WORLDWIDE PDB
PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:30 PM GMT

PDB ID	:	1XKV
Title	:	Crystal Structure Of ATP-Dependent Phosphoenolpyruvate Carboxykinase From Thermus thermophilus HB8
Authors	:	Sugahara, M.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on	:	2004-09-29
Resolution	:	2.20 Å(reported)

The following versions of software and data (see references ⓘ) 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 2.20 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	Percentile Ranks	Value
Rfree		0.218
Clashscore		7
Ramachandran outliers		0.4%
Sidechain outliers		2.6%
RSRZ outliers		2.8%

Worse .... *Better*

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution
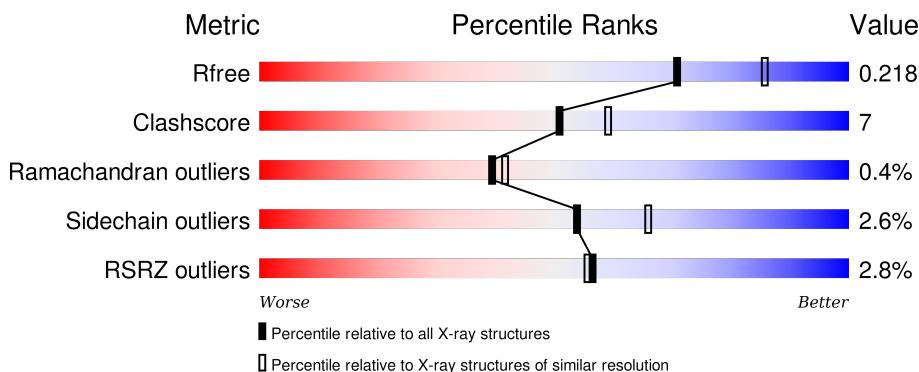
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	529	2% ▮ 84% ... 12% ..
1	B	529	3% ▮ 81% ... 15% ..
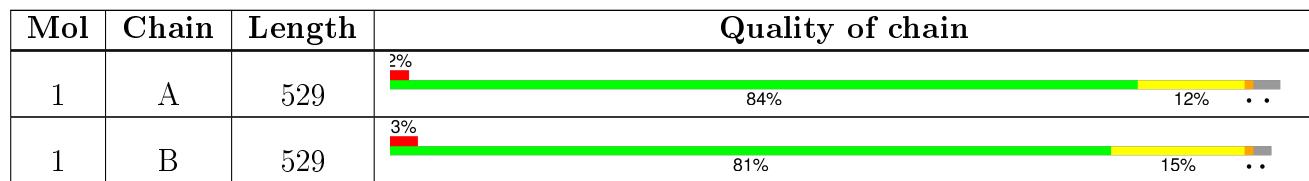
The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	B	4002	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 9038 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 ATP-dependent phosphoenolpyruvate carboxykinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	515	Total	C	N	O	S	0	0	0
			4097	2637	715	736	9			
1	B	519	Total	C	N	O	S	0	0	0
			4128	2655	720	743	10			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).
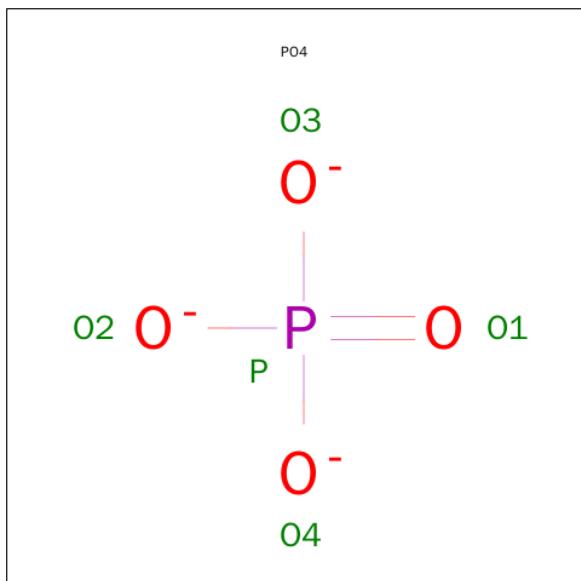
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 5	O 4	P 1	0	0
3	A	1	Total 5	O 4	P 1	0	0
3	B	1	Total 5	O 4	P 1	0	0
3	B	1	Total 5	O 4	P 1	0	0
3	B	1	Total 5	O 4	P 1	0	0
3	B	1	Total 5	O 4	P 1	0	0

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).
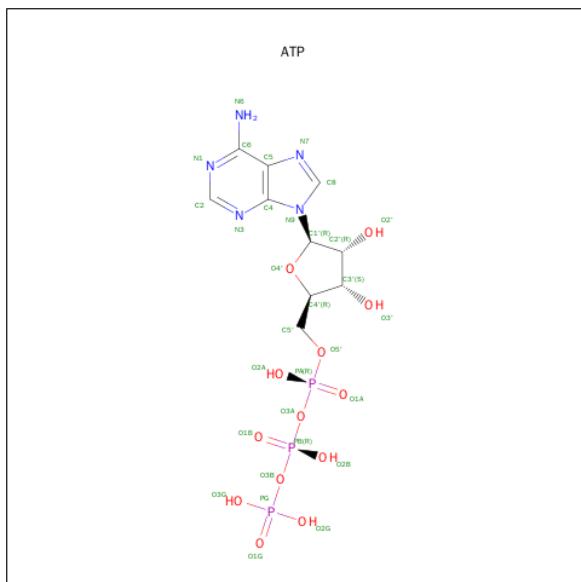
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 31	C 10	N 5	O 13	P 3	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).
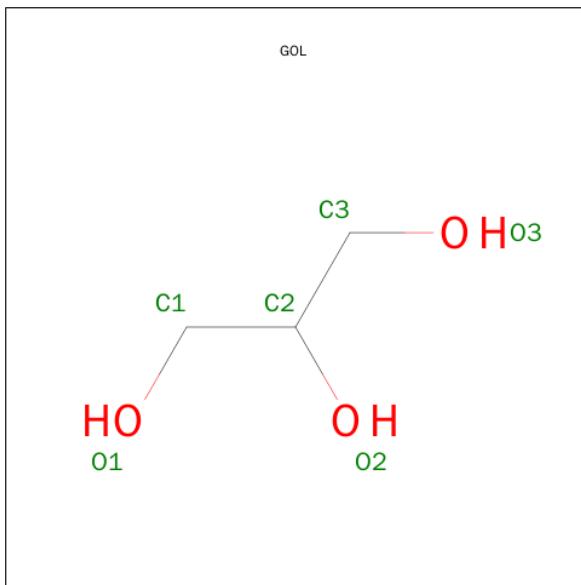
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total 6	C 3	O 3	0	0
5	B	1	Total 6	C 3	O 3	0	0

• Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	351	Total 351	O 351	0	0
6	B	387	Total 387	O 387	0	0

3 Residue-property plots ⓘ

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: ATP-dependent phosphoenolpyruvate carboxykinase
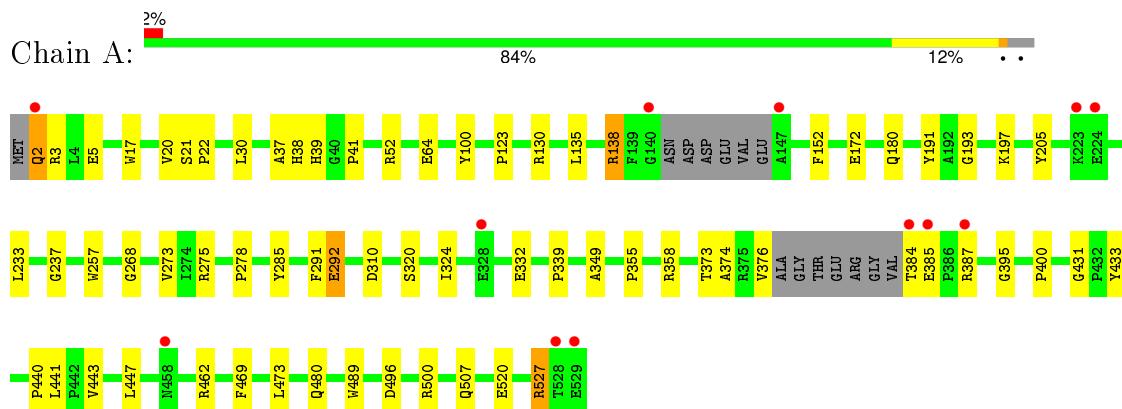
- Molecule 1: ATP-dependent phosphoenolpyruvate carboxykinase
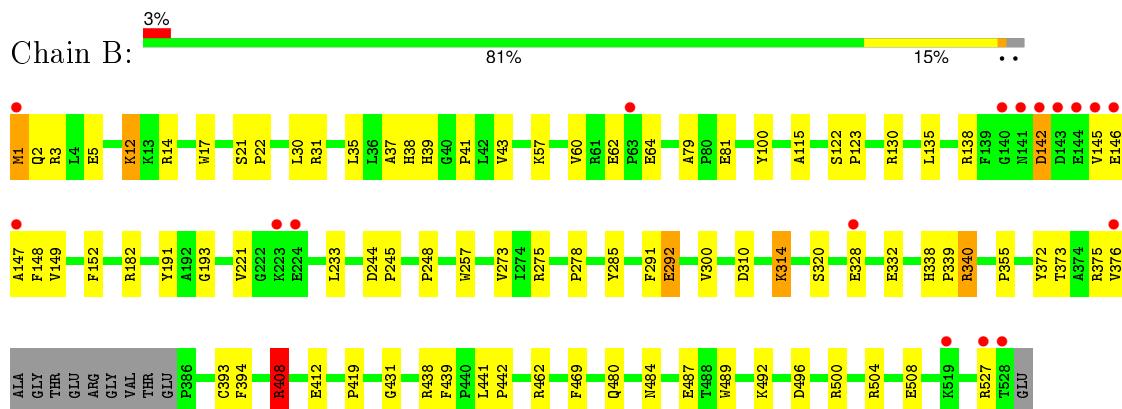
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	65.07Å 130.55Å 174.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.93 – 2.20 38.93 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (38.93-2.20) 99.7 (38.93-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$<I/\sigma(I)>$ [1]	3.45 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.186 , 0.219 0.186 , 0.218	Depositor DCC
R_{free} test set	3688 reflections (4.87%)	DCC
Wilson B-factor (Å^2)	30.6	Xtriage
Anisotropy	0.375	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 46.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$<|L|> = 0.49$, $<L^2> = 0.32$	Xtriage
Outliers	0 of 75748 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9038	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PO4, CA, ATP

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	0.35	0/4216	0.60	0/5732
1	B	0.36	0/4248	0.66	3/5777 (0.1%)
All	All	0.35	0/8464	0.63	3/11509 (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	B	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	12	LYS	CD-CE-NZ	-7.32	94.86	111.70
1	B	12	LYS	CB-CG-CD	6.59	128.74	111.60
1	B	12	LYS	CG-CD-CE	5.57	128.62	111.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	408	ARG	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	4097	0	4021	43	0
1	B	4128	0	4051	70	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	10	0	0	0	0
3	B	20	0	0	0	0
4	A	31	0	12	1	0
5	B	12	0	16	0	0
6	A	351	0	0	3	0
6	B	387	0	0	6	0
All	All	9038	0	8100	113	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 (113) 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:340:ARG:HA	1:B:340:ARG:HE	1.33	0.92
1:B:1:MET:HE2	1:B:3:ARG:HG2	1.55	0.88
1:B:1:MET:CE	1:B:3:ARG:HG2	2.13	0.78
1:B:37:ALA:HB3	1:B:41:PRO:HG2	1.71	0.70
1:B:1:MET:CE	1:B:3:ARG:CG	2.70	0.68
1:B:462:ARG:HH21	1:B:484:ASN:HA	1.59	0.67
1:A:3:ARG:HG3	1:A:5:GLU:OE2	1.95	0.67
1:B:1:MET:HE1	1:B:3:ARG:CG	2.26	0.66
1:A:123:PRO:HD2	6:A:3027:HOH:O	1.96	0.66
1:B:2:GLN:O	1:B:3:ARG:HD2	1.97	0.64
1:A:462:ARG:HD3	1:A:473:LEU:HD21	1.80	0.62
1:B:100:TYR:CZ	1:B:123:PRO:HG3	2.36	0.61
1:A:5:GLU:N	1:A:5:GLU:OE2	2.35	0.60
1:B:340:ARG:CA	1:B:340:ARG:HE	2.10	0.59
1:B:408:ARG:CZ	1:B:408:ARG:HB3	2.31	0.59
1:B:291:PHE:O	1:B:292:GLU:HB2	2.02	0.59
1:B:504:ARG:O	1:B:508:GLU:HG3	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:PRO:HB3	1:B:439:PHE:CZ	2.38	0.59
1:A:291:PHE:O	1:A:292:GLU:HB2	2.02	0.59
1:B:441:LEU:HB3	1:B:442:PRO:HD3	1.85	0.58
1:B:355:PRO:HG2	6:B:4321:HOH:O	2.03	0.58
1:A:233:LEU:HD12	1:A:373:THR:HG22	1.85	0.58
1:A:64:GLU:HG2	1:A:180:GLN:OE1	2.04	0.58
1:A:237:GLY:HA2	4:A:1001:ATP:H5'1	1.84	0.57
1:A:376:VAL:HG21	1:A:433:TYR:CD2	2.40	0.57
1:B:123:PRO:HD2	6:B:4024:HOH:O	2.04	0.57
1:A:275:ARG:HA	1:A:310:ASP:OD1	2.06	0.56
1:B:278:PRO:HG3	1:B:285:TYR:CE1	2.40	0.56
1:B:1:MET:HE1	1:B:3:ARG:HG3	1.87	0.55
1:B:462:ARG:NH2	1:B:484:ASN:HA	2.21	0.55
1:A:38:HIS:O	1:A:39:HIS:HB2	2.05	0.55
1:B:100:TYR:CE1	1:B:123:PRO:HG3	2.42	0.55
1:A:358:ARG:HD2	6:A:3145:HOH:O	2.07	0.54
1:B:81:GLU:H	1:B:81:GLU:CD	2.11	0.54
1:B:43:VAL:HG21	1:B:300:VAL:HG21	1.90	0.52
1:A:440:PRO:HG2	1:A:443:VAL:HG23	1.92	0.52
1:B:1:MET:HE2	1:B:3:ARG:CG	2.31	0.52
1:B:3:ARG:HB3	1:B:5:GLU:OE2	2.09	0.52
1:A:37:ALA:HB3	1:A:41:PRO:HG2	1.92	0.51
1:A:520:GLU:OE1	1:A:520:GLU:N	2.40	0.51
1:A:205:TYR:HB2	1:A:400:PRO:HG3	1.93	0.51
1:A:2:GLN:HB2	1:A:17:TRP:CH2	2.46	0.51
1:B:35:LEU:C	1:B:35:LEU:HD12	2.31	0.50
1:A:431:GLY:HA2	1:A:489:TRP:CE2	2.46	0.50
1:B:60:VAL:HG12	1:B:62:GLU:HG3	1.94	0.50
1:A:355:PRO:HB2	1:A:447:LEU:CD1	2.42	0.50
1:B:138:ARG:HH22	1:B:332:GLU:HG2	1.76	0.49
1:B:130:ARG:HG3	1:B:135:LEU:HD21	1.95	0.49
1:B:57:LYS:HD3	6:B:4311:HOH:O	2.12	0.49
1:B:221:VAL:O	1:B:248:PRO:HD2	2.13	0.48
1:A:291:PHE:O	1:A:292:GLU:CB	2.61	0.48
1:B:145:VAL:HG23	1:B:146:GLU:N	2.29	0.48
1:B:496:ASP:O	1:B:500:ARG:HG3	2.14	0.48
1:B:441:LEU:N	1:B:442:PRO:CD	2.77	0.47
1:B:441:LEU:N	1:B:442:PRO:HD2	2.30	0.47
1:A:100:TYR:CZ	1:A:123:PRO:HG3	2.49	0.47
1:A:257:TRP:CZ2	1:A:339:PRO:HD2	2.49	0.47
1:A:130:ARG:HG3	1:A:135:LEU:HD21	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:THR:HG22	1:A:385:GLU:N	2.30	0.47
1:A:384:THR:HG22	1:A:385:GLU:H	1.80	0.47
1:B:21:SER:HB2	1:B:22:PRO:HD3	1.96	0.46
1:B:147:ALA:O	1:B:149:VAL:HG23	2.15	0.46
1:A:349:ALA:H	1:A:374:ALA:HB3	1.80	0.46
1:B:314:LYS:HB2	6:B:4008:HOH:O	2.15	0.46
1:A:507:GLN:OE1	1:A:507:GLN:HA	2.15	0.46
1:B:79:ALA:HB1	1:B:81:GLU:OE1	2.16	0.46
1:B:291:PHE:O	1:B:292:GLU:CB	2.64	0.46
1:A:496:ASP:O	1:A:500:ARG:HG3	2.16	0.45
1:B:2:GLN:HA	1:B:17:TRP:CH2	2.50	0.45
1:B:1:MET:CE	1:B:3:ARG:HG3	2.45	0.45
1:B:527:ARG:NH1	1:B:527:ARG:HG2	2.32	0.44
1:A:355:PRO:HB2	1:A:447:LEU:HD11	1.99	0.44
1:B:275:ARG:HA	1:B:310:ASP:OD1	2.18	0.44
1:A:278:PRO:HG3	1:A:285:TYR:CE1	2.52	0.44
1:B:191:TYR:CE2	1:B:193:GLY:HA3	2.53	0.44
1:B:408:ARG:O	1:B:412:GLU:HG3	2.18	0.44
1:A:138:ARG:HB3	1:A:138:ARG:HE	1.52	0.43
1:B:182:ARG:HD2	1:B:182:ARG:HH11	1.61	0.43
1:A:268:GLY:HA2	1:A:324:ILE:N	2.33	0.43
1:B:233:LEU:HD12	1:B:373:THR:CG2	2.49	0.43
1:A:527:ARG:O	1:A:527:ARG:HG3	2.18	0.43
1:B:31:ARG:HD2	1:B:31:ARG:HH11	1.60	0.43
1:B:487:GLU:HA	1:B:492:LYS:NZ	2.33	0.43
1:B:14:ARG:HG3	1:B:14:ARG:NH1	2.34	0.43
1:B:393:CYS:O	1:B:394:PHE:HB2	2.19	0.43
1:B:14:ARG:HG3	1:B:14:ARG:HH11	1.84	0.42
1:A:52:ARG:HD3	1:A:197:LYS:HG3	2.01	0.42
1:B:338:HIS:HA	1:B:339:PRO:HD3	1.93	0.42
1:B:64:GLU:N	1:B:64:GLU:OE1	2.49	0.42
1:B:257:TRP:CZ2	1:B:339:PRO:HD2	2.55	0.42
1:B:273:VAL:HG11	1:B:320:SER:HB2	2.01	0.42
1:B:1:MET:HA	6:B:4168:HOH:O	2.19	0.42
1:A:20:VAL:HG23	1:A:22:PRO:HD2	2.02	0.42
1:B:38:HIS:O	1:B:39:HIS:HB2	2.19	0.42
1:B:372:TYR:CG	1:B:373:THR:N	2.88	0.42
1:B:35:LEU:O	1:B:35:LEU:HD12	2.20	0.41
1:A:21:SER:HB2	1:A:22:PRO:HD3	2.02	0.41
1:A:273:VAL:HG11	1:A:320:SER:HB2	2.02	0.41
1:A:191:TYR:CE2	1:A:193:GLY:HA3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:ARG:HD3	1:A:387:ARG:HH11	1.64	0.41
1:A:2:GLN:HB2	1:A:17:TRP:CZ2	2.56	0.41
1:B:339:PRO:O	1:B:419:PRO:HB3	2.21	0.41
1:B:115:ALA:HB1	1:B:148:PHE:CZ	2.55	0.41
1:A:462:ARG:HD3	1:A:473:LEU:CD2	2.48	0.41
1:B:438:ARG:NH2	6:B:4368:HOH:O	2.41	0.41
1:B:244:ASP:HA	1:B:245:PRO:HD3	1.98	0.41
1:B:527:ARG:HG2	1:B:527:ARG:HH11	1.85	0.41
1:B:81:GLU:N	1:B:81:GLU:CD	2.73	0.40
1:A:233:LEU:HD12	1:A:373:THR:CG2	2.49	0.40
1:B:431:GLY:HA2	1:B:489:TRP:CE2	2.56	0.40
1:A:30:LEU:HG	6:A:3292:HOH:O	2.21	0.40
1:B:233:LEU:HD12	1:B:373:THR:HG22	2.02	0.40
1:B:122:SER:HA	1:B:123:PRO:HD3	1.93	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	509/529 (96%)	497 (98%)	10 (2%)	2 (0%)	39	42
1	B	515/529 (97%)	495 (96%)	18 (4%)	2 (0%)	39	42
All	All	1024/1058 (97%)	992 (97%)	28 (3%)	4 (0%)	39	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	292	GLU
1	B	142	ASP
1	B	292	GLU
1	A	395	GLY

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	427/438 (98%)	418 (98%)	9 (2%)	61	74
1	B	431/438 (98%)	418 (97%)	13 (3%)	48	60
All	All	858/876 (98%)	836 (97%)	22 (3%)	54	66

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	138	ARG
1	A	152	PHE
1	A	172	GLU
1	A	332	GLU
1	A	441	LEU
1	A	469	PHE
1	A	480	GLN
1	A	527	ARG
1	B	1	MET
1	B	12	LYS
1	B	30	LEU
1	B	142	ASP
1	B	152	PHE
1	B	314	LYS
1	B	328	GLU
1	B	340	ARG
1	B	375	ARG
1	B	376	VAL
1	B	408	ARG
1	B	469	PHE
1	B	480	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	GLN
1	A	480	GLN
1	A	497	GLN
1	B	2	GLN
1	B	92	GLN
1	B	458	ASN
1	B	480	GLN
1	B	497	GLN
1	B	507	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)

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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	$\#|Z| > 2$
4	ATP	A	1001	-	24,33,33	1.35	3 (12%)	31,52,52	2.63	6 (19%)
3	PO4	A	3001	-	4,4,4	1.13	0	6,6,6	0.27	0
3	PO4	A	3002	-	4,4,4	1.13	0	6,6,6	0.27	0
3	PO4	B	3003	-	4,4,4	1.18	0	6,6,6	0.27	0
3	PO4	B	3004	-	4,4,4	1.11	0	6,6,6	0.27	0
3	PO4	B	3005	-	4,4,4	1.04	0	6,6,6	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
3	PO4	B	3006	-	4,4,4	1.07	0	6,6,6	0.27	0
5	GOL	B	4001	-	5,5,5	0.23	0	5,5,5	0.28	0
5	GOL	B	4002	-	5,5,5	0.15	0	5,5,5	0.23	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
4	ATP	A	1001	-	-	0/18/38/38	0/3/3/3
3	PO4	A	3001	-	-	0/0/0/0	0/0/0/0
3	PO4	A	3002	-	-	0/0/0/0	0/0/0/0
3	PO4	B	3003	-	-	0/0/0/0	0/0/0/0
3	PO4	B	3004	-	-	0/0/0/0	0/0/0/0
3	PO4	B	3005	-	-	0/0/0/0	0/0/0/0
3	PO4	B	3006	-	-	0/0/0/0	0/0/0/0
5	GOL	B	4001	-	-	0/4/4/4	0/0/0/0
5	GOL	B	4002	-	-	0/4/4/4	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1001	ATP	O4'-C1'	2.03	1.43	1.41
4	A	1001	ATP	C2-N1	2.35	1.38	1.33
4	A	1001	ATP	C2-N3	3.81	1.38	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1001	ATP	N3-C2-N1	-11.69	119.94	128.89
4	A	1001	ATP	C4-C5-N7	-3.95	105.84	109.48
4	A	1001	ATP	O3A-PA-O5'	-2.57	96.11	102.94
4	A	1001	ATP	O4'-C1'-N9	-2.17	103.56	108.10
4	A	1001	ATP	N6-C6-N1	-2.14	114.60	119.20
4	A	1001	ATP	C2'-C1'-N9	4.36	120.95	114.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1001	ATP	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	515/529 (97%)	-0.26	12 (2%)	64	63	20, 31, 46, 77	0
1	B	519/529 (98%)	-0.27	17 (3%)	50	49	18, 30, 50, 81	0
All	All	1034/1058 (97%)	-0.26	29 (2%)	56	55	18, 30, 50, 81	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	142	ASP	5.9
1	B	141	ASN	5.0
1	B	145	VAL	4.5
1	B	144	GLU	4.2
1	B	146	GLU	3.6
1	B	143	ASP	3.1
1	B	528	THR	3.1
1	A	2	GLN	3.0
1	A	384	THR	2.8
1	B	1	MET	2.7
1	A	140	GLY	2.7
1	A	223	LYS	2.7
1	A	385	GLU	2.7
1	B	140	GLY	2.6
1	B	328	GLU	2.5
1	A	528	THR	2.5
1	A	224	GLU	2.5
1	B	147	ALA	2.4
1	A	458	ASN	2.4
1	B	223	LYS	2.3
1	B	527	ARG	2.3
1	B	224	GLU	2.3
1	A	387	ARG	2.2
1	B	63	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	519	LYS	2.1
1	B	376	VAL	2.1
1	A	529	GLU	2.0
1	A	147	ALA	2.0
1	A	328	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

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, 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
5	GOL	B	4002	6/6	0.93	0.13	<mark>2.04</mark>	46,48,51,54	0
4	ATP	A	1001	31/31	0.79	0.21	1.68	76,82,103,103	0
3	PO4	B	3006	5/5	0.91	0.21	1.29	78,78,79,79	0
5	GOL	B	4001	6/6	0.97	0.12	0.58	31,32,33,33	0
3	PO4	B	3004	5/5	0.96	0.11	-0.18	59,59,60,60	0
3	PO4	A	3002	5/5	0.98	0.08	-0.84	52,52,52,53	0
3	PO4	B	3005	5/5	0.97	0.08	-1.11	40,40,42,43	0
3	PO4	B	3003	5/5	0.99	0.07	-1.42	23,26,26,28	0
3	PO4	A	3001	5/5	0.99	0.07	-1.89	22,23,24,25	0
2	CA	A	2001	1/1	0.98	0.07	-2.07	28,28,28,28	0
2	CA	B	2002	1/1	1.00	0.05	-2.64	31,31,31,31	0

6.5 Other polymers (i)

There are no such residues in this entry.