



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:56 AM GMT

PDB ID : 3KAR
Title : THE MOTOR DOMAIN OF KINESIN-LIKE PROTEIN KAR3, A SACCHAROMYCES CEREVISIAE KINESIN-RELATED PROTEIN
Authors : Gulick, A.M.; Song, H.; Endow, S.; Rayment, I.
Deposited on : 1997-11-26
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

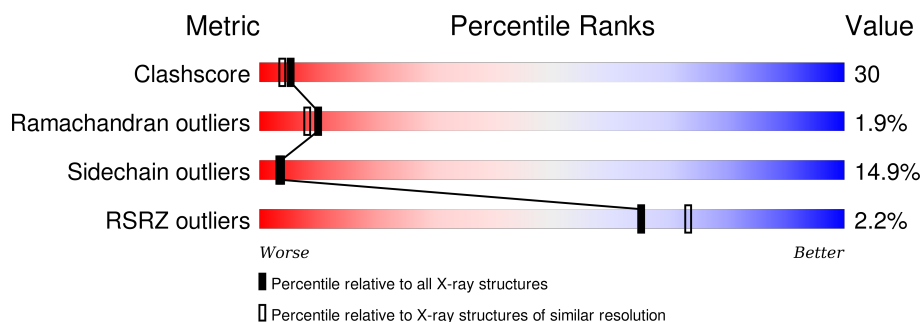
1 Overall quality at a glance

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))
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	346	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2608 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 KAR3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2475	1540	432	494	9			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
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		

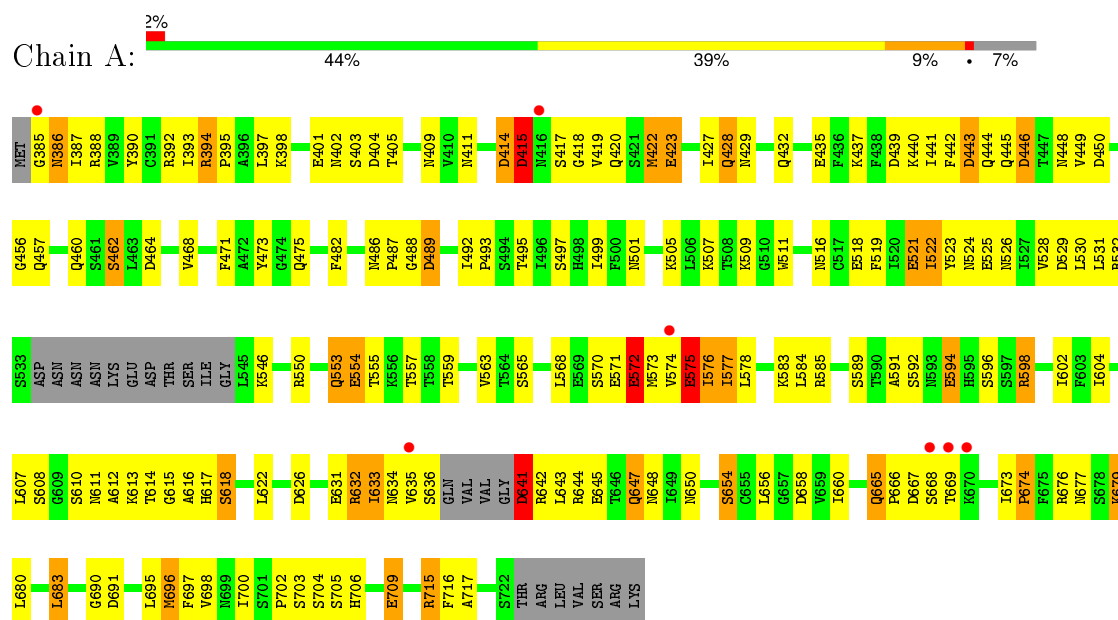
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	105	Total 105	O 105	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: KINESIN-LIKE PROTEIN KAR3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	44.13Å 81.21Å 48.30Å 90.00° 105.78° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 40.60 – 1.88	Depositor EDS
% Data completeness (in resolution range)	93.2 (30.00-2.30) 77.4 (40.60-1.88)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.56 (at 1.88Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.175 , (Not available) 0.168 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.573	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 123.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 22220 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2608	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.31% 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: 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	1.07	11/2514 (0.4%)	1.34	29/3407 (0.9%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	554	GLU	CD-OE2	7.07	1.33	1.25
1	A	594	GLU	CD-OE2	6.72	1.33	1.25
1	A	423	GLU	CD-OE1	6.53	1.32	1.25
1	A	521	GLU	CD-OE1	6.46	1.32	1.25
1	A	572	GLU	CD-OE1	6.37	1.32	1.25
1	A	518	GLU	CD-OE1	-6.13	1.19	1.25
1	A	435	GLU	CD-OE2	6.08	1.32	1.25
1	A	571	GLU	CD-OE2	5.70	1.31	1.25
1	A	709	GLU	CD-OE2	5.51	1.31	1.25
1	A	645	GLU	CD-OE2	5.15	1.31	1.25
1	A	518	GLU	CD-OE2	5.01	1.31	1.25

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	489	ASP	CB-CG-OD1	-10.76	108.62	118.30
1	A	443	ASP	CB-CG-OD2	-8.42	110.72	118.30
1	A	658	ASP	CB-CG-OD1	-8.21	110.91	118.30
1	A	691	ASP	CB-CG-OD1	-8.08	111.03	118.30
1	A	450	ASP	CB-CG-OD2	-7.96	111.14	118.30
1	A	529	ASP	CB-CG-OD1	7.93	125.44	118.30
1	A	667	ASP	CB-CG-OD1	-7.84	111.25	118.30
1	A	658	ASP	CB-CG-OD2	7.83	125.35	118.30
1	A	489	ASP	CB-CG-OD2	7.79	125.31	118.30
1	A	529	ASP	CB-CG-OD2	-7.73	111.34	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	443	ASP	CB-CG-OD1	7.40	124.96	118.30
1	A	667	ASP	CB-CG-OD2	7.01	124.61	118.30
1	A	439	ASP	CB-CG-OD1	-6.99	112.01	118.30
1	A	439	ASP	CB-CG-OD2	6.44	124.10	118.30
1	A	464	ASP	CB-CG-OD1	-6.44	112.50	118.30
1	A	450	ASP	CB-CG-OD1	6.38	124.04	118.30
1	A	414	ASP	CB-CG-OD1	-6.31	112.62	118.30
1	A	676	ARG	NE-CZ-NH1	6.31	123.45	120.30
1	A	626	ASP	CB-CG-OD1	6.27	123.94	118.30
1	A	691	ASP	CB-CG-OD2	6.09	123.78	118.30
1	A	464	ASP	CB-CG-OD2	5.85	123.56	118.30
1	A	585	ARG	CG-CD-NE	-5.84	99.55	111.80
1	A	641	ASP	CB-CG-OD2	-5.78	113.10	118.30
1	A	632	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	715	ARG	NE-CZ-NH1	5.33	122.96	120.30
1	A	446	ASP	CB-CG-OD2	-5.28	113.55	118.30
1	A	641	ASP	CB-CG-OD1	5.17	122.96	118.30
1	A	415	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	A	394	ARG	NE-CZ-NH1	5.03	122.81	120.30

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	2475	0	2360	147	1
2	A	1	0	0	0	0
3	A	27	0	12	0	0
4	A	105	0	0	14	0
All	All	2608	0	2372	147	1

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 30.

All (147) 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:572:GLU:HA	1:A:575:GLU:HG2	1.26	1.15
1:A:574:VAL:HG13	1:A:577:ILE:HD12	1.43	1.01
1:A:448:ASN:HD22	1:A:489:ASP:HB3	1.25	0.99
1:A:568:LEU:HD21	1:A:574:VAL:HG23	1.47	0.94
1:A:486:ASN:HD22	1:A:489:ASP:HB2	1.36	0.91
1:A:395:PRO:HA	1:A:444:GLN:HG2	1.51	0.90
1:A:568:LEU:CD2	1:A:574:VAL:HG23	2.05	0.86
1:A:392:ARG:NH1	4:A:823:HOH:O	2.10	0.85
1:A:486:ASN:HD22	1:A:489:ASP:CB	1.91	0.84
1:A:468:VAL:HB	1:A:622:LEU:HD23	1.59	0.83
1:A:524:ASN:OD1	1:A:648:ASN:ND2	2.12	0.81
1:A:633:ILE:HG23	1:A:635:VAL:HG23	1.63	0.80
1:A:574:VAL:HG12	1:A:578:LEU:HG	1.62	0.80
1:A:386:ASN:C	1:A:387:ILE:HD12	2.03	0.78
1:A:644:ARG:HA	1:A:647:GLN:NE2	1.99	0.78
1:A:419:VAL:HG11	1:A:437:LYS:HB3	1.64	0.77
1:A:550:ARG:HB2	1:A:559:THR:HG22	1.67	0.77
1:A:568:LEU:HD21	1:A:574:VAL:CG2	2.15	0.76
1:A:419:VAL:HG11	1:A:437:LYS:HD3	1.68	0.76
1:A:572:GLU:CA	1:A:575:GLU:HG2	2.11	0.75
1:A:475:GLN:HE22	1:A:706:HIS:HB3	1.50	0.75
1:A:419:VAL:HG12	1:A:420:GLN:N	2.02	0.75
1:A:387:ILE:O	1:A:388:ARG:HD2	1.86	0.74
1:A:486:ASN:ND2	1:A:489:ASP:HB2	2.02	0.74
1:A:387:ILE:N	1:A:387:ILE:HD12	2.03	0.74
1:A:456:GLY:O	1:A:460:GLN:HG3	1.88	0.74
1:A:643:LEU:C	1:A:647:GLN:HE21	1.91	0.73
1:A:570:SER:O	1:A:573:MET:HB2	1.89	0.72
1:A:698:VAL:HG21	1:A:717:ALA:HB2	1.71	0.71
1:A:634:ASN:ND2	1:A:636:SER:O	2.22	0.70
1:A:574:VAL:O	1:A:577:ILE:HG13	1.92	0.70
1:A:473:TYR:HB2	1:A:696:MET:CE	2.23	0.69
1:A:401:GLU:HB3	1:A:703:SER:OG	1.94	0.68
1:A:526:ASN:ND2	4:A:897:HOH:O	2.27	0.68
1:A:403:SER:HA	1:A:704:SER:OG	1.94	0.67
1:A:522:ILE:HG13	1:A:602:ILE:HD11	1.76	0.67
1:A:417:SER:O	1:A:419:VAL:HG23	1.95	0.67
1:A:394:ARG:O	1:A:702:PRO:HG3	1.96	0.65
1:A:572:GLU:O	1:A:575:GLU:HG3	1.97	0.64
1:A:574:VAL:CG1	1:A:578:LEU:HG	2.27	0.64
1:A:473:TYR:HB2	1:A:696:MET:HE2	1.78	0.64
1:A:422:MET:HG3	1:A:441:ILE:HD13	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:ASP:O	1:A:418:GLY:N	2.31	0.64
1:A:462:SER:HB3	1:A:607:LEU:HD13	1.80	0.63
1:A:419:VAL:CG1	1:A:420:GLN:N	2.62	0.62
1:A:674:PRO:HB3	1:A:677:ASN:ND2	2.14	0.62
1:A:449:VAL:HG13	4:A:824:HOH:O	2.00	0.62
1:A:555:THR:HG22	1:A:557:THR:HG23	1.81	0.62
1:A:505:LYS:O	1:A:505:LYS:HG2	1.98	0.62
1:A:501:ASN:HB2	4:A:896:HOH:O	1.99	0.62
1:A:635:VAL:HG12	1:A:635:VAL:O	1.99	0.61
1:A:473:TYR:CB	1:A:696:MET:HE2	2.31	0.60
1:A:419:VAL:HG11	1:A:437:LYS:CD	2.31	0.60
1:A:419:VAL:CG1	1:A:437:LYS:HB3	2.32	0.60
1:A:522:ILE:HG13	1:A:602:ILE:CD1	2.31	0.59
1:A:402:ASN:ND2	1:A:404:ASP:H	1.99	0.59
1:A:574:VAL:HA	1:A:577:ILE:HD11	1.83	0.59
1:A:633:ILE:HG23	1:A:635:VAL:CG2	2.32	0.59
1:A:448:ASN:ND2	1:A:489:ASP:HB3	2.09	0.58
1:A:444:GLN:NE2	4:A:843:HOH:O	2.29	0.58
1:A:492:ILE:HB	1:A:493:PRO:HD3	1.86	0.58
1:A:656:LEU:O	1:A:660:ILE:HG12	2.03	0.58
1:A:641:ASP:N	1:A:641:ASP:OD1	2.37	0.57
1:A:397:LEU:HB2	1:A:401:GLU:OE2	2.04	0.56
1:A:574:VAL:HA	1:A:577:ILE:CD1	2.36	0.56
1:A:643:LEU:O	1:A:647:GLN:HG3	2.06	0.56
1:A:665:GLN:O	1:A:668:SER:OG	2.25	0.54
1:A:495:THR:O	1:A:499:ILE:HD12	2.06	0.54
1:A:409:ASN:O	1:A:423:GLU:N	2.41	0.54
1:A:386:ASN:OD1	1:A:388:ARG:NH1	2.41	0.54
1:A:471:PHE:HB2	1:A:696:MET:HE3	1.89	0.54
1:A:427:ILE:O	1:A:429:ASN:N	2.40	0.54
1:A:482:PHE:O	1:A:486:ASN:HB2	2.08	0.53
1:A:419:VAL:HG11	1:A:437:LYS:CB	2.37	0.53
1:A:487:PRO:C	1:A:489:ASP:H	2.13	0.52
1:A:611:ASN:O	1:A:615:GLY:N	2.43	0.52
1:A:635:VAL:O	1:A:636:SER:HB3	2.10	0.52
1:A:411:ASN:ND2	4:A:902:HOH:O	2.40	0.51
1:A:635:VAL:HG13	1:A:643:LEU:HD13	1.93	0.51
1:A:519:PHE:HB3	1:A:531:LEU:HG	1.93	0.51
1:A:419:VAL:CG1	1:A:420:GLN:H	2.24	0.50
1:A:643:LEU:O	1:A:647:GLN:NE2	2.37	0.50
1:A:427:ILE:C	1:A:429:ASN:N	2.64	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:610:SER:HA	1:A:616:ALA:O	2.12	0.50
1:A:650:ASN:O	1:A:654:SER:HB3	2.12	0.50
1:A:568:LEU:HD22	1:A:574:VAL:HG23	1.93	0.49
1:A:387:ILE:N	1:A:387:ILE:CD1	2.75	0.49
1:A:607:LEU:N	1:A:607:LEU:HD12	2.28	0.49
1:A:523:TYR:O	1:A:524:ASN:HB2	2.12	0.49
1:A:511:TRP:NE1	4:A:906:HOH:O	2.23	0.49
1:A:521:GLU:HG2	1:A:528:VAL:HB	1.94	0.49
1:A:521:GLU:HB3	1:A:530:LEU:HD11	1.95	0.49
1:A:598:ARG:NH2	1:A:631:GLU:OE2	2.45	0.49
1:A:633:ILE:O	1:A:633:ILE:HG23	2.12	0.48
1:A:523:TYR:CZ	1:A:524:ASN:ND2	2.81	0.48
1:A:525:GLU:O	1:A:679:LYS:HD3	2.13	0.48
1:A:397:LEU:HD22	1:A:401:GLU:OE2	2.14	0.48
1:A:397:LEU:HB2	1:A:401:GLU:HG3	1.94	0.48
1:A:419:VAL:HG12	1:A:420:GLN:H	1.77	0.47
1:A:674:PRO:CB	1:A:677:ASN:HD22	2.26	0.47
1:A:395:PRO:HA	1:A:444:GLN:CG	2.35	0.47
1:A:611:ASN:CG	1:A:614:THR:OG1	2.54	0.47
1:A:574:VAL:HG12	1:A:578:LEU:CG	2.40	0.47
1:A:715:ARG:NE	4:A:840:HOH:O	2.28	0.47
1:A:427:ILE:CG2	1:A:428:GLN:N	2.78	0.46
1:A:521:GLU:O	1:A:521:GLU:HG3	2.14	0.46
1:A:680:LEU:HB3	4:A:877:HOH:O	2.15	0.46
1:A:565:SER:HB2	4:A:846:HOH:O	2.15	0.46
1:A:572:GLU:O	1:A:575:GLU:CG	2.63	0.45
1:A:578:LEU:HA	1:A:578:LEU:HD23	1.67	0.45
1:A:633:ILE:O	1:A:635:VAL:HG23	2.16	0.45
1:A:553:GLN:H	1:A:553:GLN:CD	2.20	0.45
1:A:574:VAL:CG1	1:A:577:ILE:HD12	2.30	0.45
1:A:644:ARG:CA	1:A:647:GLN:NE2	2.75	0.45
1:A:473:TYR:HB2	1:A:696:MET:HE1	1.98	0.44
1:A:612:ALA:C	1:A:614:THR:H	2.21	0.44
1:A:674:PRO:O	1:A:677:ASN:HB2	2.17	0.44
1:A:665:GLN:CG	1:A:666:PRO:HD2	2.48	0.44
1:A:443:ASP:C	1:A:443:ASP:OD1	2.57	0.43
1:A:402:ASN:HD22	1:A:404:ASP:H	1.62	0.43
1:A:427:ILE:C	1:A:429:ASN:H	2.22	0.43
1:A:583:LYS:O	1:A:591:ALA:HB2	2.19	0.43
1:A:584:LEU:O	1:A:591:ALA:HB1	2.18	0.43
1:A:709:GLU:HA	1:A:709:GLU:OE1	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:633:ILE:HD13	1:A:635:VAL:HG22	2.00	0.42
1:A:432:GLN:O	4:A:801:HOH:O	2.21	0.42
1:A:507:LYS:HG2	4:A:845:HOH:O	2.18	0.42
1:A:555:THR:HG22	1:A:557:THR:CG2	2.46	0.42
1:A:550:ARG:HD3	4:A:813:HOH:O	2.19	0.42
1:A:390:TYR:CD1	1:A:440:LYS:HB3	2.54	0.42
1:A:716:PHE:C	1:A:716:PHE:CD1	2.92	0.42
1:A:524:ASN:HA	1:A:648:ASN:ND2	2.35	0.42
1:A:443:ASP:OD1	1:A:445:GLN:N	2.36	0.41
1:A:574:VAL:O	1:A:576:ILE:N	2.54	0.41
1:A:471:PHE:CB	1:A:696:MET:HE3	2.50	0.41
1:A:674:PRO:CB	1:A:677:ASN:ND2	2.81	0.41
1:A:499:ILE:HG21	1:A:607:LEU:HD21	2.03	0.41
1:A:618:SER:HB3	4:A:809:HOH:O	2.19	0.41
1:A:393:ILE:HG12	1:A:700:ILE:HD11	2.01	0.41
1:A:405:THR:HB	1:A:704:SER:HB3	2.00	0.41
1:A:617:HIS:N	1:A:617:HIS:CD2	2.86	0.41
1:A:577:ILE:HG13	1:A:577:ILE:H	1.35	0.41
1:A:635:VAL:HG13	1:A:643:LEU:CD1	2.50	0.41
1:A:427:ILE:O	1:A:428:GLN:C	2.59	0.41
1:A:442:PHE:HA	1:A:446:ASP:OD2	2.20	0.41
1:A:679:LYS:O	1:A:683:LEU:HB2	2.22	0.40
1:A:385:GLY:N	1:A:690:GLY:O	2.55	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:GLN:OE1	1:A:563:VAL:O[1_454]	2.05	0.15

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	317/346 (92%)	287 (90%)	24 (8%)	6 (2%)	10 8

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	398	LYS
1	A	575	GLU
1	A	669	THR
1	A	428	GLN
1	A	488	GLY
1	A	613	LYS

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	275/316 (87%)	234 (85%)	41 (15%)	4 3

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	386	ASN
1	A	414	ASP
1	A	415	ASP
1	A	422	MET
1	A	457	GLN
1	A	462	SER
1	A	497	SER
1	A	509	LYS
1	A	516	ASN
1	A	522	ILE
1	A	532	ARG
1	A	546	LYS
1	A	553	GLN
1	A	554	GLU
1	A	572	GLU

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Mol	Chain	Res	Type
1	A	575	GLU
1	A	576	ILE
1	A	577	ILE
1	A	589	SER
1	A	592	SER
1	A	594	GLU
1	A	596	SER
1	A	598	ARG
1	A	604	ILE
1	A	608	SER
1	A	618	SER
1	A	632	ARG
1	A	633	ILE
1	A	641	ASP
1	A	642	ARG
1	A	647	GLN
1	A	654	SER
1	A	665	GLN
1	A	673	ILE
1	A	674	PRO
1	A	679	LYS
1	A	683	LEU
1	A	695	LEU
1	A	696	MET
1	A	697	PHE
1	A	705	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	402	ASN
1	A	460	GLN
1	A	475	GLN
1	A	486	ASN
1	A	526	ASN
1	A	593	ASN
1	A	647	GLN
1	A	677	ASN

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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	999	2	22,29,29	1.07	1 (4%)	27,45,45	1.94	3 (11%)

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	999	2	-	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)
3	A	999	ADP	C6-N6	-2.94	1.25	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	A	999	ADP	N3-C2-N1	-7.13	123.44	128.89
3	A	999	ADP	N6-C6-N1	2.11	123.73	119.20
3	A	999	ADP	C2-N1-C6	4.85	127.44	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	323/346 (93%)	-0.52	7 (2%) 65 73	11, 33, 79, 100	6 (1%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	385	GLY	11.4
1	A	669	THR	5.3
1	A	635	VAL	3.3
1	A	574	VAL	3.0
1	A	416	ASN	2.7
1	A	668	SER	2.7
1	A	670	LYS	2.4

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
3	ADP	A	999	27/27	0.97	0.08	-1.26	2,22,44,56	0
2	MG	A	997	1/1	0.96	0.03	-	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.