



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:14 AM UTC

PDB ID : 5JH5 / pdb_00005jh5
Title : Structural Basis for the Hierarchical Assembly of the Core of PRC1.1
Authors : Wong, S.J.; Taylor, A.B.; Hart, P.J.; Kim, C.A.
Deposited on : 2016-04-20
Resolution : 2.55 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
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

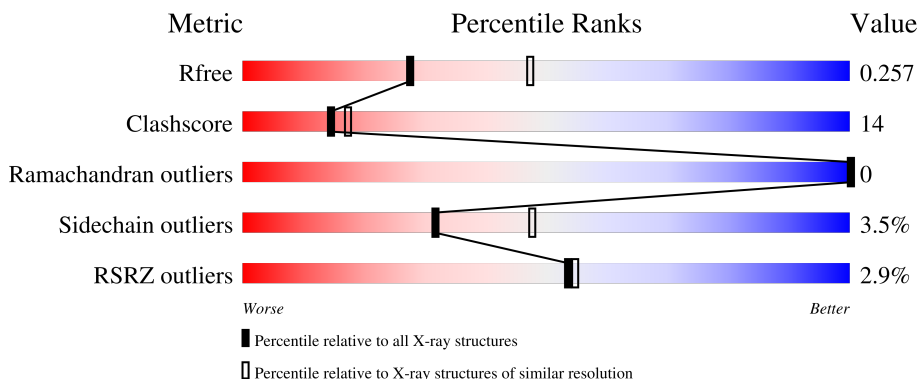
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	281	2% 59% 34% • 5%
2	B	162	6% 64% 23% • 8%
3	C	109	55% 26% 19%
4	D	122	2% 69% 15% • 14%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5015 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 Lysine-specific demethylase 2B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	Se	0	0	0
			2140	1344	389	384	16	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1056	GLY	-	expression tag	UNP Q8NHM5
A	1057	THR	-	expression tag	UNP Q8NHM5
A	1058	ARG	-	expression tag	UNP Q8NHM5

- Molecule 2 is a protein called S-phase kinase-associated protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	149	Total	C	N	O	S	Se	0	0	0
			1196	760	196	235	2	3			

- Molecule 3 is a protein called Polycomb group RING finger protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	88	Total	C	N	O	S	Se	0	0	0
			755	483	143	123	3	3			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	147	GLY	-	expression tag	UNP Q9BSM1
C	148	THR	-	expression tag	UNP Q9BSM1
C	149	ARG	-	expression tag	UNP Q9BSM1

- Molecule 4 is a protein called BCL-6 corepressor-like protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	D	105	Total	C	N	O	S	Se	0	0	0
			860	555	143	157	3	2			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1590	MSE	-	initiating methionine	UNP Q5H9F3
D	1591	GLU	-	expression tag	UNP Q5H9F3
D	1592	THR	-	expression tag	UNP Q5H9F3
D	1593	ARG	-	expression tag	UNP Q5H9F3

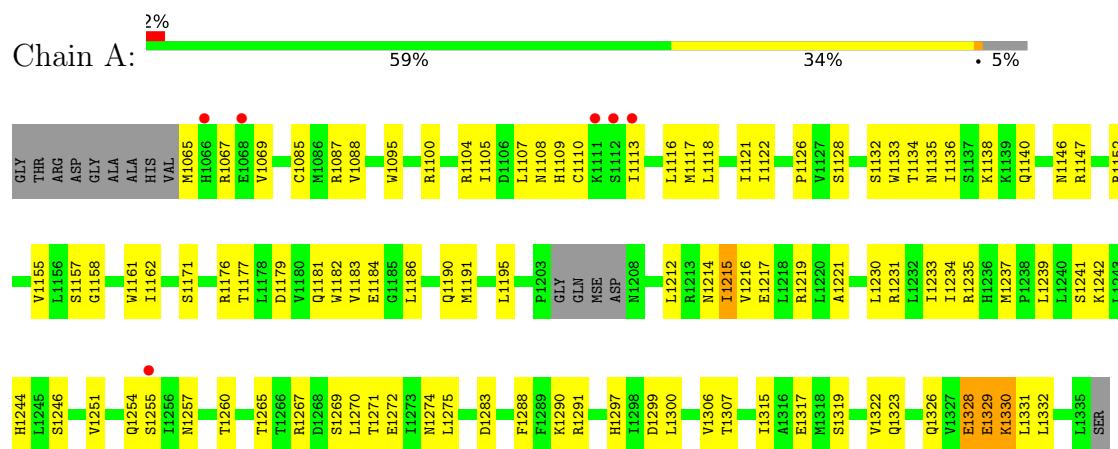
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	22	Total	O	0	0
			22	22		
5	B	10	Total	O	0	0
			10	10		
5	C	16	Total	O	0	0
			16	16		
5	D	16	Total	O	0	0
			16	16		

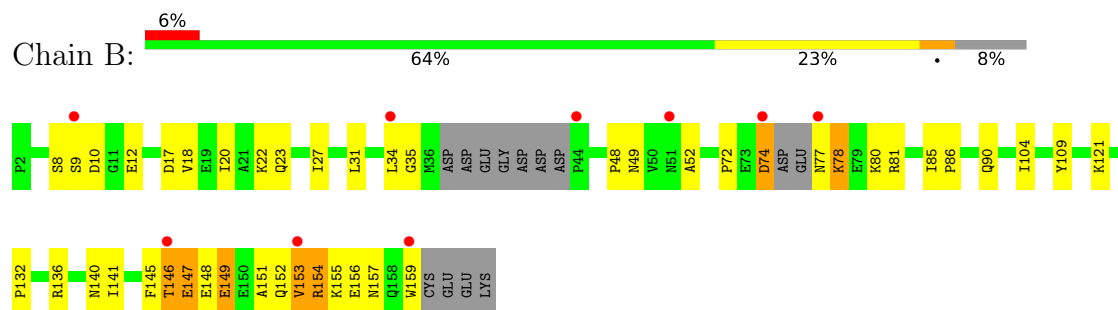
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

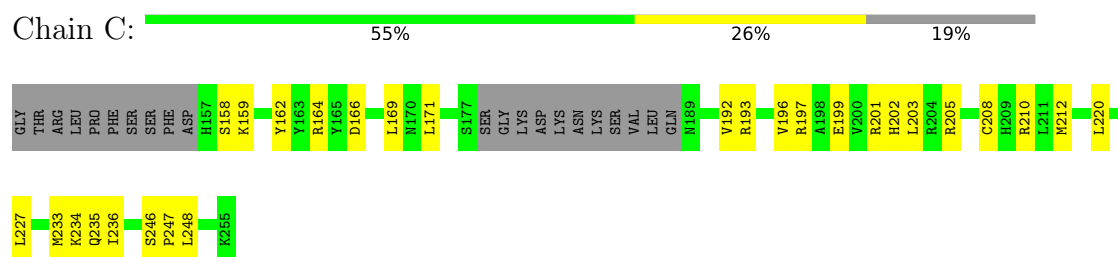
- Molecule 1: Lysine-specific demethylase 2B



- Molecule 2: S-phase kinase-associated protein 1



- Molecule 3: Polycomb group RING finger protein 1



- Molecule 4: BCL-6 corepressor-like protein 1



MSE	GLU	THR	ARG	D1594	M1597	F1598	E1599	L1606	Y1609	R1617	E1647	T1650	K1653	F1656	V1660	L1665	L1666	T1667	P1668	R1671	P1672	G1673	GLY	LEU	ASP	ASP	ARG	SER	PRO	GLY	SER	SER	E1685	T1686	P1694	E1703	V1704	E1705	S1708	C1709	ASN	SER
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.05Å 73.80Å 123.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.28 – 2.55 63.28 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.5 (63.28-2.55) 99.7 (63.28-2.55)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.55Å)	Xtriage
Refinement program	PHENIX (1.10_2155)	Depositor
R, R_{free}	0.206 , 0.257 0.207 , 0.257	Depositor DCC
R_{free} test set	1999 reflections (9.34%)	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.219	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5015	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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](#)

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	0/2171	0.45	0/2927
2	B	0.30	1/1214 (0.1%)	0.48	1/1636 (0.1%)
3	C	0.19	0/771	0.40	0/1034
4	D	0.16	0/879	0.36	0/1185
All	All	0.24	1/5035 (0.0%)	0.43	1/6782 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	153	VAL	C-N	-5.87	1.26	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	146	THR	N-CA-C	-5.72	102.44	110.50

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	2140	0	2192	76	0
2	B	1196	0	1200	39	0
3	C	755	0	764	23	0
4	D	860	0	851	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	22	0	0	0	0
5	B	10	0	0	0	0
5	C	16	0	0	0	0
5	D	16	0	0	0	0
All	All	5015	0	5007	139	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 14.

All (139) 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:1306:VAL:O	1:A:1330:LYS:NZ	1.77	1.18
2:B:152:GLN:O	2:B:156:GLU:HB2	1.62	0.98
3:C:193:ARG:NH2	4:D:1606:LEU:O	2.05	0.87
2:B:152:GLN:O	2:B:156:GLU:CB	2.25	0.84
2:B:152:GLN:O	2:B:156:GLU:N	2.13	0.82
1:A:1108:ASN:OD1	1:A:1109:HIS:ND1	2.15	0.79
2:B:34:LEU:HD12	2:B:35:GLY:N	1.97	0.78
2:B:154:ARG:HG3	2:B:155:LYS:N	2.01	0.76
1:A:1319:SER:HA	1:A:1322:VAL:O	1.87	0.75
1:A:1230:LEU:HD22	1:A:1255:SER:HB2	1.70	0.73
2:B:72:PRO:HG2	2:B:78:LYS:HE3	1.69	0.73
2:B:74:ASP:N	2:B:74:ASP:OD1	2.19	0.72
1:A:1107:LEU:O	1:A:1110:CYS:HB2	1.93	0.68
1:A:1152:ARG:HB3	1:A:1176:ARG:HG2	1.75	0.67
2:B:34:LEU:HD12	2:B:35:GLY:H	1.59	0.67
1:A:1329:GLU:HA	1:A:1329:GLU:OE2	1.93	0.66
2:B:153:VAL:O	2:B:156:GLU:HB3	1.95	0.66
2:B:81:ARG:O	2:B:121:LYS:HE2	1.95	0.66
1:A:1297:HIS:CE1	1:A:1331:LEU:HD11	2.31	0.66
2:B:74:ASP:O	2:B:78:LYS:O	2.14	0.65
4:D:1668:PRO:HA	4:D:1671:ARG:HD2	1.78	0.64
2:B:152:GLN:O	2:B:156:GLU:CA	2.45	0.64
3:C:203:LEU:HD23	3:C:220:LEU:HD21	1.80	0.64
1:A:1219:ARG:HG2	1:A:1244:HIS:CD2	2.32	0.63
1:A:1104:ARG:HG2	1:A:1128:SER:HB2	1.79	0.63
1:A:1122:ILE:HG12	1:A:1147:ARG:HB2	1.80	0.62
1:A:1283:ASP:OD2	1:A:1307:THR:OG1	2.14	0.62
1:A:1107:LEU:HD11	1:A:1121:ILE:HD11	1.82	0.62
2:B:151:ALA:O	2:B:155:LYS:HB3	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:1650:THR:CG2	4:D:1686:THR:HB	2.31	0.60
1:A:1328:GLU:O	1:A:1328:GLU:HG3	2.00	0.60
2:B:132:PRO:O	2:B:136:ARG:HG3	2.01	0.60
1:A:1326:GLN:HG2	1:A:1332:LEU:HD23	1.82	0.60
1:A:1184:GLU:HB2	3:C:197:ARG:HD2	1.83	0.60
1:A:1113:ILE:HG12	1:A:1117:MSE:HE1	1.85	0.59
1:A:1315:ILE:HD11	1:A:1326:GLN:HG3	1.84	0.59
1:A:1195:LEU:HD23	1:A:1212:LEU:HD12	1.83	0.59
4:D:1617:ARG:NH1	4:D:1647:GLU:OE2	2.37	0.57
2:B:17:ASP:HB2	2:B:20:ILE:HD12	1.85	0.57
1:A:1107:LEU:HD13	1:A:1117:MSE:HE3	1.88	0.55
3:C:202:HIS:O	3:C:205:ARG:HB2	2.06	0.55
2:B:27:ILE:O	2:B:31:LEU:HD13	2.06	0.55
2:B:18:VAL:O	2:B:22:LYS:HG3	2.07	0.55
1:A:1179:ASP:OD1	1:A:1181:GLN:HG3	2.08	0.54
1:A:1271:THR:HG22	1:A:1272:GLU:HG3	1.89	0.54
1:A:1186:LEU:HD12	1:A:1190:GLN:HB2	1.88	0.54
1:A:1110:CYS:H	1:A:1134:THR:HG22	1.74	0.53
1:A:1290:LYS:NZ	1:A:1317:GLU:OE1	2.34	0.52
1:A:1067:ARG:HH12	1:A:1100:ARG:NH2	2.08	0.52
2:B:34:LEU:CD1	2:B:35:GLY:H	2.22	0.52
1:A:1182:TRP:CH2	3:C:235:GLN:HG2	2.45	0.52
1:A:1113:ILE:CG2	1:A:1118:LEU:HD21	2.39	0.52
1:A:1067:ARG:HH12	1:A:1100:ARG:HH22	1.56	0.52
3:C:208:CYS:O	3:C:212:MSE:N	2.43	0.52
3:C:159:LYS:HG3	3:C:162:TYR:CZ	2.44	0.52
1:A:1155:VAL:HA	1:A:1179:ASP:HB3	1.92	0.51
1:A:1214:ASN:HA	1:A:1239:LEU:HD12	1.93	0.51
1:A:1085:CYS:HA	1:A:1088:VAL:HG23	1.93	0.51
4:D:1653:LYS:HZ2	4:D:1685:GLU:HG2	1.76	0.51
1:A:1133:TRP:CE3	1:A:1158:GLY:HA3	2.46	0.51
2:B:154:ARG:HG3	2:B:155:LYS:H	1.74	0.51
1:A:1190:GLN:OE1	1:A:1190:GLN:N	2.41	0.51
3:C:210:ARG:NH2	4:D:1703:GLU:OE1	2.43	0.51
1:A:1217:GLU:OE2	1:A:1219:ARG:NH2	2.44	0.50
1:A:1182:TRP:CZ2	3:C:235:GLN:HG2	2.46	0.50
1:A:1065:MSE:HE2	2:B:141:ILE:HD12	1.94	0.50
1:A:1257:ASN:HB2	1:A:1288:PHE:CE2	2.46	0.50
1:A:1219:ARG:HA	1:A:1244:HIS:HB2	1.94	0.49
3:C:166:ASP:OD2	3:C:197:ARG:HD3	2.12	0.49
1:A:1067:ARG:NH1	1:A:1100:ARG:HH22	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1138:LYS:HB2	1:A:1162:ILE:O	2.12	0.49
1:A:1231:ARG:HA	1:A:1234:ILE:HG12	1.94	0.49
3:C:202:HIS:HE2	4:D:1709:CYS:HA	1.77	0.49
1:A:1234:ILE:HG13	1:A:1235:ARG:N	2.28	0.48
1:A:1217:GLU:HG3	1:A:1242:LYS:HB2	1.96	0.48
2:B:145:PHE:HD1	2:B:149:GLU:HB3	1.78	0.48
1:A:1113:ILE:HD13	1:A:1136:ILE:HG13	1.95	0.48
1:A:1069:VAL:HG13	2:B:104:ILE:HG21	1.95	0.47
3:C:193:ARG:NH1	4:D:1599:GLU:OE1	2.47	0.47
4:D:1660:VAL:HG13	4:D:1666:LEU:HD22	1.96	0.47
1:A:1105:ILE:HD12	1:A:1126:PRO:HG3	1.97	0.47
4:D:1609:TYR:HE1	4:D:1666:LEU:HD21	1.80	0.47
1:A:1087:ARG:O	2:B:136:ARG:NH2	2.47	0.47
2:B:85:ILE:HD11	2:B:121:LYS:HB3	1.96	0.47
2:B:146:THR:O	2:B:149:GLU:N	2.47	0.47
3:C:169:LEU:HD21	3:C:233:MSE:HB3	1.96	0.47
1:A:1113:ILE:HG12	1:A:1117:MSE:CE	2.44	0.47
1:A:1272:GLU:HA	1:A:1297:HIS:HB3	1.97	0.47
2:B:86:PRO:O	2:B:90:GLN:HG3	2.15	0.47
1:A:1265:THR:O	1:A:1269:SER:HB2	2.15	0.46
3:C:169:LEU:C	3:C:169:LEU:HD12	2.40	0.46
3:C:164:ARG:NE	4:D:1597:MSE:HE2	2.30	0.46
1:A:1161:TRP:CE3	1:A:1186:LEU:HD13	2.50	0.46
1:A:1231:ARG:O	1:A:1234:ILE:HG13	2.15	0.46
2:B:148:GLU:O	2:B:152:GLN:HB2	2.15	0.46
2:B:136:ARG:O	2:B:140:ASN:N	2.48	0.46
2:B:155:LYS:O	2:B:155:LYS:HG2	2.15	0.46
1:A:1116:LEU:HD13	2:B:156:GLU:O	2.16	0.46
4:D:1597:MSE:SE	4:D:1694:PRO:HG3	2.66	0.46
3:C:158:SER:C	3:C:159:LYS:HD2	2.41	0.45
1:A:1132:SER:HB2	1:A:1157:SER:HB2	1.97	0.45
1:A:1162:ILE:H	1:A:1162:ILE:HD12	1.80	0.45
2:B:8:SER:HB2	2:B:10:ASP:OD1	2.17	0.45
3:C:210:ARG:NH1	4:D:1705:GLU:OE2	2.50	0.44
3:C:227:LEU:HD13	3:C:236:ILE:HD13	1.99	0.44
3:C:246:SER:OG	3:C:247:PRO:HA	2.17	0.44
1:A:1146:ASN:OD1	1:A:1171:SER:HB2	2.18	0.43
3:C:196:VAL:HB	3:C:234:LYS:HB2	1.99	0.43
4:D:1656:PHE:C	4:D:1656:PHE:CD1	2.97	0.43
2:B:12:GLU:OE1	2:B:52:ALA:HB1	2.18	0.43
2:B:34:LEU:CD1	2:B:35:GLY:N	2.76	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:74:ASP:HB2	2:B:77:ASN:HB2	2.01	0.43
1:A:1135:ASN:ND2	4:D:1594:ASP:OD2	2.50	0.43
1:A:1260:THR:O	1:A:1291:ARG:HD3	2.18	0.43
1:A:1329:GLU:HB3	1:A:1330:LYS:H	1.54	0.43
3:C:199:GLU:OE1	3:C:201:ARG:NE	2.49	0.43
1:A:1191:MSE:HE1	1:A:1233:ILE:HD11	2.01	0.43
1:A:1212:LEU:HD22	1:A:1215:ILE:HD12	2.02	0.42
2:B:147:GLU:OE1	2:B:147:GLU:HA	2.19	0.42
1:A:1087:ARG:NH2	2:B:132:PRO:HG2	2.34	0.42
1:A:1107:LEU:CD1	1:A:1117:MSE:HE3	2.49	0.42
1:A:1267:ARG:NH2	1:A:1291:ARG:HB3	2.35	0.41
3:C:192:VAL:HA	4:D:1599:GLU:O	2.20	0.41
1:A:1274:ASN:HA	1:A:1299:ASP:HB3	2.02	0.41
1:A:1087:ARG:NE	2:B:132:PRO:HG2	2.35	0.41
4:D:1660:VAL:CG1	4:D:1666:LEU:HD22	2.50	0.41
1:A:1067:ARG:HG2	1:A:1095:TRP:CE2	2.56	0.41
4:D:1653:LYS:HB2	4:D:1653:LYS:HE3	1.81	0.41
1:A:1254:GLN:HA	1:A:1254:GLN:OE1	2.20	0.41
1:A:1195:LEU:CD2	1:A:1237:MSE:HE3	2.51	0.41
1:A:1275:LEU:HB2	1:A:1300:LEU:HD23	2.02	0.41
3:C:171:LEU:HD23	3:C:248:LEU:HB3	2.03	0.41
1:A:1177:THR:HG22	1:A:1217:GLU:HB3	2.03	0.41
1:A:1221:ALA:HA	1:A:1246:SER:O	2.21	0.41
1:A:1216:VAL:HG13	1:A:1241:SER:HB3	2.02	0.40
2:B:9:SER:HB3	2:B:48:PRO:C	2.46	0.40
2:B:49:ASN:HB3	2:B:109:TYR:CD2	2.56	0.40
1:A:1118:LEU:HD11	1:A:1140:GLN:HG2	2.04	0.40
1:A:1323:GLN:HA	1:A:1323:GLN:NE2	2.36	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	263/281 (94%)	254 (97%)	9 (3%)	0	100	100
2	B	143/162 (88%)	139 (97%)	4 (3%)	0	100	100
3	C	84/109 (77%)	84 (100%)	0	0	100	100
4	D	101/122 (83%)	101 (100%)	0	0	100	100
All	All	591/674 (88%)	578 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	250/251 (100%)	243 (97%)	7 (3%)	38	57
2	B	137/146 (94%)	128 (93%)	9 (7%)	15	21
3	C	85/101 (84%)	85 (100%)	0	100	100
4	D	98/110 (89%)	94 (96%)	4 (4%)	27	41
All	All	570/608 (94%)	550 (96%)	20 (4%)	32	48

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1183	VAL
1	A	1215	ILE
1	A	1251	VAL
1	A	1270	LEU
1	A	1328	GLU
1	A	1329	GLU
1	A	1330	LYS
2	B	23	GLN
2	B	74	ASP
2	B	78	LYS
2	B	80	LYS
2	B	147	GLU
2	B	149	GLU

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Mol	Chain	Res	Type
2	B	154	ARG
2	B	157	ASN
2	B	159	TRP
4	D	1597	MSE
4	D	1665	LEU
4	D	1666	LEU
4	D	1709	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1214	ASN
2	B	23	GLN
2	B	64	HIS
2	B	97	GLN
3	C	224	ASN
4	D	1645	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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	260/281 (92%)	0.23	6 (2%) 61 62	27, 47, 69, 79	0
2	B	146/162 (90%)	0.50	9 (6%) 26 26	31, 55, 84, 91	0
3	C	85/109 (77%)	-0.14	0 100 100	26, 37, 54, 72	0
4	D	103/122 (84%)	-0.09	2 (1%) 66 66	26, 37, 54, 68	0
All	All	594/674 (88%)	0.19	17 (2%) 53 55	26, 45, 72, 91	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	34	LEU	4.0
4	D	1709	CYS	3.0
1	A	1066	HIS	2.7
1	A	1111	LYS	2.5
1	A	1113	ILE	2.4
1	A	1255	SER	2.4
1	A	1068	GLU	2.4
2	B	159	TRP	2.2
2	B	44	PRO	2.2
2	B	74	ASP	2.2
2	B	153	VAL	2.2
2	B	146	THR	2.2
2	B	51	ASN	2.1
4	D	1708	SER	2.1
2	B	77	ASN	2.0
1	A	1112	SER	2.0
2	B	9	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.