



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:57 pm GMT

PDB ID : 2UZD
Title : CRYSTAL STRUCTURE OF HUMAN CDK2 COMPLEXED WITH A THI-AZOLIDINONE INHIBITOR
Authors : Richardson, C.M.; Dokurno, P.; Murray, J.B.; Surgenor, A.E.
Deposited on : 2007-04-27
Resolution : 2.72 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

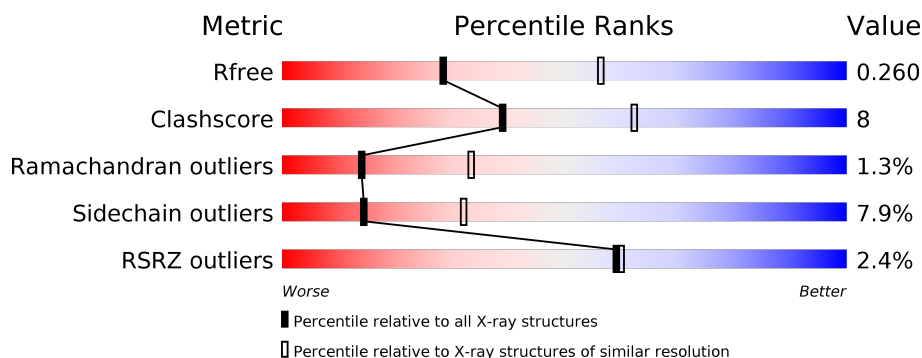
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.72 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	2649 (2.74-2.70)
Clashscore	112137	2993 (2.74-2.70)
Ramachandran outliers	110173	2946 (2.74-2.70)
Sidechain outliers	110143	2947 (2.74-2.70)
RSRZ outliers	101464	2665 (2.74-2.70)

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 ≥ 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	298	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>18%</div> <div>• •</div> </div> </div>
1	C	298	<div> <div>4%</div> <div> <div></div> <div>78%</div> <div>18%</div> <div>• •</div> </div> </div>
2	B	258	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>18%</div> <div>•</div> </div> </div>
2	D	258	<div> <div>%</div> <div> <div></div> <div>81%</div> <div>17%</div> <div>•</div> </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
1	TPO	A	160	X	-	-	-
1	TPO	C	160	X	-	-	-
3	C85	C	1297	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9177 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.

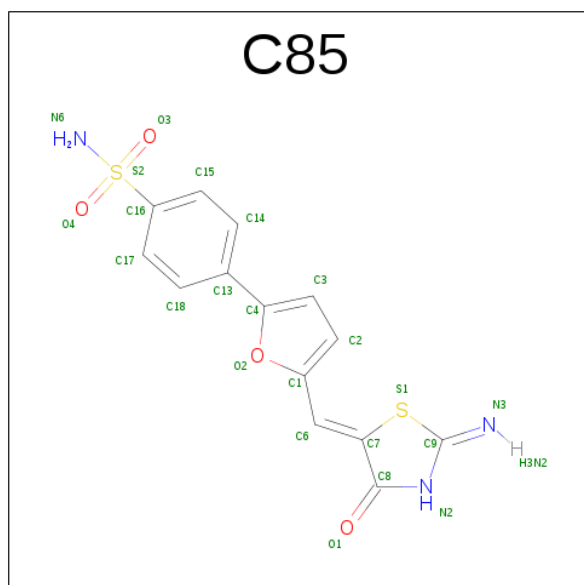
- Molecule 1 is a protein called CELL DIVISION PROTEIN KINASE 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	P	S	0	0	0
			2382	1547	403	423	1	8			
1	C	296	Total	C	N	O	P	S	0	0	0
			2382	1547	403	423	1	8			

- 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	258	Total	C	N	O	S	0	0	0
			2083	1350	339	383	11			

- Molecule 3 is 4-{5-[(Z)-(2-IMINO-4-OXO-1,3-THIAZOLIDIN-5-YLIDENE)METHYL]FURAN-2-YL}BENZENESULFONAMIDE (three-letter code: C85) (formula: C₁₄H₁₁N₃O₄S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			23	14	3	4	2		
3	C	1	Total	C	N	O	S	0	0
			23	14	3	4	2		

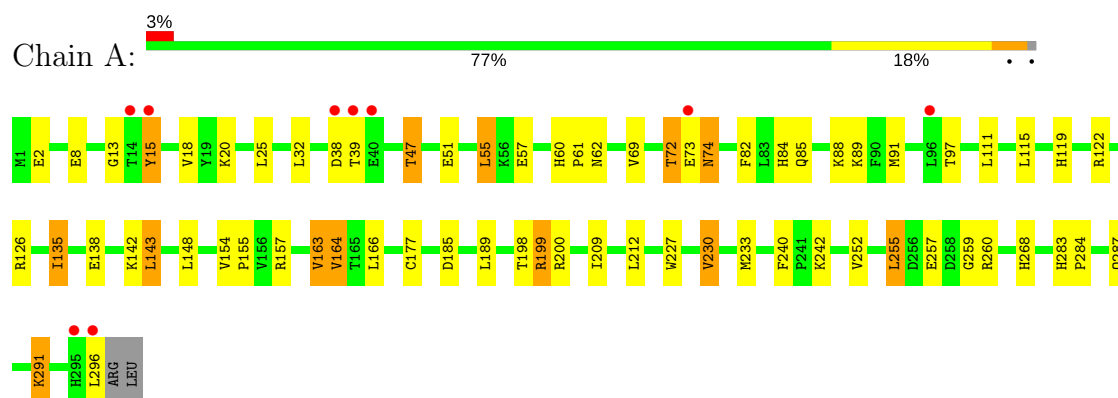
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	53	Total	O	0	0
			53	53		
4	B	42	Total	O	0	0
			42	42		
4	C	60	Total	O	0	0
			60	60		
4	D	46	Total	O	0	0
			46	46		

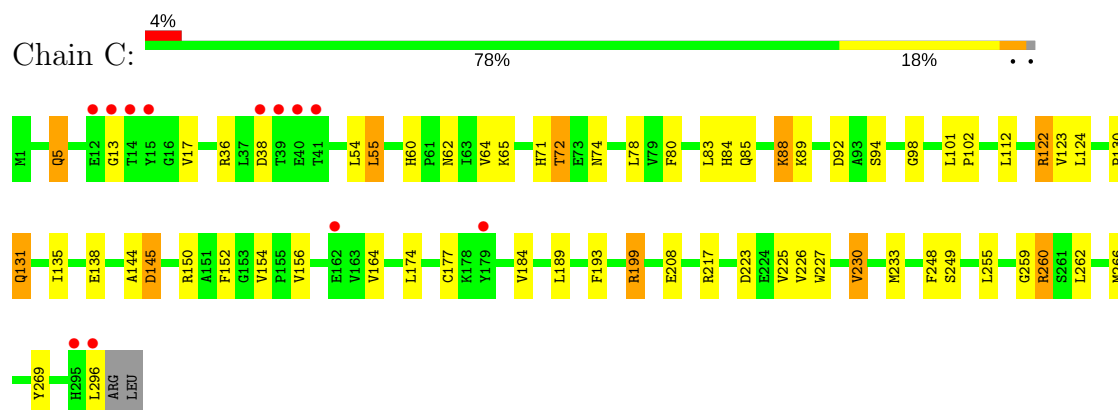
3 Residue-property plots [i](#)

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.

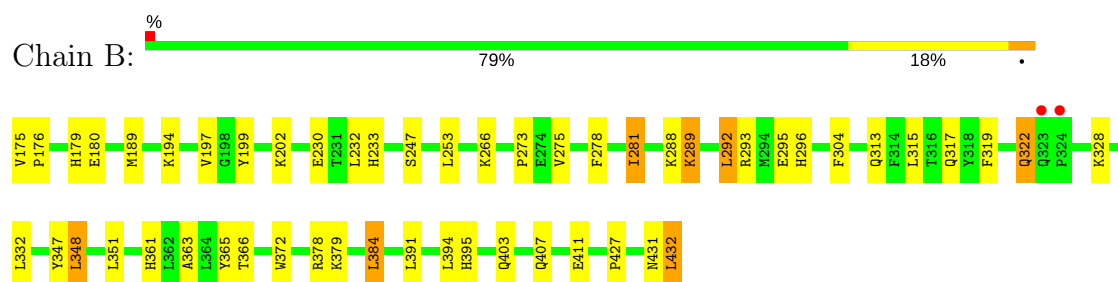
• Molecule 1: CELL DIVISION PROTEIN KINASE 2



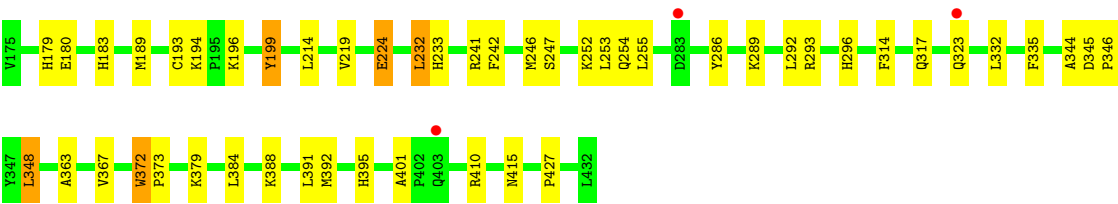
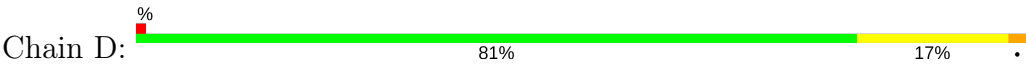
• Molecule 1: CELL DIVISION PROTEIN KINASE 2



• Molecule 2: CYCLIN A2



• Molecule 2: CYCLIN A2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	160.56Å 160.29Å 66.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.72 14.98 – 2.72	Depositor EDS
% Data completeness (in resolution range)	93.4 (15.00-2.72) 93.4 (14.98-2.72)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.208 , 0.269 0.202 , 0.260	Depositor DCC
R_{free} test set	2223 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	29.3	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9177	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, C85

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	0.61	0/2432	0.70	0/3300
1	C	0.61	0/2432	0.71	1/3300 (0.0%)
2	B	0.59	0/2133	0.69	0/2897
2	D	0.60	0/2133	0.68	0/2897
All	All	0.60	0/9130	0.69	1/12394 (0.0%)

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	1	0
1	C	1	0
All	All	2	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	296	LEU	CA-CB-CG	5.42	127.78	115.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	160	TPO	CB
1	C	160	TPO	CB

There are no planarity outliers.

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	2382	0	2425	49	0
1	C	2382	0	2425	46	0
2	B	2083	0	2107	30	0
2	D	2083	0	2107	32	0
3	A	23	0	10	1	0
3	C	23	0	10	5	0
4	A	53	0	0	5	0
4	B	42	0	0	6	0
4	C	60	0	0	9	0
4	D	46	0	0	6	0
All	All	9177	0	9084	153	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 (153) 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:296:HIS:HD2	4:B:2025:HOH:O	1.57	0.88
1:A:177:CYS:SG	1:A:233:MET:SD	2.72	0.88
1:C:260:ARG:HG3	1:C:260:ARG:HH11	1.39	0.88
2:B:296:HIS:CD2	4:B:2025:HOH:O	2.27	0.87
1:C:60:HIS:HD2	1:C:62:ASN:H	1.24	0.85
1:C:227:TRP:O	1:C:230:VAL:HG22	1.77	0.85
1:C:260:ARG:CG	1:C:260:ARG:HH11	1.92	0.82
1:A:155:PRO:HB2	4:A:2024:HOH:O	1.79	0.82
1:C:60:HIS:CD2	1:C:62:ASN:H	1.99	0.81
1:A:60:HIS:CD2	1:A:62:ASN:H	1.99	0.79
1:A:60:HIS:HD2	1:A:62:ASN:H	1.30	0.79
1:C:223:ASP:HB2	4:C:2052:HOH:O	1.82	0.79
2:D:344:ALA:HB1	2:D:348:LEU:HD22	1.65	0.78
1:A:198:THR:O	1:A:199:ARG:HG3	1.85	0.76
1:C:36:ARG:HB2	4:C:2008:HOH:O	1.87	0.74
2:D:219:VAL:HG22	2:D:232:LEU:HD11	1.68	0.74
1:A:15:TYR:HB2	4:A:2005:HOH:O	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:SER:O	1:C:199:ARG:HD2	1.91	0.70
2:D:233:HIS:HD2	4:D:2034:HOH:O	1.77	0.67
1:A:177:CYS:HB2	4:A:2033:HOH:O	1.95	0.66
2:B:289:LYS:HE3	2:B:293:ARG:HE	1.60	0.66
1:A:227:TRP:O	1:A:230:VAL:HG22	1.95	0.66
1:C:5:GLN:HA	1:C:5:GLN:HE21	1.61	0.65
2:D:332:LEU:HD23	2:D:363:ALA:HA	1.80	0.62
1:C:217:ARG:HD3	4:C:2049:HOH:O	1.99	0.62
1:C:71:HIS:CE1	2:D:296:HIS:HD2	2.17	0.61
2:B:361:HIS:HB2	2:B:372:TRP:HB2	1.82	0.60
1:A:260:ARG:HD3	4:A:2045:HOH:O	1.99	0.60
1:C:262:LEU:HG	1:C:266:MET:CE	2.31	0.60
1:A:199:ARG:CZ	1:A:200:ARG:HG2	2.30	0.60
1:A:227:TRP:HB3	1:A:230:VAL:HG22	1.84	0.59
2:B:332:LEU:HD23	2:B:363:ALA:HA	1.83	0.59
1:A:126:ARG:HD2	1:A:163:VAL:HG21	1.85	0.59
2:B:266:LYS:NZ	2:B:295:GLU:OE2	2.33	0.58
1:C:260:ARG:NH1	1:C:260:ARG:HG3	2.10	0.58
1:A:74:ASN:H	1:A:74:ASN:HD22	1.52	0.58
1:A:252:VAL:O	1:A:252:VAL:HG13	2.03	0.57
2:B:361:HIS:CE1	2:B:384:LEU:HD21	2.39	0.57
1:C:262:LEU:HG	1:C:266:MET:HE2	1.86	0.57
2:D:254:GLN:HG2	2:D:286:TYR:HE2	1.69	0.57
1:A:143:LEU:HD23	1:A:143:LEU:N	2.20	0.56
2:B:275:VAL:HG21	2:B:292:LEU:HD11	1.86	0.56
1:C:154:VAL:HG11	2:D:179:HIS:HD2	1.71	0.56
2:D:363:ALA:O	2:D:367:VAL:HG23	2.04	0.56
1:A:255:LEU:HG	1:A:259:GLY:HA3	1.87	0.56
2:B:202:LYS:HE2	4:C:2018:HOH:O	2.06	0.56
1:A:20:LYS:HD2	1:A:82:PHE:CZ	2.41	0.55
1:A:74:ASN:N	1:A:74:ASN:HD22	2.04	0.55
1:C:227:TRP:CD2	1:C:230:VAL:HG13	2.42	0.55
2:D:401:ALA:HB1	2:D:410:ARG:HD2	1.86	0.55
2:B:384:LEU:HD22	2:B:432:LEU:HD23	1.90	0.54
1:C:64:VAL:HG21	1:C:144:ALA:HB2	1.90	0.53
1:C:227:TRP:CE3	1:C:269:TYR:HB3	2.44	0.53
1:A:51:GLU:O	1:A:55:LEU:HB2	2.09	0.53
1:A:74:ASN:ND2	1:A:74:ASN:H	2.06	0.53
1:C:83:LEU:O	3:C:1297:C85:H3	2.09	0.53
1:A:119:HIS:HE1	1:A:185:ASP:OD2	1.92	0.53
1:A:119:HIS:HD2	4:B:2004:HOH:O	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:2009:HOH:O	2:B:304:PHE:HD1	1.91	0.52
2:D:194:LYS:HD3	4:D:2012:HOH:O	2.08	0.52
2:D:179:HIS:CE1	2:D:379:LYS:NZ	2.79	0.51
1:A:291:LYS:HG2	1:A:291:LYS:O	2.10	0.51
1:C:124:LEU:HG	1:C:152:PHE:CD1	2.46	0.51
1:C:260:ARG:NH1	1:C:260:ARG:CG	2.62	0.50
1:C:255:LEU:O	1:C:260:ARG:NH1	2.44	0.50
2:D:183:HIS:CE1	4:D:2009:HOH:O	2.65	0.50
2:B:289:LYS:HD3	2:B:293:ARG:NH2	2.26	0.49
2:D:388:LYS:O	2:D:392:MET:HG2	2.12	0.49
1:C:60:HIS:HD2	1:C:62:ASN:N	2.03	0.49
2:D:255:LEU:HB2	2:D:286:TYR:CE1	2.47	0.49
1:A:61:PRO:O	1:A:142:LYS:HE2	2.13	0.49
2:B:179:HIS:CE1	2:B:379:LYS:NZ	2.81	0.49
1:A:199:ARG:HD2	1:A:199:ARG:O	2.13	0.49
1:A:85:GLN:HE21	1:A:89:LYS:HB3	1.78	0.48
2:D:196:LYS:HB2	4:D:2013:HOH:O	2.13	0.48
1:C:54:LEU:HD13	1:C:123:VAL:HG13	1.95	0.48
1:A:25:LEU:O	2:D:252:LYS:HD2	2.14	0.48
1:A:163:VAL:HG13	1:A:164:VAL:HG23	1.96	0.48
2:B:319:PHE:O	2:B:322:GLN:HB2	2.14	0.48
2:D:183:HIS:HE1	4:D:2009:HOH:O	1.97	0.47
2:B:407:GLN:O	2:B:411:GLU:HG2	2.13	0.47
1:A:227:TRP:O	1:A:230:VAL:CG2	2.62	0.47
1:A:154:VAL:HG13	2:B:176:PRO:HG2	1.96	0.47
2:D:214:LEU:HD22	2:D:253:LEU:HG	1.95	0.47
2:D:289:LYS:HD3	2:D:293:ARG:HE	1.80	0.47
1:A:15:TYR:CG	1:A:47:THR:HG21	2.50	0.47
1:C:193:PHE:CD2	1:C:266:MET:HE1	2.50	0.47
1:C:101:LEU:N	1:C:102:PRO:HD2	2.31	0.46
1:C:84:HIS:O	3:C:1297:C85:H15	2.15	0.46
1:A:115:LEU:HA	1:A:115:LEU:HD12	1.76	0.46
1:C:145:ASP:HA	4:C:2015:HOH:O	2.14	0.46
2:D:395:HIS:HE1	2:D:427:PRO:O	1.99	0.46
1:A:143:LEU:HD23	1:A:143:LEU:H	1.81	0.46
1:C:78:LEU:HB2	1:C:80:PHE:CE1	2.51	0.46
1:C:156:VAL:HG22	4:C:2033:HOH:O	2.15	0.46
1:A:209:ILE:HD11	1:A:240:PHE:HE1	1.81	0.46
1:A:74:ASN:ND2	1:A:74:ASN:N	2.64	0.45
1:A:155:PRO:HD3	4:B:2032:HOH:O	2.16	0.45
1:C:223:ASP:OD1	1:C:225:VAL:HB	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:289:LYS:HD3	2:B:293:ARG:HH21	1.81	0.45
1:C:88:LYS:HB2	1:C:130:PRO:HB2	1.97	0.45
1:C:131:GLN:HG3	4:C:2025:HOH:O	2.16	0.45
2:D:179:HIS:CE1	2:D:379:LYS:HZ1	2.34	0.45
2:B:281:ILE:HG13	2:B:281:ILE:H	1.61	0.45
2:D:372:TRP:HA	2:D:373:PRO:HD3	1.66	0.44
1:A:85:GLN:O	1:A:135:ILE:HG23	2.18	0.44
2:D:224:GLU:HG3	2:D:224:GLU:O	2.17	0.44
2:B:347:TYR:OH	2:B:394:LEU:HA	2.17	0.44
1:C:55:LEU:HD12	1:C:55:LEU:HA	1.76	0.44
1:A:212:LEU:HA	1:A:212:LEU:HD23	1.63	0.44
1:C:145:ASP:HB2	3:C:1297:C85:N3	2.33	0.44
2:B:288:LYS:HD2	4:B:2023:HOH:O	2.17	0.43
1:C:255:LEU:HG	1:C:259:GLY:HA3	2.00	0.43
1:A:227:TRP:CE3	1:A:230:VAL:HG13	2.54	0.43
2:D:242:PHE:CZ	2:D:246:MET:HG3	2.53	0.43
2:B:179:HIS:CE1	2:B:379:LYS:HZ3	2.35	0.43
1:A:85:GLN:NE2	1:A:89:LYS:HB3	2.32	0.43
1:A:88:LYS:HA	1:A:91:MET:HE2	2.00	0.43
2:D:323:GLN:HA	4:D:2036:HOH:O	2.18	0.43
2:B:317:GLN:HE21	2:B:379:LYS:HZ3	1.65	0.43
1:A:84:HIS:HD2	1:A:296:LEU:HD22	1.83	0.43
2:D:345:ASP:HA	2:D:346:PRO:HA	1.80	0.43
1:C:135:ILE:HB	4:C:2027:HOH:O	2.18	0.42
2:D:314:PHE:O	2:D:317:GLN:HB3	2.19	0.42
1:A:85:GLN:HA	3:A:1297:C85:C15	2.49	0.42
2:B:315:LEU:HD23	2:B:315:LEU:HA	1.91	0.42
1:A:73:GLU:OE1	1:A:74:ASN:ND2	2.52	0.42
1:C:230:VAL:HA	1:C:233:MET:HG3	2.01	0.42
1:A:209:ILE:HD11	1:A:240:PHE:CE1	2.54	0.42
2:B:230:GLU:OE2	2:B:313:GLN:NE2	2.53	0.42
1:C:112:LEU:HA	1:C:112:LEU:HD23	1.80	0.42
1:C:174:LEU:HA	4:C:2037:HOH:O	2.18	0.42
1:C:262:LEU:HG	1:C:266:MET:HE3	2.00	0.42
1:C:85:GLN:HB2	3:C:1297:C85:O3	2.19	0.42
2:D:196:LYS:HE2	2:D:196:LYS:HB3	1.90	0.42
1:C:83:LEU:O	3:C:1297:C85:H14	2.20	0.42
2:B:175:VAL:HA	2:B:176:PRO:HD3	1.88	0.41
1:A:111:LEU:HD13	1:A:189:LEU:HD21	2.01	0.41
1:A:283:HIS:CG	1:A:284:PRO:HD2	2.55	0.41
2:B:273:PRO:HG2	2:B:278:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:GLY:HA2	1:C:199:ARG:NE	2.36	0.41
1:A:73:GLU:H	1:A:73:GLU:CD	2.23	0.41
2:B:395:HIS:HE1	2:B:427:PRO:O	2.04	0.41
2:D:348:LEU:HA	2:D:348:LEU:HD12	1.86	0.41
1:C:78:LEU:HD23	1:C:78:LEU:N	2.36	0.41
1:A:18:VAL:HA	1:A:32:LEU:O	2.20	0.41
2:D:332:LEU:O	2:D:335:PHE:HB3	2.21	0.41
2:B:233:HIS:HE1	4:B:2034:HOH:O	2.04	0.41
2:D:193:CYS:O	2:D:241:ARG:HD2	2.20	0.41
2:B:194:LYS:HE3	2:B:351:LEU:HD23	2.02	0.40
2:D:199:TYR:CD1	2:D:199:TYR:C	2.94	0.40
2:B:348:LEU:HD12	2:B:348:LEU:HA	1.88	0.40
1:C:122:ARG:HA	1:C:152:PHE:CE1	2.57	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	293/298 (98%)	276 (94%)	12 (4%)	5 (2%)	11	26
1	C	293/298 (98%)	273 (93%)	15 (5%)	5 (2%)	11	26
2	B	256/258 (99%)	248 (97%)	6 (2%)	2 (1%)	22	48
2	D	256/258 (99%)	243 (95%)	11 (4%)	2 (1%)	22	48
All	All	1098/1112 (99%)	1040 (95%)	44 (4%)	14 (1%)	14	33

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	GLY
1	A	38	ASP

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Mol	Chain	Res	Type
1	A	164	VAL
1	C	72	THR
1	C	145	ASP
1	C	164	VAL
1	A	39	THR
1	A	72	THR
2	D	415	ASN
2	B	365	TYR
2	B	366	THR
1	C	38	ASP
2	D	372	TRP
1	C	13	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	260/262 (99%)	234 (90%)	26 (10%)	9	20
1	C	260/262 (99%)	237 (91%)	23 (9%)	12	26
2	B	232/232 (100%)	213 (92%)	19 (8%)	13	30
2	D	232/232 (100%)	222 (96%)	10 (4%)	33	62
All	All	984/988 (100%)	906 (92%)	78 (8%)	14	32

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	8	GLU
1	A	15	TYR
1	A	47	THR
1	A	55	LEU
1	A	57	GLU
1	A	69	VAL
1	A	72	THR
1	A	74	ASN

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Mol	Chain	Res	Type
1	A	97	THR
1	A	122	ARG
1	A	135	ILE
1	A	138	GLU
1	A	143	LEU
1	A	148	LEU
1	A	157	ARG
1	A	163	VAL
1	A	166	LEU
1	A	199	ARG
1	A	230	VAL
1	A	242	LYS
1	A	255	LEU
1	A	257	GLU
1	A	268	HIS
1	A	287	GLN
1	A	291	LYS
2	B	180	GLU
2	B	189	MET
2	B	197	VAL
2	B	199	TYR
2	B	232	LEU
2	B	247	SER
2	B	253	LEU
2	B	281	ILE
2	B	289	LYS
2	B	292	LEU
2	B	322	GLN
2	B	328	LYS
2	B	348	LEU
2	B	378	ARG
2	B	384	LEU
2	B	391	LEU
2	B	403	GLN
2	B	431	ASN
2	B	432	LEU
1	C	5	GLN
1	C	17	VAL
1	C	55	LEU
1	C	65	LYS
1	C	72	THR
1	C	74	ASN

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Mol	Chain	Res	Type
1	C	88	LYS
1	C	89	LYS
1	C	92	ASP
1	C	122	ARG
1	C	131	GLN
1	C	138	GLU
1	C	150	ARG
1	C	177	CYS
1	C	184	VAL
1	C	189	LEU
1	C	199	ARG
1	C	208	GLU
1	C	226	VAL
1	C	230	VAL
1	C	248	PHE
1	C	249	SER
1	C	260	ARG
2	D	180	GLU
2	D	189	MET
2	D	199	TYR
2	D	224	GLU
2	D	232	LEU
2	D	247	SER
2	D	292	LEU
2	D	348	LEU
2	D	384	LEU
2	D	391	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	74	ASN
1	A	84	HIS
1	A	85	GLN
1	A	113	GLN
1	A	119	HIS
1	A	272	ASN
2	B	179	HIS
2	B	183	HIS
2	B	254	GLN
2	B	296	HIS

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Mol	Chain	Res	Type
2	B	313	GLN
2	B	395	HIS
2	B	431	ASN
1	C	5	GLN
1	C	59	ASN
1	C	60	HIS
1	C	62	ASN
1	C	84	HIS
1	C	85	GLN
1	C	119	HIS
1	C	246	GLN
2	D	179	HIS
2	D	183	HIS
2	D	233	HIS
2	D	296	HIS
2	D	313	GLN
2	D	395	HIS
2	D	396	GLN
2	D	419	HIS
2	D	431	ASN

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	0.95	1 (11%)	10,14,16	1.22	1 (10%)
1	TPO	C	160	1	9,10,11	1.33	1 (11%)	10,14,16	1.20	2 (20%)

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	1/1/3/4	0/8/11/13	0/0/0/0
1	TPO	C	160	1	1/1/3/4	0/8/11/13	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	160	TPO	P-O1P	2.15	1.58	1.50
1	C	160	TPO	P-O1P	2.53	1.59	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	160	TPO	O-C-CA	-2.01	120.46	125.15
1	C	160	TPO	CG2-CB-CA	2.77	118.35	113.22
1	A	160	TPO	CG2-CB-CA	3.03	118.85	113.22

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	160	TPO	CB
1	A	160	TPO	CB

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](#)

2 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	C85	A	1297	-	21,25,25	2.35	6 (28%)	26,37,37	1.93	4 (15%)
3	C85	C	1297	-	21,25,25	2.43	6 (28%)	26,37,37	2.00	5 (19%)

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	C85	A	1297	-	-	0/8/26/26	0/2/3/3
3	C85	C	1297	-	-	0/8/26/26	0/2/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1297	C85	C8-C7	-6.28	1.40	1.48
3	A	1297	C85	C8-C7	-6.18	1.40	1.48
3	C	1297	C85	C9-S1	-5.30	1.68	1.74
3	A	1297	C85	C13-C4	-4.36	1.38	1.46
3	A	1297	C85	C9-S1	-4.22	1.69	1.74
3	C	1297	C85	C13-C4	-3.60	1.40	1.46
3	C	1297	C85	C7-S1	-3.29	1.66	1.73
3	A	1297	C85	C7-S1	-2.24	1.69	1.73
3	C	1297	C85	O3-S2	3.08	1.50	1.43
3	A	1297	C85	O3-S2	3.41	1.50	1.43
3	C	1297	C85	O4-S2	3.56	1.51	1.43
3	A	1297	C85	O4-S2	3.58	1.51	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1297	C85	O3-S2-O4	-5.11	109.32	118.70
3	A	1297	C85	O3-S2-O4	-5.10	109.35	118.70
3	C	1297	C85	C6-C7-S1	-3.69	124.42	129.26
3	A	1297	C85	C6-C7-S1	-3.48	124.71	129.26
3	C	1297	C85	C3-C4-C13	-2.60	126.40	128.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1297	C85	C7-C8-N2	2.08	111.97	110.19
3	C	1297	C85	C3-C2-C1	2.42	108.39	106.30
3	A	1297	C85	C6-C7-C8	5.04	124.33	120.43
3	C	1297	C85	C6-C7-C8	5.15	124.42	120.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1297	C85	1	0
3	C	1297	C85	5	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	295/298 (98%)	-0.32	9 (3%) 49 50	7, 19, 42, 59	12 (4%)
1	C	295/298 (98%)	-0.30	12 (4%) 38 37	6, 18, 48, 63	8 (2%)
2	B	258/258 (100%)	-0.52	2 (0%) 86 87	6, 21, 36, 51	2 (0%)
2	D	258/258 (100%)	-0.41	3 (1%) 79 80	6, 22, 37, 50	2 (0%)
All	All	1106/1112 (99%)	-0.38	26 (2%) 59 60	6, 20, 43, 63	24 (2%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	295	HIS	6.6
1	C	39	THR	6.1
1	A	296	LEU	5.4
1	C	38	ASP	4.3
1	C	40	GLU	3.9
1	C	14	THR	3.8
1	C	296	LEU	3.8
1	A	39	THR	3.6
1	C	13	GLY	3.5
1	A	15	TYR	3.4
1	C	41	THR	3.4
1	A	40	GLU	3.1
1	A	38	ASP	3.0
2	B	324	PRO	2.9
1	A	73	GLU	2.9
1	C	12	GLU	2.9
2	D	283	ASP	2.9
2	D	323	GLN	2.8
1	C	179	TYR	2.6
1	C	15	TYR	2.5
2	B	323	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	295	HIS	2.2
1	A	14	THR	2.1
1	C	162	GLU	2.1
2	D	403	GLN	2.0
1	A	96	LEU	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. 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, 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	LLDF	B-factors(Å ²)	Q<0.9
1	TPO	C	160	11/12	0.95	0.14	-	30,31,31,32	0
1	TPO	A	160	11/12	0.98	0.10	-	4,14,19,19	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, 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	LLDF	B-factors(Å ²)	Q<0.9
3	C85	C	1297	23/23	0.75	0.34	6.83	60,61,62,62	23
3	C85	A	1297	23/23	0.78	0.25	1.46	56,58,61,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.