Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 04:23 PM GMT

PDB ID	:	4EOS
Title	:	Thr 160 phosphorylated CDK2 WT - human cyclin A3 complex with the inhibitor RO3306
Authors	:	Echalier, A.; Cot, E.; Camasses, A.; Hodimont, E.; Hoh, F.; Sheinerman, F.; Krasinska, L.; Fisher, D.
Deposited on	:	2012-04-14
Resolution	:	2.57 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

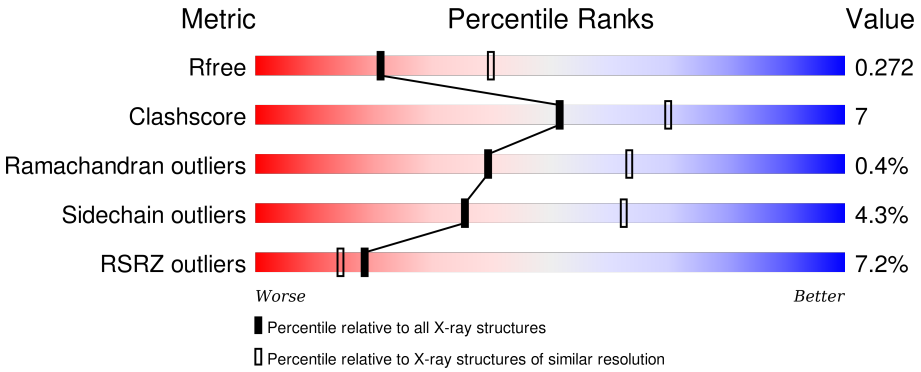
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 2.57 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}	91344	2636 (2.60-2.56)
Clashscore	102246	3003 (2.60-2.56)
Ramachandran outliers	100387	2956 (2.60-2.56)
Sidechain outliers	100360	2956 (2.60-2.56)
RSRZ outliers	91569	2642 (2.60-2.56)

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 <=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	2% 87% 11%
1	C	300	9% 78% 18%
2	B	258	6% 92% 7%
2	D	258	12% 90% 8%

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	1RO	A	301	-	-	X	-
3	1RO	C	301	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9110 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	297	Total	C	N	O	P	S	0	1	0
			2395	1555	410	421	1	8			
1	C	297	Total	C	N	O	P	S	0	2	0
			2406	1560	407	430	1	8			

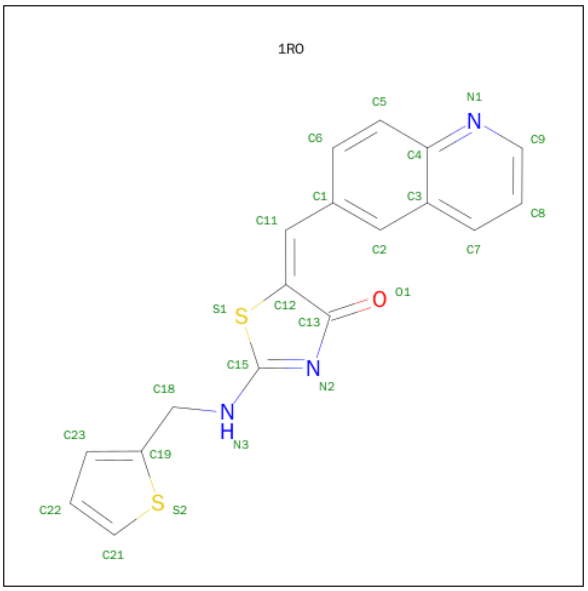
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	PRO	-	EXPRESSION TAG	UNP P24941
A	-1	GLY	-	EXPRESSION TAG	UNP P24941
A	0	SER	-	EXPRESSION TAG	UNP P24941
C	-2	PRO	-	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	257	Total	C	N	O	S	0	0	0
			2076	1345	338	382	11			
2	D	255	Total	C	N	O	S	0	0	0
			2059	1334	335	379	11			

- Molecule 3 is (5E)-5-(QUINOLIN-6-YLMETHYLIDENE)-2-[(THIOPHEN-2-YLMETHYL) AMINO]-1,3-THIAZOL-4(5H)-ONE (three-letter code: 1RO) (formula: $C_{18}H_{13}N_3OS_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			24	18	3	1	2		
3	C	1	Total	C	N	O	S	0	0
			24	18	3	1	2		

- Molecule 4 is water.

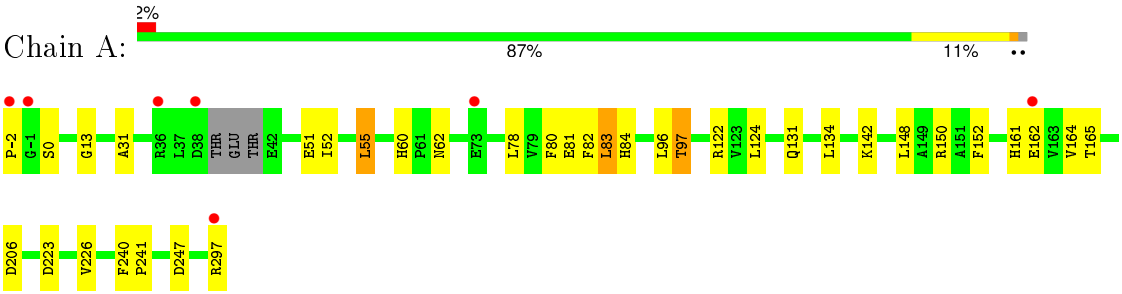
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		
4	B	25	Total	O	0	0
			25	25		
4	C	30	Total	O	0	0
			30	30		
4	D	25	Total	O	0	0
			25	25		

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 errors 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.

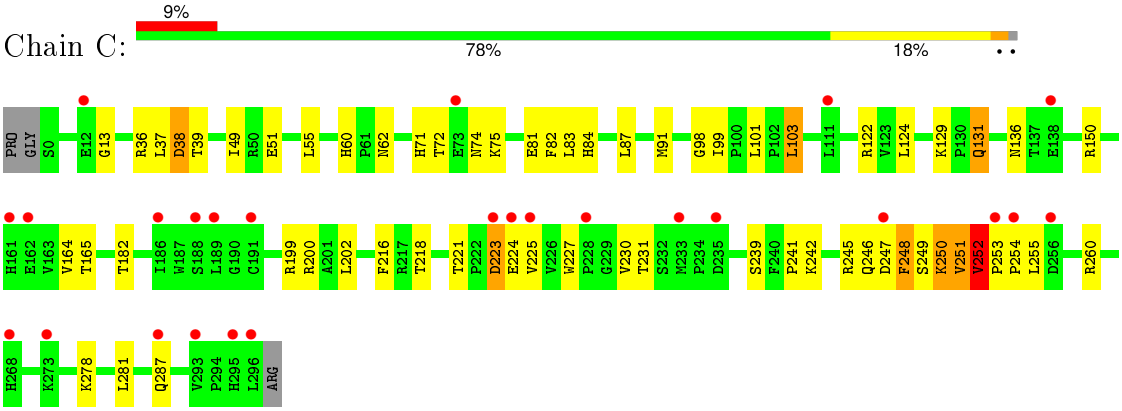
- Molecule 1: Cyclin-dependent kinase 2

Chain A:



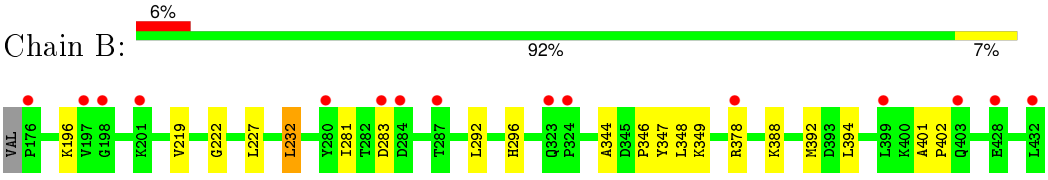
- Molecule 1: Cyclin-dependent kinase 2

Chain C:



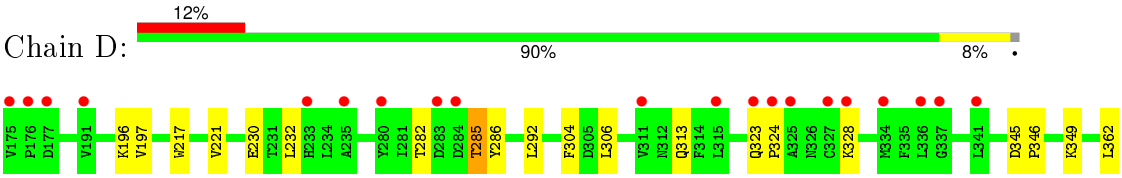
- Molecule 2: Cyclin-A2

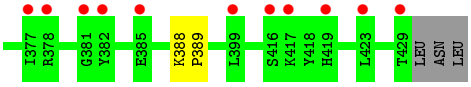
Chain B:



- Molecule 2: Cyclin-A2

Chain D:




I377
R378
G381
Y382
E385
K388
P389
L399
S416
K417
Y418
H419
L423
T429
LEU
ASN
LEU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	74.65Å 133.34Å 149.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.44 – 2.57 30.43 – 2.57	Depositor EDS
% Data completeness (in resolution range)	98.7 (30.44-2.57) 98.7 (30.43-2.57)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.88 (at 2.57Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.230 , 0.274 0.229 , 0.272	Depositor DCC
R_{free} test set	2433 reflections (5.39%)	DCC
Wilson B-factor (Å^2)	49.2	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 45.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Outliers	2 of 47493 reflections (0.004%)	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	9110	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, 1RO

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	0.46	0/2448	0.58	3/3317 (0.1%)
1	C	0.49	0/2456	0.60	3/3332 (0.1%)
2	B	0.32	0/2126	0.47	0/2886
2	D	0.33	0/2109	0.49	0/2864
All	All	0.41	0/9139	0.54	6/12399 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	84	HIS	N-CA-CB	-8.33	95.61	110.60
1	C	251	VAL	N-CA-C	-7.71	90.18	111.00
1	A	83	LEU	N-CA-C	-6.12	94.48	111.00
1	C	251	VAL	CB-CA-C	-5.68	100.60	111.40
1	C	84	HIS	N-CA-CB	-5.33	101.00	110.60
1	A	84	HIS	N-CA-C	5.13	124.84	111.00

There are no chirality outliers.

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	2395	0	2445	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2406	0	2442	74	0
2	B	2076	0	2099	8	0
2	D	2059	0	2079	14	0
3	A	24	0	13	12	0
3	C	24	0	13	9	0
4	A	46	0	0	1	0
4	B	25	0	0	0	0
4	C	30	0	0	0	0
4	D	25	0	0	1	0
All	All	9110	0	9091	123	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 7.

All (123) 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:C:252:VAL:HG13	1:C:253:PRO:CD	1.45	1.46
1:C:249:SER:HA	1:C:251:VAL:O	1.33	1.25
1:C:252:VAL:CG1	1:C:253:PRO:HD2	1.70	1.21
1:C:248:PHE:O	1:C:249:SER:OG	1.67	1.11
1:C:250:LYS:C	1:C:250:LYS:HD2	1.49	1.10
1:C:224:GLU:OE2	1:C:231:THR:OG1	1.72	1.08
1:C:252:VAL:CG1	1:C:253:PRO:CD	2.30	1.06
1:C:250:LYS:HD2	1:C:251:VAL:N	1.70	1.05
1:C:250:LYS:HG3	1:C:251:VAL:H	1.22	1.04
1:C:252:VAL:HG13	1:C:253:PRO:HD2	1.05	1.04
3:C:301:1RO:H13	3:C:301:1RO:O1	1.61	1.01
1:C:253:PRO:HA	1:C:255:LEU:H	1.27	0.99
1:C:250:LYS:CG	1:C:251:VAL:H	1.74	0.96
1:C:129:LYS:HD3	1:C:165:THR:HG21	1.48	0.94
1:C:252:VAL:HG13	1:C:253:PRO:N	1.71	0.94
1:C:250:LYS:C	1:C:250:LYS:CD	2.30	0.94
1:C:250:LYS:CD	1:C:251:VAL:N	2.30	0.93
1:C:252:VAL:CG1	1:C:253:PRO:N	2.29	0.92
1:C:249:SER:N	1:C:250:LYS:HB3	1.86	0.90
1:C:250:LYS:CG	1:C:251:VAL:N	2.30	0.90
1:C:131[A]:GLN:HE21	1:C:131[A]:GLN:N	1.70	0.90
1:C:253:PRO:HA	1:C:255:LEU:N	1.87	0.89
1:C:248:PHE:C	1:C:249:SER:HG	1.72	0.88
1:C:131[A]:GLN:HE21	1:C:131[A]:GLN:H	1.16	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:252:VAL:CB	1:C:253:PRO:HD2	2.03	0.86
1:C:246:GLN:O	1:C:247:ASP:HB2	1.74	0.85
3:A:301:1RO:O1	3:A:301:1RO:C2	2.26	0.83
1:A:82:PHE:CD2	1:A:83:LEU:O	2.32	0.81
3:C:301:1RO:C2	3:C:301:1RO:O1	2.29	0.80
1:A:206:ASP:HB3	4:A:424:HOH:O	1.81	0.79
1:C:218:THR:HG23	1:C:250:LYS:HE2	1.65	0.78
1:C:251:VAL:O	1:C:251:VAL:CG1	2.30	0.78
1:C:252:VAL:CG2	1:C:253:PRO:HD2	2.14	0.78
1:C:13:GLY:HA3	3:C:301:1RO:C18	2.16	0.75
1:C:252:VAL:HG22	1:C:253:PRO:HD2	1.69	0.75
1:C:71:HIS:HE2	2:D:304:PHE:HE2	1.35	0.74
3:A:301:1RO:N2	3:A:301:1RO:C19	2.48	0.74
1:C:131[A]:GLN:NE2	1:C:131[A]:GLN:H	1.86	0.73
3:C:301:1RO:C19	3:C:301:1RO:N2	2.49	0.73
1:C:253:PRO:CA	1:C:255:LEU:H	2.02	0.71
1:C:252:VAL:HG12	1:C:253:PRO:O	1.89	0.71
1:C:249:SER:H	1:C:250:LYS:HB3	1.55	0.70
1:C:216:PHE:HB3	1:C:221:THR:HG22	1.77	0.67
1:C:252:VAL:CB	1:C:253:PRO:CD	2.67	0.66
3:A:301:1RO:O1	3:A:301:1RO:H13	1.94	0.66
1:C:252:VAL:HG13	1:C:253:PRO:CG	2.22	0.66
1:C:13:GLY:HA3	3:C:301:1RO:H3	1.77	0.66
1:C:38:ASP:OD2	1:C:38:ASP:N	2.30	0.65
3:C:301:1RO:C13	3:C:301:1RO:H13	2.27	0.64
2:B:219:VAL:HG22	2:B:232:LEU:HD11	1.80	0.64
1:C:253:PRO:HB2	1:C:254:PRO:HA	1.80	0.63
1:A:52:ILE:HG12	1:A:78:LEU:HD21	1.80	0.63
1:C:87:LEU:O	1:C:91:MET:HG2	2.00	0.61
1:C:82:PHE:CD2	1:C:83:LEU:O	2.53	0.61
1:C:13:GLY:HA3	3:C:301:1RO:H2	1.82	0.61
1:C:251:VAL:O	1:C:251:VAL:HG12	1.99	0.60
1:C:227:TRP:O	1:C:230:VAL:HG23	2.01	0.60
1:C:37:LEU:HB2	1:C:74:ASN:O	2.01	0.60
1:C:251:VAL:O	1:C:251:VAL:HG13	2.01	0.60
1:C:218:THR:CG2	1:C:250:LYS:HE2	2.32	0.58
1:A:60:HIS:CD2	1:A:62:ASN:H	2.22	0.57
1:C:252:VAL:HG22	1:C:253:PRO:CD	2.35	0.57
1:A:51:GLU:O	1:A:55:LEU:HB2	2.05	0.55
1:C:51:GLU:O	1:C:55:LEU:HB2	2.06	0.55
1:C:221:THR:HG21	1:C:241:PRO:O	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:GLY:HA2	3:A:301:1RO:H2	1.89	0.54
2:D:346:PRO:HB2	2:D:349:LYS:HE2	1.89	0.54
1:A:97:THR:HG23	1:A:97:THR:O	2.07	0.54
2:D:388:LYS:HB3	2:D:389:PRO:HD3	1.91	0.53
1:A:82:PHE:CE2	1:A:83:LEU:O	2.61	0.52
1:C:72:THR:HB	1:C:75:LYS:H	1.74	0.52
1:C:245:ARG:HH12	1:C:248:PHE:HD2	1.56	0.52
1:C:246:GLN:O	1:C:247:ASP:CB	2.46	0.52
2:D:282:THR:HG21	2:D:286:TYR:CD2	2.45	0.52
1:C:250:LYS:O	1:C:251:VAL:C	2.44	0.51
1:C:250:LYS:HG3	1:C:251:VAL:N	2.00	0.51
3:A:301:1RO:C13	3:A:301:1RO:C2	2.90	0.49
2:D:362:LEU:HB2	4:D:514:HOH:O	2.12	0.49
1:A:60:HIS:HD2	1:A:62:ASN:H	1.61	0.49
2:B:347:TYR:OH	2:B:394:LEU:HA	2.12	0.49
3:A:301:1RO:C13	3:A:301:1RO:H13	2.44	0.48
1:C:249:SER:OG	1:C:249:SER:O	2.30	0.48
1:C:49:ILE:HG23	2:D:306:LEU:HD12	1.94	0.48
1:C:36:ARG:HG3	1:C:37:LEU:N	2.29	0.48
1:C:252:VAL:CG1	1:C:253:PRO:O	2.58	0.47
1:A:131:GLN:O	3:A:301:1RO:H4	2.15	0.47
1:C:83:LEU:HD12	1:C:136:ASN:HB3	1.96	0.47
3:C:301:1RO:N2	3:C:301:1RO:C23	2.77	0.47
2:B:401:ALA:HB3	2:B:402:PRO:HD3	1.97	0.46
1:C:253:PRO:CA	1:C:255:LEU:N	2.66	0.46
1:A:83:LEU:O	3:A:301:1RO:H10	2.16	0.45
1:A:134:LEU:HD21	3:A:301:1RO:O1	2.16	0.45
1:A:31:ALA:HB3	1:A:80:PHE:HB2	1.98	0.45
2:B:388:LYS:O	2:B:392:MET:HG2	2.17	0.45
2:B:344:ALA:HB1	2:B:348:LEU:HD22	1.99	0.45
2:D:323:GLN:HA	2:D:324:PRO:HA	1.80	0.44
2:D:282:THR:O	2:D:285:THR:OG1	2.33	0.44
1:A:223:ASP:OD1	1:A:226:VAL:HG12	2.17	0.44
1:A:13:GLY:CA	3:A:301:1RO:H2	2.47	0.44
2:D:345:ASP:HA	2:D:346:PRO:HA	1.77	0.44
3:A:301:1RO:N2	3:A:301:1RO:C23	2.81	0.43
1:C:71:HIS:NE2	2:D:304:PHE:HE2	2.08	0.43
2:D:230:GLU:OE2	2:D:313:GLN:HG3	2.19	0.43
1:C:60:HIS:CD2	1:C:62:ASN:H	2.37	0.43
1:C:98:GLY:HA2	1:C:199:ARG:HD3	2.01	0.43
1:A:124:LEU:HG	1:A:152:PHE:CD1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:PHE:HA	1:A:241:PRO:HD3	1.94	0.42
2:D:217:TRP:O	2:D:221:VAL:HG23	2.20	0.42
2:D:282:THR:HG22	2:D:285:THR:OG1	2.19	0.42
1:A:83:LEU:HD21	1:A:142:LYS:HD2	2.00	0.42
1:C:81:GLU:O	3:C:301:1RO:H9	2.19	0.42
2:B:346:PRO:O	2:B:349:LYS:HG2	2.19	0.41
1:C:249:SER:HB3	1:C:260:ARG:HD3	2.03	0.41
1:C:124:LEU:HD21	1:C:182:THR:HA	2.02	0.41
1:A:81:GLU:O	3:A:301:1RO:H9	2.21	0.41
2:D:282:THR:HG21	2:D:286:TYR:HD2	1.86	0.41
2:B:222:GLY:HA2	2:B:227:LEU:HD12	2.02	0.40
1:C:99:ILE:HG23	1:C:103:LEU:HD12	2.03	0.40
1:C:223:ASP:C	1:C:225:VAL:H	2.25	0.40
1:C:223:ASP:O	1:C:225:VAL:N	2.54	0.40
1:C:253:PRO:HB2	1:C:254:PRO:CA	2.50	0.40
1:C:250:LYS:C	1:C:251:VAL:O	2.51	0.40
2:B:281:ILE:H	2:B:281:ILE:HG13	1.69	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	293/300 (98%)	285 (97%)	7 (2%)	1 (0%)	46 70
1	C	296/300 (99%)	276 (93%)	18 (6%)	2 (1%)	26 49
2	B	255/258 (99%)	250 (98%)	5 (2%)	0	100 100
2	D	253/258 (98%)	248 (98%)	4 (2%)	1 (0%)	39 63
All	All	1097/1116 (98%)	1059 (96%)	34 (3%)	4 (0%)	39 63

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	VAL
1	C	164	VAL
1	C	252	VAL
2	D	197	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	261/263 (99%)	248 (95%)	13 (5%)	30	55
1	C	263/263 (100%)	244 (93%)	19 (7%)	18	34
2	B	231/232 (100%)	225 (97%)	6 (3%)	54	78
2	D	229/232 (99%)	224 (98%)	5 (2%)	60	82
All	All	984/990 (99%)	941 (96%)	43 (4%)	35	61

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-2	PRO
1	A	0	SER
1	A	55	LEU
1	A	96	LEU
1	A	97	THR
1	A	122	ARG
1	A	148	LEU
1	A	150	ARG
1	A	161	HIS
1	A	162	GLU
1	A	165	THR
1	A	247	ASP
1	A	297	ARG
2	B	196	LYS
2	B	232	LEU
2	B	283	ASP
2	B	292	LEU
2	B	296	HIS

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Mol	Chain	Res	Type
2	B	378	ARG
1	C	38	ASP
1	C	39	THR
1	C	101	LEU
1	C	103	LEU
1	C	122	ARG
1	C	131[A]	GLN
1	C	131[B]	GLN
1	C	150	ARG
1	C	200	ARG
1	C	202	LEU
1	C	223	ASP
1	C	239	SER
1	C	242	LYS
1	C	248	PHE
1	C	250	LYS
1	C	252	VAL
1	C	278	LYS
1	C	281	LEU
1	C	287	GLN
2	D	196	LYS
2	D	232	LEU
2	D	285	THR
2	D	292	LEU
2	D	328	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	161	HIS
1	A	246	GLN
1	A	265	GLN
2	B	233	HIS
2	B	254	GLN
2	B	425	ASN
1	C	59	ASN
1	C	60	HIS
1	C	84	HIS
1	C	85	GLN
1	C	246	GLN
1	C	265	GLN

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Mol	Chain	Res	Type
1	C	268	HIS
2	D	317	GLN
2	D	406	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)

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	2.21	4 (50%)	7,14,16	1.84	1 (14%)
1	TPO	C	160	1	8,10,11	0.63	0	7,14,16	1.27	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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	160	TPO	P-O2P	-3.31	1.42	1.54
1	A	160	TPO	P-O1P	-3.23	1.40	1.51
1	A	160	TPO	P-O3P	-2.97	1.44	1.54
1	A	160	TPO	P-OG1	-2.29	1.53	1.60

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	160	TPO	OG1-P-O1P	-2.93	99.78	107.11

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)

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	1RO	A	301	-	25,27,27	2.83	5 (20%)	26,37,37	2.78	7 (26%)
3	1RO	C	301	-	25,27,27	2.78	4 (16%)	26,37,37	2.59	9 (34%)

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	1RO	A	301	-	-	0/5/21/21	0/4/4/4
3	1RO	C	301	-	-	0/5/21/21	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	301	1RO	C13-N2	-2.01	1.33	1.37
3	A	301	1RO	C5-C6	2.01	1.40	1.36
3	A	301	1RO	C15-N2	2.28	1.34	1.31
3	C	301	1RO	O1-C13	2.32	1.28	1.24
3	C	301	1RO	C15-N2	2.64	1.35	1.31
3	A	301	1RO	C3-C4	4.59	1.48	1.42
3	C	301	1RO	C3-C4	4.60	1.48	1.42
3	C	301	1RO	C11-C12	11.69	1.49	1.34
3	A	301	1RO	C11-C12	12.06	1.50	1.34

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	301	1RO	C1-C11-C12	-10.66	116.25	130.96
3	C	301	1RO	C1-C11-C12	-7.70	120.33	130.96
3	C	301	1RO	C22-C21-S2	-7.19	105.28	113.23
3	A	301	1RO	C22-C21-S2	-6.38	106.17	113.23
3	C	301	1RO	C11-C12-S1	-4.09	124.76	129.31
3	C	301	1RO	O1-C13-N2	-2.71	117.73	124.40
3	A	301	1RO	O1-C13-N2	-2.66	117.84	124.40
3	C	301	1RO	C18-N3-C15	-2.46	118.81	122.81
3	A	301	1RO	C3-C4-N1	-2.35	118.73	122.08
3	C	301	1RO	C3-C4-N1	-2.18	118.97	122.08
3	C	301	1RO	C5-C4-N1	2.33	122.39	118.52
3	C	301	1RO	C9-N1-C4	2.42	120.97	116.87
3	A	301	1RO	C9-N1-C4	2.47	121.05	116.87
3	A	301	1RO	C5-C4-N1	2.50	122.68	118.52
3	A	301	1RO	C18-C19-C23	2.64	134.15	128.63
3	C	301	1RO	C18-C19-C23	2.76	134.40	128.63

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	301	1RO	12	0
3	C	301	1RO	9	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{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	296/300 (98%)	0.06	7 (2%) 62 58	28, 43, 68, 86	0
1	C	296/300 (98%)	0.47	26 (8%) 12 9	35, 51, 78, 86	0
2	B	257/258 (99%)	0.33	15 (5%) 26 22	33, 51, 72, 93	0
2	D	255/258 (98%)	0.61	31 (12%) 5 4	33, 57, 94, 115	0
All	All	1104/1116 (98%)	0.36	79 (7%) 18 14	28, 50, 81, 115	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	254	PRO	7.2
2	D	323	GLN	5.0
2	D	175	VAL	4.9
2	B	432	LEU	4.6
2	D	325	ALA	4.6
2	D	284	ASP	4.4
2	B	284	ASP	4.3
2	D	416	SER	4.3
1	C	295	HIS	4.3
2	D	176	PRO	4.1
2	B	283	ASP	3.9
2	D	378	ARG	3.8
1	C	189	LEU	3.8
1	C	247	ASP	3.7
1	A	-2	PRO	3.7
1	C	162	GLU	3.6
2	B	324	PRO	3.6
1	A	38	ASP	3.5
1	A	73	GLU	3.4
2	D	324	PRO	3.4
1	C	225	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
2	D	399	LEU	3.1
2	B	323	GLN	3.1
1	A	36	ARG	3.1
1	C	12	GLU	3.0
2	B	201	LYS	3.0
1	C	233	MET	2.9
2	D	334	MET	2.8
2	D	336	LEU	2.8
2	D	280	TYR	2.8
1	C	111	LEU	2.8
2	D	381	GLY	2.8
1	A	-1	GLY	2.8
2	D	429	THR	2.7
1	C	228	PRO	2.7
2	D	235	ALA	2.7
2	B	176	PRO	2.7
1	C	256	ASP	2.7
1	C	293	VAL	2.7
2	D	419	HIS	2.6
2	B	280	TYR	2.6
2	D	311	VAL	2.6
2	D	328	LYS	2.6
2	D	191	VAL	2.6
2	B	197	VAL	2.6
1	C	161	HIS	2.5
2	B	198	GLY	2.5
1	C	186	ILE	2.4
1	C	223	ASP	2.4
2	D	341	LEU	2.4
2	D	327	CYS	2.4
2	D	423	LEU	2.4
1	C	138[A]	GLU	2.4
2	D	315	LEU	2.3
1	C	273	LYS	2.3
1	C	191	CYS	2.3
2	D	385	GLU	2.3
1	C	253	PRO	2.2
2	D	417	LYS	2.2
2	B	403	GLN	2.2
2	D	377	ILE	2.2
2	D	337	GLY	2.2
1	A	162	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	188	SER	2.2
1	C	296	LEU	2.2
1	C	268	HIS	2.2
2	D	283	ASP	2.2
2	B	428	GLU	2.1
2	D	177	ASP	2.1
2	B	287	THR	2.1
2	B	378	ARG	2.1
1	C	235	ASP	2.1
2	D	233	HIS	2.1
2	B	399	LEU	2.0
1	C	287	GLN	2.0
1	C	224	GLU	2.0
1	C	73	GLU	2.0
1	A	297	ARG	2.0
2	D	382	TYR	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, 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	LLDF	B-factors(Å^2)	Q<0.9
1	TPO	A	160	11/12	0.98	0.10	-	34,38,40,40	0
1	TPO	C	160	11/12	0.97	0.12	-	41,45,46,47	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^{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	LLDF	B-factors(Å^2)	Q<0.9
3	1RO	A	301	24/24	0.94	0.15	-0.06	45,46,48,49	0
3	1RO	C	301	24/24	0.91	0.15	-0.47	46,48,49,50	0

6.5 Other polymers (i)

There are no such residues in this entry.