Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 09:15 AM GMT

PDB ID	:	3HQD
Title	:	Human kinesin Eg5 motor domain in complex with AMPPNP and Mg2+
Authors	:	Parke, C.L.; Wojcik, E.J.; Kim, S.; Worthylake, D.K.
Deposited on	:	2009-06-05
Resolution	:	2.19 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:

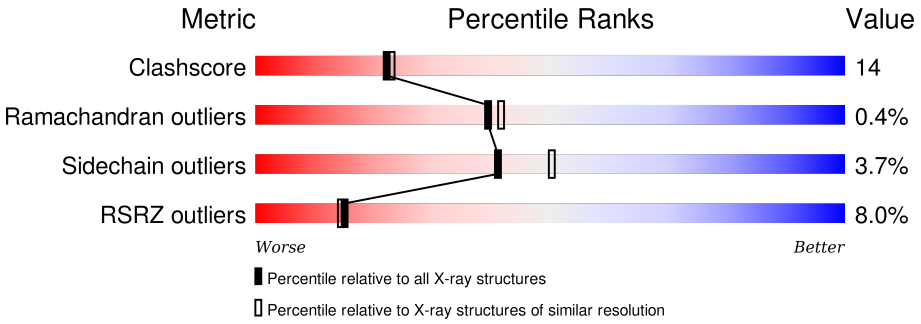
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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))
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	369	4% 72% 22% • 5%
1	B	369	11% 68% 23% • 5%

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	PO4	A	562	-	-	-	X
4	MG	B	502	-	-	-	X

2 Entry composition

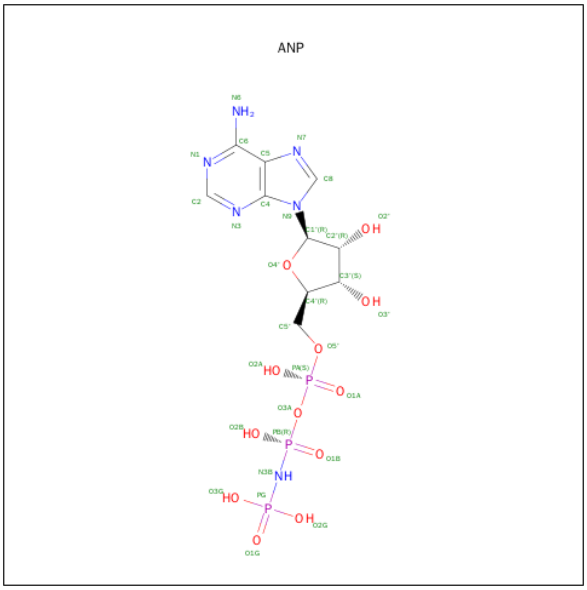
There are 5 unique types of molecules in this entry. The entry contains 5832 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Kinesin-like protein KIF11.

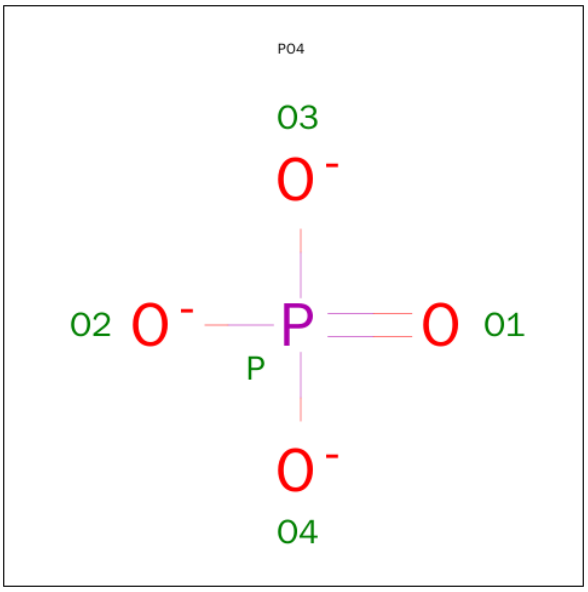
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	350	Total	C	N	O	S	0	0	0
			2741	1711	484	536	10			
1	B	350	Total	C	N	O	S	0	0	0
			2741	1711	484	536	10			

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is water.

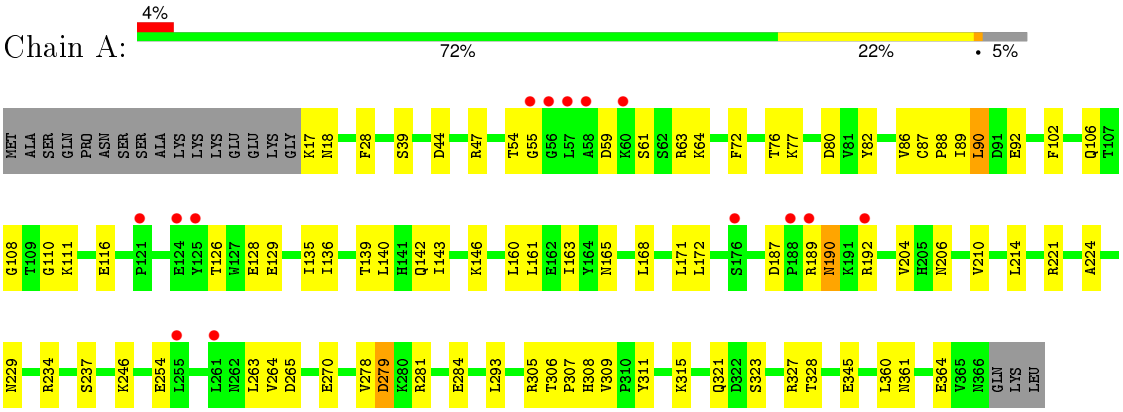
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	181	Total O 181 181	0	0
5	B	100	Total O 100 100	0	0

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.

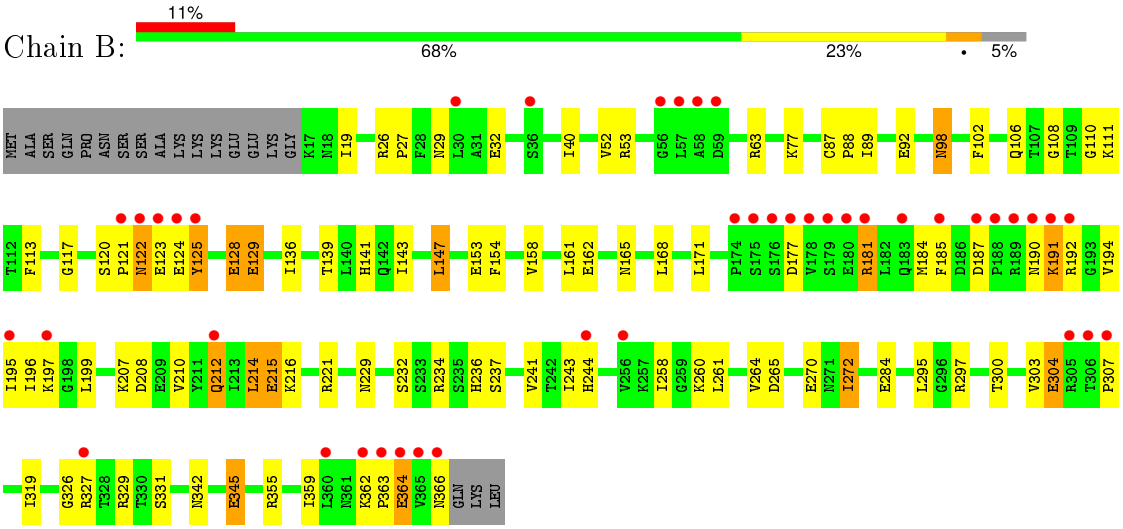
- Molecule 1: Kinesin-like protein KIF11

Chain A:



- Molecule 1: Kinesin-like protein KIF11

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	60.50Å 71.44Å 94.85Å 90.00° 98.86° 90.00°	Depositor
Resolution (Å)	25.00 – 2.19 24.78 – 2.17	Depositor EDS
% Data completeness (in resolution range)	99.6 (25.00-2.19) 97.9 (24.78-2.17)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.46 (at 2.17Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.262 0.225 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	21.7	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 49.1	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 41480 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5832	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% 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: PO4, ANP, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.34	0/2781	0.63	0/3759
1	B	0.30	0/2781	0.58	0/3759
All	All	0.32	0/5562	0.61	0/7518

There are no bond length outliers.

There are no bond angle outliers.

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	2741	0	2769	73	0
1	B	2741	0	2769	78	0
2	A	31	0	13	7	0
2	B	31	0	13	5	0
3	A	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	181	0	0	5	0
5	B	100	0	0	2	0
All	All	5832	0	5564	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 14.

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)
1:B:122:ASN:HB3	1:B:124:GLU:HG3	1.51	0.91
1:B:191:LYS:H	1:B:191:LYS:HD3	1.38	0.87
1:B:53:ARG:HD3	1:B:63:ARG:HH21	1.42	0.84
1:A:90:LEU:HD22	1:A:146:LYS:HD2	1.59	0.84
1:A:90:LEU:HD21	1:A:143:ILE:HD13	1.66	0.77
1:A:108:GLY:H	2:A:601:ANP:HNB1	1.32	0.76
1:A:192:ARG:CZ	1:A:327:ARG:HD2	2.16	0.76
1:B:110:GLY:HA2	2:B:602:ANP:O2A	1.86	0.76
1:B:272:ILE:HD13	1:B:272:ILE:H	1.52	0.74
1:B:158:VAL:HG12	1:B:241:VAL:HG12	1.72	0.71
1:A:17:LYS:HB3	1:A:360:LEU:HD11	1.73	0.70
1:A:192:ARG:HA	5:A:497:HOH:O	1.92	0.70
1:B:106:GLN:NE2	1:B:345:GLU:HG3	2.05	0.70
1:A:206:ASN:O	1:A:210:VAL:HG23	1.92	0.69
1:A:165:ASN:HD21	1:A:284:GLU:HA	1.58	0.69
1:A:110:GLY:HA2	2:A:601:ANP:O2A	1.94	0.68
1:B:125:TYR:HB3	1:B:129:GLU:HB3	1.76	0.68
1:B:108:GLY:H	2:B:602:ANP:HNB1	1.45	0.64
1:A:315:LYS:HE3	5:A:456:HOH:O	1.98	0.64
1:B:229:ASN:HD22	2:B:602:ANP:HNB1	1.47	0.61
1:B:191:LYS:CD	1:B:191:LYS:H	2.13	0.60
1:B:272:ILE:N	1:B:272:ILE:HD13	2.16	0.60
1:A:192:ARG:NE	1:A:327:ARG:HD2	2.18	0.59
1:B:258:ILE:H	1:B:366:ASN:HB3	1.68	0.59
1:A:116:GLU:HG2	5:A:503:HOH:O	2.03	0.58
1:A:55:GLY:CA	1:A:59:ASP:HB3	2.33	0.58
1:A:55:GLY:HA3	1:A:59:ASP:HB3	1.86	0.58
1:B:272:ILE:CD1	1:B:272:ILE:H	2.17	0.58
1:A:140:LEU:HD12	1:A:214:LEU:HD22	1.86	0.57
1:B:207:LYS:O	1:B:210:VAL:HG12	2.04	0.57
1:B:165:ASN:HD21	1:B:284:GLU:HA	1.68	0.57
1:B:181:ARG:H	1:B:181:ARG:HD3	1.68	0.57
1:B:98:ASN:ND2	1:B:260:LYS:HB3	2.20	0.56
1:A:54:THR:HG21	1:A:64:LYS:HD3	1.87	0.55
1:A:224:ALA:HB1	1:A:281:ARG:HE	1.71	0.55
1:A:17:LYS:HB3	1:A:360:LEU:CD1	2.36	0.55
1:B:234:ARG:HD3	1:B:284:GLU:OE1	2.05	0.55
1:B:147:LEU:HD23	1:B:154:PHE:CD1	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:LEU:HD12	1:B:161:LEU:C	2.27	0.55
1:B:92:GLU:OE2	1:B:329:ARG:HD3	2.07	0.54
1:A:102:PHE:HB3	1:A:264:VAL:HB	1.90	0.54
1:B:168:LEU:HD12	1:B:168:LEU:N	2.22	0.54
1:A:163:ILE:HG12	1:A:168:LEU:CD2	2.39	0.53
1:A:140:LEU:HD13	1:A:210:VAL:CG1	2.39	0.53
1:B:194:VAL:HG11	1:B:319:ILE:HD13	1.89	0.53
1:B:153:GLU:HB2	5:B:463:HOH:O	2.08	0.53
1:A:246:LYS:HE2	1:A:254:GLU:CD	2.28	0.52
1:A:111:LYS:NZ	2:A:601:ANP:O1G	2.40	0.52
1:B:194:VAL:HG22	1:B:195:ILE:N	2.24	0.52
1:A:306:THR:HG23	1:A:307:PRO:HD2	1.92	0.52
1:B:102:PHE:HB3	1:B:264:VAL:HB	1.92	0.52
1:A:305:ARG:NH1	1:A:305:ARG:HG3	2.24	0.52
1:B:89:ILE:HA	1:B:92:GLU:OE2	2.10	0.51
1:A:311:TYR:CG	1:A:321:GLN:HG3	2.45	0.51
1:A:82:TYR:OH	1:A:142:GLN:HG2	2.11	0.51
1:B:345:GLU:CD	1:B:345:GLU:H	2.09	0.51
1:A:305:ARG:HH11	1:A:305:ARG:HG3	1.76	0.51
1:B:40:ILE:O	1:B:52:VAL:HA	2.10	0.51
1:A:72:PHE:HB3	1:A:76:THR:OG1	2.11	0.51
1:A:140:LEU:HD13	1:A:210:VAL:HG12	1.93	0.50
1:A:192:ARG:NH1	5:A:541:HOH:O	2.39	0.50
1:B:212:GLN:HE21	1:B:212:GLN:C	2.15	0.50
1:A:161:LEU:HD12	1:A:161:LEU:C	2.32	0.50
1:B:122:ASN:HD22	1:B:124:GLU:CD	2.15	0.49
1:A:311:TYR:CD1	1:A:321:GLN:HG3	2.47	0.49
1:B:181:ARG:N	1:B:181:ARG:HD3	2.28	0.49
1:A:90:LEU:CD2	1:A:143:ILE:HD13	2.38	0.49
1:A:327:ARG:HG2	1:A:327:ARG:HH11	1.77	0.49
1:B:196:ILE:N	1:B:196:ILE:HD12	2.27	0.49
1:B:143:ILE:HD13	1:B:243:ILE:HD11	1.95	0.49
1:B:329:ARG:HG3	1:B:363:PRO:HG3	1.94	0.49
1:A:102:PHE:CB	1:A:264:VAL:HB	2.43	0.49
1:A:28:PHE:HZ	1:A:39:SER:HB2	1.76	0.49
1:B:102:PHE:CB	1:B:264:VAL:HB	2.43	0.48
1:A:86:VAL:HG21	1:A:135:ILE:HG23	1.95	0.48
1:A:136:ILE:HG12	1:A:263:LEU:HD13	1.96	0.48
1:B:207:LYS:HG3	1:B:208:ASP:N	2.28	0.48
1:B:162:GLU:HG2	1:B:171:LEU:HD21	1.95	0.48
1:A:229:ASN:HD22	2:A:601:ANP:HNB1	1.60	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:SER:O	1:A:63:ARG:HG3	2.14	0.48
1:A:108:GLY:N	2:A:601:ANP:HNB1	2.05	0.48
1:A:140:LEU:CD1	1:A:214:LEU:HD22	2.42	0.48
1:A:160:LEU:HG	1:A:171:LEU:HD12	1.96	0.47
1:A:278:VAL:HG12	1:A:279:ASP:N	2.29	0.47
1:A:234:ARG:CD	1:A:284:GLU:OE1	2.63	0.47
1:A:237:SER:HB3	1:A:265:ASP:HB3	1.95	0.47
1:B:229:ASN:HB3	1:B:232:SER:HB2	1.97	0.47
1:A:246:LYS:HE2	1:A:254:GLU:OE1	2.14	0.47
1:B:303:VAL:HG11	1:B:355:ARG:O	2.14	0.47
1:B:327:ARG:HG3	1:B:327:ARG:HH11	1.79	0.47
1:A:327:ARG:HD3	1:A:364:GLU:CD	2.35	0.47
1:B:87:CYS:N	1:B:88:PRO:HD2	2.30	0.46
1:A:234:ARG:HD3	1:A:284:GLU:OE1	2.15	0.46
1:B:329:ARG:HG2	1:B:329:ARG:HH11	1.80	0.46
1:B:185:PHE:O	1:B:194:VAL:HG23	2.16	0.45
1:A:306:THR:HG22	1:A:308:HIS:H	1.81	0.45
1:B:128:GLU:HG2	1:B:141:HIS:CD2	2.51	0.45
1:B:53:ARG:HD3	1:B:63:ARG:NH2	2.20	0.45
1:B:136:ILE:HG22	1:B:214:LEU:HD11	1.99	0.45
1:B:162:GLU:HA	1:B:236:HIS:O	2.17	0.45
1:A:87:CYS:HB2	1:A:88:PRO:HD3	1.99	0.45
2:B:602:ANP:H8	2:B:602:ANP:O2A	2.17	0.44
2:A:601:ANP:H8	2:A:601:ANP:O2A	2.18	0.44
1:A:234:ARG:NH2	1:A:270:GLU:OE1	2.45	0.44
1:B:147:LEU:HB3	1:B:154:PHE:CE2	2.53	0.44
1:B:327:ARG:HA	1:B:362:LYS:O	2.17	0.44
1:A:106:GLN:NE2	1:A:345:GLU:HG2	2.33	0.44
1:A:44:ASP:OD2	1:A:47:ARG:HD3	2.17	0.44
1:B:342:ASN:HB3	1:B:345:GLU:HG2	1.99	0.44
1:B:111:LYS:NZ	2:B:602:ANP:O1G	2.44	0.44
1:A:323:SER:HA	1:A:328:THR:HB	1.98	0.44
1:A:28:PHE:CZ	1:A:39:SER:HB2	2.53	0.43
1:B:234:ARG:CD	1:B:284:GLU:OE1	2.66	0.43
1:A:306:THR:CG2	5:A:423:HOH:O	2.67	0.43
1:B:196:ILE:HG22	1:B:199:LEU:HB2	2.01	0.43
1:A:246:LYS:NZ	1:A:254:GLU:HG3	2.33	0.43
1:A:77:LYS:HB2	1:A:80:ASP:OD2	2.19	0.43
1:A:293:LEU:C	1:A:293:LEU:HD23	2.39	0.43
1:B:139:THR:O	1:B:143:ILE:HG13	2.18	0.43
1:A:204:VAL:HB	1:A:210:VAL:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:SER:OG	1:B:265:ASP:HB3	2.19	0.43
1:B:244:HIS:CD2	5:B:448:HOH:O	2.72	0.42
1:A:140:LEU:HD22	1:A:210:VAL:HG11	2.01	0.42
1:B:234:ARG:NH2	1:B:270:GLU:OE2	2.40	0.42
1:B:184:MET:SD	1:B:194:VAL:HG21	2.59	0.42
1:A:224:ALA:HB1	1:A:281:ARG:NE	2.34	0.42
1:B:120:SER:HA	1:B:121:PRO:HD3	1.88	0.42
1:A:309:VAL:HG11	1:A:311:TYR:CZ	2.54	0.42
1:B:212:GLN:HE22	1:B:216:LYS:HD2	1.85	0.42
1:B:300:THR:O	1:B:304:GLU:HB2	2.19	0.42
1:B:192:ARG:NH1	1:B:326:GLY:HA2	2.35	0.42
1:B:297:ARG:HG2	1:B:297:ARG:HH11	1.85	0.42
1:B:241:VAL:CG2	1:B:261:LEU:HB3	2.49	0.42
1:B:19:ILE:HD12	1:B:359:ILE:HB	2.00	0.42
1:B:102:PHE:HA	1:B:264:VAL:O	2.20	0.41
1:B:77:LYS:HD3	1:B:77:LYS:HA	1.86	0.41
1:B:123:GLU:HG2	1:B:123:GLU:O	2.21	0.41
1:A:139:THR:O	1:A:143:ILE:HG12	2.21	0.41
1:B:327:ARG:HB3	1:B:364:GLU:HG3	2.03	0.41
1:A:126:THR:OG1	1:A:129:GLU:HG2	2.20	0.41
1:A:17:LYS:HA	1:A:361:ASN:O	2.21	0.41
1:A:187:ASP:HB3	1:A:190:ASN:O	2.21	0.41
1:B:215:GLU:HA	1:B:215:GLU:OE1	2.20	0.41
1:A:110:GLY:CA	2:A:601:ANP:O2A	2.67	0.41
1:A:160:LEU:HB3	1:A:172:LEU:HG	2.01	0.40
1:B:26:ARG:HB2	1:B:27:PRO:HD2	2.02	0.40
1:B:329:ARG:NH1	1:B:329:ARG:HG2	2.36	0.40
1:A:187:ASP:OD1	1:A:189:ARG:HG2	2.21	0.40
1:A:89:ILE:HA	1:A:92:GLU:OE2	2.21	0.40
1:B:89:ILE:HD11	1:B:331:SER:OG	2.21	0.40
1:B:113:PHE:O	1:B:117:GLY:HA2	2.21	0.40
1:B:29:ASN:OD1	1:B:32:GLU:HG3	2.22	0.40
1:B:187:ASP:HB3	1:B:190:ASN:HB2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	348/369 (94%)	337 (97%)	9 (3%)	2 (1%)	30	29
1	B	348/369 (94%)	333 (96%)	14 (4%)	1 (0%)	46	50
All	All	696/738 (94%)	670 (96%)	23 (3%)	3 (0%)	39	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	279	ASP
1	A	190	ASN
1	B	307	PRO

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	307/323 (95%)	303 (99%)	4 (1%)	76	87
1	B	307/323 (95%)	288 (94%)	19 (6%)	23	25
All	All	614/646 (95%)	591 (96%)	23 (4%)	41	50

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	90	LEU
1	A	128	GLU

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Mol	Chain	Res	Type
1	A	221	ARG
1	B	98	ASN
1	B	122	ASN
1	B	125	TYR
1	B	128	GLU
1	B	129	GLU
1	B	147	LEU
1	B	177	ASP
1	B	181	ARG
1	B	191	LYS
1	B	197	LYS
1	B	212	GLN
1	B	214	LEU
1	B	215	GLU
1	B	221	ARG
1	B	272	ILE
1	B	295	LEU
1	B	304	GLU
1	B	345	GLU
1	B	364	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	20	GLN
1	A	38	HIS
1	A	165	ASN
1	A	190	ASN
1	A	262	ASN
1	B	20	GLN
1	B	98	ASN
1	B	122	ASN
1	B	141	HIS
1	B	142	GLN
1	B	165	ASN
1	B	190	ASN
1	B	212	GLN
1	B	229	ASN
1	B	244	HIS
1	B	262	ASN
1	B	321	GLN

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 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	PO4	A	562	-	4,4,4	1.34	0	6,6,6	0.27	0
2	ANP	A	601	-	27,33,33	2.04	6 (22%)	30,52,52	4.19	19 (63%)
2	ANP	B	602	-	27,33,33	2.00	6 (22%)	30,52,52	4.12	15 (50%)

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	PO4	A	562	-	-	0/0/0/0	0/0/0/0
2	ANP	A	601	-	-	0/12/38/38	0/3/3/3
2	ANP	B	602	-	-	0/12/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	ANP	O5'-C5'	-2.95	1.32	1.44
2	B	602	ANP	O5'-C5'	-2.57	1.34	1.44
2	A	601	ANP	PG-O3G	-2.25	1.50	1.56
2	B	602	ANP	PB-O3A	-2.01	1.56	1.59
2	A	601	ANP	PB-N3B	2.42	1.69	1.63
2	B	602	ANP	PB-N3B	3.01	1.71	1.63
2	B	602	ANP	PA-O5'	3.14	1.73	1.59
2	A	601	ANP	PA-O5'	3.45	1.74	1.59
2	A	601	ANP	O4'-C1'	4.33	1.46	1.41
2	B	602	ANP	O4'-C1'	4.83	1.47	1.41
2	B	602	ANP	PG-O1G	5.29	1.52	1.46
2	A	601	ANP	PG-O1G	6.44	1.53	1.46

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	602	ANP	N3-C2-N1	-10.69	120.71	128.89
2	A	601	ANP	N3-C2-N1	-10.42	120.91	128.89
2	B	602	ANP	O3A-PA-O5'	-10.00	76.40	102.94
2	A	601	ANP	O3A-PA-O5'	-9.50	77.75	102.94
2	A	601	ANP	C5'-C4'-C3'	-6.84	88.06	115.21
2	A	601	ANP	PA-O3A-PB	-6.75	110.02	132.67
2	B	602	ANP	PA-O3A-PB	-6.46	111.02	132.67
2	B	602	ANP	C5'-C4'-C3'	-6.05	91.19	115.21
2	B	602	ANP	C2'-C1'-N9	-5.73	105.53	114.29
2	A	601	ANP	O3A-PB-N3B	-5.24	92.03	106.44
2	A	601	ANP	C2'-C1'-N9	-5.14	106.44	114.29
2	A	601	ANP	O5'-PA-O1A	-4.95	90.42	109.62
2	B	602	ANP	O5'-PA-O1A	-4.74	91.22	109.62
2	A	601	ANP	O2A-PA-O5'	-4.52	85.66	108.46
2	B	602	ANP	O2A-PA-O5'	-4.43	86.10	108.46
2	B	602	ANP	O3A-PB-N3B	-4.39	94.36	106.44
2	B	602	ANP	C1'-N9-C4	-2.92	122.54	126.94
2	A	601	ANP	C1'-N9-C4	-2.90	122.56	126.94
2	A	601	ANP	O4'-C4'-C5'	-2.39	100.77	109.32
2	B	602	ANP	O4'-C4'-C5'	-2.31	101.08	109.32
2	A	601	ANP	O2G-PG-O1G	-2.26	107.49	113.49
2	A	601	ANP	C4'-O4'-C1'	-2.05	107.47	109.72
2	B	602	ANP	O2B-PB-O3A	2.11	114.69	105.09
2	A	601	ANP	O2A-PA-O1A	2.32	125.11	112.53
2	A	601	ANP	O2B-PB-O3A	2.42	116.06	105.09
2	A	601	ANP	C2'-C3'-C4'	2.44	107.62	102.61
2	A	601	ANP	O1G-PG-N3B	2.45	115.65	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	602	ANP	C2'-C3'-C4'	3.28	109.35	102.61
2	A	601	ANP	O2A-PA-O3A	3.49	120.93	105.09
2	B	602	ANP	O2A-PA-O3A	3.56	121.26	105.09
2	B	602	ANP	O5'-C5'-C4'	4.41	125.38	109.12
2	A	601	ANP	O4'-C1'-N9	5.44	119.49	108.10
2	A	601	ANP	O5'-C5'-C4'	5.52	129.48	109.12
2	B	602	ANP	O4'-C1'-N9	5.84	120.33	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	ANP	7	0
2	B	602	ANP	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

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	350/369 (94%)	0.39	14 (4%) 42 41	8, 22, 43, 60	0
1	B	350/369 (94%)	0.80	42 (12%) 6 5	9, 29, 64, 90	0
All	All	700/738 (94%)	0.60	56 (8%) 15 14	8, 26, 54, 90	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	178	VAL	17.0
1	B	179	SER	15.6
1	B	57	LEU	6.6
1	B	121	PRO	6.1
1	A	57	LEU	5.9
1	A	58	ALA	5.5
1	B	185	PHE	5.4
1	B	124	GLU	5.4
1	B	175	SER	5.4
1	B	58	ALA	5.0
1	A	188	PRO	4.8
1	B	189	ARG	4.8
1	B	366	ASN	4.6
1	B	177	ASP	4.6
1	B	365	VAL	4.4
1	B	56	GLY	3.9
1	A	189	ARG	3.8
1	B	125	TYR	3.8
1	B	364	GLU	3.8
1	B	183	GLN	3.6
1	B	188	PRO	3.4
1	B	181	ARG	3.4
1	B	122	ASN	3.4
1	B	190	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	176	SER	3.3
1	A	192	ARG	3.1
1	B	360	LEU	3.1
1	A	255	LEU	2.9
1	B	327	ARG	2.9
1	B	305	ARG	2.8
1	B	180	GLU	2.8
1	B	212	GLN	2.7
1	B	59	ASP	2.7
1	B	307	PRO	2.6
1	B	197	LYS	2.5
1	A	125	TYR	2.5
1	A	124	GLU	2.4
1	A	121	PRO	2.4
1	B	195	ILE	2.4
1	A	55	GLY	2.3
1	B	191	LYS	2.3
1	A	56	GLY	2.3
1	B	362	LYS	2.2
1	A	261	LEU	2.2
1	B	306	THR	2.2
1	B	36	SER	2.1
1	B	256	VAL	2.1
1	B	174	PRO	2.1
1	B	244	HIS	2.1
1	B	30	LEU	2.1
1	B	363	PRO	2.1
1	B	187	ASP	2.1
1	A	176	SER	2.0
1	B	123	GLU	2.0
1	B	192	ARG	2.0
1	A	60	LYS	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

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
4	MG	B	502	1/1	0.83	0.25	2.85	31,31,31,31	0
3	PO4	A	562	5/5	0.90	0.17	2.14	38,39,42,43	0
2	ANP	B	602	31/31	0.91	0.19	1.62	14,21,31,35	0
2	ANP	A	601	31/31	0.91	0.18	1.11	11,17,27,31	0
4	MG	A	501	1/1	0.95	0.12	-1.26	14,14,14,14	0

6.5 Other polymers

There are no such residues in this entry.