



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:25 PM GMT

PDB ID : 4LJQ
Title : Crystal structure of the catalytic core of E3 ligase HOIP
Authors : Stieglitz, B.; Rana, R.R.; Koliopoulos, M.G.; Morris-Davies, A.C.;
Christodoulou, E.; Howell, S.; Brown, N.R.; Ritinger, K.
Deposited on : 2013-07-05
Resolution : 2.45 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

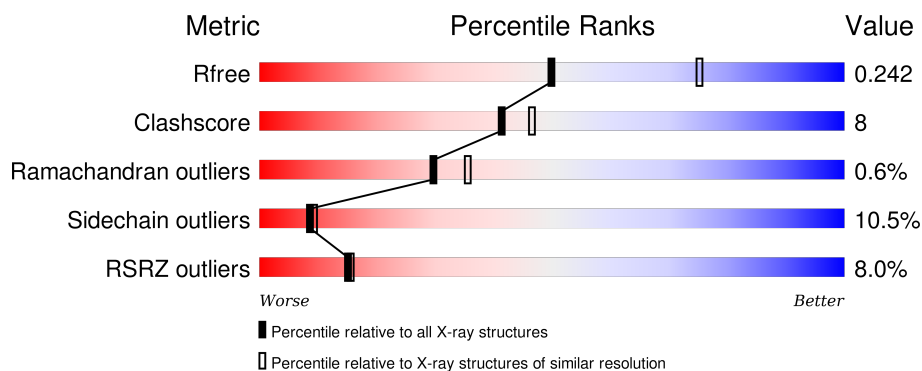
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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}	91344	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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	223	<div> <div>7%</div> <div> <div></div> <div>70%</div> <div>12%</div> <div>•</div> <div>15%</div> </div> </div>
1	B	223	<div> <div>4%</div> <div> <div></div> <div>70%</div> <div>13%</div> <div>•</div> <div>14%</div> </div> </div>
1	C	223	<div> <div>6%</div> <div> <div></div> <div>68%</div> <div>15%</div> <div>•</div> <div>14%</div> </div> </div>
1	D	223	<div> <div>9%</div> <div> <div></div> <div>70%</div> <div>12%</div> <div>•</div> <div>15%</div> </div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5913 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 E3 ubiquitin-protein ligase RNF31.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	192	Total	C	N	O	S	Se	0	0	0
			1468	939	258	255	13	3			
1	A	190	Total	C	N	O	S	Se	0	0	0
			1448	918	253	261	13	3			
1	C	192	Total	C	N	O	S	Se	0	0	0
			1479	941	254	268	13	3			
1	D	190	Total	C	N	O	S	Se	0	0	0
			1457	925	252	264	13	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	850	GLY	-	EXPRESSION TAG	UNP Q96EP0
B	851	PRO	-	EXPRESSION TAG	UNP Q96EP0
B	852	GLY	-	EXPRESSION TAG	UNP Q96EP0
A	850	GLY	-	EXPRESSION TAG	UNP Q96EP0
A	851	PRO	-	EXPRESSION TAG	UNP Q96EP0
A	852	GLY	-	EXPRESSION TAG	UNP Q96EP0
C	850	GLY	-	EXPRESSION TAG	UNP Q96EP0
C	851	PRO	-	EXPRESSION TAG	UNP Q96EP0
C	852	GLY	-	EXPRESSION TAG	UNP Q96EP0
D	850	GLY	-	EXPRESSION TAG	UNP Q96EP0
D	851	PRO	-	EXPRESSION TAG	UNP Q96EP0
D	852	GLY	-	EXPRESSION TAG	UNP Q96EP0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	5	Total	Zn	0	0
			5	5		
2	A	5	Total	Zn	0	0
			5	5		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	4	Total 4	Zn 4	0	0
2	C	4	Total 4	Zn 4	0	0

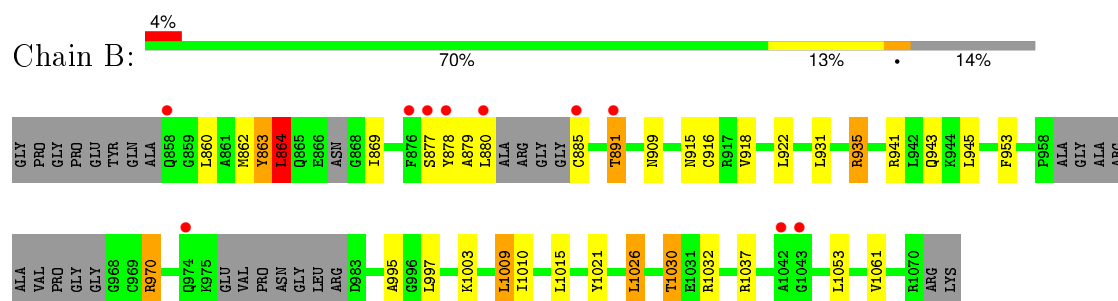
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	21	Total 21	O 21	0	0
3	A	9	Total 9	O 9	0	0
3	C	7	Total 7	O 7	0	0
3	D	6	Total 6	O 6	0	0

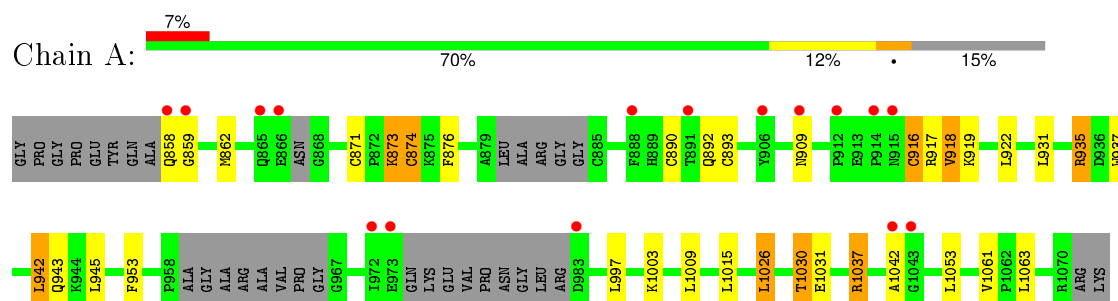
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

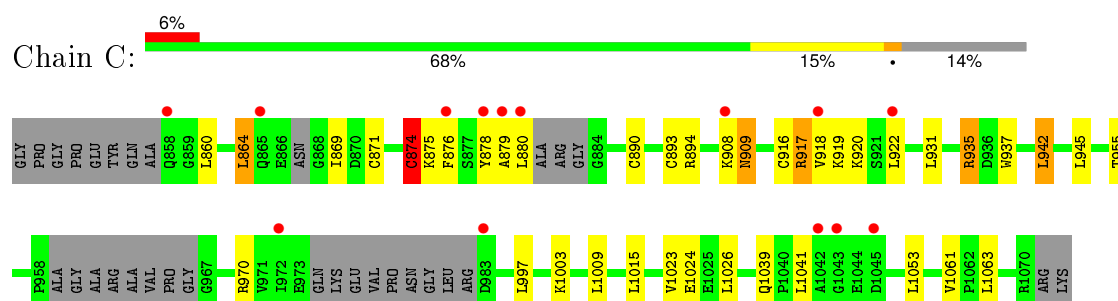
• Molecule 1: E3 ubiquitin-protein ligase RNF31



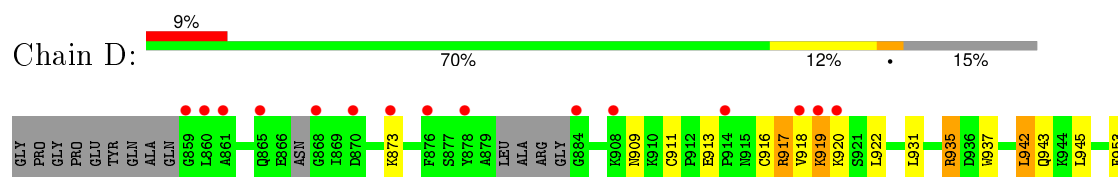
• Molecule 1: E3 ubiquitin-protein ligase RNF31

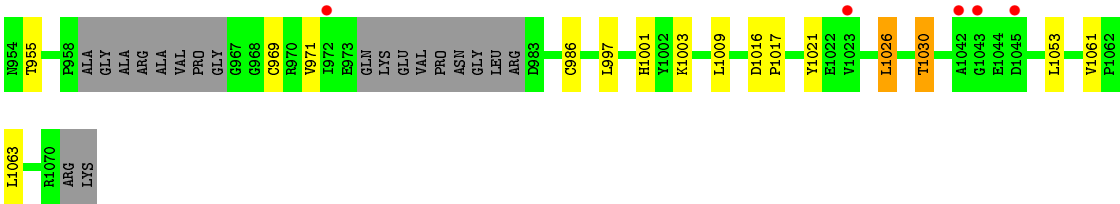


• Molecule 1: E3 ubiquitin-protein ligase RNF31



• Molecule 1: E3 ubiquitin-protein ligase RNF31





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.21Å 47.77Å 111.14Å 101.39° 90.12° 98.92°	Depositor
Resolution (Å)	29.60 – 2.45 29.61 – 2.45	Depositor EDS
% Data completeness (in resolution range)	95.0 (29.60-2.45) 90.4 (29.61-2.45)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.206 , 0.243 0.209 , 0.242	Depositor DCC
R_{free} test set	2008 reflections (6.94%)	DCC
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.9	EDS
Estimated twinning fraction	0.014 for -h,-k,k+l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 30949 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5913	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.06% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN

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	0.70	1/1481 (0.1%)	0.86	1/2005 (0.0%)
1	B	0.77	0/1500	0.92	8/2028 (0.4%)
1	C	0.67	0/1513	0.83	0/2046
1	D	0.62	0/1491	0.80	0/2017
All	All	0.69	1/5985 (0.0%)	0.85	9/8096 (0.1%)

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	0	1
1	C	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	874	CYS	CB-SG	-6.30	1.71	1.82

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	874	CYS	CB-CA-C	-9.82	90.76	110.40
1	B	878	TYR	N-CA-C	6.43	128.36	111.00
1	B	863	TYR	N-CA-C	-6.26	94.09	111.00
1	B	864	LEU	N-CA-CB	5.83	122.06	110.40
1	B	941	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	B	862	MSE	CG-SE-CE	5.25	110.44	98.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	995	ALA	C-N-CA	-5.15	111.49	122.30
1	B	1037	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	B	885	CYS	CA-CB-SG	5.03	123.05	114.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	916	CYS	Peptide
1	C	874	CYS	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	1448	0	1339	28	0
1	B	1468	0	1369	21	0
1	C	1479	0	1375	27	0
1	D	1457	0	1347	21	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	4	0	0	0	0
2	D	4	0	0	0	0
3	A	9	0	0	1	0
3	B	21	0	0	1	0
3	C	7	0	0	1	0
3	D	6	0	0	0	0
All	All	5913	0	5430	96	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 (96) 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:A:858:GLN:HB3	1:A:859:GLY:HA2	1.29	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1009:LEU:CA	1:B:1010:ILE:N	2.24	1.00
1:A:858:GLN:CB	1:A:859:GLY:HA2	1.91	1.00
1:B:1009:LEU:O	1:B:1010:ILE:N	1.98	0.97
1:A:874:CYS:HB3	1:A:876:PHE:H	1.32	0.93
1:C:874:CYS:HB3	1:C:876:PHE:H	1.34	0.91
1:D:916:CYS:HA	1:D:917:ARG:CB	2.06	0.85
1:A:1037:ARG:O	1:A:1037:ARG:HG2	1.72	0.84
1:C:916:CYS:HA	1:C:917:ARG:CB	2.05	0.84
1:B:860:LEU:O	1:B:863:TYR:O	1.98	0.82
1:D:969:CYS:SG	1:D:971:VAL:HG22	2.20	0.80
1:A:858:GLN:HB3	1:A:859:GLY:CA	2.13	0.77
1:D:971:VAL:HG21	1:D:1001:HIS:ND1	2.04	0.71
1:A:1042:ALA:O	3:A:1205:HOH:O	2.08	0.71
1:C:874:CYS:HB3	1:C:876:PHE:N	2.04	0.71
1:D:971:VAL:HG21	1:D:1001:HIS:CE1	2.25	0.71
1:B:1009:LEU:O	1:B:1009:LEU:CA	2.39	0.70
1:C:878:TYR:OH	1:C:890:CYS:O	2.09	0.70
1:C:874:CYS:HB3	1:C:875:LYS:CA	2.20	0.69
1:A:917:ARG:C	1:A:919:LYS:H	1.95	0.69
1:A:858:GLN:CG	1:A:859:GLY:HA2	2.22	0.68
1:B:891:THR:HG21	1:C:920:LYS:O	1.95	0.66
1:C:918:VAL:HG12	1:C:918:VAL:O	1.98	0.64
1:B:1030:THR:HG21	1:B:1053:LEU:HD22	1.80	0.63
1:A:871:CYS:HB3	1:A:874:CYS:HB2	1.79	0.63
1:D:918:VAL:HG12	1:D:918:VAL:O	1.99	0.62
1:C:871:CYS:HB3	1:C:874:CYS:HB2	1.80	0.62
1:A:918:VAL:HG12	1:A:918:VAL:O	1.99	0.62
1:A:858:GLN:HG2	1:A:862:MSE:HE3	1.82	0.61
1:A:874:CYS:HB3	1:A:876:PHE:N	2.10	0.61
1:B:879:ALA:O	1:B:880:LEU:C	2.40	0.60
1:A:858:GLN:CB	1:A:859:GLY:CA	2.75	0.59
1:A:890:CYS:O	1:A:893:CYS:O	2.22	0.57
1:D:913:GLU:CG	1:D:916:CYS:HB2	2.36	0.56
1:C:874:CYS:CB	1:C:875:LYS:HA	2.37	0.55
1:C:1023:VAL:HG12	1:C:1024:GLU:OE2	2.06	0.55
1:C:919:LYS:O	1:C:920:LYS:HB2	2.07	0.55
1:C:916:CYS:CA	1:C:917:ARG:CB	2.83	0.54
1:D:971:VAL:HG21	1:D:1001:HIS:CG	2.41	0.54
1:C:874:CYS:CB	1:C:875:LYS:CA	2.86	0.53
1:D:971:VAL:HG23	1:D:986:CYS:HB2	1.90	0.53
1:A:1026:LEU:O	1:A:1030:THR:HG23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:919:LYS:O	1:D:920:LYS:CB	2.57	0.52
1:D:916:CYS:CA	1:D:917:ARG:CB	2.84	0.52
1:D:937:TRP:CE3	1:D:942:LEU:HD13	2.44	0.52
1:B:880:LEU:HD23	1:B:880:LEU:H	1.75	0.52
1:A:874:CYS:CB	1:A:876:PHE:H	2.15	0.50
1:A:1037:ARG:CG	1:A:1037:ARG:O	2.51	0.50
1:C:937:TRP:CE3	1:C:942:LEU:HD13	2.46	0.50
1:C:874:CYS:CB	1:C:876:PHE:H	2.15	0.50
1:B:1026:LEU:O	1:B:1030:THR:HG23	2.12	0.50
1:D:911:CYS:HB3	1:D:919:LYS:HE2	1.94	0.49
1:D:1026:LEU:O	1:D:1030:THR:HG23	2.12	0.49
1:B:943:GLN:HG2	1:B:953:PHE:CE2	2.49	0.47
1:D:943:GLN:HG2	1:D:953:PHE:CE2	2.49	0.47
1:B:1009:LEU:O	1:B:1010:ILE:CA	2.61	0.47
1:A:917:ARG:C	1:A:919:LYS:N	2.63	0.47
1:C:1003:LYS:HE3	1:C:1003:LYS:HB2	1.75	0.47
1:A:937:TRP:CE3	1:A:942:LEU:HD13	2.50	0.46
1:A:943:GLN:HG2	1:A:953:PHE:CE2	2.50	0.46
1:C:931:LEU:O	1:C:935:ARG:HD2	2.16	0.46
1:B:1026:LEU:O	1:B:1030:THR:CG2	2.64	0.46
1:B:1003:LYS:HB2	1:B:1003:LYS:HE3	1.78	0.46
1:C:878:TYR:CZ	1:C:890:CYS:O	2.68	0.45
1:A:917:ARG:O	1:A:919:LYS:N	2.46	0.45
1:C:893:CYS:O	1:C:894:ARG:CB	2.64	0.45
1:A:876:PHE:CG	1:A:892:GLN:HG3	2.52	0.45
1:D:931:LEU:O	1:D:935:ARG:HD2	2.17	0.45
1:C:860:LEU:HD23	1:C:864:LEU:HD22	1.98	0.45
1:C:970:ARG:CG	3:C:1202:HOH:O	2.65	0.44
1:D:1003:LYS:HE3	1:D:1003:LYS:HB2	1.77	0.44
1:B:863:TYR:O	1:B:864:LEU:CB	2.64	0.44
1:D:971:VAL:HG13	1:D:997:LEU:HD22	1.99	0.44
1:D:1026:LEU:O	1:D:1030:THR:CG2	2.66	0.44
1:C:908:LYS:HG3	1:C:920:LYS:HA	2.00	0.44
1:C:919:LYS:O	1:C:920:LYS:CB	2.66	0.44
1:A:1026:LEU:O	1:A:1030:THR:CG2	2.65	0.44
1:B:1032:ARG:NH1	3:B:1206:HOH:O	2.34	0.43
1:B:869:ILE:HG23	1:B:869:ILE:O	2.18	0.43
1:C:874:CYS:HB3	1:C:875:LYS:HA	1.95	0.43
1:B:931:LEU:O	1:B:935:ARG:HD2	2.18	0.43
1:C:879:ALA:O	1:C:880:LEU:C	2.56	0.43
1:A:931:LEU:O	1:A:935:ARG:HD2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1003:LYS:HB2	1:A:1003:LYS:HE3	1.79	0.43
1:B:970:ARG:CG	1:B:970:ARG:HH11	2.32	0.43
1:A:873:LYS:HE2	1:A:873:LYS:HB3	1.74	0.43
1:A:858:GLN:HG3	1:A:859:GLY:HA2	1.98	0.43
1:B:916:CYS:SG	1:B:918:VAL:HG22	2.60	0.42
1:A:916:CYS:HB3	1:A:917:ARG:HA	2.02	0.42
1:B:863:TYR:O	1:B:864:LEU:HB2	2.20	0.42
1:B:1021:TYR:HD2	1:B:1026:LEU:HD13	1.86	0.41
1:C:937:TRP:HE3	1:C:942:LEU:HD13	1.85	0.41
1:D:937:TRP:HE3	1:D:942:LEU:HD13	1.85	0.41
1:C:909:ASN:HB2	1:C:919:LYS:HG3	2.02	0.41
1:D:1016:ASP:HA	1:D:1017:PRO:HD3	1.94	0.41
1:D:1021:TYR:HD2	1:D:1026:LEU:HD13	1.87	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	180/223 (81%)	168 (93%)	11 (6%)	1 (1%)	30	35
1	B	180/223 (81%)	173 (96%)	6 (3%)	1 (1%)	30	35
1	C	182/223 (82%)	172 (94%)	9 (5%)	1 (0%)	34	41
1	D	180/223 (81%)	172 (96%)	7 (4%)	1 (1%)	30	35
All	All	722/892 (81%)	685 (95%)	33 (5%)	4 (1%)	30	35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	917	ARG
1	D	917	ARG

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Mol	Chain	Res	Type
1	B	864	LEU
1	A	918	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	146/182 (80%)	130 (89%)	16 (11%)	8	8
1	B	145/182 (80%)	131 (90%)	14 (10%)	10	12
1	C	150/182 (82%)	132 (88%)	18 (12%)	6	6
1	D	148/182 (81%)	134 (90%)	14 (10%)	11	12
All	All	589/728 (81%)	527 (90%)	62 (10%)	8	9

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	877	SER
1	B	891	THR
1	B	909	ASN
1	B	915	ASN
1	B	922	LEU
1	B	935	ARG
1	B	945	LEU
1	B	970	ARG
1	B	997	LEU
1	B	1009	LEU
1	B	1015	LEU
1	B	1026	LEU
1	B	1030	THR
1	B	1061	VAL
1	A	873	LYS
1	A	909	ASN
1	A	922	LEU
1	A	935	ARG
1	A	942	LEU

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Mol	Chain	Res	Type
1	A	945	LEU
1	A	997	LEU
1	A	1009	LEU
1	A	1015	LEU
1	A	1026	LEU
1	A	1030	THR
1	A	1031	GLU
1	A	1037	ARG
1	A	1053	LEU
1	A	1061	VAL
1	A	1063	LEU
1	C	864	LEU
1	C	869	ILE
1	C	874	CYS
1	C	909	ASN
1	C	922	LEU
1	C	935	ARG
1	C	942	LEU
1	C	945	LEU
1	C	955	THR
1	C	997	LEU
1	C	1009	LEU
1	C	1015	LEU
1	C	1026	LEU
1	C	1039	GLN
1	C	1041	LEU
1	C	1053	LEU
1	C	1061	VAL
1	C	1063	LEU
1	D	873	LYS
1	D	909	ASN
1	D	919	LYS
1	D	922	LEU
1	D	935	ARG
1	D	942	LEU
1	D	945	LEU
1	D	955	THR
1	D	1009	LEU
1	D	1026	LEU
1	D	1030	THR
1	D	1053	LEU
1	D	1061	VAL

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Mol	Chain	Res	Type
1	D	1063	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	887	HIS
1	A	889	HIS
1	D	949	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 18 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 ⓘ

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	187/223 (83%)	0.36	16 (8%) 13 13	26, 45, 88, 107	0
1	B	189/223 (84%)	0.15	10 (5%) 30 33	17, 36, 72, 95	0
1	C	189/223 (84%)	0.46	14 (7%) 17 18	24, 47, 86, 107	0
1	D	187/223 (83%)	0.54	20 (10%) 8 7	33, 52, 91, 119	0
All	All	752/892 (84%)	0.38	60 (7%) 15 15	17, 45, 87, 119	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	876	PHE	5.6
1	D	861	ALA	5.1
1	B	876	PHE	4.3
1	D	918	VAL	4.3
1	D	1045	ASP	4.1
1	A	1043	GLY	4.0
1	A	915	ASN	3.9
1	C	880	LEU	3.9
1	D	1042	ALA	3.8
1	A	888	PHE	3.5
1	D	873	LYS	3.5
1	B	1042	ALA	3.4
1	A	914	PRO	3.4
1	C	858	GLN	3.4
1	D	920	LYS	3.3
1	D	868	GLY	3.3
1	B	1043	GLY	3.2
1	D	919	LYS	3.2
1	A	858	GLN	3.1
1	A	865	GLN	3.1
1	C	878	TYR	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	878	TYR	3.0
1	C	865	GLN	3.0
1	D	972	ILE	3.0
1	A	1042	ALA	2.9
1	A	909	ASN	2.9
1	A	983	ASP	2.9
1	D	859	GLY	2.8
1	C	908	LYS	2.8
1	C	1045	ASP	2.8
1	D	884	GLY	2.8
1	B	880	LEU	2.8
1	D	865	GLN	2.7
1	D	1023	VAL	2.7
1	C	876	PHE	2.7
1	B	891	THR	2.6
1	D	860	LEU	2.6
1	A	866	GLU	2.6
1	A	972	ILE	2.6
1	C	972	ILE	2.6
1	C	1043	GLY	2.6
1	C	1042	ALA	2.6
1	C	983	ASP	2.5
1	A	859	GLY	2.5
1	C	918	VAL	2.5
1	A	973	GLU	2.5
1	B	974	GLN	2.5
1	D	908	LYS	2.5
1	B	877	SER	2.4
1	A	912	PRO	2.3
1	B	858	GLN	2.2
1	A	891	THR	2.2
1	D	870	ASP	2.2
1	B	885	CYS	2.2
1	D	1043	GLY	2.2
1	C	922	LEU	2.2
1	D	914	PRO	2.2
1	D	878	TYR	2.1
1	C	879	ALA	2.1
1	A	906	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no 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
2	ZN	C	1102	1/1	0.99	0.16	0.68	33,33,33,33	0
2	ZN	B	1102	1/1	1.00	0.16	-0.62	22,22,22,22	0
2	ZN	A	1102	1/1	1.00	0.12	-1.09	27,27,27,27	0
2	ZN	D	1102	1/1	0.99	0.14	-1.30	35,35,35,35	0
2	ZN	A	1104	1/1	1.00	0.10	-1.47	41,41,41,41	0
2	ZN	D	1104	1/1	1.00	0.10	-1.55	38,38,38,38	0
2	ZN	C	1104	1/1	0.99	0.10	-1.58	31,31,31,31	0
2	ZN	D	1101	1/1	0.99	0.04	-1.72	69,69,69,69	0
2	ZN	B	1103	1/1	1.00	0.10	-1.75	35,35,35,35	0
2	ZN	B	1101	1/1	0.98	0.03	-1.76	49,49,49,49	0
2	ZN	C	1101	1/1	0.99	0.05	-2.04	46,46,46,46	0
2	ZN	C	1103	1/1	0.99	0.06	-2.21	51,51,51,51	0
2	ZN	B	1104	1/1	1.00	0.10	-2.43	29,29,29,29	0
2	ZN	A	1101	1/1	0.95	0.06	-2.91	60,60,60,60	0
2	ZN	A	1105	1/1	0.99	0.16	-	39,39,39,39	0
2	ZN	B	1105	1/1	1.00	0.15	-	30,30,30,30	0
2	ZN	D	1103	1/1	0.99	0.08	-	58,58,58,58	0
2	ZN	A	1103	1/1	0.98	0.07	-	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.