



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:59 PM UTC

PDB ID : 3BHT / pdb_00003bht
Title : Structure of phosphorylated Thr160 CDK2/cyclin A in complex with the inhibitor meriolin 3
Authors : Echalier, A.; Bettayeb, K.; Ferandin, Y.; Lozach, O.; Clement, M.; Valette, A.; Liger, F.; Marquet, B.; Morris, J.C.; Endicott, J.A.; Joseph, B.; Meijer, L.
Deposited on : 2007-11-29
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

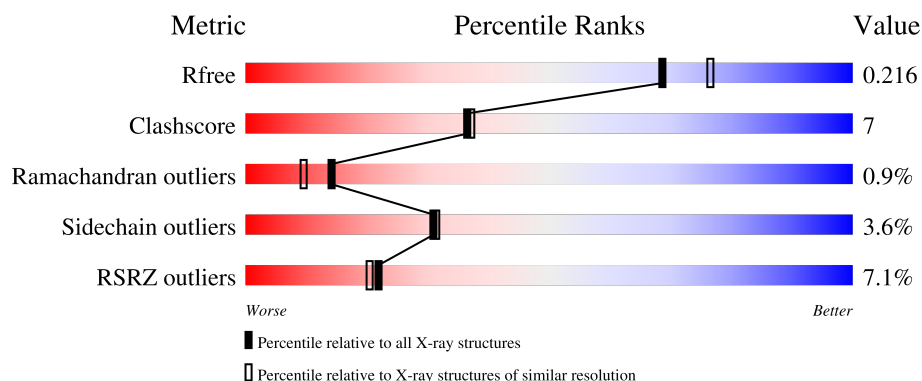
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	300	8% 83% 14% ...
1	C	300	8% 76% 12% 10%
2	B	262	5% 90% 10% .
2	D	262	6% 87% 10% .

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MFR	A	299	-	X	-	-
3	MFR	C	299	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9983 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	4	0
			2416	1568	412	426	1	9			
1	C	271	Total	C	N	O	P	S	0	2	0
			2185	1417	373	387	1	7			

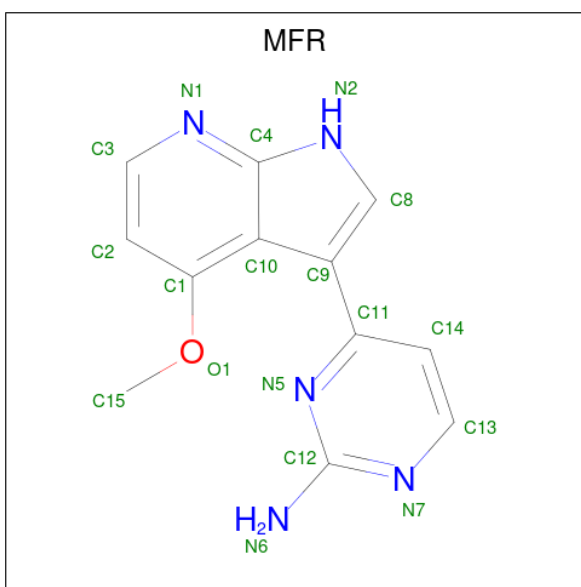
There are 4 discrepancies between the modelled and reference sequences:

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	262	Total	C	N	O	S		0	3	0
			2130	1379	346	395	10				
2	D	262	Total	C	N	O	S		0	1	0
			2118	1371	344	393	10				

- Molecule 3 is 4-(4-methoxy-1H-pyrrolo[2,3-b]pyridin-3-yl)pyrimidin-2-amine (CCD ID: MFR) (formula: C₁₂H₁₁N₅O).

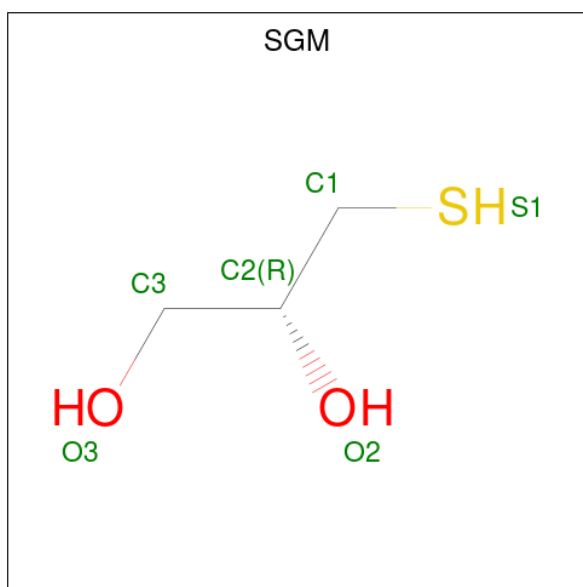


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			18	12	5	1		
3	C	1	Total	C	N	O	0	0
			18	12	5	1		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is MONOTHIOGLYCEROL (CCD ID: SGM) (formula: C₃H₈O₂S).



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

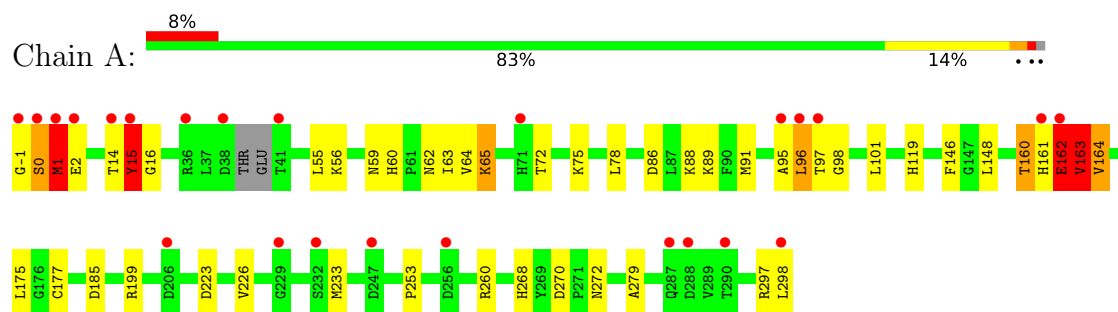
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	345	Total	O	0	0
			345	345		
6	B	292	Total	O	0	0
			292	292		
6	C	227	Total	O	0	0
			227	227		
6	D	214	Total	O	0	0
			214	214		

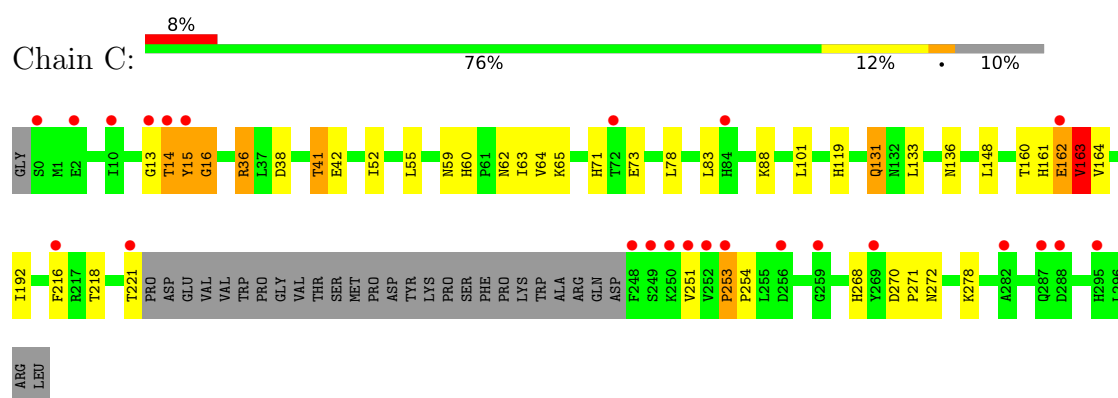
3 Residue-property plots [i](#)

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.

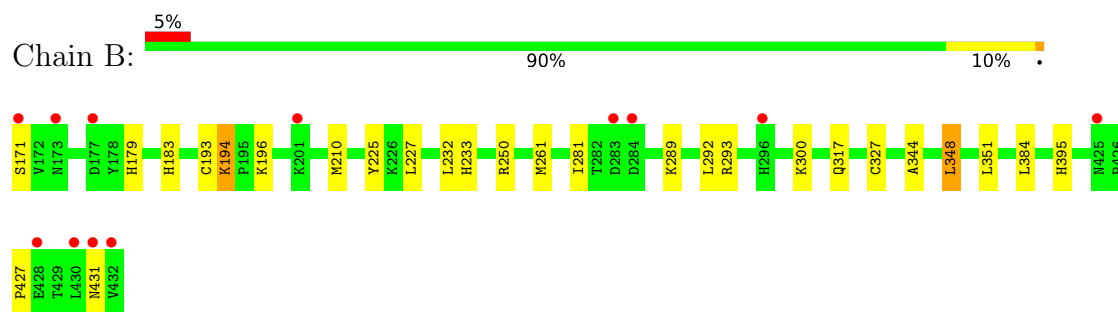
- Molecule 1: Cell division protein kinase 2



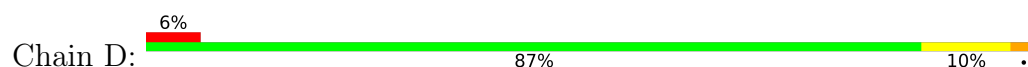
- Molecule 1: Cell division protein kinase 2

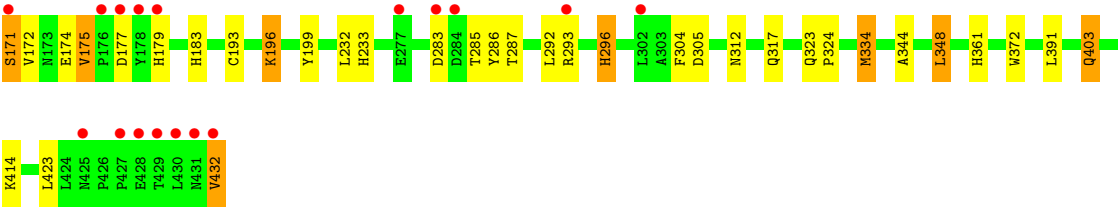


- Molecule 2: Cyclin-A2



- Molecule 2: Cyclin-A2





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.13Å 133.77Å 147.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.5 (20.00-2.00) 95.4 (20.00-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.71 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.181 , 0.220 0.181 , 0.216	Depositor DCC
R_{free} test set	4777 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	25.7	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9983	wwPDB-VP
Average B, all atoms (Å ²)	9.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	6/2477 (0.2%)	0.94	9/3357 (0.3%)
1	C	0.69	3/2229 (0.1%)	0.85	3/3016 (0.1%)
2	B	0.67	1/2190 (0.0%)	0.81	0/2978
2	D	0.64	2/2171 (0.1%)	0.97	3/2952 (0.1%)
All	All	0.73	12/9067 (0.1%)	0.90	15/12303 (0.1%)

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 (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	431	ASN	C-N	-7.14	1.23	1.33
1	C	63	ILE	C-O	-6.85	1.17	1.24
1	A	64	VAL	C-O	-6.77	1.17	1.24
2	D	171	SER	C-N	6.23	1.41	1.34
1	C	64	VAL	C-O	-5.99	1.18	1.24
1	A	56	LYS	C-O	-5.58	1.17	1.24
2	D	172	VAL	C-O	-5.56	1.17	1.24
1	A	63	ILE	C-O	-5.23	1.18	1.24
1	A	164	VAL	C-O	-5.20	1.17	1.24
1	A	15	TYR	CA-C	5.15	1.59	1.52
1	A	279	ALA	N-CA	5.09	1.52	1.46
1	C	38	ASP	C-O	-5.05	1.18	1.23

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	174	GLU	N-CA-C	20.03	138.45	112.41
1	A	162	GLU	CA-C-N	6.80	133.93	121.70
1	A	162	GLU	C-N-CA	6.80	133.93	121.70
2	D	175	VAL	N-CA-C	-6.56	92.63	111.00
1	C	163	VAL	N-CA-C	6.51	122.89	109.34
2	D	283	ASP	CB-CA-C	-6.50	109.09	116.63
1	C	253	PRO	N-CA-C	6.50	118.62	110.70
1	A	65	LYS	N-CA-C	6.24	118.79	109.25
1	A	163	VAL	N-CA-C	-6.16	93.75	111.00
1	A	15	TYR	N-CA-C	5.81	123.18	110.80
1	A	253	PRO	N-CA-C	5.75	117.72	110.70
1	A	253	PRO	CA-C-N	-5.50	114.34	120.45
1	A	253	PRO	C-N-CA	-5.50	114.34	120.45
1	C	65	LYS	N-CA-C	5.33	117.44	108.96
1	A	1	MET	N-CA-C	5.19	121.86	110.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	14	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	2416	0	2471	52	1
1	C	2185	0	2250	42	1
2	B	2130	0	2146	19	0
2	D	2118	0	2136	36	0
3	A	18	0	11	0	0
3	C	18	0	11	0	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
5	B	12	0	15	2	0
5	D	6	0	8	3	0
6	A	345	0	0	9	0
6	B	292	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	C	227	0	0	5	0
6	D	214	0	0	8	0
All	All	9983	0	9048	130	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 7.

All (130) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177[A]:CYS:SG	1:A:233:MET:HE3	1.50	1.47
2:B:193:CYS:SG	5:B:4:SGM:S1	2.31	1.27
2:D:193:CYS:SG	5:D:3:SGM:S1	2.38	1.22
2:B:171:SER:OG	6:B:595:HOH:O	1.68	1.12
1:C:272:ASN:OD1	2:D:171:SER:HB3	1.47	1.11
1:A:177[A]:CYS:SG	1:A:233:MET:CE	2.38	1.11
1:C:15:TYR:HB2	1:C:16:GLY:CA	1.82	1.08
1:C:15:TYR:HB2	1:C:16:GLY:HA3	1.32	1.06
1:A:15:TYR:CD2	1:A:16:GLY:HA3	1.90	1.05
1:A:272:ASN:OD1	2:B:171:SER:HB2	1.55	1.05
2:D:432:VAL:HG23	6:D:580:HOH:O	1.59	0.99
1:A:2:GLU:HG3	2:D:293:ARG:NH2	1.82	0.94
1:A:2:GLU:CG	2:D:293:ARG:NH2	2.31	0.93
1:C:15:TYR:CB	1:C:16:GLY:HA3	1.91	0.93
1:C:36:ARG:HG3	1:C:36:ARG:HH11	1.33	0.91
1:A:177[A]:CYS:HG	1:A:233:MET:HE3	1.14	0.91
1:A:2:GLU:HG2	1:C:73:GLU:OE2	1.70	0.91
1:A:161:HIS:HD2	6:A:353:HOH:O	1.55	0.90
2:D:171:SER:HB2	6:D:697:HOH:O	1.71	0.89
1:C:71:HIS:HD2	2:D:296:HIS:HE1	1.14	0.89
1:A:160:TPO:HG23	1:A:162:GLU:HG2	1.59	0.84
1:A:15:TYR:CG	1:A:16:GLY:HA3	2.14	0.83
1:C:71:HIS:HD2	2:D:296:HIS:CE1	1.97	0.83
1:A:268:HIS:CD2	6:A:589:HOH:O	2.33	0.81
1:A:2:GLU:HG2	2:D:293:ARG:NH2	1.99	0.78
1:A:15:TYR:CD2	1:A:16:GLY:CA	2.65	0.77
1:C:71:HIS:CD2	2:D:296:HIS:HE1	2.01	0.76
2:B:183:HIS:HB2	2:B:317:GLN:HE22	1.51	0.74
2:D:344:ALA:HB1	2:D:348:LEU:HD22	1.68	0.74
1:A:60:HIS:HD2	1:A:62:ASN:H	1.33	0.74
1:C:36:ARG:HH11	1:C:36:ARG:CG	2.00	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:ALA:O	1:A:96:LEU:HB2	1.87	0.73
1:A:161:HIS:O	1:A:161:HIS:CG	2.42	0.71
2:B:327:CYS:SG	5:B:5:SGM:S1	2.82	0.71
1:A:60:HIS:CD2	1:A:62:ASN:H	2.09	0.70
1:A:160:TPO:CG2	1:A:162:GLU:HG2	2.21	0.70
1:A:59[A]:ASN:ND2	6:A:333:HOH:O	2.26	0.69
1:A:161:HIS:O	1:A:162:GLU:C	2.36	0.69
1:A:272:ASN:CG	2:B:171:SER:HB2	2.17	0.68
2:D:305:ASP:HB3	5:D:3:SGM:H12	1.76	0.67
2:B:344:ALA:HB1	2:B:348:LEU:HD22	1.77	0.66
1:A:268:HIS:NE2	6:A:589:HOH:O	2.28	0.66
1:C:15:TYR:CB	1:C:16:GLY:CA	2.60	0.64
1:A:15:TYR:HD2	1:A:16:GLY:CA	2.12	0.62
1:A:15:TYR:CB	1:A:16:GLY:HA3	2.29	0.62
1:C:268:HIS:HD2	1:C:270:ASP:H	1.46	0.62
1:A:88:LYS:HE2	6:A:593:HOH:O	2.00	0.62
1:C:15:TYR:HB2	1:C:16:GLY:HA2	1.79	0.62
2:D:285:THR:HG22	2:D:286:TYR:CD1	2.35	0.62
1:C:52:ILE:HD11	1:C:78:LEU:HD21	1.82	0.60
1:C:71:HIS:CD2	2:D:296:HIS:CE1	2.84	0.60
1:A:2:GLU:CG	2:D:293:ARG:HH22	2.14	0.59
1:A:161:HIS:O	1:A:161:HIS:CD2	2.54	0.59
2:B:194:LYS:HD3	2:B:351:LEU:HD23	1.83	0.59
1:C:272:ASN:OD1	2:D:171:SER:CB	2.38	0.59
1:A:162:GLU:HG3	1:A:163:VAL:HG12	1.83	0.59
1:C:41:THR:HG22	1:C:42:GLU:N	2.16	0.59
1:C:60:HIS:CD2	1:C:62:ASN:H	2.21	0.58
1:C:59:ASN:ND2	6:C:458:HOH:O	2.37	0.58
1:C:278:LYS:HE3	2:D:177:ASP:O	2.04	0.58
2:D:287:THR:HB	6:D:710:HOH:O	2.04	0.57
1:C:161:HIS:O	1:C:162:GLU:C	2.46	0.57
2:B:395:HIS:HE1	2:B:427:PRO:O	1.87	0.56
2:B:183:HIS:HD2	6:B:553:HOH:O	1.87	0.56
1:C:60:HIS:HD2	1:C:62:ASN:H	1.51	0.56
1:A:0:SER:O	1:A:1:MET:CB	2.54	0.55
1:A:0:SER:O	1:A:1:MET:HB2	2.07	0.54
1:A:161:HIS:O	1:A:162:GLU:O	2.27	0.53
1:C:218:THR:HG23	1:C:251:VAL:HG21	1.91	0.52
1:A:2:GLU:HG2	2:D:293:ARG:HH21	1.75	0.52
1:A:119:HIS:HD2	6:B:543:HOH:O	1.91	0.52
1:C:15:TYR:HB3	6:C:353:HOH:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:HIS:O	1:C:162:GLU:O	2.28	0.51
1:A:2:GLU:CG	2:D:293:ARG:HH21	2.22	0.51
2:B:210:MET:HE1	2:B:250:ARG:CB	2.41	0.51
1:C:161:HIS:HD2	6:C:372:HOH:O	1.94	0.50
2:B:233:HIS:HE1	6:B:447:HOH:O	1.94	0.50
1:C:216:PHE:HB3	1:C:221:THR:HG23	1.94	0.50
1:C:71:HIS:NE2	2:D:304:PHE:HE1	2.10	0.49
2:D:233:HIS:HE1	6:D:551:HOH:O	1.94	0.49
1:C:15:TYR:CG	1:C:16:GLY:HA3	2.47	0.49
1:A:260:ARG:HD3	6:A:334:HOH:O	2.11	0.49
2:B:225:TYR:HE2	2:B:281:ILE:HG21	1.78	0.49
2:B:179[A]:HIS:CE1	6:B:492:HOH:O	2.66	0.48
1:C:161:HIS:CE1	6:C:490:HOH:O	2.67	0.48
2:B:289:LYS:HE3	2:B:293:ARG:HE	1.78	0.48
1:C:83:LEU:HD23	1:C:136:ASN:HB3	1.95	0.47
1:A:223:ASP:H	1:A:226:VAL:HG12	1.79	0.47
1:C:13:GLY:C	1:C:15:TYR:H	2.23	0.47
1:A:162:GLU:HG3	1:A:163:VAL:CG1	2.45	0.46
1:C:71:HIS:HE1	6:D:663:HOH:O	1.99	0.46
2:D:285:THR:HG22	2:D:286:TYR:CG	2.50	0.46
1:C:119:HIS:HD2	6:D:535:HOH:O	1.98	0.46
1:A:60:HIS:HE1	6:A:441:HOH:O	1.98	0.45
1:A:175:LEU:HB2	1:A:233:MET:HE2	1.97	0.45
2:D:183:HIS:HB2	2:D:317:GLN:HE22	1.81	0.45
1:A:88:LYS:HA	1:A:91:MET:CE	2.47	0.45
2:B:300:LYS:HE3	6:B:614:HOH:O	2.16	0.45
1:A:268:HIS:HD2	1:A:270:ASP:H	1.65	0.44
2:D:285:THR:HG22	2:D:286:TYR:CE1	2.52	0.44
6:C:439:HOH:O	2:D:296:HIS:HD2	2.01	0.44
1:A:98:GLY:HA2	1:A:199:ARG:NE	2.33	0.43
2:D:196:LYS:HG2	2:D:199:TYR:HB3	2.01	0.43
1:A:15:TYR:CB	1:A:16:GLY:CA	2.96	0.43
1:A:88:LYS:HA	1:A:91:MET:HE3	1.99	0.43
1:A:161:HIS:CD2	6:A:353:HOH:O	2.44	0.43
2:D:312:ASN:OD1	2:D:334:MET:HE1	2.19	0.43
1:C:88:LYS:HG3	1:C:131:GLN:NE2	2.34	0.43
2:B:194:LYS:HD3	2:B:351:LEU:CD2	2.48	0.43
1:C:272:ASN:CG	2:D:171:SER:HB3	2.32	0.43
2:B:210:MET:HE2	2:B:210:MET:HB3	1.85	0.43
1:C:36:ARG:HG3	1:C:36:ARG:NH1	2.15	0.43
1:A:65:LYS:HD2	6:A:333:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LEU:HD21	1:A:146:PHE:CD1	2.54	0.42
1:C:270:ASP:HA	1:C:271:PRO:HD3	1.92	0.42
1:A:-1:GLY:HA3	1:A:0:SER:HA	1.74	0.42
2:D:323:GLN:HA	2:D:324:PRO:HA	1.87	0.42
2:D:361:HIS:HB2	2:D:372:TRP:HB2	2.01	0.42
2:D:193:CYS:HG	5:D:3:SGM:C1	2.29	0.41
2:D:414:LYS:HG2	2:D:423:LEU:HG	2.02	0.41
1:C:41:THR:HG22	1:C:42:GLU:H	1.84	0.41
1:C:73:GLU:OE2	2:D:293:ARG:NH2	2.53	0.41
1:A:162:GLU:HG3	1:A:163:VAL:HA	2.02	0.41
1:C:133:LEU:HD11	1:C:192:ILE:HD13	2.02	0.41
2:D:403:GLN:HG3	6:D:527:HOH:O	2.21	0.41
2:D:179:HIS:HB2	6:D:716:HOH:O	2.20	0.41
1:A:119:HIS:HE1	1:A:185:ASP:OD2	2.04	0.40
1:C:253:PRO:HD2	1:C:254:PRO:HD3	2.01	0.40
1:A:86:ASP:OD2	1:A:89:LYS:HE3	2.22	0.40
2:B:227:LEU:HD12	2:B:261:MET:HE3	2.02	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)
1:A:297:ARG:NH2	1:C:251:VAL:CG2[3_454]	2.18	0.02

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	297/300 (99%)	283 (95%)	9 (3%)	5 (2%)	7	3
1	C	267/300 (89%)	254 (95%)	8 (3%)	5 (2%)	6	3
2	B	263/262 (100%)	261 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	261/262 (100%)	256 (98%)	5 (2%)	0	100	100
All	All	1088/1124 (97%)	1054 (97%)	24 (2%)	10 (1%)	14	9

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	15	TYR
1	A	162	GLU
1	C	162	GLU
1	C	163	VAL
1	A	164	VAL
1	C	164	VAL
1	A	96	LEU
1	C	14	THR
1	C	16	GLY

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	265/263 (101%)	255 (96%)	10 (4%)	29	29
1	C	239/263 (91%)	230 (96%)	9 (4%)	29	29
2	B	238/235 (101%)	232 (98%)	6 (2%)	42	45
2	D	236/235 (100%)	226 (96%)	10 (4%)	26	25
All	All	978/996 (98%)	943 (96%)	35 (4%)	31	31

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	SER
1	A	72	THR
1	A	75	LYS
1	A	78	LEU

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Mol	Chain	Res	Type
1	A	97	THR
1	A	101	LEU
1	A	148	LEU
1	A	162	GLU
1	A	163	VAL
1	A	298	LEU
2	B	194	LYS
2	B	196	LYS
2	B	232	LEU
2	B	292	LEU
2	B	348	LEU
2	B	384	LEU
1	C	14	THR
1	C	15	TYR
1	C	36	ARG
1	C	41	THR
1	C	55	LEU
1	C	101	LEU
1	C	131	GLN
1	C	148	LEU
1	C	163	VAL
2	D	175	VAL
2	D	196	LYS
2	D	232	LEU
2	D	292	LEU
2	D	296	HIS
2	D	334	MET
2	D	348	LEU
2	D	391	LEU
2	D	403	GLN
2	D	432	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	74	ASN
1	A	85	GLN
1	A	119	HIS
1	A	161	HIS
1	A	246	GLN
1	A	287	GLN

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Mol	Chain	Res	Type
2	B	183	HIS
2	B	233	HIS
2	B	254	GLN
2	B	317	GLN
2	B	395	HIS
2	B	403	GLN
1	C	59	ASN
1	C	60	HIS
1	C	71	HIS
1	C	84	HIS
1	C	85	GLN
1	C	113	GLN
1	C	119	HIS
1	C	161	HIS
1	C	265	GLN
1	C	268	HIS
1	C	287	GLN
2	D	173	ASN
2	D	228	GLN
2	D	233	HIS
2	D	254	GLN
2	D	296	HIS
2	D	313	GLN
2	D	317	GLN
2	D	396	GLN
2	D	419	HIS

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	8,10,11	2.25	3 (37%)	10,14,16	1.46	2 (20%)
1	TPO	C	160	1	8,10,11	2.35	4 (50%)	10,14,16	0.94	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	-	0/9/11/13	-
1	TPO	C	160	1	-	1/9/11/13	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	160	TPO	P-O3P	-4.25	1.39	1.54
1	C	160	TPO	P-O3P	-4.05	1.39	1.54
1	C	160	TPO	P-O2P	-3.63	1.41	1.54
1	A	160	TPO	P-O2P	-3.57	1.41	1.54
1	C	160	TPO	P-O1P	-3.03	1.41	1.50
1	A	160	TPO	P-O1P	-2.71	1.42	1.50
1	C	160	TPO	P-OG1	-2.05	1.55	1.59

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	160	TPO	OG1-P-O1P	-2.97	98.73	109.33
1	A	160	TPO	O3P-P-O2P	2.33	116.53	107.80

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	160	TPO	O-C-CA-CB

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	160	TPO	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 for Mogul analysis.

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	MFR	A	299	-	20,20,20	3.53	3 (15%)	26,28,28	3.47	17 (65%)
5	SGM	B	5	-	5,5,5	0.44	0	5,5,5	0.28	0
5	SGM	D	3	-	5,5,5	0.40	0	5,5,5	0.47	0
3	MFR	C	299	-	20,20,20	3.63	3 (15%)	26,28,28	3.70	18 (69%)
5	SGM	B	4	-	5,5,5	0.44	0	5,5,5	0.51	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	MFR	A	299	-	-	3/6/6/6	0/3/3/3
5	SGM	B	5	-	-	2/4/4/4	-
5	SGM	D	3	-	-	0/4/4/4	-
3	MFR	C	299	-	-	3/6/6/6	0/3/3/3
5	SGM	B	4	-	-	2/4/4/4	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	299	MFR	C8-C9	14.36	1.51	1.37
3	A	299	MFR	C8-C9	14.05	1.51	1.37
3	C	299	MFR	C10-C4	4.96	1.47	1.41
3	C	299	MFR	C10-C1	4.82	1.49	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	299	MFR	C10-C1	4.69	1.49	1.40
3	A	299	MFR	C10-C4	4.26	1.46	1.41

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	299	MFR	C11-N5-C12	7.79	122.12	116.37
3	C	299	MFR	C11-N5-C12	6.87	121.44	116.37
3	C	299	MFR	C4-N2-C8	6.41	111.50	108.48
3	C	299	MFR	C11-C9-C8	-6.31	116.25	127.41
3	C	299	MFR	C10-C9-C11	5.41	134.28	126.40
3	A	299	MFR	C4-N2-C8	5.34	111.00	108.48
3	A	299	MFR	N7-C12-N5	-5.19	119.92	125.70
3	A	299	MFR	C10-C9-C11	4.93	133.57	126.40
3	C	299	MFR	N7-C12-N5	-4.89	120.26	125.70
3	C	299	MFR	O1-C1-C10	4.86	122.73	115.84
3	A	299	MFR	C11-C9-C8	-4.68	119.13	127.41
3	C	299	MFR	C13-N7-C12	4.65	120.84	116.21
3	A	299	MFR	C13-N7-C12	4.34	120.53	116.21
3	A	299	MFR	N6-C12-N7	3.97	121.51	117.40
3	C	299	MFR	C3-C2-C1	3.96	122.47	118.07
3	A	299	MFR	C3-C2-C1	3.95	122.46	118.07
3	A	299	MFR	O1-C1-C10	3.88	121.34	115.84
3	A	299	MFR	C14-C11-N5	-3.83	118.46	122.92
3	A	299	MFR	C10-C4-N1	-3.77	122.62	126.64
3	C	299	MFR	C14-C13-N7	-3.64	119.51	123.97
3	C	299	MFR	C14-C11-N5	-3.59	118.74	122.92
3	C	299	MFR	C10-C4-N1	-3.55	122.86	126.64
3	A	299	MFR	C15-O1-C1	3.48	122.61	117.51
3	C	299	MFR	C13-C14-C11	3.46	119.15	117.02
3	A	299	MFR	C10-C9-C8	-3.28	103.55	105.68
3	C	299	MFR	C15-O1-C1	3.28	122.32	117.51
3	C	299	MFR	N6-C12-N7	3.27	120.78	117.40
3	C	299	MFR	C9-C8-N2	-3.19	108.41	110.46
3	C	299	MFR	C10-C9-C8	-3.16	103.63	105.68
3	C	299	MFR	C9-C11-N5	3.12	121.49	116.40
3	A	299	MFR	C14-C13-N7	-2.97	120.34	123.97
3	A	299	MFR	C9-C8-N2	-2.91	108.58	110.46
3	A	299	MFR	C2-C1-C10	-2.42	115.23	120.14
3	C	299	MFR	C2-C1-C10	-2.32	115.44	120.14
3	A	299	MFR	C13-C14-C11	2.14	118.33	117.02

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	4	SGM	S1-C1-C2-O2
3	A	299	MFR	N5-C11-C9-C10
3	C	299	MFR	N5-C11-C9-C10
5	B	5	SGM	C1-C2-C3-O3
5	B	5	SGM	O2-C2-C3-O3
3	A	299	MFR	N5-C11-C9-C8
3	C	299	MFR	N5-C11-C9-C8
3	A	299	MFR	C14-C11-C9-C8
3	C	299	MFR	C14-C11-C9-C8
5	B	4	SGM	S1-C1-C2-C3

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	5	SGM	1	0
5	D	3	SGM	3	0
5	B	4	SGM	1	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 [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	297/300 (99%)	0.34	24 (8%)	18 17	2, 5, 22, 43	4 (1%)
1	C	270/300 (90%)	0.34	24 (8%)	15 14	2, 6, 33, 59	2 (0%)
2	B	262/262 (100%)	-0.03	12 (4%)	37 36	1, 5, 15, 27	3 (1%)
2	D	262/262 (100%)	0.34	17 (6%)	25 23	2, 6, 21, 43	1 (0%)
All	All	1091/1124 (97%)	0.25	77 (7%)	22 20	1, 6, 24, 59	10 (0%)

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	428	GLU	7.2
2	D	171	SER	7.1
1	C	248	PHE	6.4
2	B	171	SER	6.1
1	A	96	LEU	5.5
1	C	249	SER	5.4
1	C	221	THR	5.3
1	A	14	THR	5.3
1	C	251	VAL	5.1
1	C	15	TYR	5.0
2	D	432	VAL	4.9
2	B	284	ASP	4.0
1	A	247	ASP	4.0
2	B	432	VAL	3.9
2	D	431	ASN	3.9
1	C	13	GLY	3.8
1	A	0	SER	3.7
1	C	295	HIS	3.7
2	D	429	THR	3.7
1	A	-1	GLY	3.6
2	D	179	HIS	3.4

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Mol	Chain	Res	Type	RSRZ
2	D	284	ASP	3.4
1	A	1	MET	3.3
1	A	2	GLU	3.3
1	C	269	TYR	3.3
1	A	162	GLU	3.2
2	B	201	LYS	3.2
2	B	431	ASN	3.2
1	C	288	ASP	3.1
1	A	41	THR	3.0
1	C	2	GLU	3.0
2	D	430	LEU	3.0
2	B	428	GLU	3.0
1	C	287	GLN	2.9
1	C	72	THR	2.8
1	C	250	LYS	2.8
1	C	84	HIS	2.8
1	C	0[A]	SER	2.8
1	C	14	THR	2.8
1	A	287	GLN	2.7
2	B	283[A]	ASP	2.7
1	C	216	PHE	2.6
1	A	38	ASP	2.6
1	A	206[A]	ASP	2.6
1	A	36	ARG	2.6
1	A	232	SER	2.6
2	B	425	ASN	2.5
2	B	173	ASN	2.5
2	D	177	ASP	2.4
2	D	283	ASP	2.4
1	C	253	PRO	2.4
2	B	177	ASP	2.4
2	B	430	LEU	2.4
1	C	162	GLU	2.3
2	D	178	TYR	2.3
1	A	229	GLY	2.3
2	D	176	PRO	2.3
1	A	161	HIS	2.2
1	A	97	THR	2.2
2	D	427	PRO	2.2
1	C	282	ALA	2.2
1	A	15	TYR	2.2
2	D	425	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	95	ALA	2.2
1	A	298	LEU	2.2
2	B	296	HIS	2.1
1	C	256	ASP	2.1
2	D	277	GLU	2.1
1	A	290	THR	2.1
1	C	252	VAL	2.1
2	D	302	LEU	2.1
1	A	288	ASP	2.1
1	C	259	GLY	2.1
1	A	71	HIS	2.1
1	A	256	ASP	2.0
2	D	293	ARG	2.0
1	C	10	ILE	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, 95th 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($\text{\AA}^2$)	Q<0.9
1	TPO	A	160	11/12	0.98	0.07	2,4,6,7	0
1	TPO	C	160	11/12	0.98	0.07	2,4,7,7	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. The B-factors column lists the minimum, median, 95th 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($\text{\AA}^2$)	Q<0.9
5	SGM	B	5	6/6	0.58	0.24	61,63,63,64	0
5	SGM	D	3	6/6	0.84	0.16	41,43,43,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SGM	B	4	6/6	0.88	0.16	30,31,33,34	0
4	MG	B	1	1/1	0.89	0.10	15,15,15,15	0
3	MFR	C	299	18/18	0.94	0.10	2,2,4,5	0
3	MFR	A	299	18/18	0.95	0.05	2,2,2,2	0
4	MG	D	2	1/1	0.98	0.18	9,9,9,9	0

6.5 Other polymers [i](#)

There are no such residues in this entry.