



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:37 PM GMT

PDB ID : 2PG2  
Title : Crystal structure of KSP in complex with ADP and thiophene containing inhibitor 15  
Authors : Lee, T.T.  
Deposited on : 2007-04-06  
Resolution : 1.85 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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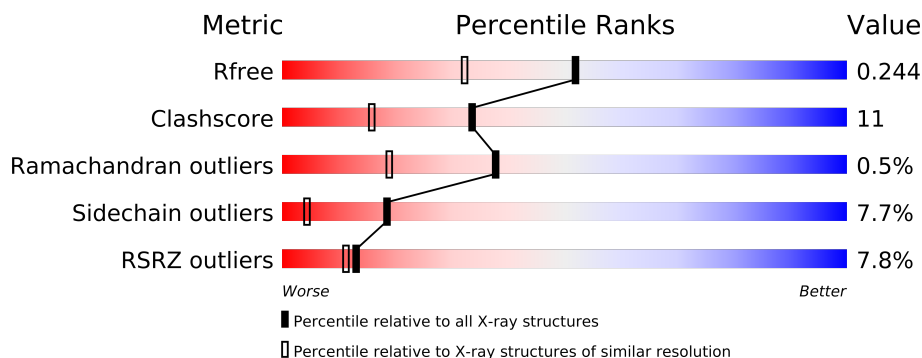
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.85 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))
$R_{free}$	66092	1857 (1.86-1.82)
Clashscore	79885	2149 (1.86-1.82)
Ramachandran outliers	78287	2124 (1.86-1.82)
Sidechain outliers	78261	2125 (1.86-1.82)
RSRZ outliers	66119	1857 (1.86-1.82)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	368	
1	B	368	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	K01	B	604	-	X

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 5776 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

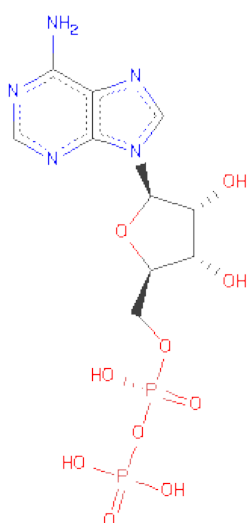
- Molecule 1 is a protein called Kinesin-like protein KIF11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	0	0
			2599	1628	453	508	10			
1	B	331	Total	C	N	O	S	0	0	0
			2599	1628	453	508	10			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

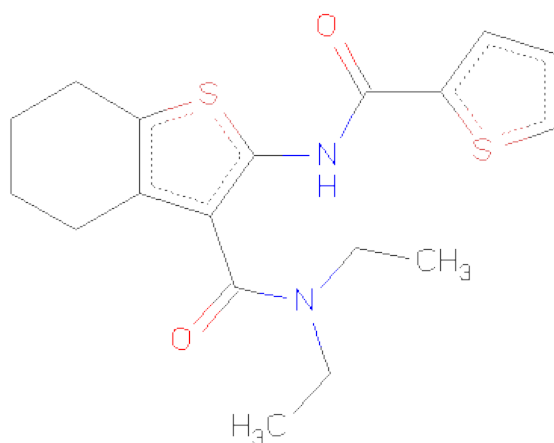
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is N,N-DIETHYL-2-[(2-THIENYLCARBONYL)AMINO]-4,5,6,7-TETRAHYDRO-1-BENZOTHIOPHENE-3-CARBOXAMIDE (three-letter code: K01) (formula:  $C_{18}H_{22}N_2O_2S_2$ ).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	247	Total	O	0	0
			247	247		
5	B	227	Total	O	0	0
			227	227		



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	161.06Å 80.35Å 69.31Å 90.00° 96.83° 90.00°	Depositor
Resolution (Å)	80.06 – 1.85 40.17 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.5 (80.06-1.85) 99.5 (40.17-1.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.97 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, $R_{free}$	0.210 , 0.248 0.207 , 0.244	Depositor DCC
$R_{free}$ test set	3760 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.5	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 42.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 74712 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5776	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K01, MG, ADP

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	0.77	0/2637	0.77	0/3564
1	B	0.70	0/2637	0.73	0/3564
All	All	0.73	0/5274	0.75	0/7128

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2599	0	2627	52	0
1	B	2599	0	2627	49	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	27	0	12	0	0
3	B	27	0	12	0	0
4	A	24	0	21	14	0
4	B	24	0	21	6	0
5	A	247	0	0	17	0
5	B	227	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5776	0	5320	117	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (117) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:A:604:K01:C2	4:A:604:K01:C1	1.75	1.54
4:B:604:K01:N14	4:B:604:K01:H21	1.54	1.20
4:A:604:K01:H21	4:A:604:K01:N14	1.57	1.20
1:A:82:TYR:OH	1:A:142:GLN:HG3	1.49	1.11
1:B:329:ARG:HG2	1:B:329:ARG:HH11	1.00	1.06
1:B:184:MET:HE2	1:B:194:VAL:HG11	1.35	1.05
4:A:604:K01:N3	4:A:604:K01:C1	2.20	1.03
1:B:329:ARG:CG	1:B:329:ARG:HH11	1.84	0.90
1:A:16:GLY:O	1:A:17:LYS:HB2	1.69	0.90
1:B:329:ARG:NH1	1:B:329:ARG:HG2	1.79	0.90
1:B:184:MET:CE	1:B:194:VAL:HG11	2.02	0.89
4:B:604:K01:H21	4:B:604:K01:C13	1.88	0.87
1:B:223:THR:HG21	5:B:698:HOH:O	1.81	0.81
1:B:173:ASN:O	1:B:220:LYS:HE3	1.81	0.81
4:B:604:K01:C2	4:B:604:K01:N14	2.44	0.79
4:A:604:K01:H21	4:A:604:K01:C13	1.95	0.78
1:B:200:GLU:HG2	5:B:769:HOH:O	1.82	0.78
1:A:186:ASP:OD1	1:A:312:ARG:NH2	2.22	0.72
1:A:299:ILE:HG23	1:A:359:ILE:HD11	1.72	0.71
1:A:223:THR:HG21	5:A:678:HOH:O	1.91	0.70
1:A:57:LEU:HB2	1:A:60:LYS:O	1.92	0.69
1:B:167:GLU:HG2	5:B:699:HOH:O	1.92	0.69
1:B:223:THR:CG2	5:B:698:HOH:O	2.40	0.68
1:A:118:GLU:O	4:A:604:K01:H19	1.94	0.67
1:A:87:CYS:SG	5:A:844:HOH:O	2.53	0.67
4:A:604:K01:O22	4:A:604:K01:C11	2.41	0.67
4:A:604:K01:C4	4:A:604:K01:C1	2.73	0.66
1:A:79:ILE:HD13	1:A:83:ARG:HD3	1.80	0.64
1:A:308:HIS:HE1	5:A:623:HOH:O	1.81	0.63
1:A:307:PRO:HG2	5:A:741:HOH:O	1.98	0.63
4:A:604:K01:C2	4:A:604:K01:N14	2.50	0.63
1:A:354:HIS:HD2	5:A:850:HOH:O	1.84	0.61
1:A:142:GLN:HB3	5:A:819:HOH:O	2.02	0.60
1:A:223:THR:HG22	5:A:691:HOH:O	2.01	0.60
1:A:248:THR:HG23	5:A:765:HOH:O	2.02	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:167:GLU:HG2	5:A:782:HOH:O	2.03	0.59
1:A:53:ARG:HD3	1:A:57:LEU:O	2.03	0.59
1:A:142:GLN:HG2	5:A:717:HOH:O	2.02	0.58
4:A:604:K01:O22	4:A:604:K01:H111	2.03	0.58
1:B:173:ASN:HD22	1:B:174:PRO:N	2.01	0.58
1:A:82:TYR:HH	1:A:142:GLN:HG3	1.64	0.58
1:B:79:ILE:HG13	1:B:83:ARG:HD3	1.85	0.58
1:A:161:LEU:HD22	1:A:196:ILE:HD13	1.85	0.57
1:A:79:ILE:CD1	1:A:83:ARG:HD3	2.34	0.57
1:B:65:THR:HG21	5:B:826:HOH:O	2.04	0.56
1:A:45:PRO:HA	1:A:71:VAL:HG13	1.87	0.56
1:A:212:GLN:HG3	5:A:827:HOH:O	2.04	0.56
1:A:48:LYS:HA	1:A:71:VAL:HG12	1.87	0.55
1:B:246:LYS:HE3	1:B:254:GLU:OE2	2.07	0.55
1:A:360:LEU:HD12	5:A:699:HOH:O	2.06	0.55
1:B:173:ASN:HD22	1:B:173:ASN:C	2.11	0.54
1:B:118:GLU:O	4:B:604:K01:H19	2.08	0.53
1:B:75:SER:O	1:B:77:LYS:HE2	2.09	0.52
1:A:244:HIS:HD2	5:A:684:HOH:O	1.91	0.52
1:B:173:ASN:HD22	1:B:174:PRO:CD	2.22	0.52
1:A:311:TYR:CD1	1:A:321:GLN:HG3	2.45	0.52
4:A:604:K01:C15	4:A:604:K01:H21	2.34	0.52
1:B:329:ARG:HD3	5:B:780:HOH:O	2.08	0.52
1:A:45:PRO:HA	1:A:71:VAL:CG1	2.40	0.51
1:B:184:MET:HE3	1:B:194:VAL:HG21	1.92	0.50
1:B:173:ASN:HD22	1:B:174:PRO:HD2	1.77	0.50
1:A:173:ASN:C	1:A:173:ASN:HD22	2.15	0.50
4:A:604:K01:C1	4:A:604:K01:C23	2.90	0.50
1:A:79:ILE:CD1	1:A:83:ARG:CD	2.90	0.49
1:A:355:ARG:NH2	5:A:711:HOH:O	2.45	0.49
1:A:87:CYS:HB2	1:A:88:PRO:HD3	1.94	0.48
1:A:149:ASP:O	1:A:150:ASN:O	2.32	0.48
1:B:38:HIS:CD2	1:B:53:ARG:HH12	2.31	0.48
1:B:246:LYS:HG2	1:B:254:GLU:HG2	1.94	0.48
1:A:241:VAL:HG13	1:A:261:LEU:HB3	1.96	0.47
1:B:234:ARG:NH1	5:B:774:HOH:O	2.34	0.47
1:B:38:HIS:HB3	5:B:751:HOH:O	2.14	0.47
1:A:173:ASN:ND2	1:A:175:SER:H	2.13	0.47
1:A:248:THR:HA	1:A:253:GLU:O	2.15	0.47
1:A:160:LEU:HB3	1:A:172:LEU:HG	1.98	0.46
4:B:604:K01:H22	4:B:604:K01:H242	1.35	0.46
1:A:321:GLN:NE2	5:A:703:HOH:O	2.46	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:191:LYS:HA	1:B:191:LYS:HE2	1.98	0.46
1:A:82:TYR:CZ	1:A:142:GLN:HG3	2.46	0.45
1:B:173:ASN:ND2	1:B:175:SER:H	2.14	0.45
1:B:171:LEU:HD13	1:B:220:LYS:HB3	1.98	0.45
1:A:191:LYS:HD2	5:A:763:HOH:O	2.16	0.45
1:A:178:VAL:CG1	1:A:220:LYS:HB3	2.47	0.44
1:A:157:LYS:HE3	1:A:157:LYS:HB2	1.72	0.44
1:A:149:ASP:CG	1:A:150:ASN:N	2.71	0.44
1:A:29:ASN:O	1:A:33:ARG:HG3	2.17	0.43
1:B:170:ASP:HB2	1:B:182:LEU:HD11	2.00	0.43
1:A:40:ILE:HD13	1:A:340:SER:HA	1.99	0.43
1:B:38:HIS:O	1:B:38:HIS:CG	2.70	0.43
1:B:167:GLU:OE1	1:B:181:ARG:NE	2.49	0.43
1:B:354:HIS:HD2	5:B:803:HOH:O	2.00	0.43
1:A:291:SER:HB3	1:A:316:LEU:CB	2.49	0.43
1:A:54:THR:HG21	1:A:64:LYS:HG3	1.99	0.43
1:B:289:ASN:ND2	1:B:292:LEU:H	2.17	0.43
1:B:119:ARG:HD2	4:B:604:K01:C19	2.49	0.43
1:B:189:ARG:HH21	1:B:195:ILE:HD11	1.83	0.42
4:A:604:K01:H22	4:A:604:K01:H242	1.30	0.42
1:B:38:HIS:HD2	1:B:53:ARG:HH22	1.66	0.42
1:B:189:ARG:NH2	1:B:195:ILE:HD11	2.35	0.42
1:B:299:ILE:HG23	1:B:359:ILE:HD11	2.02	0.42
1:A:119:ARG:HD2	4:A:604:K01:C19	2.49	0.42
1:A:178:VAL:HG13	1:A:220:LYS:HE2	2.01	0.42
1:B:325:GLY:H	1:B:361:ASN:HD21	1.67	0.42
1:A:149:ASP:OD1	1:A:150:ASN:N	2.54	0.41
1:B:27:PRO:HA	1:B:74:ALA:HB1	2.03	0.41
1:B:173:ASN:ND2	1:B:174:PRO:HD2	2.35	0.41
1:A:17:LYS:HD3	5:A:625:HOH:O	2.21	0.41
1:B:89:ILE:HG12	1:B:329:ARG:HH12	1.86	0.41
1:B:57:LEU:HD23	1:B:57:LEU:N	2.36	0.41
1:B:351:GLU:O	1:B:355:ARG:HG3	2.21	0.41
1:B:16:GLY:HA2	1:B:360:LEU:HB3	2.03	0.41
1:B:82:TYR:CD2	1:B:86:VAL:HB	2.56	0.41
1:B:40:ILE:HD13	1:B:340:SER:HA	2.02	0.40
1:A:27:PRO:HA	1:A:74:ALA:HB1	2.04	0.40
4:A:604:K01:C4	4:A:604:K01:H13	2.51	0.40
1:B:38:HIS:CE1	5:B:745:HOH:O	2.74	0.40
1:B:358:ASN:N	1:B:358:ASN:HD22	2.19	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	327/368 (89%)	321 (98%)	3 (1%)	3 (1%)	25	9
1	B	327/368 (89%)	323 (99%)	4 (1%)	0	100	100
All	All	654/736 (89%)	644 (98%)	7 (1%)	3 (0%)	38	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	149	ASP
1	A	150	ASN

### 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	293/322 (91%)	271 (92%)	22 (8%)	19	5
1	B	293/322 (91%)	270 (92%)	23 (8%)	18	4
All	All	586/644 (91%)	541 (92%)	45 (8%)	18	4

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	30	LEU
1	A	50	VAL
1	A	60	LYS
1	A	61	SER
1	A	79	ILE

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Mol	Chain	Res	Type
1	A	140	LEU
1	A	146	LYS
1	A	165	ASN
1	A	173	ASN
1	A	189	ARG
1	A	191	LYS
1	A	200	GLU
1	A	208	ASP
1	A	212	GLN
1	A	248	THR
1	A	250	ILE
1	A	254	GLU
1	A	257	LYS
1	A	289	ASN
1	A	305	ARG
1	A	341	LEU
1	B	34	LYS
1	B	50	VAL
1	B	57	LEU
1	B	61	SER
1	B	85	VAL
1	B	128	GLU
1	B	140	LEU
1	B	145	GLU
1	B	147	LEU
1	B	165	ASN
1	B	173	ASN
1	B	191	LYS
1	B	192	ARG
1	B	207	LYS
1	B	223	THR
1	B	241	VAL
1	B	254	GLU
1	B	257	LYS
1	B	271	ASN
1	B	288	ILE
1	B	289	ASN
1	B	329	ARG
1	B	341	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	173	ASN
1	A	229	ASN
1	A	244	HIS
1	A	262	ASN
1	A	271	ASN
1	A	289	ASN
1	A	308	HIS
1	A	321	GLN
1	A	361	ASN
1	B	18	ASN
1	B	38	HIS
1	B	173	ASN
1	B	229	ASN
1	B	244	HIS
1	B	262	ASN
1	B	271	ASN
1	B	289	ASN
1	B	308	HIS
1	B	342	ASN
1	B	358	ASN
1	B	361	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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
3	ADP	A	601	2	29,29,29	1.39	5 (17%)	45,45,45	1.83	8 (17%)
4	K01	A	604	-	26,26,26	1.81	8 (30%)	35,36,36	4.80	14 (40%)
3	ADP	B	601	2	29,29,29	1.33	4 (13%)	45,45,45	1.71	6 (13%)
4	K01	B	604	-	26,26,26	1.80	6 (23%)	35,36,36	5.35	16 (45%)

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	ADP	A	601	2	-	0/16/32/32	0/1/3/3
4	K01	A	604	-	-	0/17/27/27	0/1/3/3
3	ADP	B	601	2	-	0/16/32/32	0/1/3/3
4	K01	B	604	-	-	0/17/27/27	0/1/3/3

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	604	K01	C2-N3	-4.23	1.31	1.48
4	A	604	K01	C1-C2	4.09	1.75	1.49
4	A	604	K01	C2-N3	-4.07	1.31	1.48
4	B	604	K01	O22-C4	-4.06	1.14	1.22
4	A	604	K01	O22-C4	-3.65	1.15	1.22
4	B	604	K01	C1-C2	3.65	1.72	1.49
3	A	601	ADP	PB-O3A	3.35	1.66	1.60
4	A	604	K01	C15-N14	-3.20	1.27	1.35
4	B	604	K01	O21-C15	-2.91	1.17	1.23
3	A	601	ADP	C5-C4	2.90	1.47	1.40
3	B	601	ADP	C4-N9	-2.88	1.33	1.37
3	A	601	ADP	PB-O3B	-2.80	1.44	1.54
4	B	604	K01	C15-N14	-2.74	1.28	1.35
3	B	601	ADP	O4'-C1'	2.58	1.45	1.41
3	B	601	ADP	PB-O3A	2.56	1.64	1.60
4	A	604	K01	C13-S12	-2.53	1.70	1.73
3	B	601	ADP	C5-C4	2.43	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	ADP	C4-N9	-2.38	1.34	1.37
4	B	604	K01	C7-C6	2.34	1.39	1.36
4	A	604	K01	C7-C6	2.23	1.39	1.36
3	A	601	ADP	C2'-C1'	-2.22	1.50	1.53
4	A	604	K01	O21-C15	-2.01	1.19	1.23
4	A	604	K01	C4-N3	-2.01	1.29	1.34

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	604	K01	C5-C4-N3	19.32	137.88	117.99
4	A	604	K01	C5-C4-N3	18.85	137.39	117.99
4	B	604	K01	O22-C4-N3	-13.40	97.38	122.33
4	B	604	K01	C7-S12-C13	-11.57	90.38	92.36
4	A	604	K01	O22-C4-N3	-10.32	103.11	122.33
4	A	604	K01	C23-N3-C4	9.23	152.37	121.39
4	B	604	K01	C23-N3-C4	9.09	151.91	121.39
4	B	604	K01	C2-N3-C4	-8.20	93.84	121.39
4	A	604	K01	C2-N3-C4	-7.36	96.67	121.39
3	A	601	ADP	N3-C2-N1	-6.68	123.12	128.71
4	A	604	K01	C6-C5-C4	-6.26	114.65	126.30
3	B	601	ADP	N3-C2-N1	-6.03	123.67	128.71
4	B	604	K01	C9-C8-C7	-5.64	102.61	112.85
3	A	601	ADP	N3-C4-N9	5.32	135.04	125.43
4	A	604	K01	O21-C15-N14	-4.97	112.49	123.72
3	B	601	ADP	N3-C4-N9	4.97	134.40	125.43
4	A	604	K01	C9-C8-C7	-4.94	103.87	112.85
4	A	604	K01	C19-C18-S17	-4.93	107.87	113.18
4	A	604	K01	C1-C2-N3	-4.82	90.63	111.68
4	B	604	K01	C6-C5-C4	-4.42	118.07	126.30
4	B	604	K01	C5-C13-S12	4.24	113.38	111.08
4	B	604	K01	C1-C2-N3	-4.00	94.22	111.68
3	A	601	ADP	C8-N9-C4	3.97	109.93	106.90
3	B	601	ADP	C8-N9-C4	3.92	109.89	106.90
4	A	604	K01	C18-S17-C16	3.74	93.56	91.91
4	B	604	K01	C19-C18-S17	-3.64	109.26	113.18
4	B	604	K01	O21-C15-N14	-3.64	115.50	123.72
4	B	604	K01	C20-C16-S17	2.86	114.44	111.02
3	A	601	ADP	O3B-PB-O2B	2.85	118.71	107.61
3	B	601	ADP	C4-C5-N7	-2.84	107.09	109.52
4	B	604	K01	C8-C7-C6	2.77	128.30	123.61
4	A	604	K01	C8-C7-C6	2.74	128.25	123.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	604	K01	C2-N3-C23	-2.68	109.20	116.34
3	A	601	ADP	C5-C4-N3	-2.66	119.90	125.70
3	B	601	ADP	C5-C4-N3	-2.55	120.15	125.70
4	B	604	K01	C24-C23-N3	-2.38	101.27	111.68
3	B	601	ADP	C2-N1-C6	2.28	122.88	118.77
4	B	604	K01	C15-C16-S17	-2.24	112.22	120.49
3	A	601	ADP	C2-N3-C4	2.17	120.20	114.01
3	A	601	ADP	C4-C5-N7	-2.14	107.69	109.52
4	B	604	K01	O22-C4-C5	2.11	122.78	118.86
4	A	604	K01	C9-C10-C11	2.09	122.06	112.65
3	A	601	ADP	C2-N1-C6	2.08	122.53	118.77
4	A	604	K01	O21-C15-C16	2.08	130.72	118.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	331/368 (89%)	0.59	29 (8%) 10 9	13, 28, 50, 60	0
1	B	331/368 (89%)	0.39	23 (6%) 17 14	15, 30, 52, 61	0
All	All	662/736 (89%)	0.49	52 (7%) 13 11	13, 29, 51, 61	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	16	GLY	9.6
1	A	56	GLY	9.3
1	A	57	LEU	9.1
1	A	58	ALA	8.0
1	A	250	ILE	6.9
1	B	58	ALA	6.7
1	B	57	LEU	6.5
1	A	192	ARG	5.9
1	A	249	THR	5.9
1	A	16	GLY	5.3
1	A	190	ASN	4.9
1	A	60	LYS	4.6
1	B	38	HIS	4.3
1	A	36	SER	4.3
1	A	149	ASP	4.2
1	B	17	LYS	4.1
1	A	59	ASP	4.1
1	B	34	LYS	4.0
1	A	189	ARG	3.9
1	B	33	ARG	3.5
1	A	151	GLY	3.5
1	B	192	ARG	3.5
1	A	248	THR	3.5
1	A	191	LYS	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	34	LYS	3.3
1	A	360	LEU	3.3
1	A	255	LEU	3.2
1	B	361	ASN	3.2
1	B	56	GLY	3.1
1	A	253	GLU	3.1
1	B	177	ASP	2.9
1	B	31	ALA	2.9
1	B	190	ASN	2.9
1	A	256	VAL	2.8
1	B	35	ALA	2.6
1	A	251	ASP	2.6
1	A	195	ILE	2.5
1	B	195	ILE	2.4
1	B	305	ARG	2.4
1	A	150	ASN	2.4
1	B	175	SER	2.3
1	B	59	ASP	2.2
1	B	97	TYR	2.2
1	B	362	LYS	2.2
1	B	307	PRO	2.1
1	A	254	GLU	2.1
1	A	97	TYR	2.1
1	A	61	SER	2.1
1	A	153	GLU	2.1
1	A	252	GLY	2.0
1	B	191	LYS	2.0
1	B	327	ARG	2.0

## 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 ⓘ

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<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	K01	B	604	24/24	0.16	3.42	17,24,28,33	0
4	K01	A	604	24/24	0.16	1.41	14,24,28,32	0
2	MG	B	603	1/1	0.14	0.66	17,17,17,17	0
3	ADP	A	601	27/27	0.12	-0.05	13,22,25,27	0
3	ADP	B	601	27/27	0.10	-0.36	15,28,32,33	0
2	MG	A	603	1/1	0.12	-1.29	16,16,16,16	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.