Full wwPDB X-ray Structure Validation Report (i)

Dec 9, 2023 – 11:33 am GMT

PDB ID	:	1W98
Title	:	The structural basis of CDK2 activation by cyclin E
Authors	:	Lowe, E.D.; Honda, R.; Dubinina, E.; Skamnaki, V.; Cook, A.; Johnson, L.N.
Deposited on	:	2004-10-07
Resolution	:	2.15 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 (i) 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 (i)) were used in the production of this report:

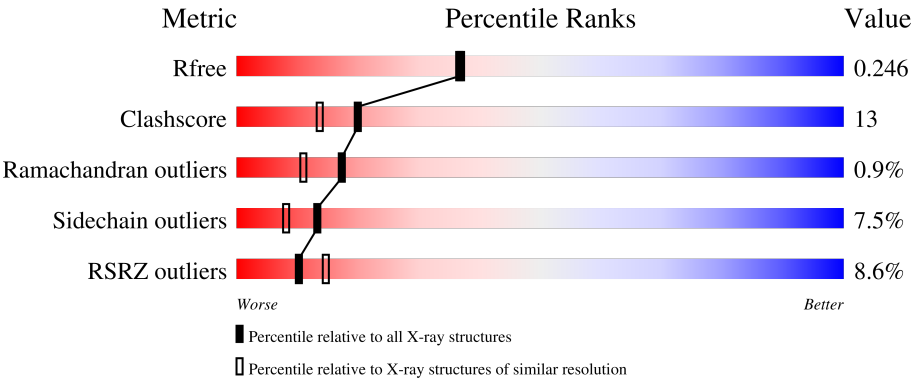
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 <=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	5% 79% 16%
2	B	283	12% 65% 26% 5% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5164 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 CELL DIVISION PROTEIN KINASE 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	P	S	0	0	0
			2399	1556	408	426	1	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP P24941

- Molecule 2 is a protein called G1/S-SPECIFIC CYCLIN E1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	270	Total	C	N	O	S	0	0	0
			2218	1444	359	398	17			

- Molecule 3 is water.

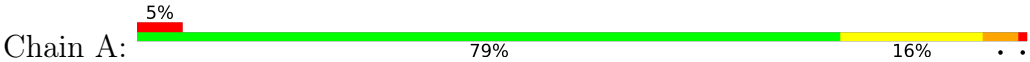
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	337	Total	O	0	0
			337	337		
3	B	210	Total	O	0	0
			210	210		

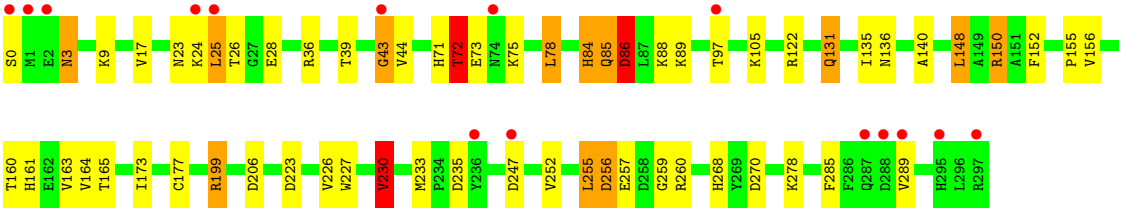
3 Residue-property plots (i)

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: CELL DIVISION PROTEIN KINASE 2

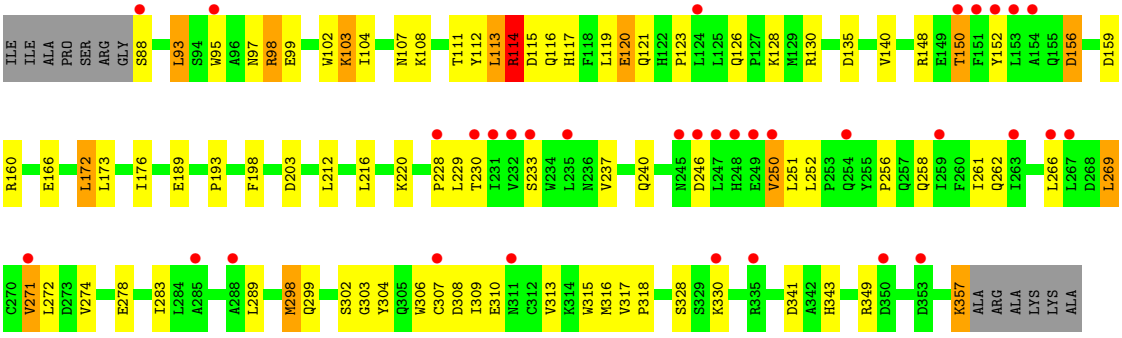
Chain A:





• Molecule 2: G1/S-SPECIFIC CYCLIN E1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	99.62Å 99.62Å 149.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	83.05 – 2.15 34.29 – 2.15	Depositor EDS
% Data completeness (in resolution range)	94.2 (83.05-2.15) 94.2 (34.29-2.15)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.43 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.181 , 0.246 0.185 , 0.246	Depositor DCC
R_{free} test set	1985 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	40.7	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 47.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5164	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

Bond lengths and bond angles in the following residue types are not validated in this section: TPO

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	0.88	1/2449 (0.0%)	0.99	11/3322 (0.3%)
2	B	0.83	2/2274 (0.1%)	0.93	13/3087 (0.4%)
All	All	0.86	3/4723 (0.1%)	0.96	24/6409 (0.4%)

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	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	43	GLY	C-N	-13.48	1.03	1.34
2	B	357	LYS	CE-NZ	12.47	1.80	1.49
2	B	357	LYS	CD-CE	10.15	1.76	1.51

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	43	GLY	O-C-N	-14.15	100.06	122.70
1	A	43	GLY	C-N-CA	9.42	145.25	121.70
2	B	160	ARG	NE-CZ-NH2	-8.63	115.98	120.30
1	A	86	ASP	N-CA-CB	-7.87	96.44	110.60
2	B	160	ARG	NE-CZ-NH1	7.37	123.98	120.30
2	B	357	LYS	CD-CE-NZ	-6.99	95.63	111.70
1	A	230	VAL	CB-CA-C	-6.83	98.43	111.40
1	A	206	ASP	CB-CG-OD2	6.78	124.40	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	43	GLY	N-CA-C	6.64	129.70	113.10
1	A	43	GLY	CA-C-N	6.53	131.57	117.20
2	B	308	ASP	CB-CG-OD2	6.35	124.02	118.30
2	B	148	ARG	NE-CZ-NH2	-6.04	117.28	120.30
2	B	341	ASP	CB-CG-OD2	6.03	123.72	118.30
2	B	135	ASP	CB-CG-OD2	5.95	123.65	118.30
2	B	203	ASP	CB-CG-OD2	5.93	123.64	118.30
2	B	156	ASP	CB-CG-OD2	5.88	123.60	118.30
1	A	235	ASP	CB-CG-OD2	5.73	123.45	118.30
2	B	114	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	256	ASP	CB-CG-OD2	5.51	123.26	118.30
2	B	115	ASP	CB-CG-OD1	5.38	123.14	118.30
1	A	78	LEU	CA-CB-CG	5.26	127.40	115.30
1	A	230	VAL	CG1-CB-CG2	5.19	119.20	110.90
2	B	271	VAL	CB-CA-C	-5.18	101.55	111.40
2	B	159	ASP	CB-CG-OD2	5.16	122.94	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	43	GLY	Mainchain,Peptide
1	A	72	THR	Peptide
1	A	85	GLN	Peptide

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	2399	0	2442	52	0
2	B	2218	0	2219	72	0
3	A	337	0	0	20	0
3	B	210	0	0	20	0
All	All	5164	0	4661	117	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 13.

All (117) 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:B:357:LYS:CD	2:B:357:LYS:CE	1.76	1.55
2:B:357:LYS:CE	2:B:357:LYS:NZ	1.80	1.42
1:A:26:THR:HG21	3:A:2054:HOH:O	1.25	1.27
1:A:177:CYS:HB2	3:A:2227:HOH:O	1.33	1.23
2:B:307:CYS:HB3	3:B:2174:HOH:O	1.73	0.86
2:B:283:ILE:HG12	3:B:2170:HOH:O	1.75	0.86
2:B:93:LEU:HD22	2:B:98:ARG:HA	1.57	0.84
1:A:268:HIS:HD2	1:A:270:ASP:H	1.27	0.83
1:A:247:ASP:HB3	3:A:2298:HOH:O	1.77	0.83
2:B:130:ARG:HD3	2:B:274:VAL:HG22	1.60	0.82
2:B:256:PRO:CB	3:B:2146:HOH:O	2.28	0.81
2:B:256:PRO:HB2	3:B:2146:HOH:O	1.80	0.80
2:B:357:LYS:CE	2:B:357:LYS:CG	2.60	0.79
1:A:233:MET:SD	3:A:2227:HOH:O	2.41	0.78
1:A:36:ARG:CD	3:A:2070:HOH:O	2.32	0.76
1:A:155:PRO:HD3	2:B:240:GLN:HG2	1.68	0.75
2:B:304:TYR:CE2	3:B:2169:HOH:O	2.40	0.75
1:A:36:ARG:HD2	3:A:2070:HOH:O	1.89	0.73
2:B:116:GLN:HG3	3:B:2155:HOH:O	1.91	0.71
1:A:85:GLN:OE1	1:A:89:LYS:HG2	1.91	0.71
2:B:117:HIS:O	2:B:120:GLU:HG3	1.90	0.70
2:B:121:GLN:O	2:B:166:GLU:O	2.10	0.69
1:A:84:HIS:HB3	1:A:135:ILE:O	1.94	0.67
2:B:126:GLN:HG2	3:B:2063:HOH:O	1.95	0.67
2:B:299:GLN:HG2	3:B:2167:HOH:O	1.94	0.66
2:B:130:ARG:HD3	2:B:274:VAL:CG2	2.24	0.66
1:A:131:GLN:H	1:A:131:GLN:HE21	1.44	0.65
1:A:25:LEU:N	3:A:2049:HOH:O	2.30	0.65
1:A:278:LYS:NZ	2:B:97:ASN:HB2	2.11	0.65
1:A:105:LYS:HE2	1:A:285:PHE:O	1.97	0.64
1:A:122:ARG:NH2	2:B:229:LEU:HD21	2.13	0.63
2:B:266:LEU:HD21	2:B:316:MET:CE	2.30	0.61
2:B:93:LEU:HD22	2:B:98:ARG:CA	2.31	0.61
2:B:357:LYS:CD	2:B:357:LYS:NZ	2.64	0.61
2:B:315:TRP:CD1	3:B:2179:HOH:O	2.52	0.60
1:A:150:ARG:HD3	3:A:2206:HOH:O	2.02	0.60
2:B:256:PRO:HB3	3:B:2146:HOH:O	1.95	0.59
2:B:298:MET:CE	2:B:302:SER:OG	2.51	0.59
2:B:302:SER:O	2:B:304:TYR:N	2.36	0.58
2:B:283:ILE:CG1	3:B:2170:HOH:O	2.44	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:88:SER:HA	3:B:2038:HOH:O	2.05	0.56
2:B:304:TYR:CE1	3:B:2171:HOH:O	2.53	0.56
2:B:298:MET:HE3	2:B:304:TYR:HB2	1.88	0.56
2:B:150:THR:HG21	2:B:189:GLU:OE2	2.05	0.56
2:B:298:MET:HE2	2:B:302:SER:OG	2.06	0.55
2:B:172:LEU:HD13	2:B:176:ILE:HD11	1.89	0.55
1:A:131:GLN:HG3	3:A:2191:HOH:O	2.08	0.54
2:B:262:GLN:HG2	2:B:328:SER:HB3	1.90	0.53
1:A:199:ARG:CZ	1:A:199:ARG:HB2	2.38	0.53
1:A:165:THR:HG22	3:A:2220:HOH:O	2.09	0.52
2:B:266:LEU:HD21	2:B:316:MET:HE3	1.91	0.52
1:A:148:LEU:HD22	3:A:2039:HOH:O	2.08	0.52
2:B:298:MET:HE2	2:B:298:MET:O	2.10	0.52
1:A:278:LYS:HZ2	2:B:97:ASN:HB2	1.75	0.51
1:A:9:LYS:HE3	1:A:17:VAL:HG13	1.92	0.51
1:A:255:LEU:HG	1:A:259:GLY:HA3	1.92	0.51
1:A:73:GLU:HG2	3:A:2135:HOH:O	2.10	0.51
2:B:103:LYS:NZ	3:B:2044:HOH:O	2.43	0.50
2:B:152:TYR:HB3	2:B:230:THR:HB	1.93	0.50
2:B:302:SER:OG	2:B:302:SER:O	2.29	0.50
1:A:23:ASN:HB3	1:A:26:THR:HB	1.94	0.50
1:A:72:THR:CG2	1:A:72:THR:O	2.60	0.49
1:A:227:TRP:O	1:A:230:VAL:HG22	2.12	0.49
1:A:84:HIS:HB2	3:A:2146:HOH:O	2.12	0.49
2:B:309:ILE:CG2	2:B:313:VAL:HG23	2.43	0.48
1:A:0:SER:HB3	3:A:2132:HOH:O	2.12	0.48
2:B:269:LEU:N	3:B:2150:HOH:O	2.46	0.48
1:A:36:ARG:HH11	1:A:75:LYS:HZ1	1.60	0.48
2:B:150:THR:CG2	2:B:189:GLU:OE2	2.63	0.47
2:B:298:MET:HE1	2:B:304:TYR:O	2.14	0.47
2:B:304:TYR:CZ	3:B:2169:HOH:O	2.56	0.47
1:A:155:PRO:HD3	2:B:240:GLN:CG	2.41	0.47
1:A:260:ARG:HD3	3:A:2304:HOH:O	2.15	0.47
1:A:3:ASN:HB3	3:A:2005:HOH:O	2.16	0.46
1:A:256:ASP:OD1	1:A:256:ASP:N	2.46	0.46
2:B:126:GLN:OE1	2:B:128:LYS:HE2	2.15	0.46
2:B:306:TRP:CH2	2:B:310:GLU:HG3	2.50	0.46
2:B:258:GLN:HE22	2:B:330:LYS:HA	1.81	0.46
2:B:298:MET:HE2	2:B:298:MET:C	2.35	0.46
2:B:156:ASP:OD1	2:B:228:PRO:HG3	2.16	0.46
2:B:261:ILE:HD11	2:B:343:HIS:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:112:TYR:HB3	2:B:114:ARG:CZ	2.46	0.46
2:B:317:VAL:HB	2:B:318:PRO:HD3	1.98	0.46
1:A:268:HIS:HD2	1:A:270:ASP:N	2.05	0.45
1:A:131:GLN:H	1:A:131:GLN:NE2	2.12	0.45
1:A:163:VAL:HG23	1:A:173:ILE:HD13	1.97	0.45
2:B:113:LEU:C	2:B:114:ARG:HD3	2.37	0.45
1:A:88:LYS:HE3	3:A:2161:HOH:O	2.17	0.45
2:B:233:SER:O	2:B:237:VAL:HG23	2.16	0.45
1:A:122:ARG:HA	1:A:152:PHE:CE1	2.51	0.45
2:B:313:VAL:O	2:B:317:VAL:HG23	2.17	0.45
1:A:268:HIS:CD2	1:A:270:ASP:H	2.18	0.44
2:B:98:ARG:HG2	2:B:99:GLU:N	2.32	0.44
2:B:298:MET:HE3	2:B:302:SER:OG	2.16	0.44
2:B:103:LYS:HG3	2:B:104:ILE:N	2.33	0.43
1:A:86:ASP:OD1	1:A:88:LYS:HB3	2.18	0.43
2:B:116:GLN:NE2	3:B:2056:HOH:O	2.17	0.43
2:B:123:PRO:HD2	3:B:2086:HOH:O	2.17	0.43
2:B:216:LEU:O	2:B:220:LYS:HG3	2.18	0.43
1:A:136:ASN:ND2	1:A:140:ALA:HB3	2.33	0.43
2:B:95:TRP:CZ3	2:B:251:LEU:HG	2.54	0.43
1:A:227:TRP:CE3	1:A:230:VAL:HG13	2.54	0.42
1:A:71:HIS:CE1	2:B:216:LEU:HD13	2.54	0.42
1:A:105:LYS:CE	1:A:285:PHE:O	2.67	0.42
2:B:261:ILE:CD1	2:B:343:HIS:HB2	2.49	0.42
1:A:89:LYS:HE3	3:A:2151:HOH:O	2.20	0.42
2:B:102:TRP:HZ2	3:B:2169:HOH:O	2.03	0.42
1:A:156:VAL:HG22	2:B:251:LEU:O	2.20	0.42
1:A:161:HIS:HD2	3:A:2104:HOH:O	2.02	0.42
2:B:261:ILE:HD11	2:B:343:HIS:CA	2.51	0.41
2:B:193:PRO:HB2	2:B:198:PHE:CE2	2.56	0.41
1:A:223:ASP:H	1:A:226:VAL:HG22	1.86	0.41
2:B:119:LEU:HD12	3:B:2059:HOH:O	2.21	0.41
1:A:97:THR:HB	3:A:2168:HOH:O	2.21	0.40
2:B:95:TRP:HZ3	2:B:250:VAL:HG23	1.84	0.40
1:A:26:THR:HG22	1:A:28:GLU:HB2	2.03	0.40
1:A:24:LYS:H	1:A:24:LYS:HG2	1.60	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	295/298 (99%)	277 (94%)	14 (5%)	4 (1%)	11 5
2	B	268/283 (95%)	258 (96%)	9 (3%)	1 (0%)	34 29
All	All	563/581 (97%)	535 (95%)	23 (4%)	5 (1%)	17 11

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	HIS
1	A	164	VAL
1	A	289	VAL
2	B	303	GLY
1	A	44	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	262/262 (100%)	248 (95%)	14 (5%)	22 19
2	B	246/254 (97%)	222 (90%)	24 (10%)	8 4
All	All	508/516 (98%)	470 (92%)	38 (8%)	13 8

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ASN

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Mol	Chain	Res	Type
1	A	25	LEU
1	A	39	THR
1	A	72	THR
1	A	78	LEU
1	A	86	ASP
1	A	131	GLN
1	A	148	LEU
1	A	150	ARG
1	A	199	ARG
1	A	230	VAL
1	A	252	VAL
1	A	255	LEU
1	A	257	GLU
2	B	93	LEU
2	B	98	ARG
2	B	103	LYS
2	B	107	ASN
2	B	108	LYS
2	B	111	THR
2	B	113	LEU
2	B	114	ARG
2	B	120	GLU
2	B	140	VAL
2	B	150	THR
2	B	172	LEU
2	B	173	LEU
2	B	212	LEU
2	B	246	ASP
2	B	250	VAL
2	B	252	LEU
2	B	269	LEU
2	B	271	VAL
2	B	272	LEU
2	B	278	GLU
2	B	289	LEU
2	B	298	MET
2	B	349	ARG

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	113	GLN

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Mol	Chain	Res	Type
1	A	131	GLN
1	A	268	HIS
2	B	174	GLN
2	B	236	ASN
2	B	258	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)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	TPO	A	160	1	8,10,11	0.88	0	10,14,16	1.07	1 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TPO	A	160	1	-	1/9/11/13	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	160	TPO	O-C-CA	-2.32	118.69	124.78

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	160	TPO	CB-OG1-P-O3P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides 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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	43:GLY	C	44:VAL	N	1.03

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, 95^{th} 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	297/298 (99%)	-0.02	15 (5%) 28 36	27, 38, 60, 69	0
2	B	270/283 (95%)	0.71	34 (12%) 3 5	33, 47, 63, 69	0
All	All	567/581 (97%)	0.33	49 (8%) 10 15	27, 43, 62, 69	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	247	LEU	7.5
2	B	246	ASP	7.2
2	B	231	ILE	5.9
1	A	25	LEU	5.2
1	A	43	GLY	4.6
2	B	267	LEU	4.6
2	B	248	HIS	4.5
2	B	88	SER	4.5
2	B	230	THR	4.4
2	B	232	VAL	4.2
2	B	153	LEU	3.8
2	B	335	ARG	3.8
2	B	235	LEU	3.7
1	A	295	HIS	3.7
2	B	245	ASN	3.7
1	A	0	SER	3.6
2	B	271	VAL	3.6
2	B	151	PHE	3.6
2	B	285	ALA	3.6
1	A	289	VAL	3.4
1	A	2	GLU	3.2
1	A	297	ARG	3.1
1	A	288	ASP	3.1
2	B	150	THR	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	249	GLU	3.0
2	B	152	TYR	2.9
1	A	236	TYR	2.9
1	A	97	THR	2.8
1	A	287	GLN	2.7
2	B	233	SER	2.6
2	B	266	LEU	2.6
2	B	288	ALA	2.5
1	A	24	LYS	2.4
1	A	1	MET	2.4
2	B	307	CYS	2.4
2	B	95	TRP	2.4
1	A	74	ASN	2.3
1	A	247	ASP	2.3
2	B	254	GLN	2.3
2	B	154	ALA	2.2
2	B	259	ILE	2.2
2	B	228	PRO	2.2
2	B	330	LYS	2.1
2	B	124	LEU	2.1
2	B	250	VAL	2.1
2	B	311	ASN	2.1
2	B	263	ILE	2.0
2	B	350	ASP	2.0
2	B	353	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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, 95^{th} 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(Å^2)	Q<0.9
1	TPO	A	160	11/12	0.98	0.07	32,35,39,40	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.