



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:52 PM GMT

PDB ID : 1YV3
Title : The structural basis of blebbistatin inhibition and specificity for myosin II
Authors : Allingham, J.S.; Smith, R.; Rayment, I.
Deposited on : 2005-02-14
Resolution : 2.00 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

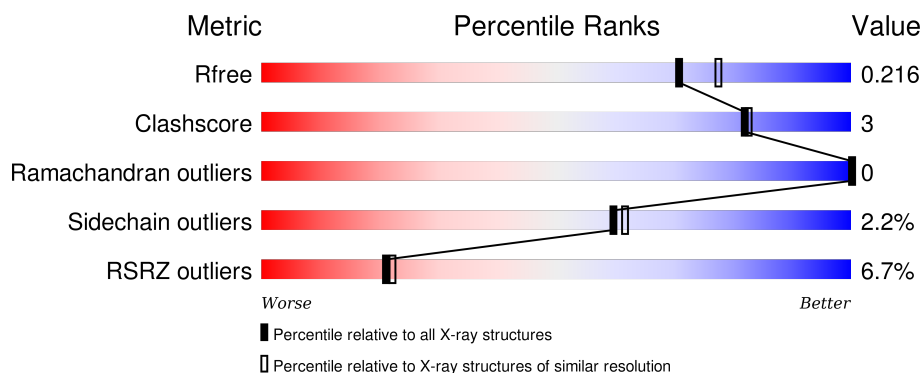
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

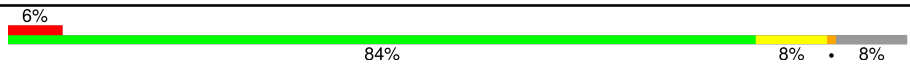
The reported resolution of this entry is 2.00 Å.

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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	762	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6219 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 Myosin II heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	703	Total	C	N	O	S	0	6	0
			5481	3509	923	1034	15			

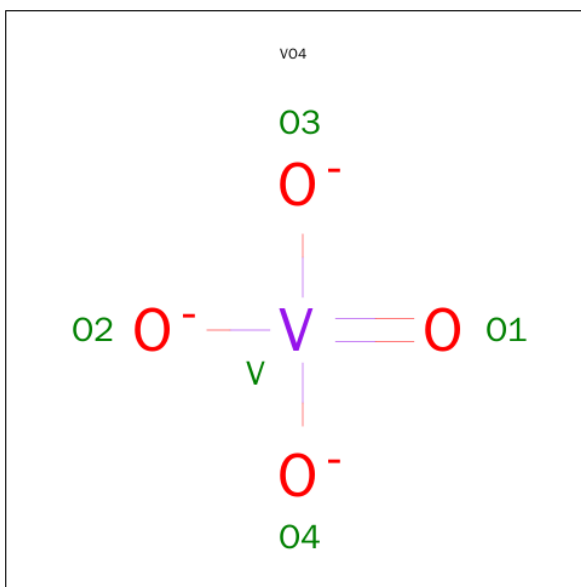
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	MET	see remark 999	UNP P08799
A	249	SER	ASN	see remark 999	UNP P08799
A	312	CYS	TYR	see remark 999	UNP P08799
A	760	LEU	GLN	see remark 999	UNP P08799
A	761	PRO	ARG	see remark 999	UNP P08799
A	762	ASN	ILE	see remark 999	UNP P08799

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

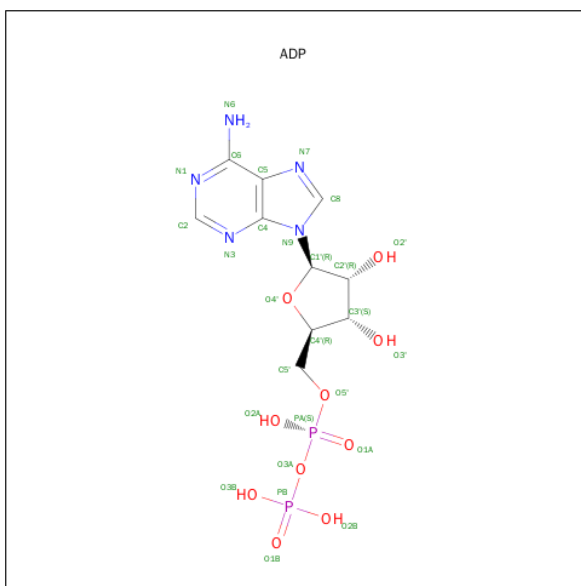
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is VANADATE ION (three-letter code: VO4) (formula: O₄V).



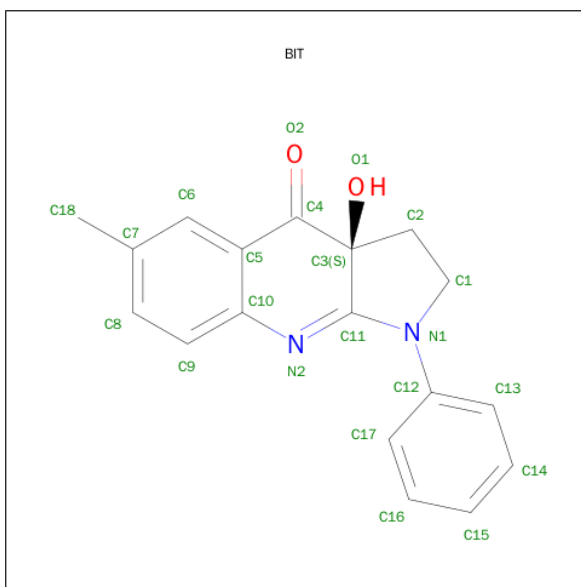
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	O	V		0	0
			5	4	1			

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



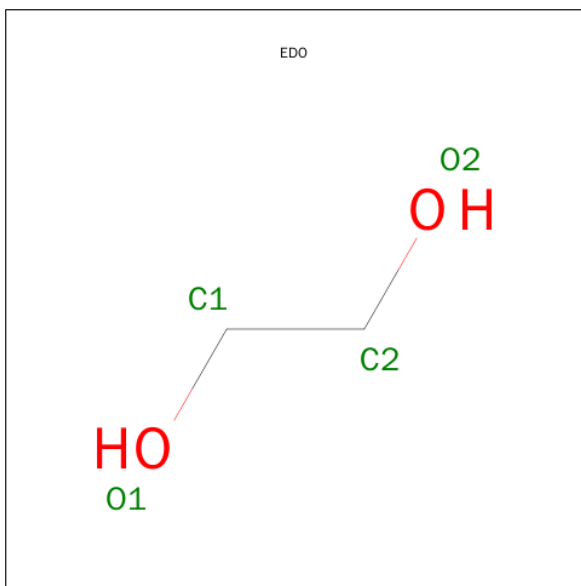
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 5 is (-)-1-PHENYL-1,2,3,4-TETRAHYDRO-4-HYDROXYPYRROLO[2,3-B]-7-M ETHYLQUINOLIN-4-ONE (three-letter code: BIT) (formula: $C_{18}H_{16}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			22	18	2	2		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		

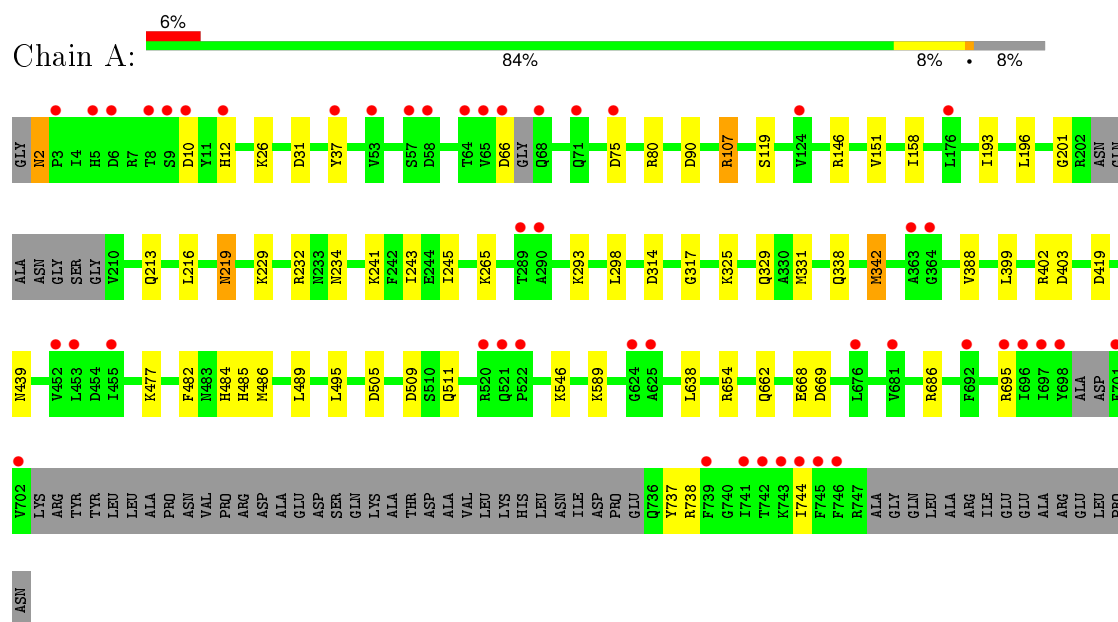
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	678	Total 678	O 678	0	0

3 Residue-property plots [i](#)

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: Myosin II heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.11Å 145.77Å 152.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.00 38.17 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.5 (40.00-2.00) 99.5 (38.17-1.99)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.56 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.173 , 0.215 0.176 , 0.216	Depositor DCC
R_{free} test set	3376 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	23.4	Xtriage
Anisotropy	0.326	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 63.0	EDS
Estimated twinning fraction	0.012 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.022 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 66866 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6219	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 4.13% 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: VO4, MG, EDO, ADP, BIT

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	0.98	5/5616 (0.1%)	0.91	13/7592 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	331[A]	MET	SD-CE	-5.77	1.45	1.77
1	A	331[B]	MET	SD-CE	-5.77	1.45	1.77
1	A	265	LYS	CE-NZ	5.62	1.63	1.49
1	A	388	VAL	CB-CG2	5.37	1.64	1.52
1	A	37	TYR	CD1-CE1	-5.28	1.31	1.39

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	314	ASP	CB-CG-OD2	8.67	126.10	118.30
1	A	419	ASP	CB-CG-OD1	7.40	124.96	118.30
1	A	509	ASP	CB-CG-OD2	6.20	123.88	118.30
1	A	75	ASP	CB-CG-OD2	6.09	123.78	118.30
1	A	402	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	A	90	ASP	CB-CG-OD2	5.89	123.60	118.30
1	A	505	ASP	CB-CG-OD2	5.71	123.44	118.30
1	A	10	ASP	CB-CG-OD2	5.64	123.38	118.30
1	A	686	ARG	NE-CZ-NH2	-5.61	117.49	120.30
1	A	314	ASP	CB-CG-OD1	-5.50	113.35	118.30
1	A	66	ASP	CB-CG-OD2	5.49	123.24	118.30
1	A	80	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	A	654	ARG	NE-CZ-NH1	5.01	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	5481	0	5206	34	0
2	A	2	0	0	0	0
3	A	5	0	0	0	0
4	A	27	0	12	0	0
5	A	22	0	16	0	0
6	A	4	0	6	0	0
7	A	678	0	0	19	2
All	All	6219	0	5240	34	2

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 3.

All (34) 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:668:GLU:OE1	7:A:1210:HOH:O	1.86	0.92
1:A:439:ASN:CB	7:A:1610:HOH:O	2.23	0.87
1:A:495:LEU:CG	7:A:1653:HOH:O	2.22	0.86
1:A:485[A]:HIS:ND1	7:A:1612:HOH:O	2.18	0.76
1:A:201:GLY:O	7:A:1455:HOH:O	2.05	0.73
1:A:669:ASP:OD1	7:A:1627:HOH:O	2.07	0.70
1:A:107:ARG:HD3	7:A:1167:HOH:O	1.98	0.63
1:A:484:HIS:CE1	1:A:489:LEU:HD11	2.34	0.63
1:A:589[A]:LYS:NZ	7:A:1492:HOH:O	2.30	0.63
1:A:325:LYS:O	1:A:329:GLN:HG3	2.00	0.61
1:A:158[A]:ILE:CG1	7:A:1502:HOH:O	2.49	0.61
1:A:241:LYS:HD2	1:A:243:ILE:HD11	1.86	0.56
1:A:232:ARG:NH1	7:A:1471:HOH:O	2.38	0.56
1:A:234:ASN:H	1:A:662:GLN:HE22	1.53	0.55
1:A:31:ASP:CB	7:A:1513:HOH:O	2.56	0.53
1:A:477:LYS:HG3	1:A:638:LEU:HD21	1.91	0.52
1:A:293:LYS:HA	1:A:298:LEU:HD12	1.93	0.50
1:A:737:TYR:O	1:A:738:ARG:HD3	2.12	0.50
1:A:485[A]:HIS:CE1	7:A:1612:HOH:O	2.63	0.49
1:A:229:LYS:HG3	1:A:234:ASN:HD22	1.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:GLY:O	7:A:1599:HOH:O	2.20	0.49
1:A:342:MET:HA	1:A:342:MET:HE2	1.97	0.47
1:A:146:ARG:NH1	7:A:1588:HOH:O	2.47	0.47
1:A:484:HIS:CE1	1:A:489:LEU:CD1	2.97	0.46
1:A:482:PHE:CZ	1:A:486:MET:HG3	2.52	0.45
1:A:2:ASN:ND2	7:A:1510:HOH:O	2.50	0.45
1:A:213:GLN:CG	7:A:1638:HOH:O	2.65	0.44
1:A:511:GLN:OE1	7:A:1488:HOH:O	2.21	0.43
1:A:193:ILE:HD12	1:A:245:ILE:HD11	2.00	0.43
1:A:26:LYS:HD2	7:A:1512:HOH:O	2.20	0.41
1:A:219:ASN:ND2	7:A:1271:HOH:O	2.29	0.41
1:A:399:LEU:HA	1:A:403:ASP:O	2.21	0.40
1:A:196:LEU:HD23	1:A:196:LEU:HA	1.95	0.40
1:A:484:HIS:CE1	1:A:489:LEU:CG	3.04	0.40

All (2) 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)
7:A:1173:HOH:O	7:A:1189:HOH:O[3_555]	2.13	0.07
7:A:1204:HOH:O	7:A:1204:HOH:O[3_555]	2.16	0.04

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	699/762 (92%)	684 (98%)	15 (2%)	0	100 100

There are no Ramachandran outliers to report.

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	555/665 (84%)	543 (98%)	12 (2%)	60	62

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	12	HIS
1	A	107	ARG
1	A	119	SER
1	A	151	VAL
1	A	216	LEU
1	A	219	ASN
1	A	338	GLN
1	A	342	MET
1	A	546	LYS
1	A	695	ARG
1	A	744	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	79	GLN
1	A	234	ASN
1	A	283	GLN
1	A	338	GLN
1	A	484	HIS
1	A	500	ASN
1	A	582	GLN
1	A	606	ASN
1	A	613	ASN
1	A	662	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	1000	3,2	22,29,29	1.15	3 (13%)	27,45,45	2.43	3 (11%)
6	EDO	A	1001	-	3,3,3	0.40	0	2,2,2	0.46	0
5	BIT	A	800	-	23,25,25	2.57	5 (21%)	27,38,38	1.57	7 (25%)
3	VO4	A	998	2,4	1,4,4	0.69	0	0,6,6	0.00	-

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	1000	3,2	-	0/12/32/32	0/3/3/3
6	EDO	A	1001	-	-	0/1/1/1	0/0/0/0
5	BIT	A	800	-	-	0/4/34/34	0/3/4/4
3	VO4	A	998	2,4	-	0/0/0/0	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	800	BIT	C5-C4	-3.81	1.42	1.48
5	A	800	BIT	C11-N1	-3.53	1.33	1.37
5	A	800	BIT	C5-C10	2.14	1.43	1.40
4	A	1000	ADP	O4'-C1'	2.20	1.44	1.41
4	A	1000	ADP	C4-N3	2.51	1.39	1.35
4	A	1000	ADP	C2-N3	3.14	1.37	1.32
5	A	800	BIT	O1-C3	5.06	1.51	1.42
5	A	800	BIT	C11-N2	8.93	1.39	1.28

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1000	ADP	N3-C2-N1	-11.12	120.38	128.89
5	A	800	BIT	C13-C12-N1	-2.81	116.01	120.17
5	A	800	BIT	C9-C10-N2	-2.13	115.46	118.55
5	A	800	BIT	C6-C5-C10	-2.09	117.94	120.07
5	A	800	BIT	C8-C7-C6	2.14	121.46	117.95
5	A	800	BIT	O2-C4-C5	2.40	126.04	122.29
5	A	800	BIT	C17-C12-N1	2.48	123.83	120.17
4	A	1000	ADP	C2-N1-C6	2.56	123.34	118.77
5	A	800	BIT	C9-C10-C5	3.19	123.00	119.19
4	A	1000	ADP	O3B-PB-O1B	3.26	121.08	110.58

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 ⓘ

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	703/762 (92%)	0.20	47 (6%) 21 22	10, 21, 43, 71	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	698	TYR	8.8
1	A	697	ILE	7.7
1	A	702	VAL	7.1
1	A	65	VAL	7.0
1	A	625	ALA	5.4
1	A	696	ILE	5.0
1	A	744	ILE	4.7
1	A	742	THR	4.5
1	A	624	GLY	4.1
1	A	363	ALA	3.9
1	A	741	ILE	3.9
1	A	6	ASP	3.8
1	A	64	THR	3.8
1	A	66	ASP	3.7
1	A	739	PHE	3.6
1	A	8	THR	3.6
1	A	5	HIS	3.5
1	A	58	ASP	3.4
1	A	745	PHE	3.2
1	A	68	GLN	3.2
1	A	701	PHE	3.0
1	A	364	GLY	2.9
1	A	9	SER	2.9
1	A	521	GLN	2.9
1	A	746	PHE	2.7
1	A	57	SER	2.6
1	A	53	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	290	ALA	2.4
1	A	10	ASP	2.4
1	A	681	VAL	2.4
1	A	522	PRO	2.4
1	A	75	ASP	2.4
1	A	695	ARG	2.4
1	A	12	HIS	2.4
1	A	71	GLN	2.3
1	A	676	LEU	2.3
1	A	692	PHE	2.3
1	A	124	VAL	2.3
1	A	520	ARG	2.2
1	A	743	LYS	2.2
1	A	176	LEU	2.1
1	A	37	TYR	2.1
1	A	452	VAL	2.1
1	A	455	ILE	2.1
1	A	3	PRO	2.0
1	A	289	THR	2.0
1	A	453	LEU	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(Å ²)	Q<0.9
6	EDO	A	1001	4/4	0.96	0.14	1.80	31,34,34,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	BIT	A	800	22/22	0.97	0.15	0.38	9,15,17,20	0
3	VO4	A	998	5/5	1.00	0.10	-0.47	9,10,11,11	0
2	MG	A	997	1/1	0.99	0.10	-0.66	10,10,10,10	0
4	ADP	A	1000	27/27	0.99	0.08	-1.67	9,12,15,18	0
2	MG	A	999	1/1	0.97	0.07	-	24,24,24,24	1

6.5 Other polymers [i](#)

There are no such residues in this entry.