



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:12 AM GMT

PDB ID : 3DSR
Title : ADP in transition binding site in the subunit B of the energy converter A1Ao ATP synthase
Authors : Kumar, A.; Manimekalai, S.M.S.; Balakrishna, A.M.; Gruber, G.
Deposited on : 2008-07-14
Resolution : 2.70 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

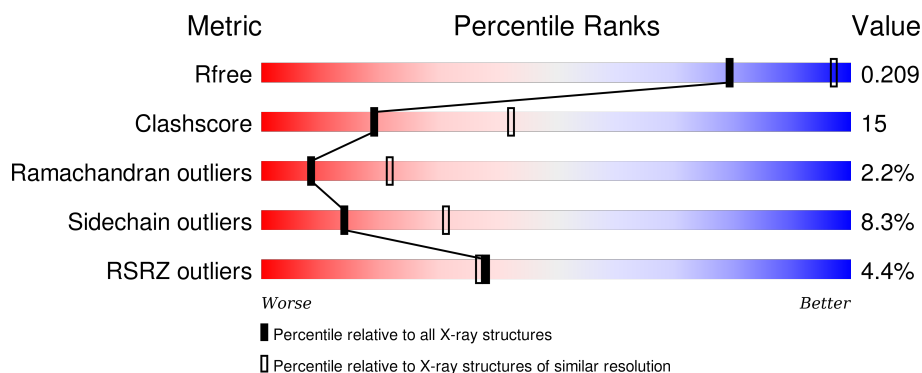
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	469	<div> <div>4%</div> <div> <div></div> <div>57%</div> <div>27%</div> <div>5%</div> <div>11%</div> </div> </div>
1	B	469	<div> <div>4%</div> <div> <div></div> <div>62%</div> <div>25%</div> <div>•</div> <div>9%</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
2	ADP	A	461	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6929 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 V-type ATP synthase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	0	0	0
			3225	2044	558	612	11			
1	B	425	Total	C	N	O	S	0	0	0
			3275	2076	569	619	11			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	MET	-	EXPRESSION TAG	UNP Q60187
A	-7	LYS	-	EXPRESSION TAG	UNP Q60187
A	-6	HIS	-	EXPRESSION TAG	UNP Q60187
A	-5	HIS	-	EXPRESSION TAG	UNP Q60187
A	-4	HIS	-	EXPRESSION TAG	UNP Q60187
A	-3	HIS	-	EXPRESSION TAG	UNP Q60187
A	-2	HIS	-	EXPRESSION TAG	UNP Q60187
A	-1	HIS	-	EXPRESSION TAG	UNP Q60187
A	0	PRO	-	EXPRESSION TAG	UNP Q60187
A	2	VAL	ALA	SEE REMARK 999	UNP Q60187
B	-8	MET	-	EXPRESSION TAG	UNP Q60187
B	-7	LYS	-	EXPRESSION TAG	UNP Q60187
B	-6	HIS	-	EXPRESSION TAG	UNP Q60187
B	-5	HIS	-	EXPRESSION TAG	UNP Q60187
B	-4	HIS	-	EXPRESSION TAG	UNP Q60187
B	-3	HIS	-	EXPRESSION TAG	UNP Q60187
B	-2	HIS	-	EXPRESSION TAG	UNP Q60187
B	-1	HIS	-	EXPRESSION TAG	UNP Q60187
B	0	PRO	-	EXPRESSION TAG	UNP Q60187
B	2	VAL	ALA	SEE REMARK 999	UNP Q60187

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

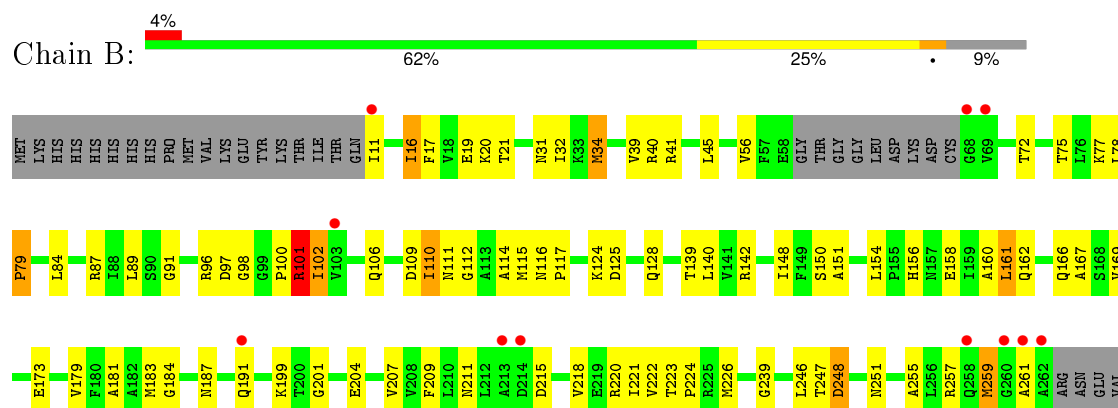


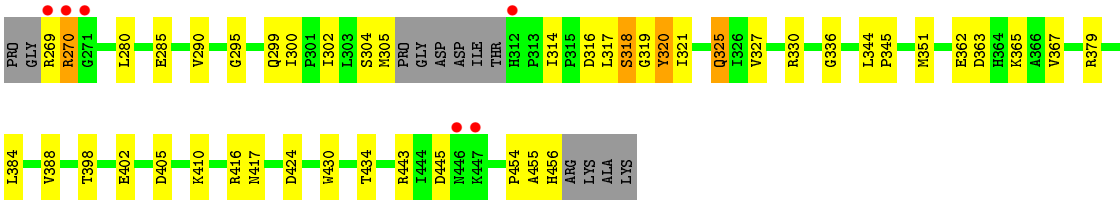
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	210	Total	O	0	0
			210	210		
3	B	192	Total	O	0	0
			192	192		

- Molecule 1: V-type ATP synthase beta chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.21Å 95.94Å 130.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.08 – 2.70 27.07 – 2.70	Depositor EDS
% Data completeness (in resolution range)	93.3 (27.08-2.70) 93.3 (27.07-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.198 , 0.267 0.205 , 0.209	Depositor DCC
R_{free} test set	1245 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	45.9	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 65.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 24439 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6929	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.54	0/3284	0.70	1/4448 (0.0%)
1	B	0.53	0/3335	0.66	1/4518 (0.0%)
All	All	0.53	0/6619	0.68	2/8966 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	ASP	N-CA-C	7.30	130.71	111.00
1	B	161	LEU	CA-CB-CG	7.22	131.90	115.30

There are no chirality outliers.

There are no planarity outliers.

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	3225	0	3239	114	0
1	B	3275	0	3287	88	0
2	A	27	0	12	8	0
3	A	210	0	0	8	0
3	B	192	0	0	8	0
All	All	6929	0	6538	199	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 15.

All (199) 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:79:PRO:HG2	1:A:102:ILE:HG12	1.38	1.06
1:B:101:ARG:H	1:B:101:ARG:HE	1.17	0.89
1:A:365:LYS:HD3	3:A:559:HOH:O	1.73	0.89
1:B:218:VAL:HA	1:B:221:ILE:HD12	1.53	0.88
1:A:184:GLY:H	1:A:211:ASN:HD22	1.22	0.86
1:B:84:LEU:HD23	1:B:102:ILE:HD12	1.58	0.85
1:B:336:GLY:HA2	3:B:637:HOH:O	1.75	0.84
1:A:106:GLN:HE21	1:A:106:GLN:HA	1.42	0.83
1:A:396:ARG:HD2	1:A:437:PRO:HG2	1.59	0.82
1:A:48:SER:HB2	1:A:218:VAL:HG21	1.60	0.81
1:B:162:GLN:HE21	1:B:166:GLN:HE22	1.29	0.80
1:B:139:THR:HB	1:B:351:MET:HG3	1.64	0.79
1:B:247:THR:O	1:B:302:ILE:HB	1.82	0.79
1:B:151:ALA:HA	1:B:305:MET:H	1.47	0.79
1:B:110:ILE:HD11	1:B:226:MET:HG2	1.64	0.78
1:A:383:GLY:O	1:A:387:ILE:HG22	1.88	0.74
1:A:119:ALA:HB2	1:A:291:LYS:HA	1.69	0.73
1:B:16:ILE:HG13	1:B:17:PHE:N	2.04	0.72
1:B:184:GLY:HA3	1:B:251:ASN:HD22	1.55	0.71
1:B:187:ASN:O	1:B:191:GLN:HG2	1.89	0.71
1:B:211:ASN:OD1	1:B:220:ARG:HG3	1.92	0.69
1:A:275:TYR:C	1:A:275:TYR:HD2	1.95	0.69
2:A:461:ADP:C2	3:A:660:HOH:O	2.46	0.68
1:A:90:SER:HB2	1:A:96:ARG:HG3	1.76	0.67
2:A:461:ADP:O2'	1:B:345:PRO:HB2	1.94	0.67
1:B:150:SER:OG	1:B:156:HIS:HD2	1.77	0.66
1:A:275:TYR:C	1:A:275:TYR:CD2	2.66	0.66
1:B:87:ARG:NH2	1:B:100:PRO:O	2.29	0.66
1:A:342:ASN:HD22	1:B:318:SER:HB3	1.59	0.66
1:B:218:VAL:HA	1:B:221:ILE:CD1	2.26	0.66
1:A:347:LEU:HD21	1:A:349:ARG:HG3	1.79	0.65
1:A:275:TYR:HD2	1:A:276:MET:N	1.94	0.65
1:B:398:THR:O	1:B:402:GLU:HG3	1.98	0.64
1:B:101:ARG:H	1:B:101:ARG:NE	1.94	0.63
1:A:416:ARG:NH1	1:A:417:ASN:OD1	2.32	0.63
1:A:115:MET:CE	1:A:237:GLU:HG3	2.28	0.63
1:A:396:ARG:HE	1:A:440:GLN:CG	2.12	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:GLN:NE2	1:A:422:ILE:H	1.97	0.62
1:A:184:GLY:H	1:A:211:ASN:ND2	1.94	0.62
1:A:324:GLY:HA2	1:A:347:LEU:HD23	1.80	0.62
1:A:25:GLY:N	1:A:28:GLU:OE2	2.30	0.62
1:A:115:MET:HE3	1:A:237:GLU:HG3	1.82	0.61
1:A:404:ALA:O	1:A:408:GLU:HG3	2.00	0.61
1:A:278:THR:O	1:A:281:ALA:HB3	2.00	0.61
1:A:188:GLU:H	1:A:188:GLU:CD	2.04	0.60
1:A:314:ILE:HD12	1:A:315:PRO:HB3	1.84	0.60
1:B:150:SER:OG	1:B:156:HIS:CD2	2.55	0.59
1:B:31:ASN:HB3	1:B:39:VAL:CG1	2.33	0.59
1:A:80:ALA:HB3	1:A:106:GLN:HB2	1.83	0.59
1:A:387:ILE:HD11	1:B:112:GLY:C	2.22	0.59
1:B:220:ARG:HD2	1:B:255:ALA:HB2	1.85	0.59
1:A:88:ILE:HD13	1:A:194:MET:HG2	1.84	0.59
1:B:248:ASP:HB3	1:B:251:ASN:HB2	1.84	0.59
1:A:22:GLU:HB3	3:A:608:HOH:O	2.03	0.58
1:B:257:ARG:HH21	1:B:270:ARG:HA	1.68	0.58
1:A:162:GLN:HE21	1:A:166:GLN:HE22	1.50	0.58
1:A:151:ALA:HA	1:A:305:MET:H	1.69	0.58
1:A:450:GLN:HA	1:A:456:HIS:CD2	2.38	0.58
1:A:90:SER:O	1:A:92:SER:N	2.30	0.57
1:A:18:VAL:HG22	1:A:52:VAL:HB	1.86	0.57
1:A:101:ARG:HG3	1:A:102:ILE:N	2.19	0.57
1:B:31:ASN:HB3	1:B:39:VAL:HG11	1.85	0.57
1:A:48:SER:HB2	1:A:218:VAL:CG2	2.32	0.57
1:A:48:SER:OG	1:A:49:ALA:N	2.37	0.57
1:A:184:GLY:HA2	1:A:213:ALA:HA	1.87	0.57
1:A:128:GLN:HE21	1:A:422:ILE:H	1.52	0.57
1:B:173:GLU:HG2	3:B:730:HOH:O	2.04	0.57
2:A:461:ADP:O2'	1:B:327:VAL:HG21	2.05	0.56
1:A:29:ILE:HA	1:A:42:GLY:O	2.05	0.56
1:B:179:VAL:HG22	1:B:207:VAL:CG1	2.34	0.56
1:A:119:ALA:O	1:A:121:LEU:N	2.39	0.56
2:A:461:ADP:H2	3:A:660:HOH:O	1.86	0.55
1:A:147:PRO:HA	1:A:301:PRO:HD2	1.87	0.55
1:A:396:ARG:CD	1:A:437:PRO:HG2	2.32	0.55
1:A:145:LYS:HD3	1:A:284:TYR:O	2.07	0.55
1:A:396:ARG:HE	1:A:440:GLN:HG3	1.72	0.54
1:A:150:SER:O	1:A:304:SER:HA	2.07	0.54
1:B:321:ILE:CG2	1:B:325:GLN:HE21	2.20	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:430:TRP:O	1:B:434:THR:HG23	2.07	0.54
1:A:106:GLN:NE2	1:A:106:GLN:HA	2.17	0.54
1:A:342:ASN:ND2	1:B:318:SER:HB3	2.22	0.54
1:B:201:GLY:O	1:B:204:GLU:HG2	2.08	0.54
1:B:162:GLN:HE21	1:B:166:GLN:NE2	1.99	0.54
1:A:424:ASP:O	1:A:428:ILE:HG23	2.08	0.53
1:A:347:LEU:HG	1:A:348:SER:N	2.22	0.53
1:A:333:HIS:HD2	3:A:493:HOH:O	1.92	0.52
1:B:34:MET:SD	1:B:40:ARG:HD3	2.50	0.52
1:B:125:ASP:HB2	1:B:142:ARG:CZ	2.40	0.52
1:B:181:ALA:HB3	1:B:246:LEU:HD23	1.91	0.52
1:B:128:GLN:O	1:B:167:ALA:HA	2.11	0.51
1:A:246:LEU:CD1	1:A:299:GLN:HG2	2.41	0.51
1:B:17:PHE:H	1:B:111:ASN:HD21	1.58	0.51
1:A:314:ILE:HB	1:A:315:PRO:HA	1.91	0.51
1:B:321:ILE:HG21	1:B:325:GLN:HE21	1.75	0.51
1:A:321:ILE:HG23	1:A:324:GLY:O	2.10	0.51
1:B:96:ARG:NH2	1:B:215:ASP:OD1	2.44	0.50
1:A:120:ARG:O	1:A:120:ARG:HG3	2.12	0.49
1:A:396:ARG:HH21	1:A:440:GLN:NE2	2.10	0.49
1:A:28:GLU:O	1:A:43:GLN:HA	2.13	0.49
1:A:32:ILE:HD11	1:A:54:VAL:HG11	1.95	0.49
1:A:41:ARG:N	1:A:41:ARG:HD2	2.27	0.49
1:A:130:GLY:HA3	1:A:419:ASN:ND2	2.27	0.49
1:B:384:LEU:O	1:B:388:VAL:HG22	2.13	0.49
1:B:101:ARG:HE	1:B:101:ARG:N	1.99	0.48
1:A:18:VAL:CG2	1:A:52:VAL:HB	2.43	0.48
1:A:280:LEU:O	1:A:283:LEU:HB2	2.13	0.48
1:A:275:TYR:O	1:A:277:TYR:N	2.47	0.48
1:A:223:THR:HB	1:A:224:PRO:CD	2.44	0.48
1:A:217:ALA:C	1:A:219:GLU:N	2.65	0.48
1:A:387:ILE:HG13	1:B:111:ASN:O	2.14	0.48
1:A:438:GLU:OE2	1:A:456:HIS:HE1	1.97	0.48
1:B:417:ASN:HA	3:B:602:HOH:O	2.13	0.47
1:A:416:ARG:HH11	1:A:417:ASN:HD21	1.62	0.47
1:A:360:THR:OG1	1:A:361:ARG:N	2.47	0.47
1:A:166:GLN:HG2	1:A:419:ASN:CG	2.34	0.47
1:B:290:VAL:HG22	1:B:295:GLY:O	2.13	0.47
1:B:75:THR:O	1:B:77:LYS:HG3	2.13	0.47
1:A:97:ASP:C	1:A:99:GLY:H	2.17	0.47
1:B:299:GLN:NE2	1:B:320:TYR:OH	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:PHE:O	1:A:58:GLU:HG2	2.14	0.47
1:A:119:ALA:HB2	1:A:291:LYS:CA	2.41	0.46
1:A:416:ARG:HH11	1:A:417:ASN:ND2	2.13	0.46
1:B:98:GLY:HA2	3:B:650:HOH:O	2.13	0.46
1:B:17:PHE:CD1	1:B:222:VAL:HG12	2.49	0.46
1:A:156:HIS:CD2	1:A:302:ILE:HG21	2.50	0.46
1:A:173:GLU:OE2	1:A:294:LYS:HE3	2.15	0.46
1:A:92:SER:HB3	1:A:94:GLU:OE1	2.15	0.46
1:B:115:MET:HG2	3:B:684:HOH:O	2.15	0.46
1:B:89:LEU:HD12	1:B:209:PHE:CE1	2.50	0.46
1:B:109:ASP:HB3	1:B:114:ALA:HB2	1.97	0.46
1:A:16:ILE:HD11	1:A:32:ILE:HD13	1.97	0.46
1:A:305:MET:HB3	1:A:306:PRO:HA	1.98	0.46
1:A:33:LYS:HB2	1:A:74:GLU:HG3	1.98	0.46
1:A:430:TRP:O	1:A:434:THR:HG23	2.16	0.46
1:B:31:ASN:O	1:B:72:THR:HA	2.16	0.45
1:A:345:PRO:HB2	2:A:461:ADP:N6	2.31	0.45
1:B:221:ILE:HD11	1:B:259:MET:SD	2.57	0.45
1:B:363:ASP:O	1:B:367:VAL:HG23	2.17	0.45
1:B:91:GLY:O	1:B:226:MET:HG3	2.17	0.45
1:A:80:ALA:O	1:A:105:ASP:HB3	2.17	0.45
1:A:212:LEU:C	1:A:214:ASP:H	2.19	0.45
1:A:114:ALA:HB3	3:A:575:HOH:O	2.17	0.44
1:A:244:VAL:HB	1:A:299:GLN:HG3	1.98	0.44
1:A:78:LEU:HB2	1:A:226:MET:SD	2.57	0.44
1:A:228:LEU:HD21	1:A:299:GLN:CD	2.38	0.44
1:A:407:PHE:O	1:A:412:VAL:HG23	2.17	0.44
1:A:32:ILE:HA	1:A:73:GLY:O	2.18	0.44
1:B:416:ARG:HG2	1:B:416:ARG:HH11	1.82	0.43
1:B:11:ILE:HD11	1:B:261:ALA:CB	2.48	0.43
1:A:41:ARG:O	1:A:57:PHE:N	2.43	0.43
1:A:223:THR:HB	1:A:224:PRO:HD3	2.00	0.43
1:B:158:GLU:OE1	1:B:416:ARG:NE	2.51	0.43
1:A:32:ILE:HB	1:A:34:MET:HE2	2.01	0.43
1:A:87:ARG:NH2	1:A:97:ASP:OD1	2.46	0.43
1:B:154:LEU:HD21	1:B:330:ARG:HG3	2.00	0.43
1:B:280:LEU:HD11	1:B:317:LEU:CD2	2.49	0.43
1:A:399:LYS:NZ	3:A:593:HOH:O	2.51	0.43
1:A:331:GLU:HB3	1:B:316:ASP:OD2	2.19	0.43
1:B:115:MET:O	1:B:117:PRO:HD3	2.19	0.43
1:B:416:ARG:HH12	1:B:417:ASN:HD21	1.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:MET:CE	1:B:259:MET:HA	2.50	0.42
1:A:118:TYR:HE1	1:A:286:ARG:HA	1.84	0.42
1:B:87:ARG:NH1	1:B:97:ASP:OD2	2.52	0.42
1:A:294:LYS:O	1:A:294:LYS:HG3	2.19	0.42
1:A:152:SER:O	1:A:330:ARG:NH2	2.53	0.42
1:B:455:ALA:O	1:B:456:HIS:ND1	2.52	0.42
1:A:148:ILE:HG22	1:A:300:ILE:CG2	2.50	0.42
1:A:305:MET:SD	2:A:461:ADP:O3B	2.78	0.42
1:B:124:LYS:HD2	3:B:613:HOH:O	2.19	0.42
1:A:339:PRO:HA	1:A:340:PRO:HD3	1.91	0.41
1:B:344:LEU:HD12	1:B:379:ARG:HD3	2.03	0.41
1:B:223:THR:HB	1:B:224:PRO:HD3	2.01	0.41
1:B:148:ILE:HB	1:B:302:ILE:HD13	2.03	0.41
1:B:140:LEU:HD22	1:B:300:ILE:HD11	2.01	0.41
1:B:443:ARG:NH1	3:B:744:HOH:O	2.52	0.41
1:A:116:ASN:N	1:A:117:PRO:HD3	2.36	0.41
1:A:78:LEU:HA	1:A:79:PRO:HD3	1.75	0.41
1:B:410:LYS:HA	1:B:410:LYS:HD3	1.91	0.41
1:A:149:PHE:HE1	2:A:461:ADP:C2	2.38	0.41
1:B:204:GLU:H	1:B:204:GLU:HG2	1.60	0.41
1:B:78:LEU:HA	1:B:79:PRO:HD2	1.87	0.41
1:B:148:ILE:O	1:B:148:ILE:HG22	2.20	0.41
3:A:546:HOH:O	1:B:319:GLY:HA2	2.21	0.41
1:B:239:GLY:HA2	3:B:674:HOH:O	2.21	0.41
1:A:384:LEU:O	1:A:387:ILE:HG23	2.21	0.41
1:B:257:ARG:NH2	1:B:270:ARG:HA	2.35	0.41
1:A:130:GLY:O	1:A:414:GLN:NE2	2.52	0.41
1:B:201:GLY:O	1:B:204:GLU:CG	2.69	0.41
1:A:332:LEU:HA	1:A:332:LEU:HD23	1.86	0.41
1:A:396:ARG:NH2	1:A:440:GLN:NE2	2.69	0.40
1:A:31:ASN:C	1:A:32:ILE:HG13	2.40	0.40
1:A:80:ALA:HB1	1:A:233:TYR:CG	2.56	0.40
2:A:461:ADP:HO2'	1:B:345:PRO:HB2	1.85	0.40
1:B:32:ILE:O	1:B:39:VAL:HA	2.20	0.40
1:A:228:LEU:HD21	1:A:299:GLN:OE1	2.21	0.40
1:A:248:ASP:HB3	1:A:251:ASN:ND2	2.37	0.40
1:B:19:GLU:O	1:B:21:THR:N	2.54	0.40
1:B:221:ILE:O	1:B:224:PRO:HD2	2.21	0.40
1:B:160:ALA:HB2	1:B:302:ILE:HD11	2.03	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	411/469 (88%)	362 (88%)	35 (8%)	14 (3%)	5	10
1	B	417/469 (89%)	390 (94%)	23 (6%)	4 (1%)	19	45
All	All	828/938 (88%)	752 (91%)	58 (7%)	18 (2%)	8	22

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	ARG
1	A	216	PRO
1	A	276	MET
1	A	314	ILE
1	A	102	ILE
1	B	101	ARG
1	A	248	ASP
1	A	357	ALA
1	A	91	GLY
1	A	213	ALA
1	A	274	GLY
1	A	438	GLU
1	B	20	LYS
1	B	454	PRO
1	A	325	GLN
1	A	215	ASP
1	B	79	PRO
1	A	315	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	342/384 (89%)	314 (92%)	28 (8%)	14	32
1	B	346/384 (90%)	317 (92%)	29 (8%)	14	30
All	All	688/768 (90%)	631 (92%)	57 (8%)	14	31

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	38	THR
1	A	40	ARG
1	A	45	LEU
1	A	47	SER
1	A	56	VAL
1	A	78	LEU
1	A	94	GLU
1	A	103	VAL
1	A	106	GLN
1	A	148	ILE
1	A	173	GLU
1	A	188	GLU
1	A	204	GLU
1	A	248	ASP
1	A	250	THR
1	A	275	TYR
1	A	278	THR
1	A	299	GLN
1	A	320	TYR
1	A	331	GLU
1	A	347	LEU
1	A	365	LYS
1	A	387	ILE
1	A	396	ARG
1	A	419	ASN
1	A	428	ILE
1	A	431	GLN
1	B	16	ILE
1	B	34	MET
1	B	41	ARG
1	B	45	LEU
1	B	56	VAL

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Mol	Chain	Res	Type
1	B	101	ARG
1	B	102	ILE
1	B	106	GLN
1	B	110	ILE
1	B	116	ASN
1	B	161	LEU
1	B	169	VAL
1	B	183	MET
1	B	199	LYS
1	B	248	ASP
1	B	259	MET
1	B	269	ARG
1	B	270	ARG
1	B	285	GLU
1	B	304	SER
1	B	314	ILE
1	B	318	SER
1	B	320	TYR
1	B	325	GLN
1	B	362	GLU
1	B	365	LYS
1	B	405	ASP
1	B	424	ASP
1	B	445	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	116	ASN
1	A	128	GLN
1	A	138	ASN
1	A	162	GLN
1	A	211	ASN
1	A	299	GLN
1	A	325	GLN
1	A	333	HIS
1	A	342	ASN
1	A	419	ASN
1	A	440	GLN
1	A	450	GLN
1	A	456	HIS

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Mol	Chain	Res	Type
1	B	27	ASN
1	B	31	ASN
1	B	55	GLN
1	B	106	GLN
1	B	111	ASN
1	B	116	ASN
1	B	128	GLN
1	B	156	HIS
1	B	162	GLN
1	B	238	HIS
1	B	251	ASN
1	B	299	GLN
1	B	325	GLN
1	B	440	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](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ADP	A	461	-	22,29,29	0.94	1 (4%)	27,45,45	1.97	5 (18%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	A	461	-	-	0/12/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	461	ADP	C5-C4	2.79	1.46	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	461	ADP	N3-C2-N1	-6.50	123.92	128.89
2	A	461	ADP	C2'-C1'-N9	-3.79	108.50	114.29
2	A	461	ADP	PA-O3A-PB	-3.48	120.99	132.67
2	A	461	ADP	C4-C5-N7	-3.32	106.42	109.48
2	A	461	ADP	O4'-C1'-N9	2.62	113.58	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	461	ADP	8	0

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, 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	419/469 (89%)	0.08	20 (4%) 34 33	27, 40, 62, 74	0
1	B	425/469 (90%)	0.02	17 (4%) 42 41	27, 40, 61, 73	0
All	All	844/938 (89%)	0.05	37 (4%) 38 37	27, 40, 62, 74	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	314	ILE	5.7
1	B	11	ILE	4.7
1	B	262	ALA	4.5
1	B	269	ARG	4.2
1	A	102	ILE	4.1
1	A	273	PRO	3.6
1	A	72	THR	3.5
1	B	270	ARG	3.4
1	B	214	ASP	3.1
1	B	446	ASN	3.1
1	B	258	GLN	3.0
1	B	69	VAL	3.0
1	A	100	PRO	2.9
1	A	446	ASN	2.9
1	B	260	GLY	2.9
1	B	312	HIS	2.9
1	A	244	VAL	2.8
1	A	455	ALA	2.8
1	A	300	ILE	2.7
1	A	261	ALA	2.7
1	B	271	GLY	2.6
1	A	315	PRO	2.5
1	B	447	LYS	2.4
1	B	261	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	173	GLU	2.3
1	A	29	ILE	2.2
1	B	68	GLY	2.2
1	A	299	GLN	2.2
1	A	171	GLY	2.2
1	A	179	VAL	2.1
1	B	213	ALA	2.1
1	B	191	GLN	2.1
1	A	101	ARG	2.1
1	A	70	ILE	2.0
1	A	181	ALA	2.0
1	A	450	GLN	2.0
1	B	103	VAL	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(\AA^2)	Q<0.9
2	ADP	A	461	27/27	0.76	0.42	2.34	58,60,64,64	27

6.5 Other polymers [i](#)

There are no such residues in this entry.