



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 04:47 pm GMT

PDB ID : 4BCO  
Title : Structure of CDK2 in complex with cyclin A and a 2-amino-4-heteroaryl-pyrimidine inhibitor  
Authors : Hole, A.J.; Baumli, S.; Wang, S.; Endicott, J.A.; Noble, M.E.M.  
Deposited on : 2012-10-02  
Resolution : 2.05 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

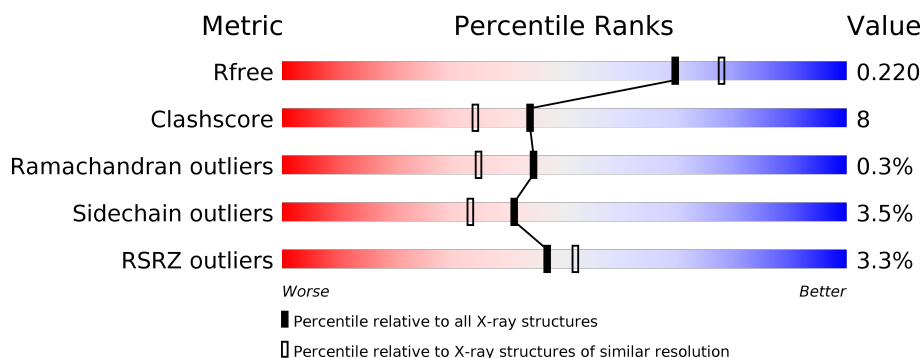
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.05 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))
$R_{free}$	100719	1316 (2.04-2.04)
Clashscore	112137	1394 (2.04-2.04)
Ramachandran outliers	110173	1383 (2.04-2.04)
Sidechain outliers	110143	1383 (2.04-2.04)
RSRZ outliers	101464	1319 (2.04-2.04)

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  $\geq 3$ , 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	300	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, orange 1%, orange 86%, yellow 86%, yellow 97%, green 97%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>86%</span> <span>11% •</span> </div> </div>
1	C	300	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 5%, orange 5%, orange 75%, yellow 75%, yellow 89%, green 89%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>5%</span> <span>75%</span> <span>14% • 10%</span> </div> </div>
2	B	262	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 2%, orange 2%, orange 85%, yellow 85%, yellow 96%, green 96%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>2%</span> <span>85%</span> <span>11% • •</span> </div> </div>
2	D	262	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 5%, orange 5%, orange 83%, yellow 83%, yellow 95%, green 95%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>5%</span> <span>83%</span> <span>12% • •</span> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SGM	B	1434	-	-	-	X
4	SGM	D	1433	-	-	X	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9356 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	299	Total	C	N	O	P	S	0	0	0
			2407	1562	409	427	1	8			
1	C	270	Total	C	N	O	P	S	0	2	0
			2178	1411	374	385	1	7			

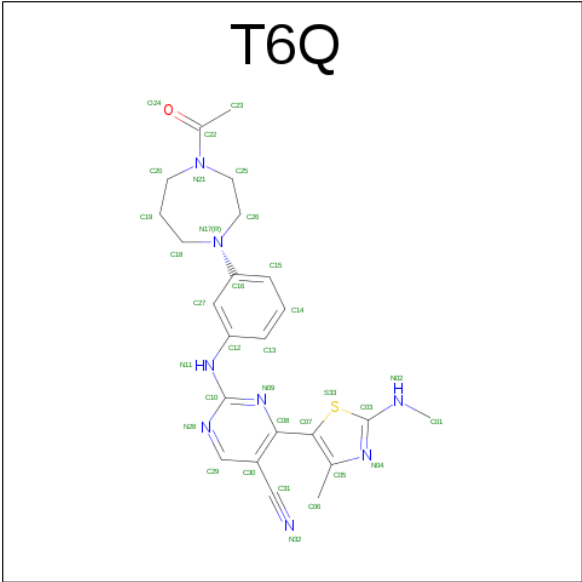
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP P24941
A	0	SER	-	EXPRESSION TAG	UNP P24941
C	-1	GLY	-	EXPRESSION TAG	UNP P24941
C	0	SER	-	EXPRESSION TAG	UNP P24941

- Molecule 2 is a protein called CYCLIN-A2.

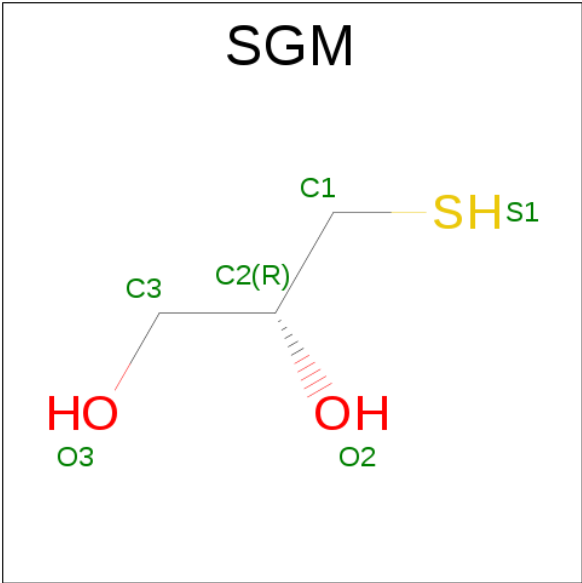
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	258	Total	C	N	O	S		0	0	0
			2083	1350	339	383	11				
2	D	256	Total	C	N	O	S		0	0	0
			2069	1340	337	381	11				

- Molecule 3 is 2-[[3-(4-ETHANOYL-1,4-DIAZEPAN-1-YL)PHENYL]AMINO]-4-[4-METHYL-2-(METHYLAMINO)-1,3-THIAZOL-5-YL]PYRIMIDINE-5-CARBONITRILE (three-letter code: T6Q) (formula: C<sub>23</sub>H<sub>26</sub>N<sub>8</sub>OS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			33	23	8	1	1		
3	C	1	Total	C	N	O	S	0	0
			33	23	8	1	1		

- Molecule 4 is MONOTHIOGLYCEROL (three-letter code: SGM) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>2</sub>S).



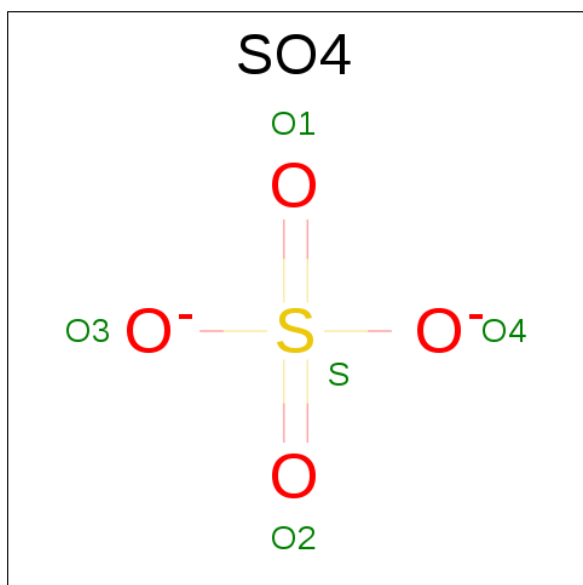
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	O	S	0	0
			6	3	2	1		
4	B	1	Total	C	O	S	0	0
			6	3	2	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	D	1	Total	C	O	S	0	0
			6	3	2	1		
4	D	1	Total	C	O	S	0	0
			6	3	2	1		

- Molecule 5 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		

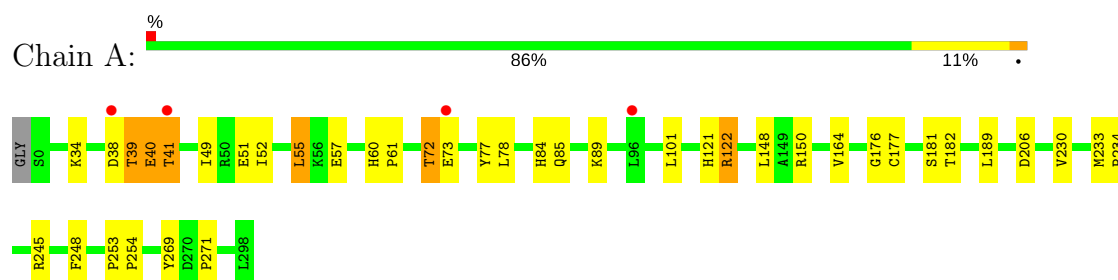
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	187	Total	O	0	0
			187	187		
6	B	140	Total	O	0	0
			140	140		
6	C	102	Total	O	0	0
			102	102		
6	D	90	Total	O	0	0
			90	90		

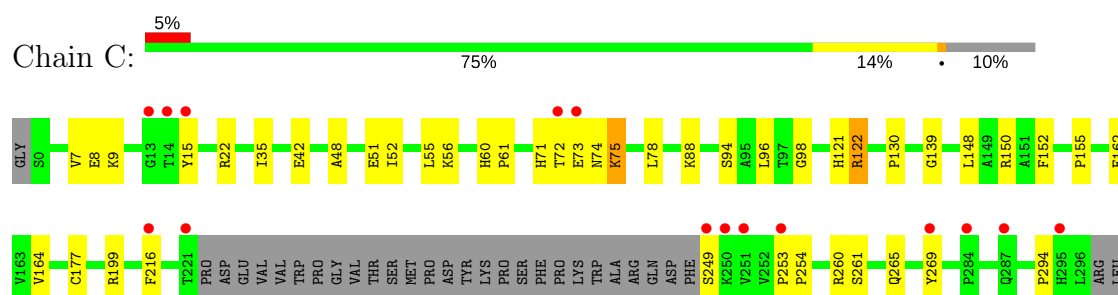
### 3 Residue-property plots [i](#)

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

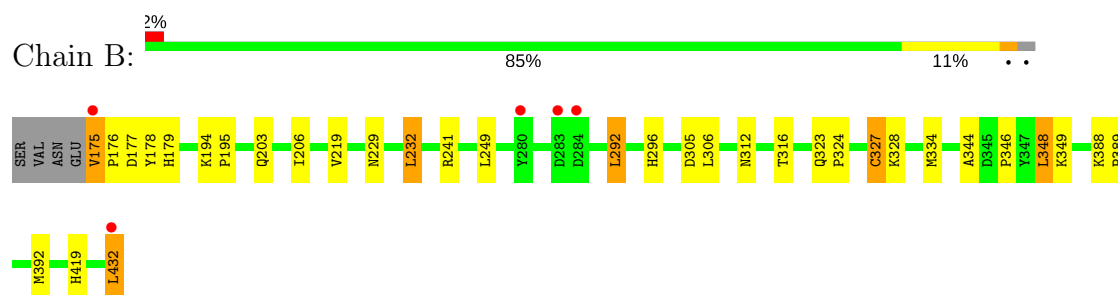
#### • Molecule 1: CYCLIN-DEPENDENT KINASE 2



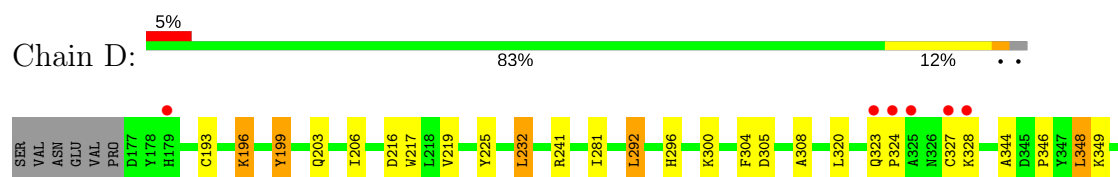
#### • Molecule 1: CYCLIN-DEPENDENT KINASE 2

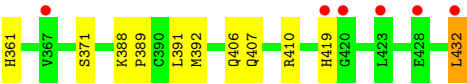


#### • Molecule 2: CYCLIN-A2



#### • Molecule 2: CYCLIN-A2







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.90Å 133.81Å 148.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.60 – 2.05 49.60 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.60-2.05) 98.8 (49.60-2.05)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 2.05Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.189 , 0.226 0.181 , 0.220	Depositor DCC
$R_{free}$ test set	4598 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.6	Xtriage
Anisotropy	0.008	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 54.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9356	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, T6Q, SGM, SO4

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	0.31	0/2457	0.48	0/3333
1	C	0.27	0/2222	0.46	0/3007
2	B	0.29	0/2133	0.46	0/2897
2	D	0.27	0/2118	0.44	0/2875
All	All	0.29	0/8930	0.46	0/12112

There are no bond length outliers.

There are no bond angle outliers.

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	2407	0	2454	29	0
1	C	2178	0	2241	34	0
2	B	2083	0	2104	34	0
2	D	2069	0	2089	37	0
3	A	33	0	26	6	0
3	C	33	0	26	5	0
4	B	12	0	15	5	0
4	D	12	0	16	9	0
5	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	5	0	0	0	0
6	A	187	0	0	1	0
6	B	140	0	0	0	0
6	C	102	0	0	1	0
6	D	90	0	0	1	0
All	All	9356	0	8971	135	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 (135) 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:388:LYS:HE3	2:B:432:LEU:HD22	1.42	1.01
2:D:327:CYS:SG	2:D:419:HIS:NE2	2.34	0.99
1:A:40:GLU:HA	1:A:41:THR:HB	1.39	0.99
1:C:177:CYS:SG	6:C:2081:HOH:O	2.29	0.89
2:D:327:CYS:HG	4:D:1434:SGM:HS1	0.96	0.89
2:D:327:CYS:SG	2:D:419:HIS:CE1	2.71	0.84
1:A:40:GLU:HA	1:A:41:THR:CB	2.07	0.82
2:B:327:CYS:SG	4:B:1433:SGM:S1	2.78	0.82
1:C:15:TYR:OH	1:C:51:GLU:OE1	1.97	0.81
2:D:193:CYS:SG	4:D:1433:SGM:S1	2.41	0.78
2:B:344:ALA:HB1	2:B:348:LEU:HD22	1.64	0.78
2:D:327:CYS:SG	4:D:1434:SGM:S1	2.75	0.76
3:A:1299:T6Q:H27	3:A:1299:T6Q:H251	1.66	0.75
2:D:432:LEU:H	2:D:432:LEU:HD12	1.54	0.73
1:C:71:HIS:CE1	2:D:304:PHE:HE1	2.07	0.72
1:A:84:HIS:O	3:A:1299:T6Q:H262	1.90	0.71
1:C:155:PRO:HG3	2:D:320:LEU:HD21	1.72	0.71
2:D:241:ARG:NH2	4:D:1433:SGM:H12	2.07	0.70
1:A:52:ILE:HD11	1:A:78:LEU:HD21	1.74	0.69
2:B:327:CYS:SG	2:B:419:HIS:NE2	2.64	0.69
2:D:305:ASP:HB3	4:D:1433:SGM:H11	1.73	0.69
2:B:327:CYS:SG	4:B:1433:SGM:O2	2.47	0.68
2:B:327:CYS:SG	2:B:419:HIS:CE1	2.88	0.67
1:C:52:ILE:HD11	1:C:78:LEU:HD21	1.76	0.66
1:A:85:GLN:OE1	1:A:89:LYS:HE2	1.98	0.64
1:C:98:GLY:HA2	1:C:199:ARG:HD3	1.79	0.64
2:D:203:GLN:O	6:D:2022:HOH:O	2.14	0.64
2:B:305:ASP:HB3	4:B:1434:SGM:H11	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:193:CYS:SG	4:D:1433:SGM:H2	2.38	0.63
2:B:327:CYS:HG	2:B:419:HIS:CE1	2.17	0.62
2:B:346:PRO:O	2:B:349:LYS:HG2	1.99	0.62
1:A:72:THR:HG23	1:A:73:GLU:N	2.15	0.61
3:A:1299:T6Q:C31	3:A:1299:T6Q:H061	2.31	0.61
2:D:344:ALA:HB1	2:D:348:LEU:HD22	1.83	0.60
2:B:229:ASN:HD22	2:B:334:MET:CE	2.14	0.60
1:C:71:HIS:CE1	2:D:304:PHE:CE1	2.90	0.59
3:A:1299:T6Q:H13	3:A:1299:T6Q:N09	2.18	0.58
2:B:175:VAL:N	2:B:176:PRO:CD	2.67	0.57
2:D:388:LYS:HB3	2:D:389:PRO:HD3	1.86	0.57
1:A:177:CYS:HB2	1:A:233:MET:CE	2.36	0.56
2:B:229:ASN:HD22	2:B:334:MET:HE2	1.71	0.55
3:C:1298:T6Q:H13	3:C:1298:T6Q:N09	2.20	0.55
1:C:216:PHE:HA	1:C:269:TYR:OH	2.07	0.55
1:A:121:HIS:O	1:A:122:ARG:HG3	2.07	0.55
1:C:121:HIS:O	1:C:122:ARG:HG3	2.07	0.54
2:B:203:GLN:HB3	2:B:206:ILE:HG12	1.90	0.54
2:D:361:HIS:HE1	2:D:371:SER:HB2	1.73	0.54
1:A:269:TYR:O	1:A:271:PRO:HD3	2.08	0.54
1:C:15:TYR:HE2	1:C:35:ILE:HD13	1.73	0.53
2:B:249:LEU:CD1	1:C:22[B]:ARG:HH22	2.21	0.53
2:D:346:PRO:O	2:D:349:LYS:HG2	2.08	0.53
2:B:392:MET:SD	2:B:432:LEU:HD23	2.48	0.53
3:A:1299:T6Q:C31	3:A:1299:T6Q:C06	2.87	0.53
2:D:361:HIS:CE1	2:D:371:SER:HB2	2.44	0.53
2:B:177:ASP:O	2:B:178:TYR:HB2	2.08	0.52
2:B:177:ASP:OD1	2:B:178:TYR:N	2.42	0.52
2:B:388:LYS:HB3	2:B:389:PRO:HD3	1.92	0.52
1:A:40:GLU:CA	1:A:41:THR:CB	2.85	0.52
2:B:219:VAL:HG22	2:B:232:LEU:HD11	1.90	0.52
1:A:121:HIS:C	1:A:122:ARG:HG3	2.30	0.51
3:C:1298:T6Q:C31	3:C:1298:T6Q:H061	2.41	0.51
2:B:175:VAL:N	2:B:176:PRO:HD2	2.26	0.51
1:C:71:HIS:NE2	2:D:304:PHE:HE1	2.08	0.51
1:A:73:GLU:H	1:A:73:GLU:CD	2.14	0.50
1:C:261:SER:O	1:C:265:GLN:HG3	2.11	0.50
2:D:296:HIS:CE1	2:D:300:LYS:HZ1	2.29	0.50
1:C:8:GLU:OE2	3:C:1298:T6Q:H231	2.13	0.49
2:D:241:ARG:CZ	4:D:1433:SGM:H12	2.41	0.49
2:D:219:VAL:HG22	2:D:232:LEU:HD11	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:327:CYS:HG	4:D:1434:SGM:C1	2.24	0.48
1:C:249:SER:HA	1:C:260:ARG:HD2	1.94	0.48
1:C:72:THR:HB	1:C:75:LYS:O	2.13	0.48
2:D:361:HIS:CD2	2:D:391:LEU:HD21	2.48	0.48
1:C:15:TYR:CZ	1:C:51:GLU:OE1	2.67	0.47
2:D:432:LEU:H	2:D:432:LEU:CD1	2.20	0.47
1:A:245:ARG:NH2	1:A:248:PHE:HE1	2.12	0.47
1:C:60:HIS:CG	1:C:61:PRO:HD2	2.50	0.47
1:A:230:VAL:HA	1:A:233:MET:HG3	1.97	0.47
1:A:72:THR:CG2	1:A:73:GLU:N	2.76	0.47
1:C:122:ARG:HA	1:C:152:PHE:CE2	2.49	0.47
2:D:308:ALA:HB2	4:D:1433:SGM:S1	2.55	0.47
1:A:39:THR:C	1:A:40:GLU:HG2	2.34	0.47
2:B:305:ASP:CB	4:B:1434:SGM:H11	2.44	0.47
1:C:73:GLU:HG2	1:C:74:ASN:N	2.30	0.47
2:D:407:GLN:OE1	2:D:410:ARG:HD3	2.15	0.47
1:A:51:GLU:O	1:A:55:LEU:HB2	2.14	0.46
2:D:432:LEU:HD12	2:D:432:LEU:N	2.25	0.46
3:C:1298:T6Q:C06	3:C:1298:T6Q:C31	2.93	0.46
2:D:296:HIS:CD2	2:D:300:LYS:HZ1	2.33	0.46
1:C:15:TYR:HE2	1:C:35:ILE:CD1	2.29	0.46
1:C:73:GLU:HG2	1:C:74:ASN:OD1	2.16	0.45
1:A:49:ILE:HG23	2:B:306:LEU:HD12	1.98	0.45
1:C:72:THR:HG22	1:C:73:GLU:N	2.32	0.45
2:D:217:TRP:CZ2	2:D:281:ILE:HD13	2.51	0.45
2:B:388:LYS:O	2:B:392:MET:HG2	2.16	0.45
2:D:196:LYS:HB2	2:D:196:LYS:NZ	2.30	0.45
1:A:181:SER:OG	1:A:182:THR:N	2.49	0.45
2:B:176:PRO:O	2:B:179:HIS:HB3	2.16	0.45
2:B:432:LEU:HD12	2:B:432:LEU:O	2.16	0.45
1:C:7:VAL:HG12	1:C:8:GLU:HG2	1.97	0.45
1:C:15:TYR:OH	1:C:48:ALA:HA	2.17	0.45
1:C:121:HIS:C	1:C:122:ARG:HG3	2.36	0.44
2:B:249:LEU:HD12	1:C:22[B]:ARG:NH2	2.33	0.44
2:B:323:GLN:HA	2:B:324:PRO:HA	1.72	0.44
2:B:348:LEU:HD12	2:B:348:LEU:HA	1.89	0.44
2:D:323:GLN:HA	2:D:324:PRO:HA	1.79	0.44
1:C:253:PRO:HB2	1:C:254:PRO:HD3	2.00	0.44
1:A:176:GLY:O	1:A:234:PRO:HG2	2.18	0.43
3:C:1298:T6Q:C13	3:C:1298:T6Q:N09	2.81	0.43
1:A:34:LYS:HE2	1:A:77:TYR:OH	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:229:ASN:HD22	2:B:334:MET:HE1	1.83	0.43
1:A:60:HIS:CG	1:A:61:PRO:HD2	2.54	0.43
1:A:230:VAL:O	1:A:233:MET:HG3	2.18	0.43
1:C:94:SER:O	1:C:199:ARG:HD2	2.17	0.43
1:C:72:THR:HG22	1:C:73:GLU:CD	2.39	0.43
2:B:241:ARG:NH2	4:B:1434:SGM:H12	2.34	0.43
2:D:216:ASP:HB2	2:D:406:GLN:HG2	2.01	0.42
1:A:57:GLU:OE2	6:A:2048:HOH:O	2.21	0.42
1:A:177:CYS:HB2	1:A:233:MET:HE3	2.01	0.42
1:A:72:THR:O	1:A:73:GLU:C	2.57	0.42
2:B:194:LYS:HA	2:B:195:PRO:HD3	1.95	0.42
1:A:49:ILE:CG2	2:B:306:LEU:HD12	2.50	0.42
2:D:225:TYR:HE1	2:D:281:ILE:HG21	1.84	0.41
2:D:292:LEU:HD12	2:D:292:LEU:HA	1.89	0.41
1:C:88:LYS:HB2	1:C:130:PRO:HB2	2.02	0.41
2:B:312:ASN:O	2:B:316:THR:HG23	2.20	0.41
1:A:189:LEU:HA	1:A:189:LEU:HD23	1.91	0.41
2:D:203:GLN:HB3	2:D:206:ILE:HG12	2.03	0.41
1:A:253:PRO:N	1:A:254:PRO:CD	2.83	0.41
1:C:51:GLU:O	1:C:55:LEU:HB2	2.20	0.41
2:D:199:TYR:CD1	2:D:199:TYR:C	2.93	0.41
1:C:139:GLY:HA2	1:C:294:PRO:HD3	2.03	0.40
2:B:292:LEU:HD12	2:B:292:LEU:HA	1.87	0.40
1:C:56:LYS:HB3	1:C:56:LYS:HE2	1.78	0.40
3:A:1299:T6Q:N09	3:A:1299:T6Q:C13	2.82	0.40

There are no symmetry-related clashes.

## 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	296/300 (99%)	287 (97%)	7 (2%)	2 (1%)	25 14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	267/300 (89%)	260 (97%)	6 (2%)	1 (0%)	38	27
2	B	256/262 (98%)	254 (99%)	2 (1%)	0	100	100
2	D	254/262 (97%)	252 (99%)	2 (1%)	0	100	100
All	All	1073/1124 (96%)	1053 (98%)	17 (2%)	3 (0%)	44	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	41	THR
1	A	164	VAL
1	C	164	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	263/263 (100%)	253 (96%)	10 (4%)	38	30
1	C	238/263 (90%)	230 (97%)	8 (3%)	42	34
2	B	232/236 (98%)	224 (97%)	8 (3%)	42	34
2	D	230/236 (98%)	222 (96%)	8 (4%)	41	33
All	All	963/998 (96%)	929 (96%)	34 (4%)	41	33

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ASP
1	A	39	THR
1	A	40	GLU
1	A	55	LEU
1	A	72	THR
1	A	101	LEU
1	A	122	ARG
1	A	148	LEU

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Mol	Chain	Res	Type
1	A	150	ARG
1	A	206	ASP
2	B	175	VAL
2	B	232	LEU
2	B	292	LEU
2	B	296	HIS
2	B	327	CYS
2	B	328	LYS
2	B	348	LEU
2	B	432	LEU
1	C	9	LYS
1	C	42	GLU
1	C	75	LYS
1	C	96	LEU
1	C	122	ARG
1	C	148	LEU
1	C	150	ARG
1	C	162	GLU
2	D	196	LYS
2	D	199	TYR
2	D	232	LEU
2	D	292	LEU
2	D	328	LYS
2	D	348	LEU
2	D	392	MET
2	D	432	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are 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	9,10,11	0.74	0	10,14,16	1.01	0
1	TPO	C	160	1	9,10,11	0.71	0	10,14,16	1.13	0

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	-	0/8/11/13	0/0/0/0
1	TPO	C	160	1	-	0/8/11/13	0/0/0/0

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.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

8 ligands are 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
3	T6Q	A	1299	-	29,36,36	1.80	6 (20%)	30,50,50	2.29	8 (26%)
4	SGM	B	1433	-	5,5,5	0.47	0	5,5,5	0.26	0
4	SGM	B	1434	2	5,5,5	0.89	0	5,5,5	0.40	0
5	SO4	B	1435	-	4,4,4	0.15	0	6,6,6	0.09	0
3	T6Q	C	1298	-	29,36,36	1.82	7 (24%)	30,50,50	2.49	7 (23%)
4	SGM	D	1433	-	5,5,5	0.48	0	5,5,5	0.67	0
4	SGM	D	1434	-	5,5,5	0.38	0	5,5,5	0.21	0
5	SO4	D	1435	-	4,4,4	0.13	0	6,6,6	0.08	0

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
3	T6Q	A	1299	-	-	0/13/31/31	0/3/4/4
4	SGM	B	1433	-	-	0/4/4/4	0/0/0/0
4	SGM	B	1434	2	-	0/4/4/4	0/0/0/0
5	SO4	B	1435	-	-	0/0/0/0	0/0/0/0
3	T6Q	C	1298	-	-	0/13/31/31	0/3/4/4
4	SGM	D	1433	-	-	0/4/4/4	0/0/0/0
4	SGM	D	1434	-	-	0/4/4/4	0/0/0/0
5	SO4	D	1435	-	-	0/0/0/0	0/0/0/0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1298	T6Q	C08-C07	-3.01	1.42	1.49
3	A	1299	T6Q	C08-C07	-2.70	1.42	1.49
3	C	1298	T6Q	C20-N21	-2.06	1.42	1.47
3	C	1298	T6Q	C14-C15	2.08	1.42	1.38
3	A	1299	T6Q	C14-C15	2.26	1.43	1.38
3	C	1298	T6Q	C12-N11	2.61	1.46	1.40
3	A	1299	T6Q	C12-N11	2.63	1.46	1.40
3	C	1298	T6Q	C22-N21	3.18	1.45	1.35
3	C	1298	T6Q	C16-N17	3.28	1.47	1.38
3	A	1299	T6Q	C22-N21	3.32	1.46	1.35
3	A	1299	T6Q	C16-N17	3.37	1.48	1.38
3	A	1299	T6Q	C10-N11	5.38	1.46	1.36
3	C	1298	T6Q	C10-N11	5.49	1.46	1.36

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1298	T6Q	C01-N02-C03	-6.74	117.08	122.85
3	C	1298	T6Q	N28-C10-N09	-6.17	120.74	126.68
3	A	1299	T6Q	N28-C10-N09	-5.73	121.17	126.68
3	A	1299	T6Q	C01-N02-C03	-4.46	119.04	122.85
3	C	1298	T6Q	C12-N11-C10	-2.89	121.65	129.17
3	A	1299	T6Q	C30-C29-N28	-2.77	119.16	123.39
3	A	1299	T6Q	C12-N11-C10	-2.72	122.09	129.17
3	C	1298	T6Q	C30-C29-N28	-2.54	119.51	123.39
3	A	1299	T6Q	C30-C08-N09	-2.05	118.22	121.29
3	A	1299	T6Q	C29-N28-C10	2.68	120.32	115.88
3	C	1298	T6Q	C29-N28-C10	2.84	120.59	115.88
3	C	1298	T6Q	C07-C08-N09	3.16	119.01	114.97
3	A	1299	T6Q	C07-C08-N09	3.89	119.94	114.97
3	A	1299	T6Q	C25-N21-C20	6.71	127.56	116.45
3	C	1298	T6Q	C25-N21-C20	6.89	127.87	116.45

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1299	T6Q	6	0
4	B	1433	SGM	2	0
4	B	1434	SGM	3	0
3	C	1298	T6Q	5	0
4	D	1433	SGM	6	0
4	D	1434	SGM	3	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	298/300 (99%)	-0.16	4 (1%) 77 80	17, 28, 62, 92	0
1	C	269/300 (89%)	0.05	15 (5%) 25 28	26, 38, 75, 96	0
2	B	258/262 (98%)	-0.27	5 (1%) 67 72	17, 32, 54, 86	0
2	D	256/262 (97%)	0.16	12 (4%) 32 35	22, 40, 71, 105	0
All	All	1081/1124 (96%)	-0.06	36 (3%) 47 52	17, 34, 68, 105	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	432	LEU	6.8
1	C	251	VAL	5.6
2	D	432	LEU	5.0
1	C	295	HIS	4.9
1	C	15	TYR	4.9
2	B	175	VAL	4.8
2	D	327	CYS	4.5
1	C	14	THR	4.4
2	B	284	ASP	4.1
1	C	73	GLU	4.0
1	A	96	LEU	3.8
1	C	221	THR	3.7
2	D	324	PRO	3.5
2	D	428	GLU	3.2
1	A	41	THR	3.0
2	D	179	HIS	3.0
2	D	367	VAL	2.9
1	C	13	GLY	2.9
2	D	325	ALA	2.8
2	D	423	LEU	2.8
2	D	323	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	328	LYS	2.7
2	B	283	ASP	2.6
1	C	269	TYR	2.6
2	B	280	TYR	2.6
1	C	287	GLN	2.5
1	A	73	GLU	2.5
1	C	249	SER	2.4
1	A	38	ASP	2.2
2	D	420	GLY	2.2
1	C	72	THR	2.2
1	C	284	PRO	2.2
1	C	216	PHE	2.2
1	C	253	PRO	2.2
1	C	250	LYS	2.1
2	D	419	HIS	2.1

## 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. 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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
1	TPO	A	160	11/12	0.99	0.09	-	21,25,27,31	0
1	TPO	C	160	11/12	0.99	0.07	-	24,28,31,32	0

## 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
4	SGM	B	1434	6/6	0.70	0.27	8.36	48,66,69,78	0
4	SGM	D	1433	6/6	0.73	0.32	7.31	68,70,71,157	0
5	SO4	B	1435	5/5	0.97	0.15	1.18	73,75,77,81	0
3	T6Q	A	1299	33/33	0.93	0.14	0.33	20,31,70,76	0
3	T6Q	C	1298	33/33	0.87	0.15	0.32	42,46,92,95	0
5	SO4	D	1435	5/5	0.93	0.16	-0.12	80,80,82,82	0
4	SGM	B	1433	6/6	0.83	0.21	-	74,80,81,146	0
4	SGM	D	1434	6/6	0.77	0.36	-	76,79,81,160	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.