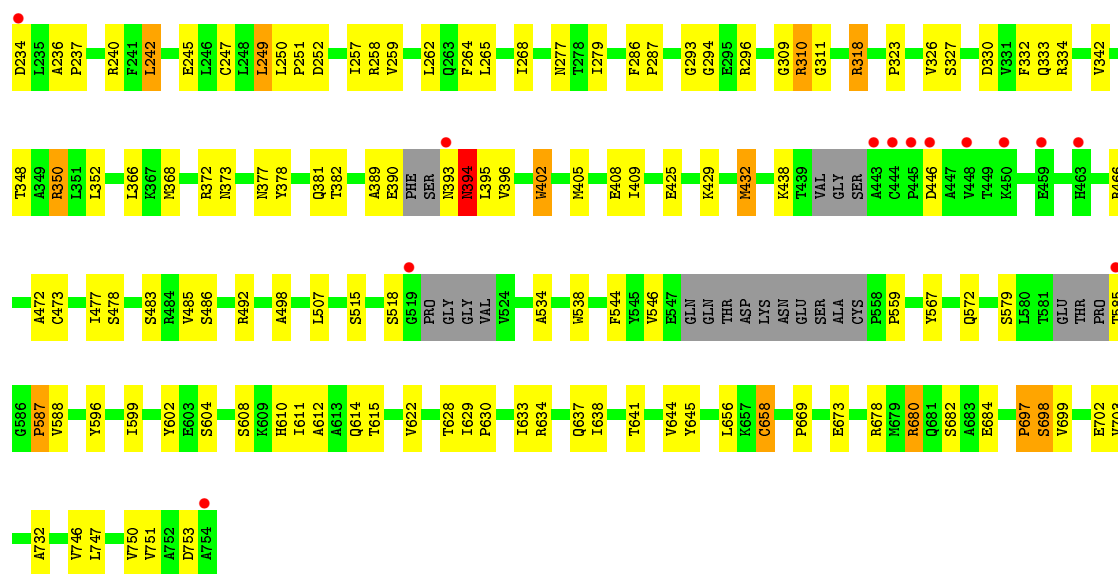
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

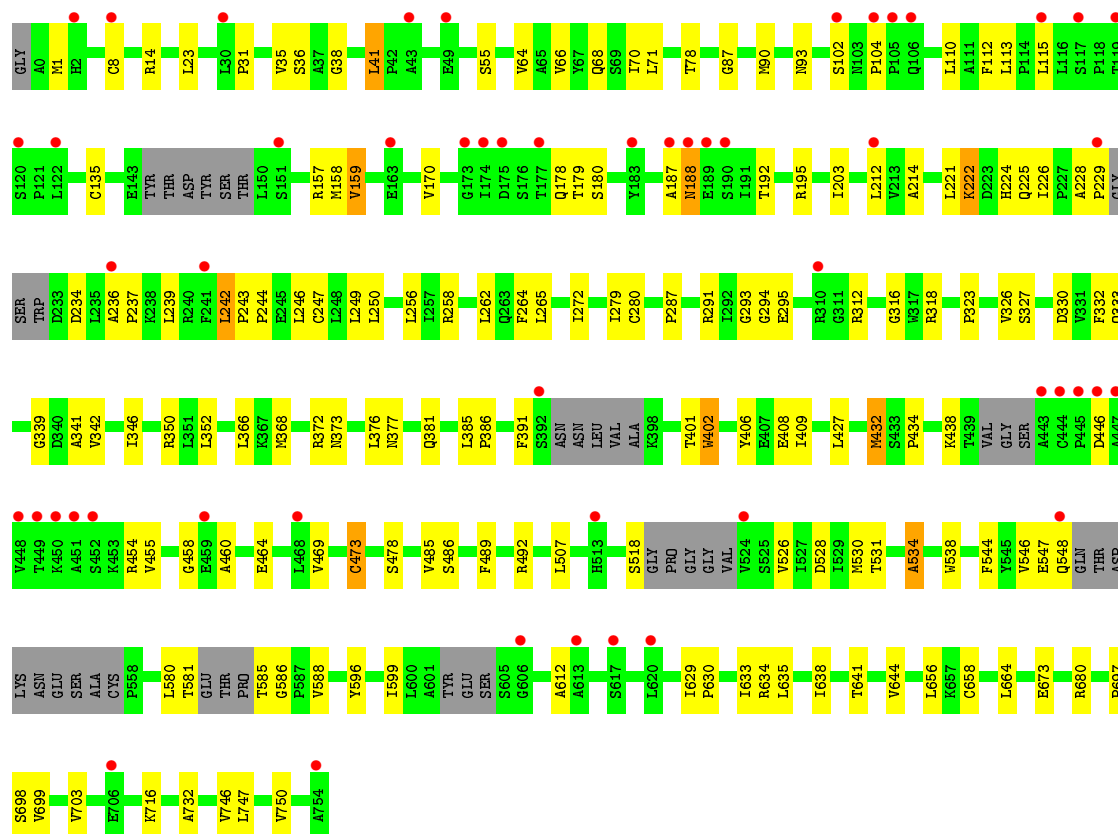
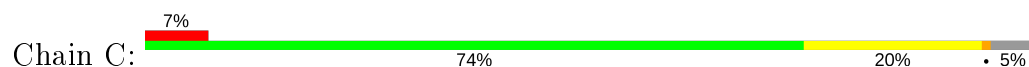


• Molecule 1: Dicer



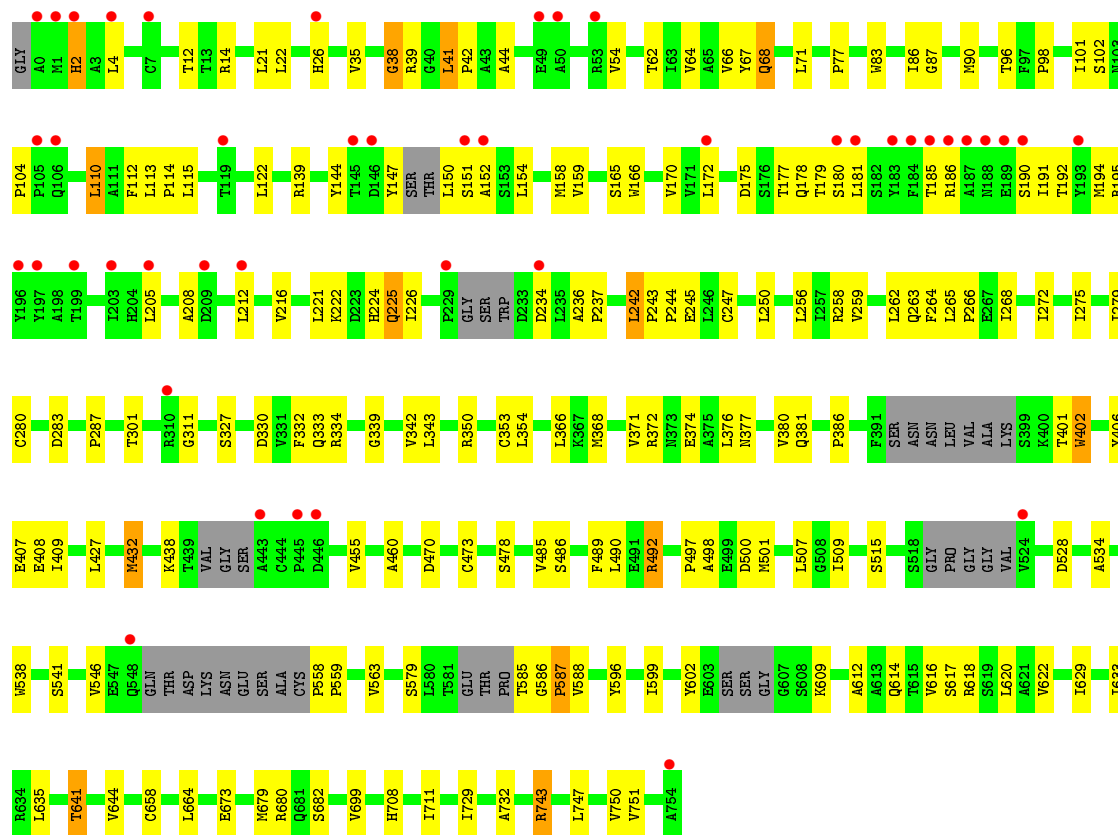


• Molecule 1: Dicer



• Molecule 1: Dicer





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	155.51Å 174.10Å 152.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.33 49.68 – 3.33	Depositor EDS
% Data completeness (in resolution range)	99.5 (20.00-3.33) 98.6 (49.68-3.33)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, R_{free}	0.243 , 0.275 0.223 , 0.259	Depositor DCC
R_{free} test set	3003 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	71.5	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for l,-k,h	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	22426	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 5.04% 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: MN

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	0.48	0/5761	0.59	1/7836 (0.0%)
1	B	0.48	0/5718	0.60	1/7777 (0.0%)
1	C	0.42	0/5666	0.55	1/7704 (0.0%)
1	D	0.44	0/5704	0.56	1/7759 (0.0%)
All	All	0.46	0/22849	0.58	4/31076 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	242	LEU	CA-CB-CG	5.36	127.62	115.30
1	C	242	LEU	CA-CB-CG	5.24	127.34	115.30
1	B	242	LEU	CA-CB-CG	5.09	127.01	115.30
1	D	38	GLY	N-CA-C	5.03	125.69	113.10

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	5630	0	5674	130	0
1	B	5590	0	5640	121	1
1	C	5539	0	5592	92	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	5574	0	5611	123	1
2	A	12	0	0	0	0
2	B	9	0	0	0	0
2	C	10	0	0	0	0
2	D	8	0	0	0	0
3	A	17	0	0	1	0
3	B	21	0	0	0	0
3	C	7	0	0	0	0
3	D	9	0	0	0	0
All	All	22426	0	22517	462	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 10.

The worst 5 of 462 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:C:455:VAL:HA	1:C:460:ALA:HB2	1.37	1.03
1:A:38:GLY:HA3	1:A:41:LEU:HD12	1.48	0.96
1:D:342:VAL:HG12	1:D:658:CYS:SG	2.09	0.92
1:C:432:MET:CE	1:C:538:TRP:HB2	2.04	0.88
1:D:38:GLY:HA3	1:D:41:LEU:HD12	1.57	0.87

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:B:466:ARG:NH1	1:D:470:ASP:OD2[4_455]	2.07	0.13

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	718/756 (95%)	656 (91%)	50 (7%)	12 (2%)	9	37
1	B	710/756 (94%)	661 (93%)	38 (5%)	11 (2%)	10	40
1	C	700/756 (93%)	644 (92%)	50 (7%)	6 (1%)	17	51
1	D	702/756 (93%)	638 (91%)	55 (8%)	9 (1%)	12	43
All	All	2830/3024 (94%)	2599 (92%)	193 (7%)	38 (1%)	12	43

5 of 38 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	0	ALA
1	A	54	VAL
1	A	229	PRO
1	A	586	GLY
1	A	606	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	611/632 (97%)	583 (95%)	28 (5%)	27	60
1	B	607/632 (96%)	574 (95%)	33 (5%)	22	55
1	C	602/632 (95%)	581 (96%)	21 (4%)	36	66
1	D	605/632 (96%)	583 (96%)	22 (4%)	35	65
All	All	2425/2528 (96%)	2321 (96%)	104 (4%)	29	62

5 of 104 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	446	ASP
1	B	658	CYS
1	D	515	SER
1	B	473	CYS
1	B	518	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 41 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	572	GLN
1	C	178	GLN
1	D	548	GLN
1	B	614	GLN
1	B	637	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 39 ligands modelled in this entry, 39 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	732/756 (96%)	-0.00	5 (0%) 87 90	21, 37, 67, 94	1 (0%)
1	B	726/756 (96%)	0.09	25 (3%) 45 44	22, 37, 69, 95	1 (0%)
1	C	718/756 (94%)	0.40	52 (7%) 15 16	20, 38, 69, 94	1 (0%)
1	D	720/756 (95%)	0.28	44 (6%) 21 22	22, 37, 68, 94	1 (0%)
All	All	2896/3024 (95%)	0.19	126 (4%) 34 35	20, 38, 68, 95	4 (0%)

The worst 5 of 126 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	190	SER	5.8
1	C	188	ASN	5.8
1	D	187	ALA	5.3
1	D	188	ASN	5.2
1	D	180	SER	5.1

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 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. 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(\AA^2)	Q<0.9
2	MN	A	5033	1/1	0.47	0.18	97,97,97,97	0
2	MN	D	5023	1/1	0.62	0.46	101,101,101,101	0
2	MN	C	5017	1/1	0.66	0.32	86,86,86,86	0
2	MN	B	5009	1/1	0.67	0.30	80,80,80,80	0
2	MN	A	5003	1/1	0.71	0.21	79,79,79,79	0
2	MN	B	5010	1/1	0.75	0.40	99,99,99,99	0
2	MN	D	5024	1/1	0.77	0.08	109,109,109,109	0
2	MN	C	5018	1/1	0.77	0.20	115,115,115,115	0
2	MN	D	5031	1/1	0.80	0.18	83,83,83,83	0
2	MN	A	5013	1/1	0.80	0.14	74,74,74,74	0
2	MN	A	5034	1/1	0.85	0.26	85,85,85,85	0
2	MN	C	5016	1/1	0.86	0.29	101,101,101,101	0
2	MN	C	5025	1/1	0.86	0.13	109,109,109,109	0
2	MN	A	5004	1/1	0.86	0.34	70,70,70,70	0
2	MN	A	5029	1/1	0.86	0.28	67,67,67,67	0
2	MN	C	5019	1/1	0.87	0.10	94,94,94,94	0
2	MN	C	5030	1/1	0.88	0.23	80,80,80,80	0
2	MN	A	5026	1/1	0.88	0.08	64,64,64,64	0
2	MN	D	5035	1/1	0.88	0.09	116,116,116,116	0
2	MN	A	5001	1/1	0.89	0.09	34,34,34,34	0
2	MN	A	5027	1/1	0.89	0.12	52,52,52,52	0
2	MN	D	5022	1/1	0.90	0.33	72,72,72,72	0
2	MN	B	5012	1/1	0.91	0.06	51,51,51,51	0
2	MN	B	5039	1/1	0.92	0.13	50,50,50,50	0
2	MN	C	5032	1/1	0.93	0.33	95,95,95,95	0
2	MN	B	5011	1/1	0.93	0.06	51,51,51,51	0
2	MN	C	5014	1/1	0.94	0.10	50,50,50,50	0
2	MN	B	5028	1/1	0.94	0.06	63,63,63,63	0
2	MN	A	5005	1/1	0.94	0.09	51,51,51,51	0
2	MN	D	5038	1/1	0.95	0.10	88,88,88,88	0
2	MN	B	5037	1/1	0.95	0.15	78,78,78,78	0
2	MN	A	5006	1/1	0.95	0.11	36,36,36,36	0
2	MN	C	5015	1/1	0.96	0.24	40,40,40,40	0
2	MN	C	5036	1/1	0.96	0.09	116,116,116,116	0
2	MN	B	5007	1/1	0.96	0.24	40,40,40,40	0
2	MN	D	5021	1/1	0.97	0.19	36,36,36,36	0
2	MN	B	5008	1/1	0.98	0.22	27,27,27,27	0
2	MN	A	5002	1/1	0.99	0.18	21,21,21,21	0
2	MN	D	5020	1/1	0.99	0.16	41,41,41,41	0

6.5 Other polymers ⓘ

There are no such residues in this entry.