



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:05 PM GMT

PDB ID : 4YIX
Title : Structure of MRB1590 bound to ADP
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Deposited on : 2015-03-02
Resolution : 2.60 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

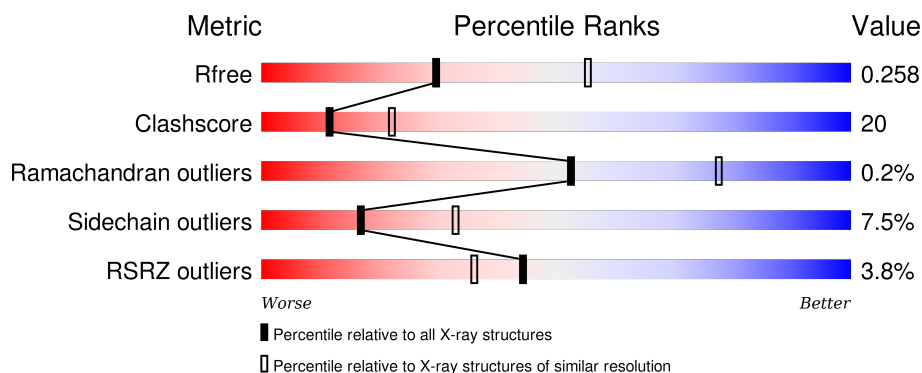
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	668	<div> <div>3%</div> <div>56%</div> <div>26%</div> <div>• •</div> <div>13%</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	MG	A	703	-	-	X	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	HG	A	705	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4585 atoms, of which 6 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	579	Total	C	N	O	S	0	0	0
			4429	2770	785	848	26			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	242	GLU	VAL	conflict	UNP Q57ZF2

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Mg	0	0
			3	3		

- Molecule 3 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	Hg	0	0
			11	11		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

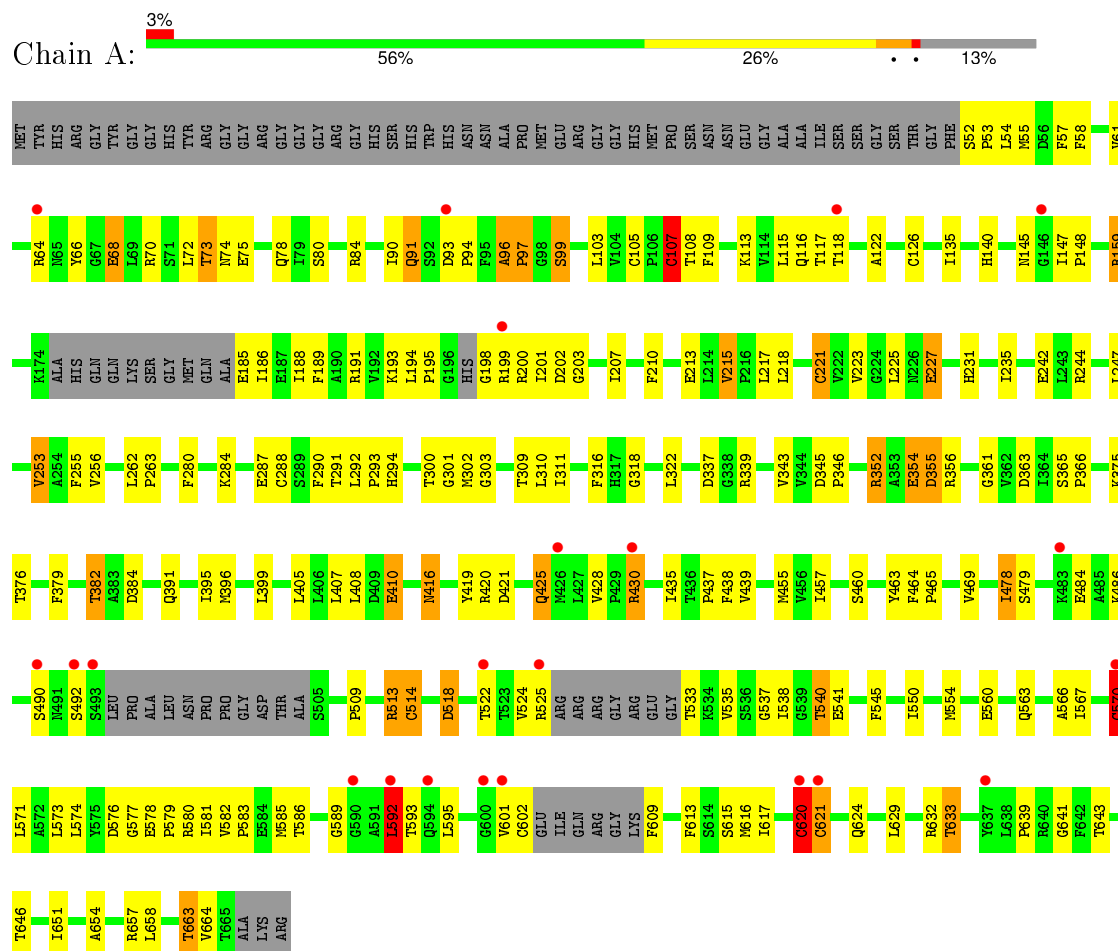
- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	109	Total	H	O	0	0
			115	6	109		

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: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	105.58Å 184.71Å 73.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.60 – 2.60 25.60 – 2.60	Depositor EDS
% Data completeness (in resolution range)	91.3 (25.60-2.60) 91.3 (25.60-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.26 (at 2.60Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.204 , 0.262 0.210 , 0.258	Depositor DCC
R_{free} test set	1987 reflections (9.65%)	DCC
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.2	EDS
Estimated twinning fraction	0.021 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.038 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 20596 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4585	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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: MG, HG, 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	0.83	7/4508 (0.2%)	0.79	15/6111 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	514	CYS	CA-CB	7.51	1.70	1.53
1	A	620	CYS	CB-SG	7.20	1.94	1.82
1	A	514	CYS	CB-SG	6.97	1.94	1.82
1	A	354	GLU	CD-OE1	-6.70	1.18	1.25
1	A	354	GLU	CD-OE2	-5.71	1.19	1.25
1	A	244	ARG	CZ-NH2	-5.14	1.26	1.33
1	A	221	CYS	CA-CB	5.08	1.65	1.53

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	514	CYS	CA-CB-SG	11.70	135.06	114.00
1	A	97	PRO	CB-CA-C	-9.14	89.14	112.00
1	A	97	PRO	N-CA-C	-8.74	89.37	112.10
1	A	107	CYS	CA-CB-SG	8.63	129.53	114.00
1	A	221	CYS	CA-CB-SG	7.50	127.49	114.00
1	A	514	CYS	N-CA-CB	6.96	123.12	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	352	ARG	O-C-N	-6.84	111.75	122.70
1	A	513	ARG	NE-CZ-NH2	-6.65	116.98	120.30
1	A	96	ALA	C-N-CD	6.12	141.26	128.40
1	A	592	LEU	CA-CB-CG	6.08	129.29	115.30
1	A	513	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	A	570	CYS	N-CA-CB	5.83	121.10	110.60
1	A	159	ARG	C-N-CD	5.43	139.81	128.40
1	A	620	CYS	CA-CB-SG	5.34	123.61	114.00
1	A	215	VAL	C-N-CD	5.33	139.59	128.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	198	GLY	Peptide

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	4429	0	4387	181	0
2	A	3	0	0	2	0
3	A	11	0	0	2	0
4	A	27	0	12	2	0
5	A	109	6	0	9	0
All	All	4579	6	4399	181	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 20.

All (181) 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:310:LEU:HD22	1:A:457:ILE:HD11	1.46	0.96
1:A:66:TYR:CE1	1:A:94:PRO:HG3	2.01	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:CYS:HG	2:A:703:MG:MG	0.67	0.95
1:A:126:CYS:SG	5:A:856:HOH:O	2.27	0.93
1:A:107:CYS:SG	3:A:705:HG:HG	0.12	0.92
1:A:96:ALA:HB1	1:A:97:PRO:HD3	1.48	0.92
1:A:573:LEU:HD13	1:A:633:THR:HG21	1.53	0.90
1:A:620:CYS:SG	2:A:703:MG:MG	1.55	0.90
1:A:107:CYS:CB	3:A:705:HG:HG	1.85	0.84
1:A:91:GLN:HG2	1:A:99:SER:OG	1.79	0.83
1:A:55:MET:HG3	1:A:207:ILE:HD13	1.62	0.82
1:A:573:LEU:HD13	1:A:633:THR:CG2	2.10	0.81
1:A:96:ALA:HB1	1:A:97:PRO:CD	2.12	0.79
1:A:210:PHE:O	1:A:215:VAL:HG23	1.86	0.75
1:A:185:GLU:N	5:A:802:HOH:O	2.20	0.75
1:A:280:PHE:CE1	1:A:478:ILE:HG13	2.23	0.73
1:A:621:CYS:SG	1:A:651:ILE:HD11	2.28	0.73
1:A:621:CYS:HB2	1:A:651:ILE:HD12	1.70	0.72
1:A:61:VAL:HB	1:A:201:ILE:HD12	1.72	0.71
1:A:195:PRO:HB2	1:A:202:ASP:HB3	1.73	0.70
1:A:117:THR:HG21	1:A:122:ALA:HB3	1.73	0.70
1:A:567:ILE:HD11	1:A:646:THR:HG21	1.74	0.70
1:A:416:ASN:OD1	1:A:435:ILE:CD1	2.41	0.69
1:A:70:ARG:NH1	5:A:804:HOH:O	2.26	0.69
1:A:107:CYS:SG	1:A:109:PHE:O	2.48	0.68
1:A:97:PRO:O	1:A:97:PRO:CD	2.41	0.67
1:A:573:LEU:CD1	1:A:633:THR:HG21	2.22	0.67
1:A:621:CYS:HB2	1:A:651:ILE:CD1	2.25	0.67
1:A:52:SER:HB3	1:A:53:PRO:HD3	1.77	0.67
1:A:203:GLY:O	1:A:207:ILE:HG13	1.94	0.66
1:A:589:GLY:HA2	1:A:592:LEU:HD22	1.76	0.66
1:A:97:PRO:O	1:A:97:PRO:HD2	1.95	0.66
1:A:242:GLU:HG3	1:A:293:PRO:CG	2.25	0.66
1:A:525:ARG:O	1:A:533:THR:HG22	1.97	0.65
1:A:524:VAL:CG1	1:A:533:THR:HG21	2.27	0.65
1:A:524:VAL:HG12	1:A:533:THR:HG21	1.77	0.65
1:A:420:ARG:NH2	1:A:428:VAL:O	2.29	0.65
1:A:518:ASP:O	1:A:522:THR:HG23	1.95	0.65
1:A:352:ARG:NH1	1:A:554:MET:CG	2.61	0.64
1:A:576:ASP:OD2	1:A:633:THR:HB	1.98	0.63
1:A:567:ILE:HD12	1:A:651:ILE:HG12	1.80	0.63
1:A:70:ARG:O	1:A:73:THR:OG1	2.16	0.63
1:A:310:LEU:HD22	1:A:457:ILE:CD1	2.25	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:583:PRO:HA	1:A:586:THR:HG22	1.82	0.62
1:A:396:MET:HE2	1:A:399:LEU:HD12	1.80	0.62
1:A:61:VAL:HA	1:A:64:ARG:HD2	1.82	0.61
1:A:361:GLY:HA3	1:A:663:THR:HG21	1.82	0.61
1:A:438:PHE:HE2	1:A:455:MET:HE1	1.65	0.61
1:A:117:THR:HG21	1:A:122:ALA:CB	2.30	0.61
1:A:117:THR:HG22	1:A:118:THR:N	2.16	0.61
1:A:438:PHE:CE2	1:A:455:MET:HE1	2.35	0.61
1:A:293:PRO:HG2	1:A:294:HIS:CD2	2.36	0.60
1:A:61:VAL:HB	1:A:201:ILE:CD1	2.31	0.60
1:A:78:GLN:NE2	1:A:84:ARG:HB3	2.15	0.60
1:A:135:ILE:HD13	1:A:186:ILE:HD13	1.82	0.60
1:A:242:GLU:HG3	1:A:293:PRO:HG3	1.83	0.59
1:A:91:GLN:OE1	1:A:194:LEU:N	2.32	0.59
1:A:437:PRO:HB2	1:A:439:VAL:HG12	1.83	0.59
1:A:589:GLY:HA3	1:A:615:SER:O	2.03	0.59
1:A:509:PRO:O	1:A:609:PHE:HA	2.02	0.59
1:A:421:ASP:O	1:A:425:GLN:HG3	2.04	0.58
1:A:318:GLY:HA2	4:A:715:ADP:O2A	2.03	0.57
1:A:352:ARG:HH12	1:A:554:MET:CG	2.17	0.57
1:A:73:THR:O	1:A:74:ASN:HB2	2.04	0.57
1:A:657:ARG:HG3	1:A:657:ARG:HH11	1.68	0.57
1:A:57:PHE:O	1:A:61:VAL:HG22	2.04	0.57
1:A:284:LYS:HE2	1:A:287:GLU:OE2	2.05	0.56
1:A:560:GLU:HB3	5:A:881:HOH:O	2.05	0.56
1:A:117:THR:HG22	1:A:118:THR:H	1.71	0.56
1:A:356:ARG:HD2	1:A:384:ASP:OD1	2.06	0.56
1:A:126:CYS:SG	1:A:601:VAL:HG22	2.45	0.56
1:A:632:ARG:NH2	5:A:815:HOH:O	2.39	0.55
1:A:54:LEU:HD11	1:A:58:PHE:CE2	2.41	0.55
1:A:382:THR:HB	1:A:384:ASP:H	1.72	0.55
1:A:438:PHE:HE2	1:A:455:MET:CE	2.19	0.54
1:A:145:ASN:HB2	1:A:147:ILE:CD1	2.38	0.54
1:A:567:ILE:CD1	1:A:651:ILE:HG12	2.38	0.54
1:A:247:LEU:HD11	1:A:343:VAL:HG23	1.90	0.54
1:A:288:CYS:SG	1:A:303:GLY:HA3	2.48	0.54
1:A:352:ARG:NH1	1:A:554:MET:SD	2.71	0.53
1:A:609:PHE:N	5:A:816:HOH:O	2.41	0.53
1:A:464:PHE:HB2	1:A:465:PRO:HD3	1.90	0.53
1:A:290:PHE:CE1	1:A:292:LEU:HD23	2.43	0.53
1:A:352:ARG:NH1	1:A:554:MET:HG2	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:395:ILE:HG22	1:A:396:MET:CE	2.38	0.53
1:A:231:HIS:O	1:A:235:ILE:HG13	2.09	0.53
1:A:140:HIS:CE1	1:A:159:ARG:HG2	2.43	0.53
1:A:287:GLU:HG2	1:A:300:THR:CG2	2.40	0.52
1:A:355:ASP:O	1:A:356:ARG:HB2	2.08	0.52
1:A:291:THR:O	5:A:801:HOH:O	2.18	0.52
1:A:96:ALA:O	1:A:193:LYS:HE2	2.10	0.52
1:A:583:PRO:HA	1:A:586:THR:CG2	2.39	0.52
1:A:91:GLN:CG	1:A:99:SER:OG	2.54	0.52
1:A:291:THR:HG23	5:A:801:HOH:O	2.10	0.52
1:A:262:LEU:N	1:A:263:PRO:HD2	2.25	0.51
1:A:352:ARG:HH11	1:A:554:MET:HB3	1.75	0.51
1:A:115:LEU:O	1:A:116:GLN:HB2	2.11	0.50
1:A:280:PHE:HE1	1:A:478:ILE:HG13	1.72	0.49
1:A:580:ARG:O	1:A:583:PRO:HD2	2.13	0.49
1:A:570:CYS:O	1:A:574:LEU:HB2	2.13	0.49
1:A:352:ARG:NH1	1:A:554:MET:HB3	2.29	0.48
1:A:90:ILE:HA	1:A:99:SER:HB3	1.95	0.48
1:A:290:PHE:CD1	1:A:292:LEU:HD23	2.49	0.48
1:A:550:ILE:HD11	1:A:658:LEU:HD13	1.94	0.48
1:A:64:ARG:HD3	1:A:68:GLU:OE2	2.14	0.48
1:A:188:ILE:HD12	1:A:218:LEU:HD21	1.96	0.48
1:A:419:TYR:CE2	1:A:460:SER:HB3	2.48	0.48
1:A:379:PHE:CZ	1:A:663:THR:HG22	2.49	0.48
1:A:253:VAL:HG13	1:A:405:LEU:HD22	1.96	0.48
1:A:567:ILE:CD1	1:A:646:THR:HG21	2.42	0.47
1:A:573:LEU:HD13	1:A:633:THR:HG23	1.95	0.47
1:A:66:TYR:CE1	1:A:94:PRO:CG	2.89	0.47
1:A:93:ASP:CG	1:A:94:PRO:HD2	2.34	0.47
1:A:416:ASN:H	1:A:416:ASN:HD22	1.62	0.47
1:A:318:GLY:CA	4:A:715:ADP:O2A	2.62	0.47
1:A:225:LEU:HD12	1:A:225:LEU:C	2.35	0.47
1:A:396:MET:HE2	1:A:396:MET:HA	1.97	0.47
1:A:577:GLY:O	1:A:581:ILE:HG22	2.15	0.47
1:A:191:ARG:HD3	1:A:639:PRO:CG	2.44	0.46
1:A:354:GLU:HG3	1:A:554:MET:CE	2.45	0.46
1:A:581:ILE:O	1:A:585:MET:HG2	2.15	0.46
1:A:563:GLN:O	1:A:566:ALA:HB3	2.14	0.46
1:A:255:PHE:HA	1:A:302:MET:O	2.16	0.46
1:A:191:ARG:HD3	1:A:639:PRO:HG2	1.97	0.46
1:A:416:ASN:N	1:A:416:ASN:HD22	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:LEU:HA	1:A:195:PRO:HD3	1.82	0.46
1:A:583:PRO:CA	1:A:586:THR:HG22	2.45	0.45
1:A:379:PHE:HZ	1:A:663:THR:HG22	1.81	0.45
1:A:586:THR:HA	1:A:616:MET:HA	1.99	0.45
1:A:407:LEU:C	1:A:408:LEU:HD12	2.37	0.45
1:A:72:LEU:O	1:A:75:GLU:HB2	2.16	0.45
1:A:578:GLU:N	1:A:579:PRO:HD2	2.31	0.45
1:A:213:GLU:O	1:A:217:LEU:HB2	2.16	0.45
1:A:108:THR:HG21	1:A:223:VAL:CG1	2.46	0.45
1:A:189:PHE:HB2	1:A:641:GLY:HA3	1.99	0.45
1:A:311:ILE:HD13	1:A:322:LEU:HD23	1.99	0.44
1:A:361:GLY:HA3	1:A:663:THR:CG2	2.47	0.44
1:A:621:CYS:SG	1:A:651:ILE:CD1	3.02	0.44
1:A:578:GLU:HB3	1:A:579:PRO:HD3	2.00	0.43
1:A:365:SER:N	1:A:366:PRO:CD	2.80	0.43
1:A:218:LEU:O	1:A:223:VAL:HG23	2.19	0.43
1:A:84:ARG:NH2	1:A:185:GLU:OE2	2.51	0.43
1:A:52:SER:CB	1:A:53:PRO:HD3	2.47	0.43
1:A:309:THR:HG23	1:A:469:VAL:HB	2.01	0.43
1:A:256:VAL:HB	1:A:280:PHE:CE2	2.53	0.43
1:A:291:THR:HG23	1:A:291:THR:O	2.19	0.43
1:A:287:GLU:CG	1:A:300:THR:CG2	2.97	0.43
1:A:537:GLY:C	1:A:538:ILE:HD13	2.39	0.43
1:A:541:GLU:HG2	5:A:872:HOH:O	2.19	0.43
1:A:363:ASP:O	1:A:396:MET:HG3	2.18	0.42
1:A:96:ALA:CB	1:A:97:PRO:CD	2.92	0.42
1:A:61:VAL:O	1:A:201:ILE:HD12	2.19	0.42
1:A:148:PRO:HG3	1:A:217:LEU:HD13	2.00	0.42
1:A:613:PHE:O	1:A:617:ILE:HG12	2.19	0.42
1:A:629:LEU:O	1:A:643:THR:HG21	2.18	0.42
1:A:535:VAL:HG22	1:A:545:PHE:CD1	2.54	0.42
1:A:540:THR:HG23	1:A:540:THR:O	2.18	0.42
1:A:621:CYS:CB	1:A:651:ILE:CD1	2.96	0.42
1:A:227:GLU:O	1:A:227:GLU:HG3	2.20	0.42
1:A:601:VAL:HG12	1:A:602:CYS:N	2.35	0.42
1:A:337:ASP:OD2	1:A:339:ARG:HB2	2.20	0.42
1:A:345:ASP:HA	1:A:346:PRO:HD3	1.92	0.42
1:A:188:ILE:HD12	1:A:218:LEU:CD2	2.49	0.41
1:A:490:SER:C	1:A:492:SER:H	2.22	0.41
1:A:465:PRO:O	1:A:486:LYS:NZ	2.54	0.41
1:A:567:ILE:HG21	1:A:654:ALA:CB	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:621:CYS:SG	1:A:624:GLN:HB3	2.60	0.41
1:A:108:THR:HG21	1:A:223:VAL:HG12	2.02	0.41
1:A:375:LYS:HA	1:A:375:LYS:HD2	1.83	0.41
1:A:117:THR:CG2	1:A:118:THR:N	2.82	0.41
1:A:464:PHE:N	1:A:465:PRO:CD	2.83	0.41
1:A:439:VAL:HA	1:A:463:TYR:CE2	2.55	0.41
1:A:410:GLU:OE2	1:A:457:ILE:HG23	2.19	0.41
1:A:595:LEU:HD23	1:A:595:LEU:HA	1.81	0.41
1:A:430:ARG:HD2	1:A:430:ARG:H	1.86	0.41
1:A:524:VAL:HG12	1:A:533:THR:CG2	2.48	0.41
1:A:582:VAL:O	1:A:586:THR:HG22	2.21	0.40
1:A:391:GLN:O	1:A:395:ILE:HG13	2.21	0.40
1:A:396:MET:CE	1:A:399:LEU:HD12	2.49	0.40
1:A:287:GLU:HG2	1:A:300:THR:HG23	2.04	0.40
1:A:300:THR:HG22	1:A:301:GLY:N	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	567/668 (85%)	544 (96%)	22 (4%)	1 (0%)	52 77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	664	VAL

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	492/554 (89%)	455 (92%)	37 (8%)	17	33

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	68	GLU
1	A	73	THR
1	A	80	SER
1	A	91	GLN
1	A	99	SER
1	A	103	LEU
1	A	105	CYS
1	A	107	CYS
1	A	113	LYS
1	A	199	ARG
1	A	200	ARG
1	A	221	CYS
1	A	227	GLU
1	A	253	VAL
1	A	316	PHE
1	A	355	ASP
1	A	376	THR
1	A	382	THR
1	A	410	GLU
1	A	416	ASN
1	A	425	GLN
1	A	430	ARG
1	A	478	ILE
1	A	479	SER
1	A	484	GLU
1	A	513	ARG
1	A	514	CYS
1	A	518	ASP
1	A	540	THR
1	A	570	CYS

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Mol	Chain	Res	Type
1	A	571	LEU
1	A	592	LEU
1	A	593	THR
1	A	620	CYS
1	A	621	CYS
1	A	633	THR
1	A	663	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	462	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](#)

Of 15 ligands modelled in this entry, 14 are 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	$\# Z > 2$
4	ADP	A	715	-	24,29,29	0.99	1 (4%)	23,45,45	1.76	2 (8%)

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
4	ADP	A	715	-	-	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)
4	A	715	ADP	C5-C4	3.10	1.47	1.40

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	715	ADP	N3-C2-N1	-6.57	123.71	128.87
4	A	715	ADP	C4'-O4'-C1'	2.01	111.77	109.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	715	ADP	2	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	579/668 (86%)	-0.15	22 (3%) 44 36	11, 23, 43, 57	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	594	GLN	4.1
1	A	64	ARG	3.8
1	A	525	ARG	3.5
1	A	493	SER	3.1
1	A	637	TYR	3.0
1	A	570	CYS	3.0
1	A	118	THR	2.8
1	A	620	CYS	2.7
1	A	621	CYS	2.6
1	A	430	ARG	2.5
1	A	590	GLY	2.5
1	A	146	GLY	2.5
1	A	483	LYS	2.4
1	A	426	MET	2.3
1	A	199	ARG	2.3
1	A	522	THR	2.2
1	A	600	GLY	2.2
1	A	492	SER	2.2
1	A	490	SER	2.1
1	A	592	LEU	2.1
1	A	93	ASP	2.1
1	A	601	VAL	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
2	MG	A	703	1/1	0.99	0.54	4.17	2,2,2,2	0
4	ADP	A	715	27/27	0.94	0.16	0.76	12,14,15,16	0
2	MG	A	702	1/1	0.96	0.25	0.37	11,11,11,11	0
3	HG	A	712	1/1	0.96	0.10	-1.24	99,99,99,99	0
3	HG	A	711	1/1	1.00	0.04	-1.52	33,33,33,33	0
3	HG	A	710	1/1	0.99	0.03	-2.15	39,39,39,39	0
3	HG	A	707	1/1	0.99	0.06	-3.07	55,55,55,55	0
3	HG	A	713	1/1	0.99	0.06	-6.21	63,63,63,63	0
3	HG	A	714	1/1	1.00	0.04	-	47,47,47,47	0
2	MG	A	701	1/1	0.93	0.15	-	22,22,22,22	0
3	HG	A	705	1/1	0.98	0.06	-	102,102,102,102	0
3	HG	A	704	1/1	0.97	0.07	-	55,55,55,55	0
3	HG	A	708	1/1	0.98	0.04	-	62,62,62,62	0
3	HG	A	706	1/1	0.99	0.11	-	61,61,61,61	0
3	HG	A	709	1/1	0.99	0.27	-	61,61,61,61	0

6.5 Other polymers ⓘ

There are no such residues in this entry.