



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:03 AM UTC

PDB ID : 3KAR / pdb_00003kar
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

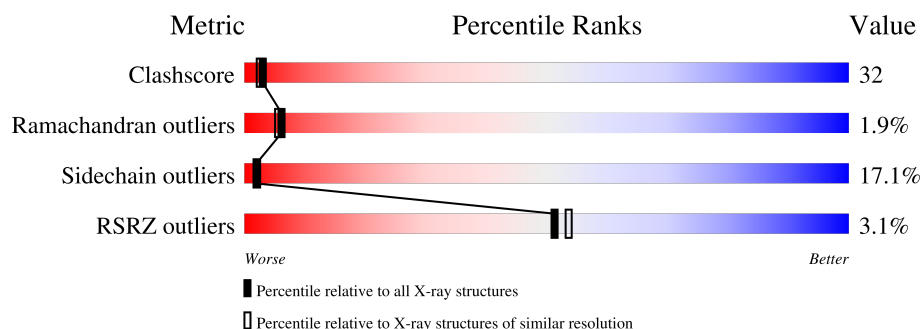
1 Overall quality at a glance

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	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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.

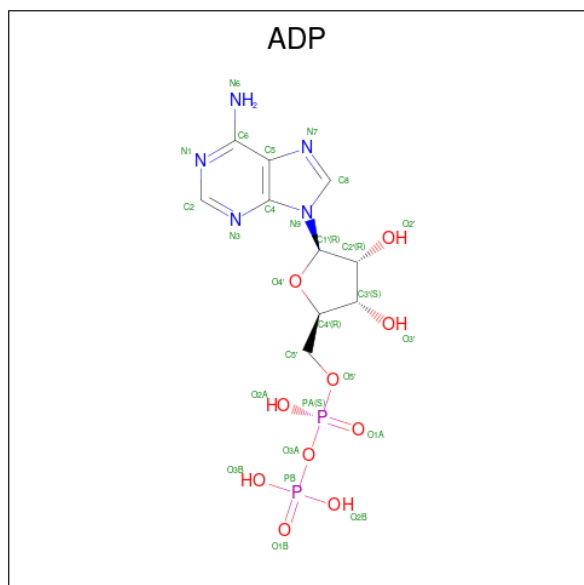
- Molecule 1 is a protein called KINESIN-LIKE PROTEIN KAR3.

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 (CCD ID: MG) (formula: Mg).

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

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

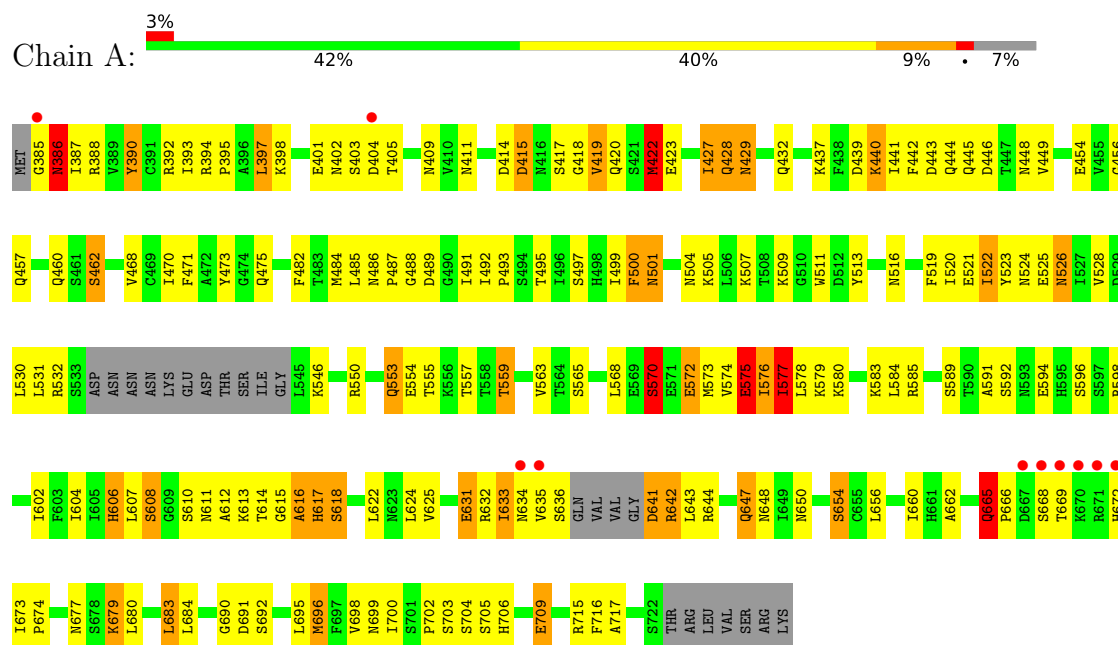
- Molecule 4 is water.

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

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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 30.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.2 (30.00-2.30) 93.3 (30.00-2.30)	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.	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.573	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 118.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	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, 0.333 respectively for untwinned datasets, and 0.375, 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: ADP, MG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.37	5/2514 (0.2%)	1.63	28/3407 (0.8%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	429	ASN	CA-C	-5.38	1.48	1.53
1	A	606	HIS	CG-ND1	5.36	1.44	1.38
1	A	470	ILE	CA-CB	-5.19	1.47	1.54
1	A	390	TYR	CA-C	-5.18	1.46	1.52
1	A	513	TYR	CA-C	-5.12	1.46	1.52

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	625	VAL	N-CA-C	7.46	119.02	108.36
1	A	500	PHE	CA-CB-CG	-7.25	106.56	113.80
1	A	616	ALA	N-CA-C	7.22	120.06	110.53
1	A	699	ASN	CA-CB-CG	-6.72	105.88	112.60
1	A	684	LEU	N-CA-C	6.70	121.47	113.16
1	A	617	HIS	CA-CB-CG	-6.67	107.13	113.80
1	A	419	VAL	N-CA-CB	-6.42	105.27	112.45
1	A	501	ASN	CA-CB-CG	-6.03	106.57	112.60
1	A	422	MET	N-CA-CB	-5.92	102.03	111.62
1	A	691	ASP	CA-CB-CG	-5.82	106.78	112.60
1	A	608	SER	CA-C-N	-5.82	118.08	122.33
1	A	608	SER	C-N-CA	-5.82	118.08	122.33
1	A	709	GLU	N-CA-C	-5.67	105.10	111.28
1	A	585	ARG	CG-CD-NE	-5.66	99.55	112.00
1	A	577	ILE	CB-CA-C	-5.66	104.73	111.97
1	A	642	ARG	N-CA-C	-5.61	105.17	112.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	665	GLN	CA-C-N	5.60	125.96	119.47
1	A	665	GLN	C-N-CA	5.60	125.96	119.47
1	A	504	ASN	CA-CB-CG	-5.43	107.17	112.60
1	A	526	ASN	CA-CB-CG	-5.41	107.19	112.60
1	A	439	ASP	CA-CB-CG	-5.29	107.31	112.60
1	A	631	GLU	N-CA-CB	5.27	117.64	109.48
1	A	485	LEU	N-CA-C	5.25	120.33	113.30
1	A	386	ASN	N-CA-C	5.21	118.08	109.85
1	A	672	HIS	CA-CB-CG	-5.18	108.62	113.80
1	A	692	SER	N-CA-C	5.11	117.44	110.24
1	A	570	SER	CA-C-N	5.06	127.37	120.54
1	A	570	SER	C-N-CA	5.06	127.37	120.54

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	2475	0	2360	155	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	155	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 32.

All (155) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:386:ASN:C	1:A:387:ILE:HD12	2.03	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:643:LEU:C	1:A:647:GLN:HE21	1.91	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:HD12	1:A:387:ILE:N	2.03	0.74
1:A:456:GLY:O	1:A:460:GLN:HG3	1.88	0.74
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:422:MET:HG3	1:A:441:ILE:HD13	1.79	0.64
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:O	1:A:635:VAL:HG12	1.99	0.61
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:402:ASN:ND2	1:A:404:ASP:H	1.99	0.59
1:A:522:ILE:HG13	1:A:602:ILE:CD1	2.31	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:471:PHE:HB2	1:A:696:MET:HE3	1.85	0.58
1:A:444:GLN:NE2	4:A:843:HOH:O	2.29	0.58
1:A:473:TYR:HB2	1:A:696:MET:HE2	1.84	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:487:PRO:C	1:A:489:ASP:H	2.13	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:473:TYR:CB	1:A:696:MET:HE2	2.37	0.55
1:A:427:ILE:C	1:A:429:ASN:N	2.64	0.55
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:427:ILE:O	1:A:429:ASN:N	2.40	0.54
1:A:419:VAL:HG11	1:A:437:LYS:CB	2.37	0.53
1:A:482:PHE:O	1:A:486:ASN:HB2	2.08	0.53
1:A:611:ASN:O	1:A:615:GLY:N	2.43	0.52
1:A:611:ASN:CG	1:A:614:THR:OG1	2.54	0.51
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:473:TYR:HB2	1:A:696:MET:HE1	1.92	0.51
1:A:519:PHE:HB3	1:A:531:LEU:HG	1.93	0.51
1:A:635:VAL:O	1:A:636:SER:HB3	2.10	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:610:SER:HA	1:A:616:ALA:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:650:ASN:O	1:A:654:SER:HB3	2.12	0.50
1:A:387:ILE:N	1:A:387:ILE:CD1	2.75	0.49
1:A:553:GLN:H	1:A:553:GLN:CD	2.20	0.49
1:A:568:LEU:HD22	1:A:574:VAL:HG23	1.93	0.49
1:A:607:LEU:N	1:A:607:LEU:HD12	2.28	0.49
1:A:612:ALA:C	1:A:614:THR:H	2.21	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:523:TYR:O	1:A:524:ASN:HB2	2.12	0.48
1:A:633:ILE:HG23	1:A:633:ILE:O	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:443:ASP:OD1	1:A:443:ASP:C	2.57	0.48
1:A:427:ILE:C	1:A:429:ASN:H	2.22	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:574:VAL:HG12	1:A:578:LEU:CG	2.40	0.47
1:A:716:PHE:C	1:A:716:PHE:CD1	2.92	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:471:PHE:CB	1:A:696:MET:HE3	2.44	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:427:ILE:O	1:A:428:GLN:C	2.59	0.45
1:A:572:GLU:O	1:A:575:GLU:CG	2.63	0.45
1:A:578:LEU:HD23	1:A:578:LEU:HA	1.67	0.45
1:A:633:ILE:O	1:A:635:VAL:HG23	2.16	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:674:PRO:O	1:A:677:ASN:HB2	2.17	0.44
1:A:579:LYS:O	1:A:580:LYS:C	2.60	0.44
1:A:665:GLN:CG	1:A:666:PRO:HD2	2.48	0.44
1:A:521:GLU:O	1:A:521:GLU:HG3	2.14	0.43
1:A:402:ASN:HD22	1:A:404:ASP:H	1.62	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:611:ASN:OD1	1:A:611:ASN:C	2.62	0.42
1:A:709:GLU:OE1	1:A:709:GLU:HA	2.18	0.42
1:A:432:GLN:O	4:A:801:HOH:O	2.21	0.42
1:A:633:ILE:HD13	1:A:635:VAL:HG22	2.00	0.42
1:A:415:ASP:C	1:A:417:SER:N	2.76	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:390:TYR:CD1	1:A:440:LYS:HB3	2.54	0.42
1:A:550:ARG:HD3	4:A:813:HOH:O	2.19	0.42
1:A:524:ASN:HA	1:A:648:ASN:ND2	2.35	0.42
1:A:662:ALA:O	1:A:668:SER:OG	2.23	0.42
1:A:612:ALA:C	1:A:614:THR:N	2.78	0.41
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:674:PRO:CB	1:A:677:ASN:ND2	2.81	0.41
1:A:393:ILE:HG12	1:A:700:ILE:HD11	2.01	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:405:THR:HB	1:A:704:SER:HB3	2.00	0.41
1:A:617:HIS:CD2	1:A:617:HIS:N	2.86	0.41
1:A:415:ASP:C	1:A:418:GLY:H	2.28	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: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:484:MET:HE1	1:A:624:LEU:HB3	2.02	0.40
1:A:385:GLY:N	1:A:690:GLY:O	2.55	0.40
1:A:500:PHE:O	1:A:501:ASN:C	2.63	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%)	6 5

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%)	228 (83%)	47 (17%)	2 2

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	386	ASN
1	A	397	LEU
1	A	414	ASP
1	A	415	ASP
1	A	422	MET

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Mol	Chain	Res	Type
1	A	427	ILE
1	A	440	LYS
1	A	454	GLU
1	A	457	GLN
1	A	462	SER
1	A	491	ILE
1	A	497	SER
1	A	509	LYS
1	A	516	ASN
1	A	520	ILE
1	A	522	ILE
1	A	532	ARG
1	A	546	LYS
1	A	553	GLN
1	A	554	GLU
1	A	559	THR
1	A	570	SER
1	A	572	GLU
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	604	ILE
1	A	606	HIS
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	679	LYS
1	A	683	LEU
1	A	695	LEU
1	A	696	MET
1	A	705	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) 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
1	A	699	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	28,29,29	1.29	3 (10%)	43,45,45	1.40	4 (9%)

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	-	2/16/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	999	ADP	PA-O3A	-3.85	1.55	1.59
3	A	999	ADP	C6-N6	-3.26	1.25	1.34
3	A	999	ADP	C2-N3	-2.84	1.28	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	999	ADP	C2-N1-C6	5.30	127.44	118.73
3	A	999	ADP	N3-C2-N1	-3.40	123.44	128.58
3	A	999	ADP	C5-C6-N1	-3.06	109.75	117.51
3	A	999	ADP	N6-C6-N1	2.41	123.73	118.38

There are no chirality outliers.

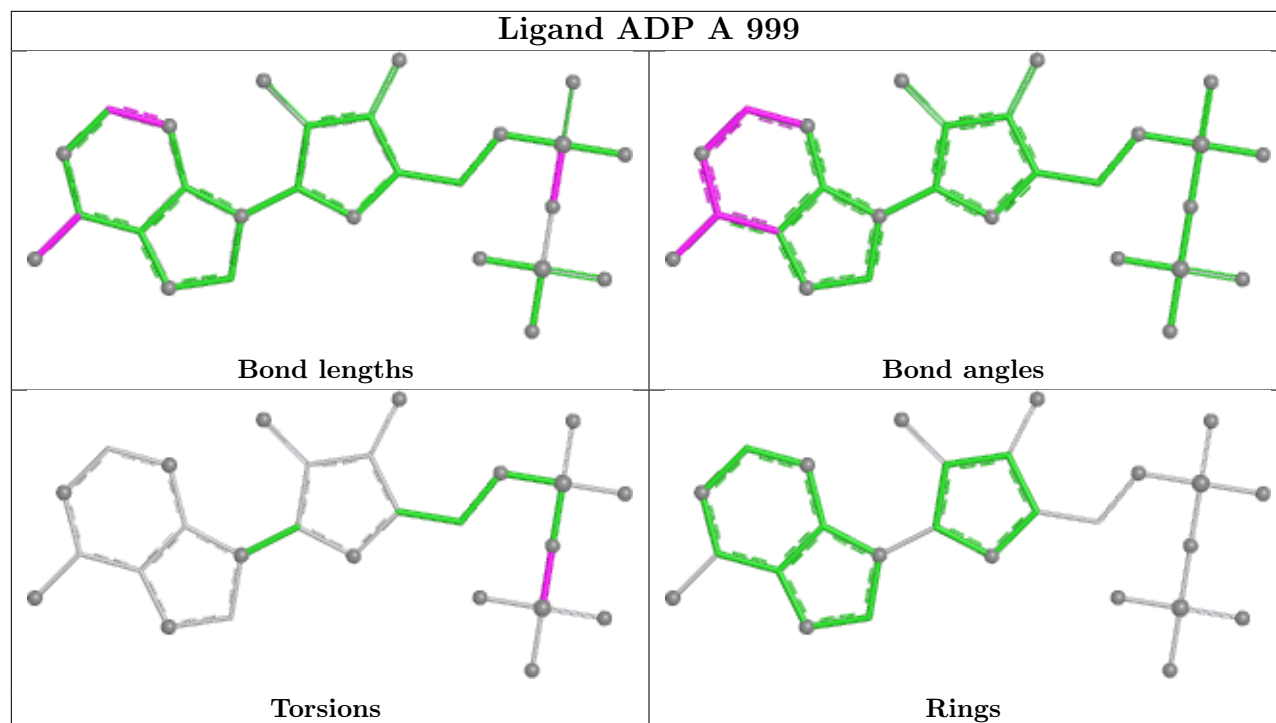
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	999	ADP	PA-O3A-PB-O2B
3	A	999	ADP	PA-O3A-PB-O1B

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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%)	-0.17	10 (3%) 51 53	11, 33, 78, 100	6 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	671	ARG	4.9
1	A	385	GLY	4.0
1	A	668	SER	3.6
1	A	672	HIS	3.6
1	A	669	THR	3.5
1	A	635	VAL	2.6
1	A	404	ASP	2.6
1	A	670	LYS	2.5
1	A	667	ASP	2.5
1	A	634	ASN	2.2

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](#)

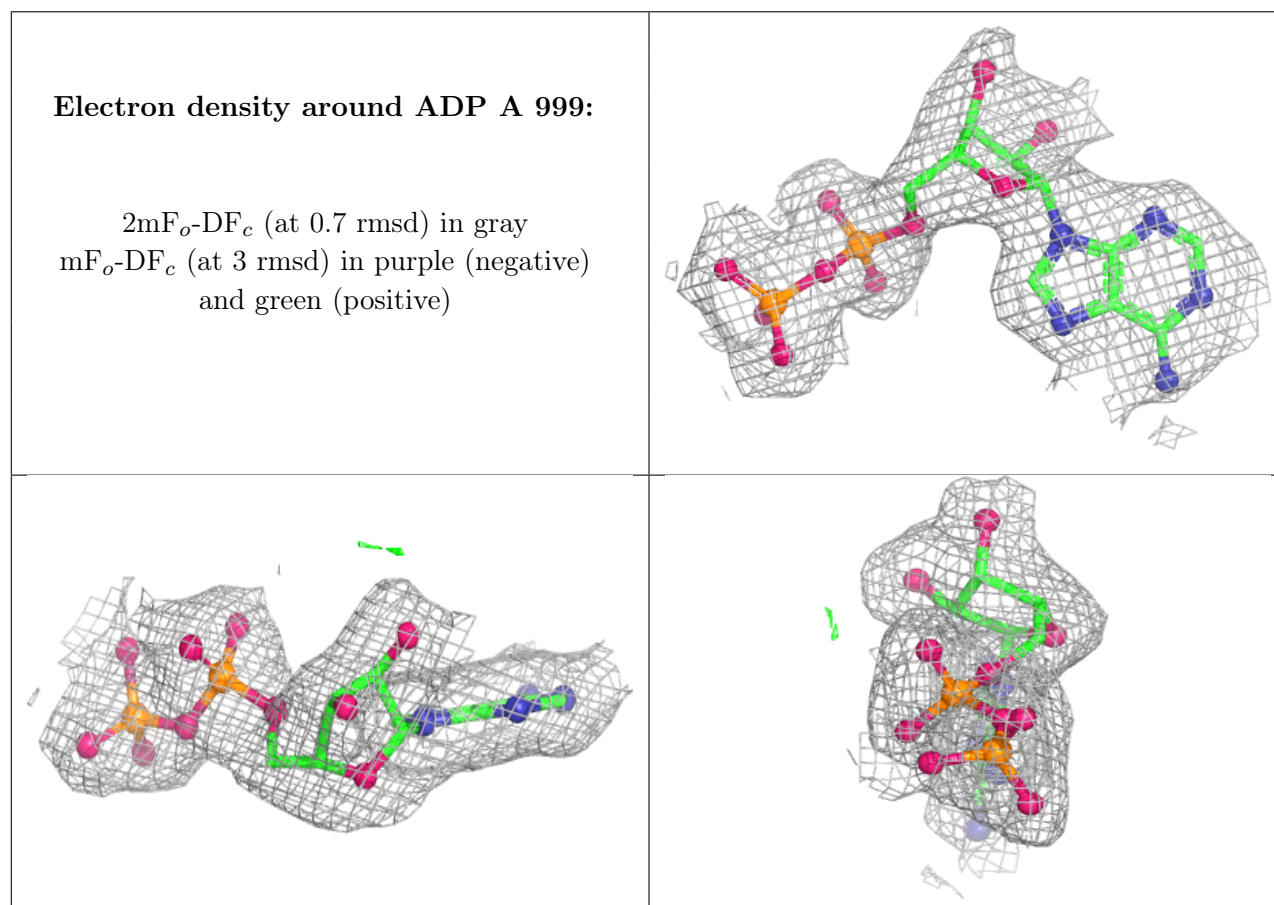
There are no oligosaccharides in this entry.

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. 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	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	A	997	1/1	0.99	0.02	21,21,21,21	0
3	ADP	A	999	27/27	0.99	0.04	2,22,44,56	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers ⓘ

There are no such residues in this entry.