



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:33 PM UTC

PDB ID : 5UQ1 / pdb_00005uq1
Title : Crystal structure of human Cdk2-Spy1 complex
Authors : McGrath, D.A.; Tripathi, S.M.; Rubin, S.M.
Deposited on : 2017-02-06
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

<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
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

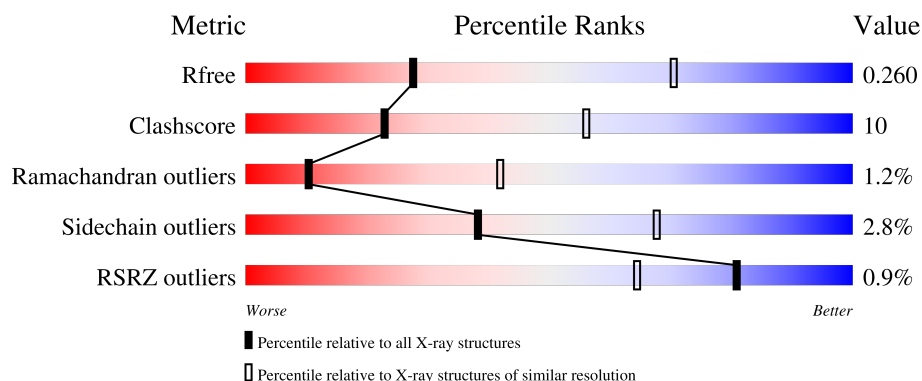
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	301	70% 22% • •
1	C	301	71% 22% • •
2	B	160	56% 26% • 16%
2	D	160	69% 16% • 13%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6913 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 Cyclin-dependent kinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2294	1497	387	403	7			
1	C	289	Total	C	N	O	S	0	0	0
			2293	1491	386	408	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P24941
A	-1	GLU	-	expression tag	UNP P24941
A	0	PHE	-	expression tag	UNP P24941
C	-2	GLY	-	expression tag	UNP P24941
C	-1	GLU	-	expression tag	UNP P24941
C	0	PHE	-	expression tag	UNP P24941

- Molecule 2 is a protein called Speedy protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	134	Total	C	N	O	S	0	0	0
			1143	744	192	198	9			
2	D	139	Total	C	N	O	S	0	0	0
			1183	771	198	204	10			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	54	GLY	-	expression tag	UNP Q5MJ70
B	55	ALA	-	expression tag	UNP Q5MJ70
B	56	MET	-	expression tag	UNP Q5MJ70
B	57	ASP	-	expression tag	UNP Q5MJ70
B	58	PRO	-	expression tag	UNP Q5MJ70
B	59	GLU	-	expression tag	UNP Q5MJ70

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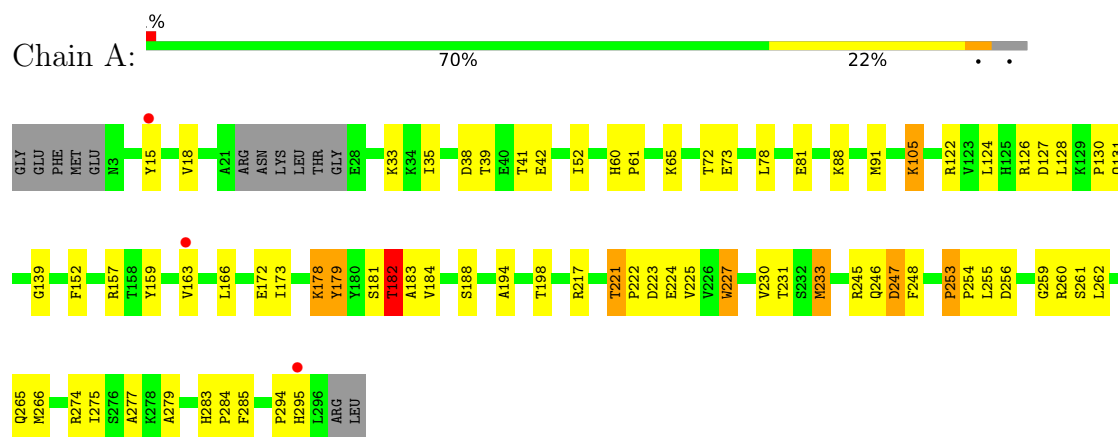
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Chain	Residue	Modelled	Actual	Comment	Reference
B	60	PHE	-	expression tag	UNP Q5MJ70
D	54	GLY	-	expression tag	UNP Q5MJ70
D	55	ALA	-	expression tag	UNP Q5MJ70
D	56	MET	-	expression tag	UNP Q5MJ70
D	57	ASP	-	expression tag	UNP Q5MJ70
D	58	PRO	-	expression tag	UNP Q5MJ70
D	59	GLU	-	expression tag	UNP Q5MJ70
D	60	PHE	-	expression tag	UNP Q5MJ70

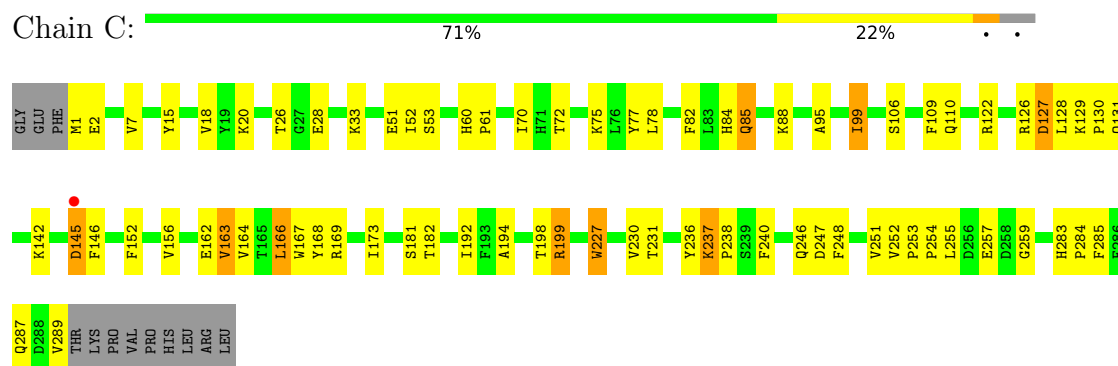
3 Residue-property plots [i](#)

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.

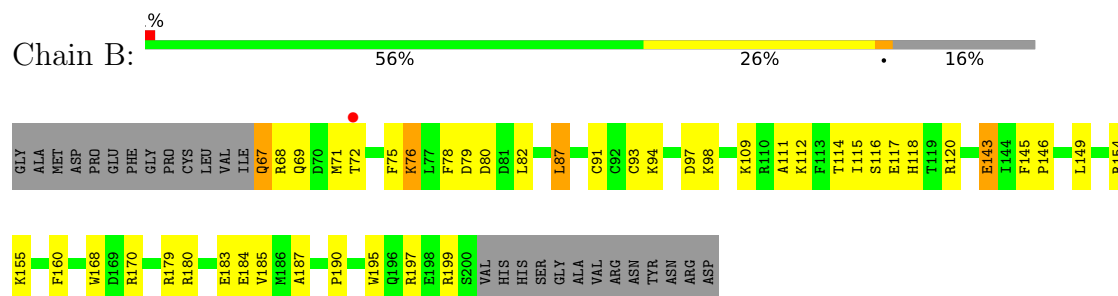
• Molecule 1: Cyclin-dependent kinase 2



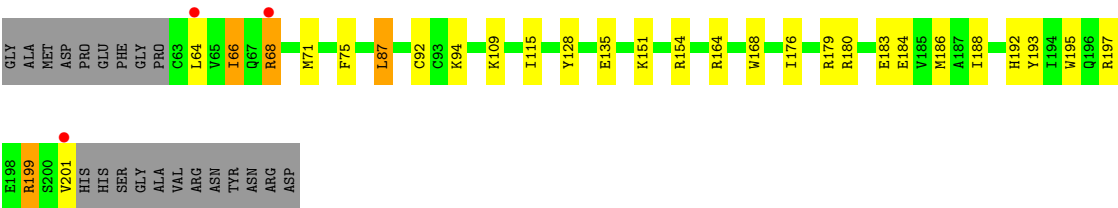
• Molecule 1: Cyclin-dependent kinase 2



• Molecule 2: Speedy protein A



● Molecule 2: Speedy protein A



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	75.36Å 75.36Å 325.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.11 – 3.20 46.11 – 3.20	Depositor EDS
% Data completeness (in resolution range)	94.4 (46.11-3.20) 92.2 (46.11-3.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.205 , 0.261 0.210 , 0.260	Depositor DCC
R_{free} test set	866 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	62.9	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 60.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.057 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6913	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.43	1/2355 (0.0%)	0.93	11/3203 (0.3%)
1	C	0.38	0/2350	0.92	8/3192 (0.3%)
2	B	0.42	0/1173	0.81	1/1582 (0.1%)
2	D	0.42	0/1213	0.83	0/1637
All	All	0.41	1/7091 (0.0%)	0.89	20/9614 (0.2%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	279	ALA	CA-CB	5.48	1.62	1.53

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	227	TRP	CA-C-N	6.19	126.37	119.32
1	A	227	TRP	C-N-CA	6.19	126.37	119.32
1	A	233	MET	CA-C-N	5.84	125.98	119.32
1	A	233	MET	C-N-CA	5.84	125.98	119.32
1	C	145	ASP	CA-C-N	-5.74	113.30	122.66
1	C	145	ASP	C-N-CA	-5.74	113.30	122.66
1	C	246	GLN	N-CA-C	-5.51	103.11	110.55
1	C	85	GLN	N-CA-C	5.44	117.11	108.79
1	C	227	TRP	CA-C-N	5.39	126.57	119.84
1	C	227	TRP	C-N-CA	5.39	126.57	119.84
1	A	253	PRO	CA-C-N	-5.37	114.16	120.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	PRO	C-N-CA	-5.37	114.16	120.12
1	A	221	THR	CA-C-N	5.35	125.27	119.76
1	A	221	THR	C-N-CA	5.35	125.27	119.76
1	C	129	LYS	CA-C-N	5.25	125.31	119.32
1	C	129	LYS	C-N-CA	5.25	125.31	119.32
1	A	277	ALA	N-CA-C	5.21	118.09	111.69
2	B	79	ASP	N-CA-C	-5.14	106.97	114.12
1	A	178	LYS	N-CA-C	-5.13	103.15	110.59
1	A	253	PRO	N-CA-C	5.02	116.83	110.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	247	ASP	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	2294	0	2318	49	0
1	C	2293	0	2321	48	0
2	B	1143	0	1114	33	1
2	D	1183	0	1165	19	1
All	All	6913	0	6918	144	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.

All (144) 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:247:ASP:OD1	1:C:248:PHE:N	2.08	0.86
2:D:135:GLU:OE1	2:D:199:ARG:NH2	2.09	0.84
1:A:178:LYS:HG3	1:A:179:TYR:HD1	1.43	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:ALA:HA	1:C:199:ARG:HD2	1.65	0.78
1:A:178:LYS:HG3	1:A:179:TYR:CD1	2.19	0.77
2:D:94:LYS:HA	2:D:199:ARG:NH1	2.00	0.75
2:B:68:ARG:HD3	2:B:69:GLN:H	1.52	0.74
1:A:181:SER:O	1:A:183:ALA:N	2.19	0.74
1:A:247:ASP:OD1	1:A:248:PHE:N	2.19	0.72
1:C:248:PHE:HA	1:C:251:VAL:HG12	1.72	0.71
1:A:159:TYR:O	1:A:178:LYS:NZ	2.25	0.69
1:C:52:ILE:HD11	1:C:78:LEU:HD11	1.77	0.66
1:C:166:LEU:O	1:C:168:TYR:N	2.30	0.65
1:C:166:LEU:O	1:C:169:ARG:N	2.28	0.64
2:B:82:LEU:HD11	2:B:120:ARG:HD2	1.79	0.64
1:A:52:ILE:HD12	2:B:168:TRP:HH2	1.63	0.64
1:C:88:LYS:HB2	1:C:130:PRO:HB2	1.81	0.63
1:C:52:ILE:HD12	2:D:168:TRP:HH2	1.63	0.63
1:C:127:ASP:HB3	1:C:145:ASP:OD1	1.99	0.62
1:C:163:VAL:HG21	1:C:173:ILE:HD13	1.82	0.61
2:D:201:VAL:HG12	2:D:201:VAL:O	2.01	0.61
1:A:265:GLN:HB2	1:A:275:ILE:HD12	1.83	0.60
1:A:157:ARG:NH1	2:B:97:ASP:OD2	2.23	0.60
1:C:70:ILE:HB	1:C:77:TYR:HB2	1.84	0.59
1:A:72:THR:HG22	1:A:73:GLU:H	1.66	0.59
2:D:135:GLU:CD	2:D:199:ARG:HH22	2.10	0.59
2:D:68:ARG:HH21	2:D:188:ILE:HG23	1.68	0.59
1:C:257:GLU:N	1:C:257:GLU:OE1	2.36	0.58
1:C:251:VAL:HG13	1:C:252:VAL:HG12	1.85	0.58
2:B:72:THR:O	2:B:76:LYS:HD2	2.04	0.58
2:D:66:ILE:O	2:D:109:LYS:NZ	2.34	0.57
1:A:38:ASP:OD1	1:A:39:THR:N	2.37	0.57
1:C:26:THR:HG23	1:C:28:GLU:HG2	1.86	0.56
1:C:88:LYS:HD3	1:C:131:GLN:HG3	1.87	0.56
1:A:256:ASP:O	1:A:260:ARG:HG3	2.06	0.55
2:B:72:THR:HG22	2:B:76:LYS:HG3	1.88	0.55
1:A:172:GLU:OE1	1:A:274:ARG:NH1	2.33	0.55
2:B:71:MET:HE1	2:B:185:VAL:O	2.07	0.55
2:B:93:CYS:HB3	2:B:97:ASP:HA	1.88	0.54
2:B:111:ALA:O	2:B:112:LYS:HG2	2.08	0.54
1:A:52:ILE:HD11	1:A:78:LEU:HD11	1.88	0.54
1:A:124:LEU:HD21	1:A:182:THR:H	1.71	0.54
1:C:52:ILE:HD12	2:D:168:TRP:CH2	2.43	0.54
1:A:227:TRP:HB3	1:A:230:VAL:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:284:PRO:C	1:C:287:GLN:HE22	2.16	0.53
1:C:1:MET:HG2	1:C:2:GLU:H	1.74	0.53
1:A:194:ALA:O	1:A:198:THR:HG22	2.09	0.53
1:A:15:TYR:HE1	1:A:18:VAL:HG23	1.74	0.53
1:A:230:VAL:HA	1:A:233:MET:SD	2.50	0.52
1:A:224:GLU:OE2	1:A:231:THR:OG1	2.25	0.52
1:C:72:THR:HG23	1:C:75:LYS:H	1.75	0.52
1:C:164:VAL:HG21	1:C:169:ARG:HA	1.92	0.51
2:B:117:GLU:OE1	2:B:170:ARG:NH1	2.43	0.51
1:C:99:ILE:HG22	1:C:199:ARG:HE	1.76	0.51
1:A:126:ARG:HD2	1:A:163:VAL:HG12	1.94	0.50
1:A:159:TYR:HB2	1:A:178:LYS:HZ2	1.77	0.50
1:C:227:TRP:HB3	1:C:230:VAL:HB	1.94	0.50
2:B:94:LYS:HA	2:B:199:ARG:HH11	1.76	0.49
1:C:109:PHE:CD2	1:C:289:VAL:HG21	2.47	0.49
1:A:223:ASP:HB3	1:A:225:VAL:HG22	1.94	0.49
2:B:80:ASP:OD2	2:B:120:ARG:NH1	2.37	0.49
1:C:194:ALA:O	1:C:198:THR:HG22	2.12	0.49
1:C:128:LEU:O	1:C:192:ILE:HD11	2.13	0.49
1:A:253:PRO:HG2	1:A:254:PRO:HD3	1.95	0.49
1:C:15:TYR:HE1	1:C:18:VAL:HG23	1.77	0.48
1:A:283:HIS:C	1:A:285:PHE:H	2.21	0.48
2:B:71:MET:HE2	2:B:75:PHE:CE2	2.48	0.48
2:D:186:MET:HG2	2:D:195:TRP:CE3	2.49	0.48
2:D:94:LYS:HA	2:D:199:ARG:HH11	1.77	0.48
1:A:255:LEU:HD22	1:A:259:GLY:HA3	1.96	0.48
1:C:253:PRO:HG2	1:C:254:PRO:HD3	1.95	0.47
2:D:68:ARG:NH2	2:D:188:ILE:HG23	2.29	0.47
2:B:109:LYS:HD3	2:B:109:LYS:HA	1.63	0.47
2:B:98:LYS:HD3	2:B:195:TRP:O	2.14	0.47
1:C:283:HIS:O	1:C:285:PHE:N	2.44	0.47
1:C:126:ARG:O	1:C:127:ASP:HB2	2.15	0.47
2:D:151:LYS:NZ	2:D:154:ARG:HH21	2.13	0.47
2:D:71:MET:HE3	2:D:75:PHE:CZ	2.50	0.47
2:B:115:ILE:HA	2:B:118:HIS:CD2	2.50	0.46
1:C:122:ARG:HA	1:C:152:PHE:CZ	2.50	0.46
1:A:105:LYS:HD3	1:A:285:PHE:O	2.14	0.46
1:C:287:GLN:N	1:C:287:GLN:OE1	2.46	0.46
1:A:262:LEU:O	1:A:266:MET:HG3	2.14	0.46
2:B:75:PHE:HA	2:B:78:PHE:HD1	1.80	0.46
2:B:179:ARG:O	2:B:183:GLU:HG2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:67:GLN:HG3	2:B:68:ARG:N	2.31	0.46
2:B:114:THR:HG1	2:B:116:SER:HG	1.61	0.46
2:B:143:GLU:O	2:B:146:PRO:HD2	2.16	0.46
1:C:15:TYR:CD1	1:C:33:LYS:HE3	2.51	0.46
1:C:7:VAL:HG22	1:C:20:LYS:O	2.15	0.46
1:A:139:GLY:HA2	1:A:294:PRO:HD3	1.98	0.45
1:A:173:ILE:HD11	1:A:184:VAL:HG11	1.98	0.45
2:D:192:HIS:CD2	2:D:193:TYR:H	2.34	0.45
1:A:245:ARG:HG3	1:A:246:GLN:O	2.17	0.45
1:C:84:HIS:CE1	1:C:85:GLN:HG2	2.51	0.45
1:A:60:HIS:CG	1:A:61:PRO:HD2	2.52	0.45
1:C:248:PHE:HA	1:C:251:VAL:CG1	2.46	0.45
2:D:87:LEU:HD12	2:D:197:ARG:CZ	2.47	0.45
1:A:33:LYS:HE2	1:A:35:ILE:HD11	1.99	0.45
1:A:15:TYR:CD1	1:A:33:LYS:HE3	2.52	0.45
2:B:115:ILE:HA	2:B:118:HIS:HD2	1.83	0.44
2:B:187:ALA:HA	2:B:190:PRO:HG3	1.98	0.44
1:A:217:ARG:HG3	1:A:217:ARG:HH11	1.82	0.44
2:B:87:LEU:HD12	2:B:197:ARG:CZ	2.48	0.44
2:D:179:ARG:O	2:D:183:GLU:HG2	2.17	0.43
2:D:128:TYR:HD1	2:D:164:ARG:HD3	1.84	0.43
1:C:60:HIS:CG	1:C:61:PRO:HD2	2.52	0.43
1:C:230:VAL:HG13	1:C:231:THR:HG23	2.00	0.43
1:A:266:MET:HA	1:A:275:ILE:HG22	2.01	0.43
1:C:51:GLU:HG3	1:C:146:PHE:HB2	2.01	0.43
1:C:20:LYS:HD3	1:C:82:PHE:CE1	2.54	0.43
2:D:180:ARG:O	2:D:184:GLU:HG3	2.19	0.43
1:A:88:LYS:HA	1:A:91:MET:HE2	2.01	0.43
2:B:180:ARG:O	2:B:184:GLU:HG3	2.19	0.43
1:C:163:VAL:HG21	1:C:173:ILE:HG21	1.99	0.43
1:A:261:SER:O	1:A:265:GLN:HG3	2.19	0.43
1:C:126:ARG:HD2	1:C:163:VAL:HG12	2.01	0.42
1:C:255:LEU:HD22	1:C:259:GLY:HA3	2.00	0.42
1:A:128:LEU:HB2	1:A:188:SER:CB	2.49	0.42
2:B:115:ILE:HD12	2:B:115:ILE:H	1.84	0.42
1:A:248:PHE:N	1:A:248:PHE:CD1	2.87	0.42
1:C:61:PRO:O	1:C:142:LYS:HE2	2.20	0.42
1:A:248:PHE:N	1:A:248:PHE:HD1	2.18	0.42
1:A:159:TYR:C	1:A:178:LYS:NZ	2.78	0.42
1:C:237:LYS:HE2	1:C:237:LYS:HB2	1.81	0.42
1:A:41:THR:O	1:A:42:GLU:HG2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:LYS:HG3	1:C:238:PRO:HD2	2.01	0.42
1:A:65:LYS:H	1:A:81:GLU:HG2	1.86	0.41
1:C:106:SER:O	1:C:110:GLN:HG3	2.20	0.41
1:A:88:LYS:HB2	1:A:130:PRO:HB2	2.01	0.41
2:B:93:CYS:SG	2:B:197:ARG:HB3	2.60	0.41
1:A:221:THR:HA	1:A:222:PRO:HD3	1.81	0.41
2:B:120:ARG:H	2:B:120:ARG:HG2	1.68	0.41
1:A:255:LEU:HA	1:A:255:LEU:HD23	1.88	0.41
1:A:159:TYR:H	1:A:178:LYS:HZ1	1.68	0.41
2:B:145:PHE:HB2	2:B:146:PRO:HD3	2.03	0.41
2:B:94:LYS:HA	2:B:199:ARG:NH1	2.36	0.41
2:B:93:CYS:O	2:B:199:ARG:NH1	2.54	0.41
1:A:131:GLN:H	1:A:131:GLN:HG3	1.57	0.40
2:B:111:ALA:C	2:B:112:LYS:HG2	2.46	0.40
1:A:122:ARG:HA	1:A:152:PHE:CZ	2.57	0.40
2:B:149:LEU:HD21	2:B:160:PHE:HB2	2.03	0.40
1:C:53:SER:HB3	2:D:176:ILE:HG12	2.03	0.40
1:C:236:TYR:CE1	1:C:240:PHE:HD2	2.40	0.40

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)
2:B:91:CYS:CB	2:D:92:CYS:SG[6_655]	2.04	0.16

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	284/301 (94%)	270 (95%)	10 (4%)	4 (1%)	9	39
1	C	287/301 (95%)	272 (95%)	9 (3%)	6 (2%)	5	31

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	132/160 (82%)	129 (98%)	3 (2%)	0	100	100
2	D	137/160 (86%)	130 (95%)	7 (5%)	0	100	100
All	All	840/922 (91%)	801 (95%)	29 (4%)	10 (1%)	10	42

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	182	THR
1	C	127	ASP
1	C	166	LEU
1	C	167	TRP
1	A	127	ASP
1	C	163	VAL
1	A	166	LEU
1	C	181	SER
1	A	284	PRO
1	C	156	VAL

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	249/265 (94%)	245 (98%)	4 (2%)	55	75
1	C	248/265 (94%)	243 (98%)	5 (2%)	48	72
2	B	120/144 (83%)	114 (95%)	6 (5%)	22	55
2	D	126/144 (88%)	120 (95%)	6 (5%)	23	56
All	All	743/818 (91%)	722 (97%)	21 (3%)	38	68

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	LYS
1	A	179	TYR

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Mol	Chain	Res	Type
1	A	182	THR
1	A	295	HIS
2	B	67	GLN
2	B	76	LYS
2	B	87	LEU
2	B	143	GLU
2	B	154	ARG
2	B	155	LYS
1	C	99	ILE
1	C	162	GLU
1	C	182	THR
1	C	199	ARG
1	C	237	LYS
2	D	64	LEU
2	D	66	ILE
2	D	68	ARG
2	D	87	LEU
2	D	115	ILE
2	D	199	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	121	HIS
2	B	118	HIS
1	C	74	ASN
1	C	85	GLN
1	C	113	GLN
1	C	121	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 [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	288/301 (95%)	-0.16	3 (1%) 79 63	36, 61, 101, 155	0
1	C	289/301 (96%)	-0.18	1 (0%) 90 81	29, 63, 109, 151	0
2	B	134/160 (83%)	-0.15	1 (0%) 84 69	40, 69, 104, 128	0
2	D	139/160 (86%)	-0.17	3 (2%) 62 42	33, 60, 96, 143	0
All	All	850/922 (92%)	-0.17	8 (0%) 81 64	29, 63, 104, 155	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	145	ASP	4.0
1	A	295	HIS	2.7
2	D	68	ARG	2.5
1	A	163	VAL	2.4
2	D	64	LEU	2.4
2	D	201	VAL	2.1
1	A	15	TYR	2.1
2	B	72	THR	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.