



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 06:19 PM EDT

PDB ID : 5IF1 / pdb_00005if1
Title : Crystal structure apo CDK2/cyclin A
Authors : Ayaz, P.; Andres, D.; Kwiatkowski, D.A.; Kolbe, C.; Lienau, P.; Siemeister, G.; Luecking, U.; Stegmann, C.M.
Deposited on : 2016-02-25
Resolution : 2.61 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

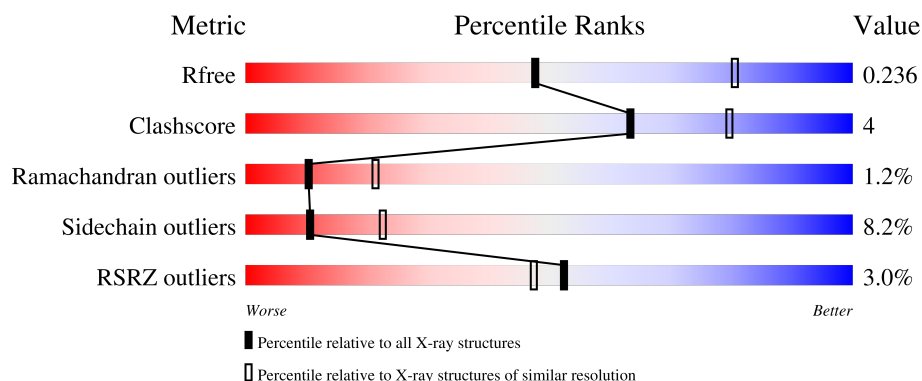
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

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	298	4% 83% 13% .
1	C	298	5% 81% 15% ..
2	B	265	% 83% 12% ..
2	D	265	% 82% 14% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9059 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	298	Total	C	N	O	S	0	1	0
			2400	1561	411	420	8			
1	C	296	Total	C	N	O	S	0	0	0
			2372	1544	403	417	8			

- Molecule 2 is a protein called Cyclin-A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	257	Total	C	N	O	S	0	1	0
			2083	1349	338	385	11			
2	D	257	Total	C	N	O	S	0	1	0
			2083	1349	339	384	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	168	HIS	-	expression tag	UNP P20248
B	169	HIS	-	expression tag	UNP P20248
B	170	HIS	-	expression tag	UNP P20248
B	171	HIS	-	expression tag	UNP P20248
B	172	HIS	-	expression tag	UNP P20248
B	173	HIS	-	expression tag	UNP P20248
D	168	HIS	-	expression tag	UNP P20248
D	169	HIS	-	expression tag	UNP P20248
D	170	HIS	-	expression tag	UNP P20248
D	171	HIS	-	expression tag	UNP P20248
D	172	HIS	-	expression tag	UNP P20248
D	173	HIS	-	expression tag	UNP P20248

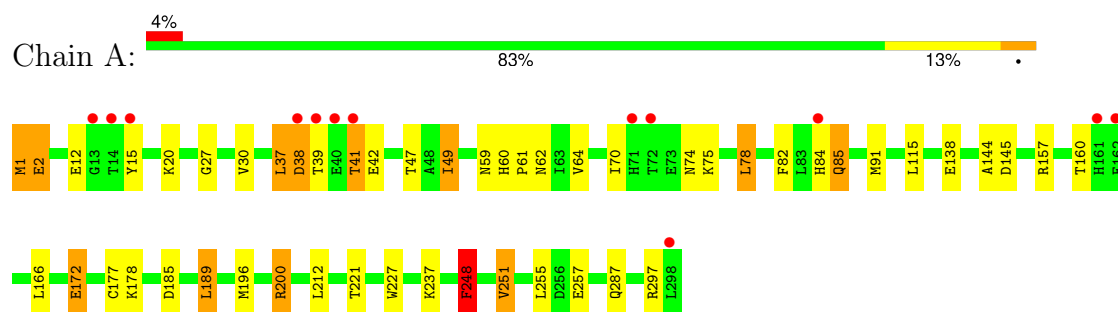
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	55	Total 55	O 55	0	0
3	B	20	Total 20	O 20	0	0
3	C	17	Total 17	O 17	0	0
3	D	29	Total 29	O 29	0	0

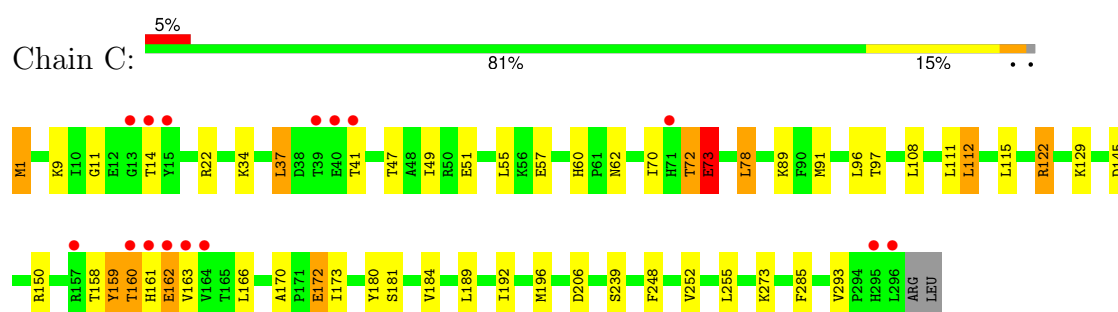
3 Residue-property plots [i](#)

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.

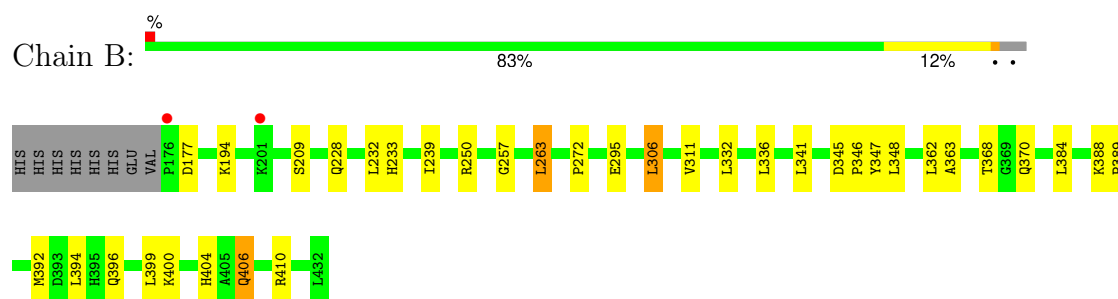
• Molecule 1: Cyclin-dependent kinase 2



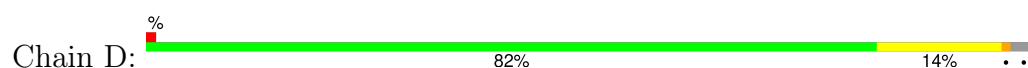
• Molecule 1: Cyclin-dependent kinase 2

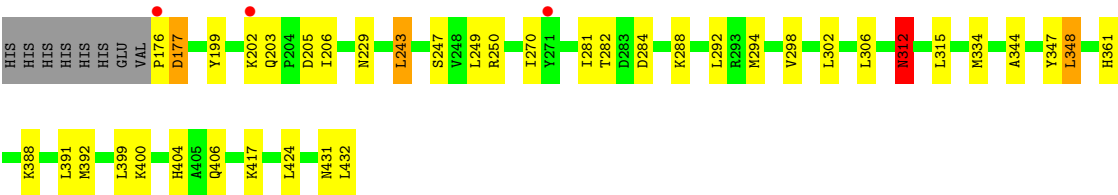


• Molecule 2: Cyclin-A2



• Molecule 2: Cyclin-A2





4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	186.57Å 186.57Å 213.07Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.07 – 2.61 48.07 – 2.61	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.07-2.61) 100.0 (48.07-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.187 , 0.236 0.194 , 0.236	Depositor DCC
R_{free} test set	3312 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	56.2	Xtriage
Anisotropy	0.250	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9059	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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	1.10	2/2465 (0.1%)	1.24	14/3344 (0.4%)
1	C	1.03	0/2434	1.15	5/3305 (0.2%)
2	B	0.98	1/2136 (0.0%)	1.11	2/2898 (0.1%)
2	D	1.00	1/2136 (0.0%)	1.14	3/2898 (0.1%)
All	All	1.03	4/9171 (0.0%)	1.16	24/12445 (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	C	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	37	LEU	CA-C	8.16	1.63	1.53
2	B	177	ASP	N-CA	6.18	1.53	1.46
2	D	177	ASP	N-CA	5.51	1.52	1.46
1	A	59	ASN	N-CA	-5.01	1.40	1.46

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	85	GLN	N-CA-C	8.75	121.38	108.60
1	A	37	LEU	N-CA-C	7.35	121.20	111.28
1	C	11	GLY	N-CA-C	6.50	120.05	110.63
1	A	221	THR	CA-C-N	-6.23	113.54	120.14
1	A	221	THR	C-N-CA	-6.23	113.54	120.14
1	A	12	GLU	N-CA-C	-6.08	99.75	109.96
1	C	159	TYR	CA-CB-CG	5.88	124.49	113.90
2	D	312	ASN	CB-CA-C	-5.86	101.44	110.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	237	LYS	CA-C-N	-5.55	114.05	120.04
1	A	237	LYS	C-N-CA	-5.55	114.05	120.04
1	A	200	ARG	CB-CA-C	5.52	119.31	109.65
1	A	227	TRP	CA-C-N	5.52	126.74	119.84
1	A	227	TRP	C-N-CA	5.52	126.74	119.84
1	A	172	GLU	CB-CA-C	-5.44	101.71	110.74
2	D	312	ASN	N-CA-C	5.42	117.61	111.11
2	B	272	PRO	CA-C-N	-5.38	114.16	119.92
2	B	272	PRO	C-N-CA	-5.38	114.16	119.92
2	D	347	TYR	N-CA-C	5.37	118.05	111.82
1	A	37	LEU	O-C-N	-5.15	116.78	122.91
1	A	248	PHE	N-CA-C	5.12	117.60	111.71
1	C	172	GLU	CB-CA-C	-5.12	101.98	110.68
1	A	84	HIS	N-CA-C	-5.12	105.78	111.36
1	C	37	LEU	CA-C-N	5.01	129.47	122.36
1	C	37	LEU	C-N-CA	5.01	129.47	122.36

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	37	LEU	Peptide
1	C	72	THR	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	2400	0	2454	24	0
1	C	2372	0	2417	21	0
2	B	2083	0	2105	13	0
2	D	2083	0	2107	17	0
3	A	55	0	0	2	1
3	B	20	0	0	2	0
3	C	17	0	0	0	0
3	D	29	0	0	2	0
All	All	9059	0	9083	72	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 4.

All (72) 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)
2:D:391:LEU:HD23	2:D:432:LEU:HD11	1.72	0.69
1:C:122:ARG:O	1:C:122:ARG:HG2	1.93	0.67
1:A:1:MET:HE1	1:A:70:ILE:HD13	1.80	0.63
2:D:229:ASN:HD22	2:D:334:MET:HE2	1.64	0.62
2:B:332:LEU:HD23	2:B:363:ALA:HA	1.80	0.62
2:B:239:ILE:HD11	2:B:257:GLY:HA2	1.82	0.61
2:B:404:HIS:HD2	2:B:406:GLN:H	1.46	0.61
1:A:60:HIS:HD2	1:A:62:ASN:H	1.49	0.60
1:A:60:HIS:CD2	1:A:62:ASN:H	2.19	0.60
2:D:203:GLN:NE2	2:D:247:SER:HA	2.17	0.59
1:C:1:MET:HE1	1:C:70:ILE:HD13	1.84	0.59
1:A:37:LEU:HA	1:A:38:ASP:C	2.29	0.58
2:D:205:ASP:OD2	2:D:250:ARG:HG2	2.05	0.56
2:B:368:THR:HG22	2:B:368:THR:O	2.04	0.56
2:D:344:ALA:HB1	2:D:348:LEU:HD22	1.86	0.56
1:A:248:PHE:HA	1:A:251:VAL:HG12	1.87	0.56
1:C:160:THR:HA	1:C:161:HIS:CG	2.40	0.56
2:D:404:HIS:HD2	2:D:406:GLN:H	1.53	0.56
2:B:336:LEU:HD13	2:B:362:LEU:HD23	1.87	0.56
1:C:163:VAL:HG22	1:C:180:TYR:OH	2.07	0.54
1:C:129:LYS:HA	1:C:192:ILE:HD11	1.91	0.53
1:A:91:MET:HE3	1:A:196:MET:HG2	1.91	0.52
1:C:161:HIS:CB	1:C:162:GLU:HA	2.39	0.52
2:D:404:HIS:CD2	2:D:406:GLN:H	2.29	0.51
1:C:161:HIS:HB3	1:C:162:GLU:HA	1.93	0.50
1:C:150:ARG:HD2	3:D:510:HOH:O	2.12	0.50
1:A:41:THR:OG1	1:A:42:GLU:N	2.45	0.49
2:B:228:GLN:HG3	3:B:508:HOH:O	2.12	0.49
2:D:294:MET:O	2:D:298:VAL:HG23	2.12	0.49
2:D:203:GLN:HG2	2:D:206:ILE:CG1	2.43	0.49
2:D:205:ASP:OD1	2:D:250:ARG:NH1	2.46	0.49
2:B:263:LEU:HD13	2:B:295:GLU:HG3	1.94	0.48
2:B:368:THR:HG21	2:B:370:GLN:OE1	2.13	0.48
2:D:361:HIS:CD2	2:D:391:LEU:HD21	2.49	0.48
1:C:57:GLU:OE1	1:C:122:ARG:NH1	2.41	0.48
1:A:115:LEU:HD21	1:A:185:ASP:HB3	1.95	0.47
2:D:312:ASN:OD1	2:D:334:MET:HE1	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:GLU:O	1:C:55:LEU:HB2	2.15	0.47
1:A:78:LEU:N	1:A:78:LEU:HD23	2.30	0.46
1:A:1:MET:HA	1:A:1:MET:HE2	1.96	0.46
1:C:49:ILE:HG23	2:D:306:LEU:HD12	1.97	0.46
2:D:199:TYR:CE1	2:D:243:LEU:HB3	2.50	0.46
2:D:203:GLN:HE22	2:D:247:SER:HA	1.81	0.46
1:A:157:ARG:NH2	3:A:303:HOH:O	2.49	0.45
1:A:60:HIS:CG	1:A:61:PRO:HD2	2.51	0.45
1:A:49:ILE:HG23	2:B:306:LEU:HD23	1.99	0.45
1:C:72:THR:HB	1:C:73:GLU:HG3	1.98	0.45
2:B:233:HIS:HD2	3:B:509:HOH:O	1.99	0.45
1:A:115:LEU:HD22	1:A:189:LEU:HD22	1.98	0.44
1:C:91:MET:HE3	1:C:196:MET:CG	2.47	0.44
1:C:60:HIS:HD2	1:C:62:ASN:H	1.65	0.44
1:A:172:GLU:HB3	1:A:177:CYS:SG	2.58	0.43
1:C:158:THR:HA	1:C:180:TYR:CE1	2.54	0.43
1:A:20:LYS:HG3	1:A:82:PHE:CE1	2.54	0.43
1:C:91:MET:HE3	1:C:196:MET:HG2	2.01	0.43
1:C:1:MET:HE3	1:C:70:ILE:HG21	2.01	0.43
1:A:1:MET:HE2	1:A:1:MET:CA	2.49	0.42
1:A:64:VAL:HG21	1:A:144:ALA:HB2	2.01	0.42
1:C:163:VAL:HG21	1:C:173:ILE:HG21	2.02	0.42
2:D:400:LYS:NZ	3:D:503:HOH:O	2.52	0.42
1:C:78:LEU:HD23	1:C:78:LEU:N	2.35	0.42
2:B:347:TYR:OH	2:B:394:LEU:HA	2.20	0.42
1:A:27:GLY:HA3	2:D:249:LEU:HD22	2.02	0.41
1:C:108:LEU:O	1:C:112:LEU:HB2	2.20	0.41
1:A:85:GLN:HE21	1:A:85:GLN:HB2	1.60	0.41
2:B:388:LYS:HB3	2:B:389:PRO:HD3	2.03	0.41
1:A:160:THR:HG22	3:A:318:HOH:O	2.20	0.41
1:A:2:GLU:CD	1:A:2:GLU:H	2.29	0.41
2:B:345:ASP:HA	2:B:346:PRO:HA	1.91	0.41
1:A:248:PHE:HA	1:A:251:VAL:CG1	2.51	0.40
1:A:91:MET:HE3	1:A:196:MET:CG	2.51	0.40
1:C:170:ALA:HB1	1:C:172:GLU:OE2	2.22	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)
3:A:343:HOH:O	3:A:343:HOH:O[11_455]	0.83	1.37

5.3 Torsion angles

5.3.1 Protein backbone

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	297/298 (100%)	277 (93%)	14 (5%)	6 (2%)	6	11
1	C	294/298 (99%)	270 (92%)	18 (6%)	6 (2%)	6	11
2	B	256/265 (97%)	251 (98%)	4 (2%)	1 (0%)	30	49
2	D	256/265 (97%)	252 (98%)	4 (2%)	0	100	100
All	All	1103/1126 (98%)	1050 (95%)	40 (4%)	13 (1%)	10	21

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	TYR
2	B	250	ARG
1	C	14	THR
1	C	73	GLU
1	A	39	THR
1	A	41	THR
1	A	38	ASP
1	A	145	ASP
1	C	239	SER
1	C	160	THR
1	A	287	GLN
1	C	41	THR
1	C	285	PHE

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	262/263 (100%)	243 (93%)	19 (7%)	13	27
1	C	259/263 (98%)	232 (90%)	27 (10%)	7	13
2	B	232/239 (97%)	217 (94%)	15 (6%)	15	33
2	D	232/239 (97%)	212 (91%)	20 (9%)	10	20
All	All	985/1004 (98%)	904 (92%)	81 (8%)	10	22

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	GLU
1	A	30	VAL
1	A	47	THR
1	A	49	ILE
1	A	74	ASN
1	A	75	LYS
1	A	78	LEU
1	A	138	GLU
1	A	166	LEU
1	A	178	LYS
1	A	189	LEU
1	A	200	ARG
1	A	212	LEU
1	A	248	PHE
1	A	251	VAL
1	A	255	LEU
1	A	257	GLU
1	A	297	ARG
2	B	194	LYS
2	B	209	SER
2	B	232	LEU
2	B	263	LEU
2	B	306	LEU
2	B	311	VAL
2	B	341	LEU
2	B	348	LEU
2	B	384	LEU
2	B	392	MET
2	B	396	GLN
2	B	399	LEU
2	B	400	LYS
2	B	406	GLN

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Mol	Chain	Res	Type
2	B	410	ARG
1	C	1	MET
1	C	9	LYS
1	C	22	ARG
1	C	34	LYS
1	C	47	THR
1	C	73	GLU
1	C	78	LEU
1	C	89	LYS
1	C	96	LEU
1	C	97	THR
1	C	111	LEU
1	C	112	LEU
1	C	115	LEU
1	C	122	ARG
1	C	145	ASP
1	C	159	TYR
1	C	162	GLU
1	C	166	LEU
1	C	181	SER
1	C	184	VAL
1	C	189	LEU
1	C	206	ASP
1	C	248	PHE
1	C	252	VAL
1	C	255	LEU
1	C	273	LYS
1	C	293	VAL
2	D	176	PRO
2	D	177	ASP
2	D	202	LYS
2	D	243	LEU
2	D	270	ILE
2	D	281	ILE
2	D	282	THR
2	D	284	ASP
2	D	288	LYS
2	D	292	LEU
2	D	302	LEU
2	D	312	ASN
2	D	315	LEU
2	D	348	LEU

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Mol	Chain	Res	Type
2	D	388	LYS
2	D	392	MET
2	D	399	LEU
2	D	417	LYS
2	D	424	LEU
2	D	431	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	60	HIS
1	A	85	GLN
1	A	131	GLN
1	A	161	HIS
1	A	265	GLN
1	A	287	GLN
1	A	295	HIS
2	B	179	HIS
2	B	203	GLN
2	B	208	ASN
2	B	313	GLN
2	B	317	GLN
2	B	323	GLN
2	B	404	HIS
2	B	431	ASN
1	C	60	HIS
1	C	84	HIS
1	C	85	GLN
1	C	113	GLN
1	C	268	HIS
2	D	203	GLN
2	D	208	ASN
2	D	361	HIS
2	D	370	GLN
2	D	396	GLN
2	D	404	HIS
2	D	415	ASN
2	D	419	HIS
2	D	425	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 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

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	298/298 (100%)	-0.29	13 (4%)	39 33	22, 50, 92, 143	1 (0%)
1	C	296/298 (99%)	0.22	15 (5%)	33 28	41, 67, 107, 151	0
2	B	257/265 (96%)	-0.21	2 (0%)	82 80	29, 58, 86, 110	1 (0%)
2	D	257/265 (96%)	-0.43	3 (1%)	76 73	29, 49, 83, 102	1 (0%)
All	All	1108/1126 (98%)	-0.17	33 (2%)	52 47	22, 56, 94, 151	3 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	164	VAL	7.4
2	D	176	PRO	6.0
1	A	298	LEU	5.2
2	B	176	PRO	5.0
1	C	160	THR	4.8
1	C	163	VAL	4.4
1	A	14	THR	4.0
1	C	14	THR	3.8
1	C	40	GLU	3.7
1	C	161	HIS	3.5
1	A	40	GLU	3.2
1	A	15	TYR	3.1
1	C	296	LEU	3.0
1	C	39	THR	3.0
2	B	201	LYS	2.9
1	C	41	THR	2.9
1	C	71	HIS	2.8
1	C	162	GLU	2.7
1	A	13	GLY	2.6
1	A	72	THR	2.6
1	C	295	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	41	THR	2.5
1	A	162	GLU	2.5
1	A	161	HIS	2.4
1	A	84	HIS	2.3
1	C	13	GLY	2.2
1	C	15	TYR	2.2
1	A	71	HIS	2.2
1	A	39	THR	2.1
2	D	202	LYS	2.1
2	D	271	TYR	2.1
1	A	38	ASP	2.1
1	C	157	ARG	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.