Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:07 AM GMT

PDB ID	:	2W27
Title	:	CRYSTAL STRUCTURE OF THE BACILLUS SUBTILIS YKUI PROTEIN, WITH AN EAL DOMAIN, IN COMPLEX WITH SUBSTRATE C-DI-GMP AND CALCIUM
Authors	:	Padavattan, S.; Anderson, W.F.; Schirmer, T.
Deposited on	:	2008-10-24
Resolution	:	2.80 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:

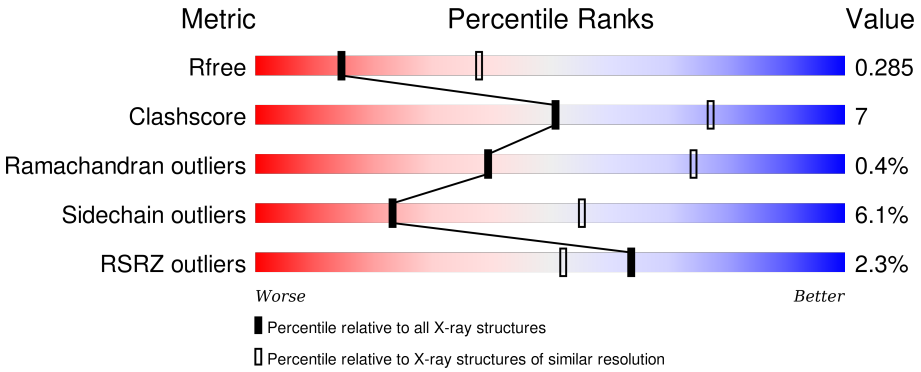
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	431	2% 73% 18% • 7%
1	B	431	2% 71% 18% • 8%

2 Entry composition

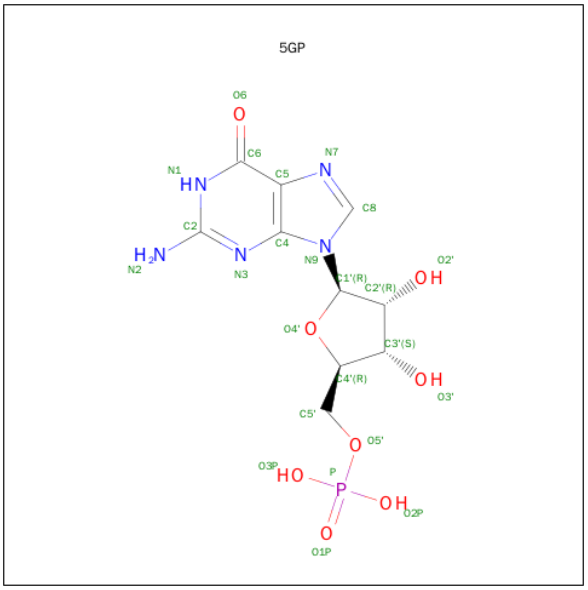
There are 4 unique types of molecules in this entry. The entry contains 6735 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 YKUI PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	54	0	0
			3342	2152	537	642	11			
1	B	395	Total	C	N	O	S	33	0	0
			3292	2120	531	629	12			

- Molecule 2 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: 5GP) (formula: $C_{10}H_{14}N_5O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Ca 1 1	0	0
3	A	1	Total Ca 1 1	0	0

- Molecule 4 is water.

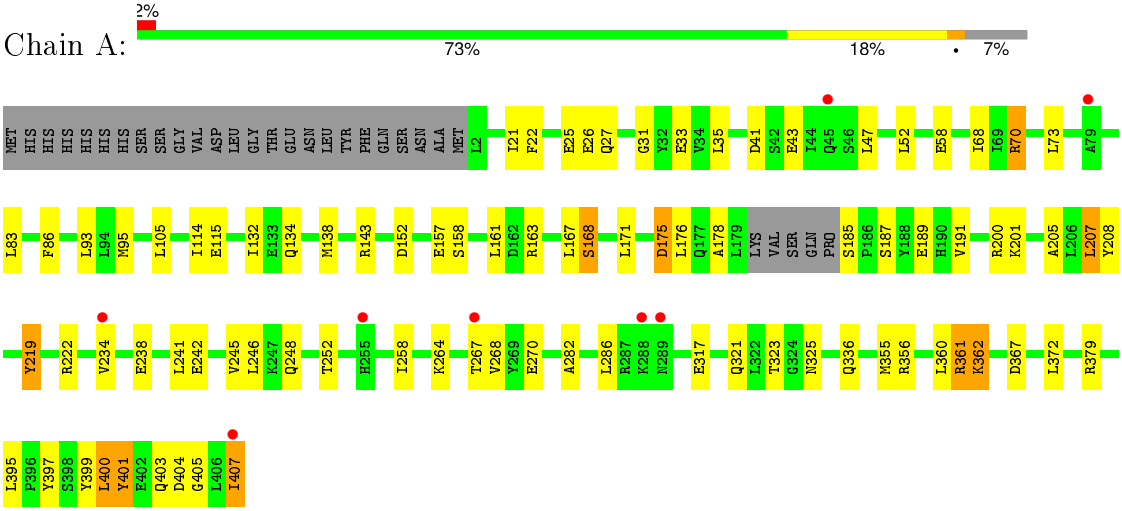
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total O 4 4	0	0
4	B	3	Total O 3 3	0	0

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.

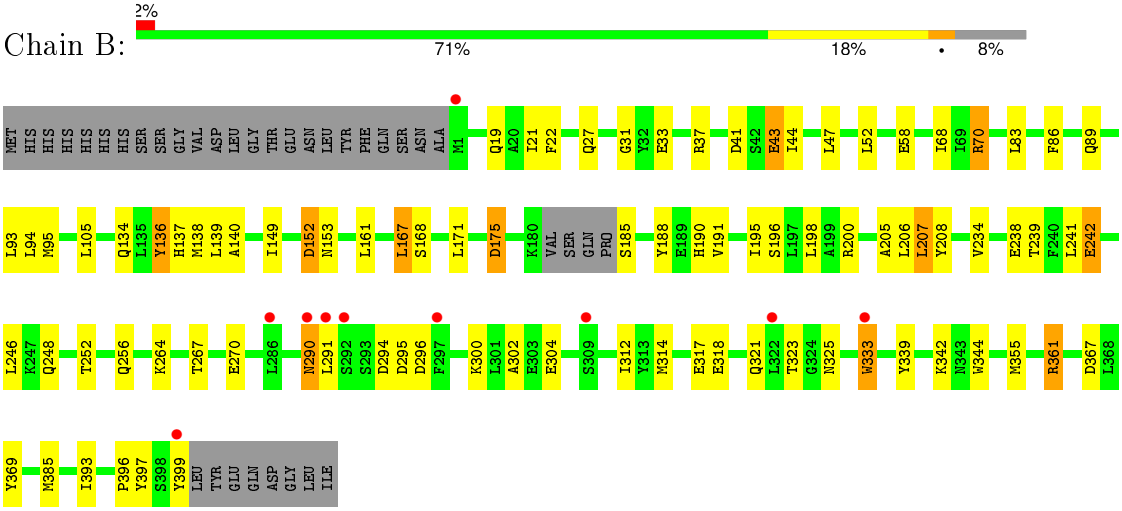
- Molecule 1: YKUI PROTEIN

Chain A:



- Molecule 1: YKUI PROTEIN

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	46.17Å 124.52Å 168.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 38.51 – 2.80	Depositor EDS
% Data completeness (in resolution range)	86.3 (30.00-2.80) 86.3 (38.51-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.57 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.226 , 0.275 0.234 , 0.285	Depositor DCC
R_{free} test set	1046 reflections (5.14%)	DCC
Wilson B-factor (Å^2)	68.8	Xtriage
Anisotropy	0.352	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 67.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 21414 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6735	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% 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: CA, 5GP

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	2.24	8/3418 (0.2%)	2.14	15/4617 (0.3%)
1	B	0.74	4/3367 (0.1%)	0.66	6/4547 (0.1%)
All	All	1.67	12/6785 (0.2%)	1.59	21/9164 (0.2%)

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)
1	A	219	TYR	CG-CD2	113.26	2.86	1.39
1	A	270	GLU	CG-CD	25.36	1.90	1.51
1	A	219	TYR	CG-CD1	25.25	1.72	1.39
1	A	401	TYR	CG-CD1	-24.18	1.07	1.39
1	B	242	GLU	CG-CD	23.93	1.87	1.51
1	A	356	ARG	CD-NE	-22.39	1.08	1.46
1	A	401	TYR	CG-CD2	19.86	1.65	1.39
1	A	362	LYS	CB-CG	-15.74	1.10	1.52
1	A	222	ARG	CD-NE	-14.35	1.22	1.46
1	B	136	TYR	CB-CG	13.09	1.71	1.51
1	B	361	ARG	CD-NE	11.73	1.66	1.46
1	B	43	GLU	CB-CG	8.32	1.68	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	219	TYR	CB-CG-CD2	-103.03	59.18	121.00
1	A	219	TYR	CD1-CG-CD2	-54.02	58.47	117.90
1	A	219	TYR	CG-CD1-CE1	46.34	158.37	121.30
1	A	401	TYR	CB-CG-CD2	-38.54	97.87	121.00
1	A	401	TYR	CB-CG-CD1	28.06	137.84	121.00
1	A	219	TYR	CB-CG-CD1	-19.29	109.42	121.00
1	A	356	ARG	CD-NE-CZ	-16.35	100.70	123.60
1	A	361	ARG	CA-CB-CG	-13.44	83.84	113.40
1	A	362	LYS	CA-CB-CG	12.43	140.74	113.40
1	A	401	TYR	CG-CD2-CE2	-11.51	112.09	121.30
1	B	242	GLU	CB-CG-CD	-8.50	91.25	114.20
1	A	270	GLU	CB-CG-CD	-7.28	94.55	114.20
1	B	242	GLU	CG-CD-OE1	6.78	131.85	118.30
1	B	242	GLU	CG-CD-OE2	-6.43	105.45	118.30
1	B	361	ARG	CD-NE-CZ	-6.36	114.70	123.60
1	A	361	ARG	NE-CZ-NH2	-6.18	117.21	120.30
1	B	304	GLU	CB-CA-C	5.49	121.37	110.40
1	A	219	TYR	CG-CD2-CE2	5.26	125.51	121.30
1	A	401	TYR	CG-CD1-CE1	5.12	125.39	121.30
1	A	361	ARG	CG-CD-NE	-5.10	101.09	111.80
1	B	136	TYR	CA-CB-CG	-5.04	103.82	113.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	219	TYR	Sidechain

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	3342	0	3231	49	1
1	B	3292	0	3193	51	1
2	A	46	0	22	1	0
2	B	46	0	22	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	4	0	0	0	0
4	B	3	0	0	0	0
All	All	6735	0	6468	90	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 (90) 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:B:291:LEU:HD11	1:B:300:LYS:HE2	1.56	0.88
1:A:161:LEU:HD12	1:B:195:ILE:CD1	2.12	0.79
1:B:234:VAL:HG11	1:B:241:LEU:HD11	1.67	0.75
1:A:399:TYR:O	1:A:403:GLN:HG2	1.90	0.71
1:A:234:VAL:HG11	1:A:241:LEU:HD11	1.73	0.70
1:A:161:LEU:HD12	1:B:195:ILE:HD13	1.74	0.67
1:A:161:LEU:HD12	1:B:195:ILE:HD11	1.78	0.64
1:B:302:ALA:HB1	1:B:333:TRP:CE2	2.32	0.64
1:B:19:GLN:OE1	2:B:501:5GP:H8	1.98	0.63
1:B:264:LYS:O	1:B:267:THR:HG22	2.00	0.61
1:B:152:ASP:HB3	1:B:153:ASN:HD22	1.65	0.61
1:B:248:GLN:O	1:B:252:THR:HG23	2.02	0.60
1:B:321:GLN:HG2	1:B:323:THR:O	2.02	0.60
1:A:248:GLN:O	1:A:252:THR:HG23	2.04	0.58
1:B:70:ARG:HH21	1:B:105:LEU:HD13	1.69	0.57
1:B:291:LEU:HD22	1:B:296:ASP:HB3	1.87	0.57
1:A:207:LEU:HD12	1:A:208:TYR:N	2.20	0.56
1:B:302:ALA:HB1	1:B:333:TRP:CZ2	2.41	0.56
1:A:200:ARG:NH2	1:B:367:ASP:OD1	2.38	0.56
1:A:321:GLN:HG2	1:A:323:THR:O	2.06	0.56
1:A:264:LYS:O	1:A:267:THR:HG22	2.07	0.55
1:B:207:LEU:HD12	1:B:208:TYR:N	2.20	0.55
1:B:21:ILE:HD12	1:B:86:PHE:CE1	2.42	0.54
1:A:397:TYR:CE2	1:A:407:ILE:HD13	2.43	0.53
1:A:258:ILE:HD13	1:B:318:GLU:O	2.08	0.52
1:A:70:ARG:HH21	1:A:105:LEU:HD13	1.75	0.52
1:B:22:PHE:CE2	1:B:246:LEU:HD12	2.46	0.51
1:B:152:ASP:HB3	1:B:153:ASN:ND2	2.26	0.51
1:B:207:LEU:C	1:B:207:LEU:HD12	2.31	0.51
1:B:175:ASP:C	1:B:175:ASP:OD1	2.48	0.51
1:A:201:LYS:HB3	1:B:190:HIS:CD2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:GLU:HA	1:A:86:PHE:HB2	1.94	0.50
1:B:137:HIS:O	1:B:140:ALA:HB3	2.12	0.50
1:B:33:GLU:HA	1:B:86:PHE:HB2	1.94	0.50
1:A:245:VAL:HG12	1:A:246:LEU:HG	1.94	0.50
1:A:171:LEU:HD23	1:A:205:ALA:O	2.12	0.49
1:B:161:LEU:HD22	1:B:198:LEU:CD2	2.43	0.49
1:A:22:PHE:CE2	1:A:246:LEU:HD12	2.47	0.49
1:A:21:ILE:HD12	1:A:86:PHE:CE1	2.47	0.49
1:A:207:LEU:HD12	1:A:207:LEU:C	2.33	0.49
1:A:367:ASP:OD1	1:B:200:ARG:NH2	2.45	0.49
1:B:314:MET:HG2	1:B:385:MET:HE1	1.95	0.48
1:A:95:MET:HE2	1:A:134:GLN:HG2	1.96	0.48
1:B:31:GLY:HA2	1:B:83:LEU:HD22	1.96	0.47
1:B:70:ARG:NH2	1:B:105:LEU:HD13	2.28	0.47
1:A:176:LEU:HD11	1:A:208:TYR:CD2	2.50	0.47
1:A:372:LEU:H	1:A:372:LEU:HD12	1.80	0.47
1:B:37:ARG:HG3	1:B:44:ILE:HG23	1.95	0.47
1:B:312:ILE:HG23	1:B:393:ILE:HG12	1.96	0.46
1:A:70:ARG:NH2	1:A:105:LEU:HD13	2.31	0.46
1:A:26:GLU:HG3	1:B:342:LYS:HZ2	1.79	0.46
1:B:171:LEU:HD23	1:B:205:ALA:O	2.16	0.45
1:B:234:VAL:HG11	1:B:241:LEU:CD1	2.43	0.45
1:A:395:LEU:N	1:A:395:LEU:HD23	2.32	0.45
1:A:268:VAL:HG22	1:A:405:GLY:O	2.16	0.45
1:A:31:GLY:HA2	1:A:83:LEU:HD22	1.99	0.45
1:A:114:ILE:O	1:A:115:GLU:C	2.55	0.44
1:A:175:ASP:C	1:A:175:ASP:OD1	2.55	0.44
1:B:95:MET:HE2	1:B:134:GLN:HG2	1.99	0.44
1:A:325:ASN:HB2	1:A:336:GLN:HB2	1.98	0.44
1:B:188:TYR:O	1:B:191:VAL:HB	2.17	0.44
1:A:25:GLU:HG2	1:B:344:TRP:CZ3	2.53	0.44
1:B:21:ILE:HD12	1:B:86:PHE:CD1	2.53	0.44
1:A:132:ILE:HD13	1:A:163:ARG:HH11	1.83	0.44
1:A:95:MET:CE	1:A:134:GLN:HG2	2.48	0.43
1:B:58:GLU:HG2	1:B:93:LEU:CD2	2.48	0.43
1:A:35:LEU:HD22	2:A:502:5GP:O1P	2.18	0.43
1:A:187:SER:O	1:A:191:VAL:HG23	2.18	0.43
1:A:379:ARG:HG3	1:A:400:LEU:HD21	2.01	0.42
1:A:58:GLU:HG2	1:A:93:LEU:CD2	2.50	0.42
1:B:238:GLU:HG3	1:B:239:THR:HG23	2.01	0.42
1:A:242:GLU:O	1:A:245:VAL:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ARG:NH2	1:A:168:SER:O	2.52	0.42
1:B:89:GLN:OE1	1:B:94:LEU:HD13	2.20	0.42
1:B:95:MET:CE	1:B:134:GLN:HG2	2.50	0.41
1:A:21:ILE:HD12	1:A:86:PHE:CD1	2.55	0.41
1:B:47:LEU:HD12	1:B:68:ILE:HD11	2.03	0.41
1:B:290:ASN:OD1	1:B:290:ASN:N	2.53	0.41
1:B:196:SER:HA	1:B:206:LEU:HD11	2.02	0.41
1:B:161:LEU:HD22	1:B:198:LEU:HD21	2.03	0.41
1:B:136:TYR:HB2	1:B:167:LEU:CD2	2.51	0.41
1:A:47:LEU:HD12	1:A:68:ILE:HD11	2.01	0.41
1:A:200:ARG:NH1	1:B:369:TYR:CG	2.89	0.41
1:A:264:LYS:O	1:A:268:VAL:HG23	2.21	0.41
1:B:139:LEU:HD22	1:B:149:ILE:HD12	2.01	0.41
1:A:282:ALA:O	1:A:286:LEU:HD12	2.21	0.41
1:A:73:LEU:HD23	1:A:73:LEU:N	2.36	0.41
1:B:325:ASN:ND2	1:B:339:TYR:CD2	2.89	0.41
1:A:185:SER:HB3	1:A:189:GLU:HB3	2.02	0.41
1:A:58:GLU:HG2	1:A:93:LEU:HD21	2.03	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:401:TYR:CE2	1:B:256:GLN:NE2[1_455]	2.12	0.08

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	397/431 (92%)	377 (95%)	19 (5%)	1 (0%)	46 79
1	B	391/431 (91%)	367 (94%)	22 (6%)	2 (0%)	34 69

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	788/862 (91%)	744 (94%)	41 (5%)	3 (0%)	39 74

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	397	TYR
1	A	178	ALA
1	B	396	PRO

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	362/389 (93%)	340 (94%)	22 (6%)	23 55
1	B	357/389 (92%)	335 (94%)	22 (6%)	23 54
All	All	719/778 (92%)	675 (94%)	44 (6%)	23 55

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	41	ASP
1	A	43	GLU
1	A	52	LEU
1	A	70	ARG
1	A	138	MET
1	A	152	ASP
1	A	157	GLU
1	A	158	SER
1	A	167	LEU
1	A	168	SER
1	A	175	ASP
1	A	207	LEU
1	A	238	GLU
1	A	317	GLU

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Mol	Chain	Res	Type
1	A	355	MET
1	A	360	LEU
1	A	361	ARG
1	A	362	LYS
1	A	400	LEU
1	A	404	ASP
1	A	407	ILE
1	B	27	GLN
1	B	41	ASP
1	B	43	GLU
1	B	52	LEU
1	B	70	ARG
1	B	138	MET
1	B	152	ASP
1	B	167	LEU
1	B	168	SER
1	B	175	ASP
1	B	185	SER
1	B	207	LEU
1	B	242	GLU
1	B	270	GLU
1	B	290	ASN
1	B	294	ASP
1	B	295	ASP
1	B	317	GLU
1	B	333	TRP
1	B	355	MET
1	B	361	ARG
1	B	399	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	27	GLN
1	A	117	HIS
1	A	126	HIS
1	A	134	GLN
1	A	214	ASN
1	A	289	ASN
1	B	27	GLN
1	B	134	GLN

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Mol	Chain	Res	Type
1	B	190	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\#\|Z\| > 2$
2	5GP	A	501	3,2	16,25,26	0.70	0	19,37,40	2.09	6 (31%)
2	5GP	A	502	2	16,25,26	0.74	0	19,37,40	1.62	3 (15%)
2	5GP	B	501	3,2	16,25,26	0.71	0	19,37,40	2.06	6 (31%)
2	5GP	B	502	2	16,25,26	0.61	0	19,37,40	1.71	5 (26%)

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
2	5GP	A	501	3,2	-	0/3/25/26	0/3/3/3
2	5GP	A	502	2	-	0/3/25/26	0/3/3/3
2	5GP	B	501	3,2	-	0/3/25/26	0/3/3/3
2	5GP	B	502	2	-	0/3/25/26	0/3/3/3

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	5GP	N3-C2-N1	-3.92	121.47	127.44
2	B	502	5GP	N3-C2-N1	-3.73	121.76	127.44
2	A	501	5GP	C1'-N9-C4	-3.66	121.42	126.94
2	A	501	5GP	C5-C6-N1	-3.66	118.59	123.59
2	B	501	5GP	C5-C6-N1	-3.57	118.70	123.59
2	B	501	5GP	N3-C2-N1	-3.53	122.07	127.44
2	A	502	5GP	N3-C2-N1	-3.51	122.09	127.44
2	B	502	5GP	C4-C5-N7	-3.41	106.35	109.48
2	A	502	5GP	C5-C6-N1	-3.27	119.12	123.59
2	B	501	5GP	C1'-N9-C4	-3.18	122.15	126.94
2	B	501	5GP	C4-C5-N7	-3.09	106.64	109.48
2	B	502	5GP	C5-C6-N1	-2.80	119.77	123.59
2	A	501	5GP	C4-C5-N7	-2.55	107.13	109.48
2	B	502	5GP	O4'-C4'-C3'	2.07	109.31	105.15
2	A	501	5GP	C2'-C1'-N9	2.49	118.09	114.29
2	B	502	5GP	C6-N1-C2	2.89	119.95	115.94
2	B	501	5GP	C2'-C1'-N9	3.08	119.00	114.29
2	A	502	5GP	C6-N1-C2	3.37	120.62	115.94
2	B	501	5GP	C6-N1-C2	3.60	120.94	115.94
2	A	501	5GP	C6-N1-C2	3.65	121.00	115.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	5GP	1	0
2	B	501	5GP	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

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	401/431 (93%)	0.15	8 (1%)	68	58	26, 38, 46, 54	13 (3%)
1	B	395/431 (91%)	0.24	10 (2%)	61	48	32, 38, 45, 77	8 (2%)
All	All	796/862 (92%)	0.20	18 (2%)	64	52	26, 38, 46, 77	21 (2%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	290	ASN	5.5
1	A	79	ALA	4.8
1	B	292	SER	3.4
1	B	322	LEU	3.0
1	B	286	LEU	2.9
1	B	399	TYR	2.8
1	B	309	SER	2.8
1	A	288	LYS	2.8
1	A	407	ILE	2.7
1	B	291	LEU	2.6
1	A	45	GLN	2.6
1	A	289	ASN	2.6
1	A	267	THR	2.5
1	B	1	MET	2.4
1	B	297	PHE	2.3
1	A	234	VAL	2.1
1	B	333	TRP	2.1
1	A	255	HIS	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

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	CA	B	503	1/1	0.88	0.21	0.47	50,50,50,50	0
2	5GP	A	501	23/24	0.92	0.15	-0.84	35,40,41,46	0
2	5GP	B	501	23/24	0.95	0.16	-1.14	35,40,41,45	0
3	CA	A	503	1/1	0.74	0.14	-1.44	50,50,50,50	0
2	5GP	B	502	23/24	0.95	0.15	-1.50	30,40,44,45	0
2	5GP	A	502	23/24	0.96	0.14	-1.74	30,41,44,45	0

6.5 Other polymers (i)

There are no such residues in this entry.