



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:50 PM GMT

PDB ID : 4JN4
Title : Allosteric opening of the polypeptide-binding site when an Hsp70 binds ATP
Authors : Qi, R.; Sarbeng, E.B.; Liu, Q.; Le, K.Q.; Xu, X.; Xu, H.; Yang, J.; Wong, J.L.;
Vorvis, C.; Hendrickson, W.A.; Zhou, L.; Liu, Q.
Deposited on : 2013-03-14
Resolution : 2.30 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 ⓘ symbol.

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

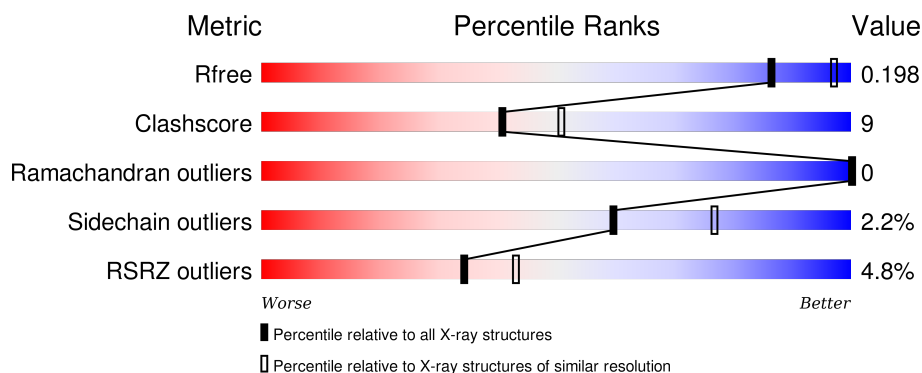
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 $\leq 5\%$. 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	608	<div> <div>4%</div> <div>85%</div> <div>13%</div> <div>..</div> </div>
1	B	608	<div> <div>6%</div> <div>83%</div> <div>14%</div> <div>..</div> </div>

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
4	SO4	A	1003	-	-	-	X
4	SO4	A	1006	-	-	-	X
4	SO4	B	1006	-	-	-	X
5	GOL	A	1009	-	-	-	X
5	GOL	B	1010	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10366 atoms, of which 32 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 Chaperone protein DnaK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	600	Total	C	N	O	S	0	0	0
			4550	2828	791	915	16			
1	B	600	Total	C	N	O	S	0	0	0
			4550	2828	791	915	16			

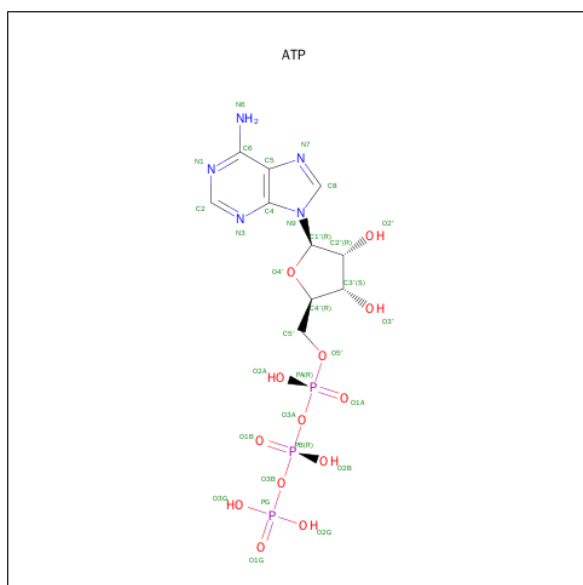
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	EXPRESSION TAG	UNP P0A6Y8
A	199	ALA	THR	CONFLICT	UNP P0A6Y8
A	?	-	THR	DELETION	UNP P0A6Y8
A	?	-	ALA	DELETION	UNP P0A6Y8
A	?	-	GLU	DELETION	UNP P0A6Y8
A	?	-	ASP	DELETION	UNP P0A6Y8
A	428	MET	ASN	ENGINEERED MUTATION	UNP P0A6Y8
A	429	GLY	GLN	ENGINEERED MUTATION	UNP P0A6Y8
A	430	GLY	SER	ENGINEERED MUTATION	UNP P0A6Y8
A	611	HIS	-	EXPRESSION TAG	UNP P0A6Y8
A	612	ALA	-	EXPRESSION TAG	UNP P0A6Y8
B	1	SER	-	EXPRESSION TAG	UNP P0A6Y8
B	199	ALA	THR	CONFLICT	UNP P0A6Y8
B	?	-	THR	DELETION	UNP P0A6Y8
B	?	-	ALA	DELETION	UNP P0A6Y8
B	?	-	GLU	DELETION	UNP P0A6Y8
B	?	-	ASP	DELETION	UNP P0A6Y8
B	428	MET	ASN	ENGINEERED MUTATION	UNP P0A6Y8
B	429	GLY	GLN	ENGINEERED MUTATION	UNP P0A6Y8
B	430	GLY	SER	ENGINEERED MUTATION	UNP P0A6Y8
B	611	HIS	-	EXPRESSION TAG	UNP P0A6Y8
B	612	ALA	-	EXPRESSION TAG	UNP P0A6Y8

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

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

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O P 31 10 5 13 3	0	0
3	B	1	Total C N O P 31 10 5 13 3	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	1
			10	8	2		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			14	3	8	3		
5	A	1	Total	C	H	O	0	0
			14	3	8	3		
5	B	1	Total	C	H	O	0	0
			14	3	8	3		
5	B	1	Total	C	H	O	0	0
			14	3	8	3		

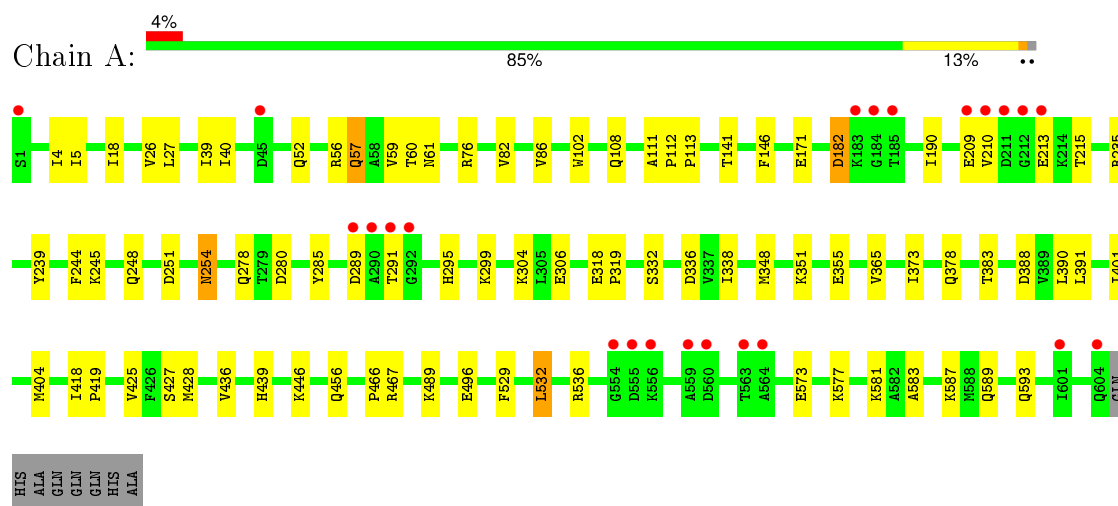
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	526	Total	O	0	0
			526	526		
6	B	560	Total	O	0	0
			560	560		

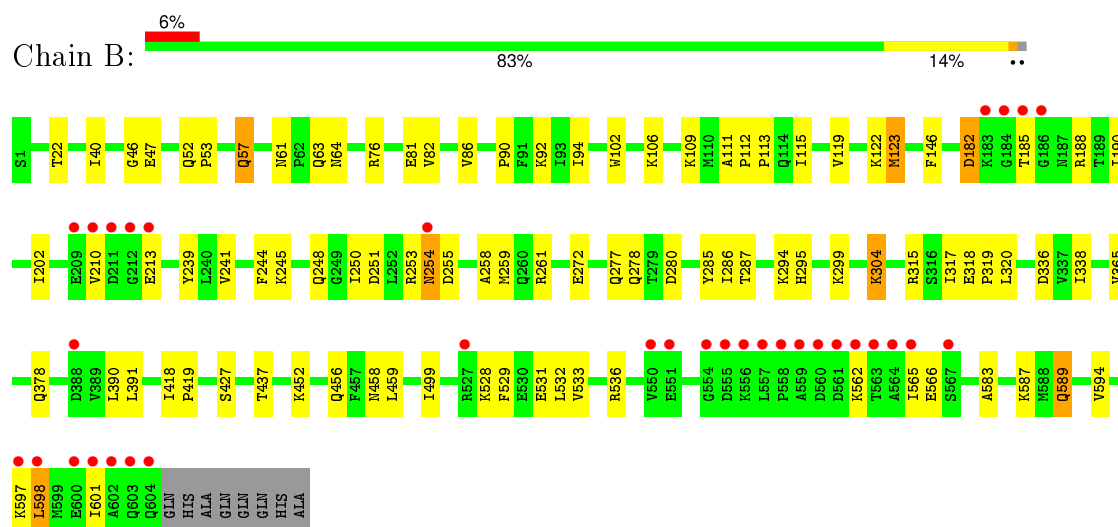
3 Residue-property plots

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: Chaperone protein DnaK



• Molecule 1: Chaperone protein DnaK



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	291.74Å 291.74Å 99.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.85 – 2.30 39.85 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.85-2.30) 100.0 (39.85-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.06 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.162 , 0.197 0.165 , 0.198	Depositor DCC
R_{free} test set	4740 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.382	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 94523 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10366	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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: GOL, MG, SO4, 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	0.29	0/4602	0.45	0/6217
1	B	0.29	0/4602	0.46	0/6217
All	All	0.29	0/9204	0.45	0/12434

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	4550	0	4610	67	0
1	B	4550	0	4610	94	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	31	0	12	0	0
3	B	31	0	12	0	0
4	A	25	0	0	1	0
4	B	35	0	0	0	0
5	A	12	16	16	3	0
5	B	12	16	16	4	0
6	A	526	0	0	6	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	560	0	0	12	0
All	All	10334	32	9276	159	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 9.

All (159) 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:A:285:TYR:H	1:A:295:HIS:HD2	1.14	0.95
1:A:278:GLN:OE1	1:B:533:VAL:HG22	1.72	0.88
1:A:254:ASN:HD22	1:A:254:ASN:H	1.26	0.84
1:A:285:TYR:H	1:A:295:HIS:CD2	1.96	0.83
1:B:57:GLN:HE22	1:B:61:ASN:HD22	1.22	0.83
1:B:280:ASP:OD1	1:B:299:LYS:HG2	1.79	0.82
1:B:531:GLU:HG3	6:B:1606:HOH:O	1.77	0.82
1:B:64:ASN:HD22	1:B:106:LYS:HE2	1.47	0.79
1:B:123:MET:SD	6:B:1208:HOH:O	2.40	0.79
1:B:255:ASP:HB3	6:B:1572:HOH:O	1.82	0.77
1:B:123:MET:HA	1:B:123:MET:HE2	1.67	0.76
1:B:63:GLN:HA	6:B:1539:HOH:O	1.85	0.76
1:B:285:TYR:H	1:B:295:HIS:CD2	2.05	0.74
1:A:489:LYS:HE2	6:A:1549:HOH:O	1.87	0.74
1:B:123:MET:HE3	1:B:123:MET:N	2.04	0.73
1:B:437:THR:HG23	1:B:458:ASN:HD21	1.52	0.73
1:A:57:GLN:HE22	1:A:61:ASN:HD22	1.36	0.72
1:B:64:ASN:ND2	1:B:106:LYS:HE2	2.03	0.72
1:A:59:VAL:HG13	1:A:60:THR:HG23	1.74	0.69
1:A:52:GLN:HG2	6:A:1591:HOH:O	1.92	0.68
1:B:210:VAL:O	1:B:213:GLU:HG2	1.94	0.68
1:B:562:LYS:O	1:B:566:GLU:HG3	1.93	0.68
1:B:123:MET:HA	1:B:123:MET:CE	2.24	0.67
1:A:304:LYS:HA	1:A:304:LYS:HE3	1.76	0.67
1:A:280:ASP:OD1	1:A:299:LYS:HG2	1.96	0.65
1:A:27:LEU:HA	5:A:1008:GOL:H12	1.79	0.64
1:B:57:GLN:NE2	1:B:61:ASN:HD22	1.95	0.63
1:A:40:ILE:HD12	1:A:40:ILE:N	2.14	0.63
1:B:255:ASP:O	1:B:259:MET:HG2	2.01	0.61
1:B:532:LEU:CD1	1:B:536:ARG:HE	2.14	0.61
1:A:209:GLU:O	1:A:209:GLU:HG3	1.99	0.60
1:B:123:MET:CE	1:B:123:MET:CA	2.79	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:258:ALA:HB3	6:B:1572:HOH:O	2.02	0.59
1:A:285:TYR:N	1:A:295:HIS:HD2	1.93	0.59
1:B:40:ILE:N	1:B:40:ILE:HD12	2.17	0.59
1:A:425:VAL:HG11	1:A:428:MET:HE1	1.85	0.58
1:A:278:GLN:NE2	1:A:299:LYS:HD2	2.18	0.58
1:B:437:THR:HG23	1:B:458:ASN:ND2	2.18	0.58
1:B:532:LEU:HD11	1:B:536:ARG:NH2	2.19	0.58
1:B:278:GLN:NE2	1:B:299:LYS:HD2	2.19	0.57
1:B:315:ARG:HH22	5:B:1010:GOL:H12	1.69	0.57
1:B:598:LEU:HD12	1:B:598:LEU:O	2.04	0.57
1:B:248:GLN:HE21	1:B:294:LYS:HD2	1.70	0.57
1:B:94:ILE:CD1	1:B:109:LYS:HG2	2.35	0.57
1:B:81:GLU:HG2	6:B:1537:HOH:O	2.05	0.56
1:B:122:LYS:C	1:B:123:MET:HE3	2.26	0.55
1:B:437:THR:HA	1:B:458:ASN:HD22	1.72	0.55
1:B:532:LEU:CD1	1:B:536:ARG:NE	2.70	0.55
1:B:239:TYR:CZ	1:B:304:LYS:HD3	2.41	0.55
1:B:299:LYS:HE2	6:B:1559:HOH:O	2.06	0.54
1:A:108:GLN:HG2	6:A:1571:HOH:O	2.08	0.54
1:B:245:LYS:HE2	1:B:251:ASP:CG	2.29	0.54
1:A:235:ARG:HH21	5:A:1009:GOL:H2	1.72	0.53
1:A:573:GLU:O	1:A:577:LYS:HD3	2.08	0.53
1:B:244:PHE:CE1	1:B:248:GLN:HG3	2.44	0.53
1:B:90:PRO:C	6:B:1539:HOH:O	2.48	0.52
1:A:210:VAL:O	1:A:213:GLU:HG2	2.09	0.52
1:A:338:ILE:HD12	1:A:365:VAL:HG21	1.92	0.52
1:A:244:PHE:CE1	1:A:248:GLN:HG3	2.44	0.52
1:A:254:ASN:N	1:A:254:ASN:HD22	1.96	0.51
1:B:64:ASN:HD21	1:B:106:LYS:HG3	1.75	0.51
1:B:90:PRO:HB2	6:B:1539:HOH:O	2.09	0.51
1:A:56:ARG:NH1	1:B:272:GLU:OE1	2.43	0.51
1:B:532:LEU:HD11	1:B:536:ARG:HH21	1.75	0.51
1:A:141:THR:HB	1:A:171:GLU:HG3	1.91	0.51
1:B:528:LYS:C	6:B:1606:HOH:O	2.49	0.51
1:A:306:GLU:HB3	1:A:348:MET:CE	2.40	0.51
1:B:245:LYS:HE2	1:B:251:ASP:CB	2.41	0.51
1:A:427:SER:HB2	1:A:436:VAL:CG2	2.41	0.51
1:B:285:TYR:H	1:B:295:HIS:HD2	1.56	0.51
1:B:46:GLY:HA2	6:B:1596:HOH:O	2.11	0.50
1:B:315:ARG:HH12	5:B:1010:GOL:H32	1.76	0.50
1:A:351:LYS:O	1:A:355:GLU:HG3	2.10	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:ILE:HD12	1:B:294:LYS:HG3	1.93	0.50
1:B:94:ILE:HD11	1:B:109:LYS:HG2	1.94	0.50
1:B:315:ARG:HH22	5:B:1010:GOL:C1	2.25	0.50
1:A:245:LYS:NZ	1:A:251:ASP:HB2	2.27	0.50
1:A:112:PRO:HB2	1:A:113:PRO:HD3	1.93	0.50
1:B:241:VAL:HG11	1:B:253:ARG:HG3	1.94	0.50
1:B:52:GLN:HB3	1:B:53:PRO:HD3	1.93	0.49
1:B:112:PRO:HB2	1:B:113:PRO:HD3	1.95	0.49
1:B:188:ARG:HD3	1:B:336:ASP:OD2	2.13	0.49
1:B:254:ASN:HD22	1:B:255:ASP:N	2.11	0.49
1:A:57:GLN:NE2	1:A:61:ASN:HD22	2.06	0.49
1:B:565:ILE:HD13	1:B:594:VAL:HG12	1.95	0.49
1:A:76:ARG:NH1	6:A:1303:HOH:O	2.45	0.48
1:B:82:VAL:O	1:B:86:VAL:HG23	2.13	0.48
1:A:532:LEU:CD1	1:A:536:ARG:NE	2.77	0.48
1:B:589:GLN:HA	1:B:589:GLN:HE21	1.78	0.48
1:A:532:LEU:HD13	1:A:532:LEU:O	2.13	0.47
1:A:391:LEU:O	1:A:391:LEU:HD12	2.14	0.47
1:B:598:LEU:HD12	1:B:598:LEU:C	2.34	0.47
1:A:26:VAL:HG23	1:A:373:ILE:HD11	1.95	0.47
1:A:254:ASN:ND2	1:A:254:ASN:H	2.04	0.47
1:B:40:ILE:HD11	1:B:119:VAL:HG23	1.95	0.47
1:B:123:MET:N	1:B:123:MET:CE	2.77	0.47
1:A:306:GLU:HB3	1:A:348:MET:HE1	1.96	0.47
1:A:190:ILE:HG22	1:A:336:ASP:HB2	1.97	0.47
1:B:261:ARG:NH1	1:B:286:ILE:HA	2.29	0.47
1:B:102:TRP:CE2	1:B:111:ALA:HB2	2.51	0.46
1:A:418:ILE:HA	1:A:419:PRO:C	2.36	0.46
1:A:5:ILE:HG22	1:A:18:ILE:HG22	1.98	0.46
1:A:536:ARG:HD3	6:A:1446:HOH:O	2.14	0.46
1:A:532:LEU:HD23	1:A:581:LYS:HB2	1.97	0.46
1:B:459:LEU:HD21	1:B:499:ILE:HG12	1.98	0.46
1:A:39:ILE:C	1:A:40:ILE:HD12	2.37	0.46
1:B:255:ASP:C	1:B:255:ASP:OD1	2.55	0.45
1:A:589:GLN:HE21	1:A:593:GLN:HG3	1.80	0.45
1:A:245:LYS:HZ2	1:A:251:ASP:HB2	1.82	0.45
1:B:102:TRP:CZ2	1:B:111:ALA:HB2	2.51	0.45
1:A:4:ILE:HD12	1:A:383:THR:HG23	1.99	0.45
1:B:254:ASN:C	1:B:254:ASN:ND2	2.69	0.45
1:B:245:LYS:HE2	1:B:251:ASP:HB2	1.99	0.45
1:B:532:LEU:O	1:B:536:ARG:HG3	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:390:LEU:HD23	1:B:391:LEU:N	2.31	0.45
1:B:378:GLN:OE1	1:B:391:LEU:HD21	2.16	0.45
1:A:583:ALA:O	1:A:587:LYS:HG2	2.17	0.45
1:B:315:ARG:HH22	5:B:1010:GOL:H2	1.82	0.45
1:B:202:ILE:HG12	1:B:320:LEU:HD23	1.99	0.45
1:A:239:TYR:CZ	1:A:304:LYS:HD3	2.52	0.44
1:B:64:ASN:ND2	1:B:106:LYS:CE	2.77	0.44
1:A:235:ARG:HH21	5:A:1009:GOL:C2	2.31	0.44
1:A:82:VAL:O	1:A:86:VAL:HG23	2.18	0.44
1:B:182:ASP:OD1	1:B:182:ASP:N	2.51	0.44
1:B:64:ASN:ND2	6:B:1531:HOH:O	2.51	0.43
1:B:533:VAL:HA	1:B:536:ARG:HD2	2.00	0.43
1:B:589:GLN:NE2	1:B:589:GLN:HA	2.33	0.43
1:B:583:ALA:O	1:B:587:LYS:HG2	2.19	0.43
1:B:57:GLN:HE22	1:B:61:ASN:ND2	2.01	0.43
1:A:532:LEU:HD12	1:A:536:ARG:HE	1.84	0.43
1:A:278:GLN:HE22	1:A:299:LYS:HD2	1.81	0.43
1:A:391:LEU:HD12	1:A:391:LEU:C	2.39	0.43
1:A:532:LEU:CD1	1:A:536:ARG:HE	2.31	0.43
1:A:378:GLN:OE1	1:A:391:LEU:HD21	2.19	0.43
1:B:76:ARG:NH2	1:B:452:LYS:HZ2	2.16	0.43
1:B:338:ILE:HD12	1:B:365:VAL:HG21	2.01	0.42
1:A:182:ASP:N	1:A:182:ASP:OD1	2.52	0.42
1:B:182:ASP:O	1:B:185:THR:HG23	2.19	0.42
1:A:318:GLU:HB2	1:A:319:PRO:HD3	2.01	0.42
1:A:102:TRP:CZ2	1:A:111:ALA:HB2	2.55	0.42
1:B:418:ILE:HA	1:B:419:PRO:C	2.39	0.42
1:A:215:THR:HA	1:A:390:LEU:O	2.20	0.42
1:B:119:VAL:O	1:B:123:MET:HG2	2.20	0.42
1:B:190:ILE:HG22	1:B:336:ASP:HB2	2.02	0.42
1:A:289:ASP:HB3	1:A:291:THR:OG1	2.20	0.42
1:A:141:THR:HB	1:A:171:GLU:CG	2.50	0.41
1:B:254:ASN:ND2	1:B:255:ASP:N	2.68	0.41
1:A:213:GLU:HA	1:A:388:ASP:HB3	2.02	0.41
1:B:112:PRO:N	1:B:113:PRO:CD	2.83	0.41
1:A:466:PRO:O	1:A:467:ARG:HB2	2.20	0.41
1:B:286:ILE:HG23	1:B:287:THR:HG22	2.03	0.41
1:A:446:LYS:HG3	4:A:1004:SO4:O3	2.21	0.41
1:A:496:GLU:HG2	6:A:1452:HOH:O	2.19	0.41
1:B:318:GLU:HB2	1:B:319:PRO:HD3	2.03	0.41
1:B:597:LYS:O	1:B:601:ILE:HG13	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:GLU:HB2	1:B:529:PHE:CZ	2.56	0.41
1:B:317:ILE:O	1:B:320:LEU:HB2	2.22	0.40
1:B:40:ILE:HD13	1:B:115:ILE:HG22	2.03	0.40
1:A:401:ILE:HA	1:A:439:HIS:O	2.21	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	598/608 (98%)	591 (99%)	7 (1%)	0	100	100
1	B	598/608 (98%)	592 (99%)	6 (1%)	0	100	100
All	All	1196/1216 (98%)	1183 (99%)	13 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	489/495 (99%)	480 (98%)	9 (2%)	66	82
1	B	489/495 (99%)	476 (97%)	13 (3%)	52	70
All	All	978/990 (99%)	956 (98%)	22 (2%)	60	77

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

Mol	Chain	Res	Type
1	A	57	GLN
1	A	146	PHE
1	A	182	ASP
1	A	254	ASN
1	A	332	SER
1	A	404	MET
1	A	456	GLN
1	A	529	PHE
1	A	532	LEU
1	B	22	THR
1	B	57	GLN
1	B	92	LYS
1	B	123	MET
1	B	146	PHE
1	B	182	ASP
1	B	254	ASN
1	B	277	GLN
1	B	304	LYS
1	B	427	SER
1	B	456	GLN
1	B	589	GLN
1	B	598	LEU

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

Mol	Chain	Res	Type
1	A	44	GLN
1	A	57	GLN
1	A	254	ASN
1	A	277	GLN
1	A	295	HIS
1	A	458	ASN
1	A	589	GLN
1	B	57	GLN
1	B	64	ASN
1	B	248	GLN
1	B	254	ASN
1	B	277	GLN
1	B	278	GLN
1	B	295	HIS
1	B	458	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	589	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 20 ligands modelled in this entry, 2 are monoatomic - leaving 18 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	ATP	A	1002	2	24,33,33	0.96	1 (4%)	31,52,52	1.82	5 (16%)
4	SO4	A	1003	-	4,4,4	0.31	0	6,6,6	0.08	0
4	SO4	A	1004	-	4,4,4	0.30	0	6,6,6	0.14	0
4	SO4	A	1005	-	4,4,4	0.18	0	6,6,6	0.14	0
4	SO4	A	1006	-	4,4,4	0.23	0	6,6,6	0.08	0
4	SO4	A	1007	-	4,4,4	0.18	0	6,6,6	0.09	0
5	GOL	A	1008	-	5,5,5	0.33	0	5,5,5	0.29	0
5	GOL	A	1009	-	5,5,5	0.39	0	5,5,5	0.15	0
3	ATP	B	1002	2	24,33,33	0.93	1 (4%)	31,52,52	1.79	6 (19%)
4	SO4	B	1003	-	4,4,4	0.17	0	6,6,6	0.14	0
4	SO4	B	1004	-	4,4,4	0.24	0	6,6,6	0.14	0
4	SO4	B	1005	-	4,4,4	0.27	0	6,6,6	0.13	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	B	1006	-	4,4,4	0.22	0	6,6,6	0.13	0
4	SO4	B	1007	-	4,4,4	0.21	0	6,6,6	0.06	0
4	SO4	B	1008[A]	-	4,4,4	0.31	0	6,6,6	0.14	0
4	SO4	B	1008[B]	-	4,4,4	0.23	0	6,6,6	0.07	0
5	GOL	B	1009	-	5,5,5	0.36	0	5,5,5	0.22	0
5	GOL	B	1010	-	5,5,5	0.36	0	5,5,5	0.20	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
3	ATP	A	1002	2	-	0/18/38/38	0/3/3/3
4	SO4	A	1003	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1004	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1005	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1006	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1007	-	-	0/0/0/0	0/0/0/0
5	GOL	A	1008	-	-	0/4/4/4	0/0/0/0
5	GOL	A	1009	-	-	0/4/4/4	0/0/0/0
3	ATP	B	1002	2	-	0/18/38/38	0/3/3/3
4	SO4	B	1003	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1004	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1005	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1006	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1007	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1008[A]	-	-	0/0/0/0	0/0/0/0
4	SO4	B	1008[B]	-	-	0/0/0/0	0/0/0/0
5	GOL	B	1009	-	-	0/4/4/4	0/0/0/0
5	GOL	B	1010	-	-	0/4/4/4	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1002	ATP	C5-C4	2.85	1.46	1.40
3	A	1002	ATP	C5-C4	2.97	1.47	1.40

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1002	ATP	N3-C2-N1	-7.40	123.23	128.89

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1002	ATP	N3-C2-N1	-7.27	123.33	128.89
3	A	1002	ATP	C1'-N9-C4	-2.85	122.64	126.94
3	A	1002	ATP	C4-C5-N7	-2.76	106.94	109.48
3	B	1002	ATP	C4-C5-N7	-2.74	106.96	109.48
3	B	1002	ATP	C1'-N9-C4	-2.72	122.83	126.94
3	A	1002	ATP	PA-O3A-PB	-2.47	125.79	132.73
3	B	1002	ATP	PA-O3A-PB	-2.15	126.70	132.73
3	B	1002	ATP	O3G-PG-O2G	2.03	115.10	107.38
3	A	1002	ATP	C2-N1-C6	2.10	122.53	118.77
3	B	1002	ATP	C2-N1-C6	2.11	122.53	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1004	SO4	1	0
5	A	1008	GOL	1	0
5	A	1009	GOL	2	0
5	B	1010	GOL	4	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, 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	600/608 (98%)	-0.11	23 (3%) 44 53	10, 26, 63, 145	0
1	B	600/608 (98%)	0.00	34 (5%) 27 36	10, 25, 70, 145	0
All	All	1200/1216 (98%)	-0.05	57 (4%) 34 43	10, 26, 65, 145	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	210	VAL	17.4
1	A	210	VAL	15.7
1	B	211	ASP	10.2
1	A	211	ASP	7.5
1	A	183	LYS	7.1
1	B	601	ILE	7.1
1	B	183	LYS	6.8
1	A	212	GLY	5.2
1	B	560	ASP	5.2
1	B	557	LEU	5.1
1	B	556	LYS	5.0
1	B	604	GLN	4.8
1	B	555	ASP	4.8
1	B	559	ALA	4.5
1	A	185	THR	4.4
1	A	601	ILE	4.4
1	B	558	PRO	4.2
1	B	527	ARG	4.0
1	B	212	GLY	3.9
1	B	554	GLY	3.8
1	A	555	ASP	3.7
1	B	597	LYS	3.5
1	B	563	THR	3.5
1	B	564	ALA	3.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	209	GLU	3.4
1	B	184	GLY	3.3
1	A	1	SER	3.3
1	A	559	ALA	3.3
1	A	184	GLY	3.3
1	B	598	LEU	3.2
1	B	185	THR	3.2
1	B	561	ASP	3.2
1	B	186	GLY	3.1
1	B	551	GLU	3.1
1	A	213	GLU	3.0
1	A	556	LYS	2.9
1	B	565	ILE	2.9
1	B	213	GLU	2.8
1	B	562	LYS	2.7
1	B	567	SER	2.6
1	B	603	GLN	2.5
1	A	290	ALA	2.5
1	B	209	GLU	2.5
1	B	388	ASP	2.5
1	B	254	ASN	2.4
1	A	292	GLY	2.4
1	A	564	ALA	2.4
1	B	602	ALA	2.4
1	A	560	ASP	2.4
1	B	600	GLU	2.4
1	A	563	THR	2.4
1	A	45	ASP	2.3
1	A	604	GLN	2.3
1	A	554	GLY	2.3
1	A	289	ASP	2.2
1	B	550	VAL	2.1
1	A	291	THR	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, 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
4	SO4	A	1006	5/5	0.96	0.29	15.03	2,10,28,31	5
4	SO4	B	1006	5/5	0.96	0.20	9.43	11,15,35,41	5
5	GOL	A	1009	6/6	0.90	0.29	4.45	68,87,101,105	0
4	SO4	A	1003	5/5	0.99	0.18	3.12	19,26,34,37	5
5	GOL	B	1010	6/6	0.94	0.15	1.48	43,59,79,89	0
5	GOL	A	1008	6/6	0.94	0.14	1.18	25,49,59,67	0
4	SO4	A	1007	5/5	0.98	0.15	0.72	33,33,37,48	5
4	SO4	B	1007	5/5	0.98	0.16	0.70	30,35,43,49	5
5	GOL	B	1009	6/6	0.96	0.12	0.65	24,36,48,48	0
4	SO4	B	1005	5/5	0.91	0.24	0.62	26,37,37,44	5
3	ATP	A	1002	31/31	0.99	0.14	-0.21	6,15,23,25	0
3	ATP	B	1002	31/31	0.99	0.15	-0.21	6,12,17,18	0
4	SO4	A	1005	5/5	0.92	0.16	-0.41	48,55,58,66	5
4	SO4	B	1003	5/5	0.99	0.09	-0.86	30,34,38,42	0
4	SO4	B	1008[B]	5/5	0.96	0.15	-	28,32,38,39	5
4	SO4	B	1004	5/5	0.97	0.22	-	39,41,47,48	5
2	MG	B	1001	1/1	0.97	0.32	-	9,9,9,9	1
4	SO4	B	1008[A]	5/5	0.96	0.15	-	17,23,27,41	5
4	SO4	A	1004	5/5	0.92	0.29	-	15,16,25,34	5
2	MG	A	1001	1/1	0.99	0.31	-	13,13,13,13	1

6.5 Other polymers ⓘ

There are no such residues in this entry.