



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:17 PM GMT

PDB ID : 4E7Z
Title : Myosin VI (MD) pre-powerstroke state, P21 crystal form
Authors : Isabet, T.; Sweeney, H.L.; Houdusse, A.
Deposited on : 2012-03-19
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

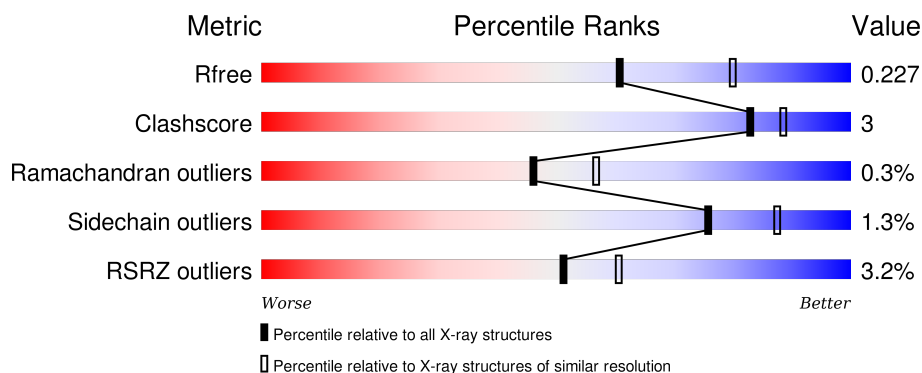
1 Overall quality at a glance

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))
R_{free}	91344	3852 (2.30-2.30)
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	798	
1	B	798	

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
4	MG	A	803	-	-	-	X
4	MG	B	803	-	-	-	X
5	GOL	A	804	-	-	X	-
5	GOL	A	806	-	-	-	X
5	GOL	B	804	-	-	-	X
5	GOL	B	805	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12444 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-VI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	759	Total	C	N	O	S	0	3	0
			6037	3845	1033	1128	31			
1	B	745	Total	C	N	O	S	0	1	0
			5849	3728	1001	1093	27			

There are 20 discrepancies between the modelled and reference sequences:

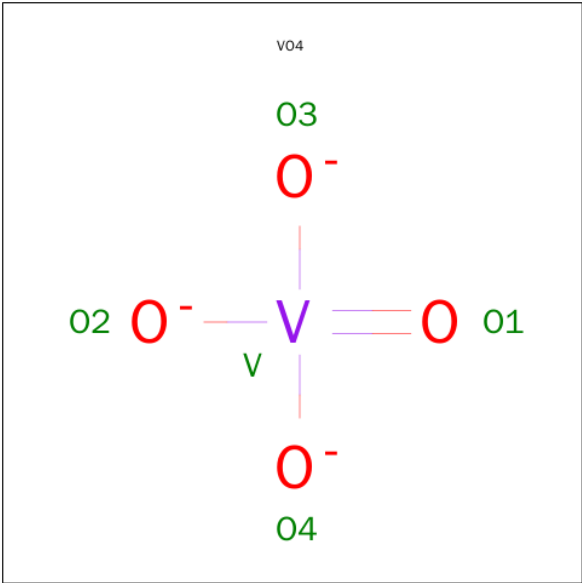
Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	MET	-	EXPRESSION TAG	UNP F1RQI7
A	-7	ASP	-	EXPRESSION TAG	UNP F1RQI7
A	-6	TYR	-	EXPRESSION TAG	UNP F1RQI7
A	-5	LYS	-	EXPRESSION TAG	UNP F1RQI7
A	-4	ASP	-	EXPRESSION TAG	UNP F1RQI7
A	-3	ASP	-	EXPRESSION TAG	UNP F1RQI7
A	-2	ASP	-	EXPRESSION TAG	UNP F1RQI7
A	-1	ASP	-	EXPRESSION TAG	UNP F1RQI7
A	0	LYS	-	EXPRESSION TAG	UNP F1RQI7
A	1	GLY	-	EXPRESSION TAG	UNP F1RQI7
B	-8	MET	-	EXPRESSION TAG	UNP F1RQI7
B	-7	ASP	-	EXPRESSION TAG	UNP F1RQI7
B	-6	TYR	-	EXPRESSION TAG	UNP F1RQI7
B	-5	LYS	-	EXPRESSION TAG	UNP F1RQI7
B	-4	ASP	-	EXPRESSION TAG	UNP F1RQI7
B	-3	ASP	-	EXPRESSION TAG	UNP F1RQI7
B	-2	ASP	-	EXPRESSION TAG	UNP F1RQI7
B	-1	ASP	-	EXPRESSION TAG	UNP F1RQI7
B	0	LYS	-	EXPRESSION TAG	UNP F1RQI7
B	1	GLY	-	EXPRESSION TAG	UNP F1RQI7

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

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

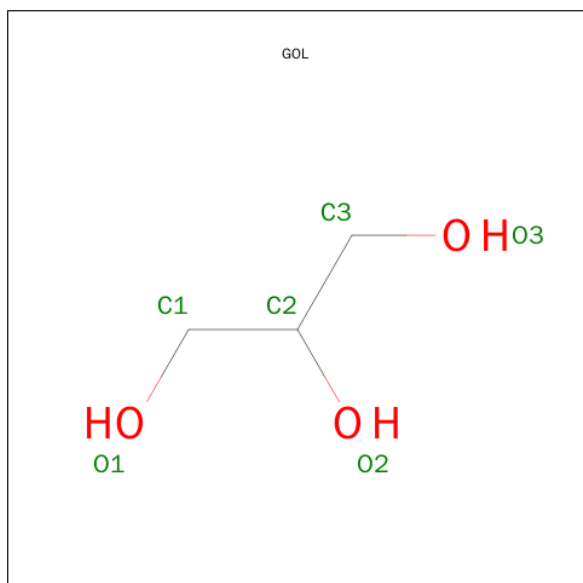


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O V	0	0
			5	4 1		
3	B	1	Total	O V	0	0
			5	4 1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	255	Total	O	0	0
			255	255		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	201	Total	O	0	0
			201	201		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.15Å 93.32Å 101.88Å 90.00° 90.60° 90.00°	Depositor
Resolution (Å)	29.65 – 2.30 29.65 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.4 (29.65-2.30) 98.4 (29.65-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.43 (at 2.31Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.176 , 0.231 0.169 , 0.227	Depositor DCC
R_{free} test set	4028 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.6	EDS
Estimated twinning fraction	0.001 for l,k,-h 0.022 for h,-k,-l 0.015 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 80428 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12444	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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: VO4, GOL, 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	0.49	0/6164	0.59	0/8323
1	B	0.47	1/5967 (0.0%)	0.58	0/8066
All	All	0.48	1/12131 (0.0%)	0.59	0/16389

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	278	CYS	CB-SG	-5.04	1.73	1.81

There are no bond angle outliers.

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	6037	0	5873	35	0
1	B	5849	0	5619	39	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	24	0	32	6	0
5	B	12	0	16	3	0
6	A	255	0	0	5	0
6	B	201	0	0	6	0
All	All	12444	0	11564	74	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 3.

All (74) 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:692:SER:HB3	5:A:805:GOL:H12	1.57	0.87
1:B:618:ARG:O	6:B:1060:HOH:O	2.05	0.72
1:B:68:LEU:HD13	1:B:702:GLN:HG3	1.74	0.69
1:B:618:ARG:NH2	5:B:805:GOL:O2	2.27	0.67
1:B:142:LYS:NZ	1:B:450:TYR:OH	2.28	0.64
1:A:68:LEU:HD13	1:A:702:GLN:HB2	1.80	0.63
1:A:169:THR:HB	1:A:179:ASP:HB3	1.80	0.63
1:A:177:ASP:N	6:A:1094:HOH:O	2.32	0.62
1:B:181:ARG:HD2	1:B:327:ILE:HA	1.82	0.62
1:B:200:ASN:HB3	1:B:203:SER:HB2	1.83	0.60
1:A:457:ILE:HD13	1:A:480:LEU:HD13	1.84	0.59
1:A:250:ARG:HH22	5:A:804:GOL:H31	1.68	0.58
1:A:305:SER:H	5:A:806:GOL:H2	1.70	0.56
1:A:174:THR:HG21	1:A:178:ILE:HD12	1.87	0.56
1:B:618:ARG:HH21	5:B:805:GOL:HO2	1.53	0.55
1:A:723:PRO:HD2	1:A:726:LEU:HD12	1.89	0.55
1:B:119:SER:OG	1:B:767:ASP:OD2	2.15	0.53
1:B:618:ARG:NE	5:B:805:GOL:O2	2.42	0.53
1:A:190:GLU:OE1	5:A:804:GOL:H32	2.10	0.51
1:A:565:LEU:HD12	1:A:568:HIS:CE1	2.46	0.50
1:A:327:ILE:HB	1:A:442[A]:CYS:SG	2.51	0.50
1:B:118:LYS:O	1:B:136:ARG:NH1	2.45	0.50
1:A:531:GLU:HG3	6:A:1145:HOH:O	2.11	0.50
1:B:614:ASP:OD2	1:B:617:ILE:HG12	2.11	0.49
1:A:310:LEU:HD21	5:A:804:GOL:O1	2.12	0.49
1:A:671:PRO:HA	1:A:685:ILE:HD11	1.95	0.49
1:A:260:ARG:HG3	1:A:265:LEU:HB2	1.94	0.49
1:A:250:ARG:NH2	5:A:804:GOL:H31	2.27	0.48
1:A:755:LYS:HB2	1:A:757:PHE:CE1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:671:PRO:HA	1:B:685:ILE:HD11	1.96	0.48
1:A:603:MET:N	1:A:603:MET:SD	2.84	0.48
1:A:233:SER:HA	1:A:425:LYS:HE2	1.95	0.47
1:B:615:LYS:HE2	1:B:619:GLU:OE1	2.14	0.47
1:A:406:VAL:HG12	1:A:407:ILE:HG12	1.96	0.47
1:B:180:ASP:O	1:B:184:GLU:HG2	2.14	0.47
1:B:5:LYS:HA	1:B:6:PRO:HD3	1.71	0.47
1:B:719:LYS:HA	6:B:1066:HOH:O	2.16	0.46
1:A:780:LEU:O	1:A:784:VAL:HG23	2.14	0.46
1:A:364:LEU:HD13	1:A:387:ARG:HG3	1.96	0.46
1:B:563:SER:HB2	1:B:578:ILE:HD11	1.96	0.46
1:A:177:ASP:O	1:A:181:ARG:HG3	2.17	0.45
1:A:232:LYS:HE3	1:A:429:SER:HB2	1.97	0.45
1:B:61:ASP:OD2	6:B:1097:HOH:O	2.21	0.45
1:A:742:LEU:HD21	1:A:780:LEU:HD22	1.98	0.45
1:B:372:LEU:HD12	1:B:386:LEU:HD23	1.99	0.44
1:A:5:LYS:HA	1:A:6:PRO:HD3	1.86	0.44
1:B:729:LEU:HB3	1:B:730:ASP:H	1.41	0.44
1:A:166:ARG:NH1	6:A:1155:HOH:O	2.48	0.44
1:B:382:ASP:HB3	1:B:385:ASP:OD2	2.17	0.44
1:A:765:GLU:N	1:A:765:GLU:OE2	2.33	0.44
1:B:98:ASN:HB2	1:B:155:ALA:O	2.17	0.44
1:B:487:ARG:NH1	6:B:1041:HOH:O	2.44	0.44
1:A:680:PHE:CE1	1:A:685:ILE:HD13	2.53	0.43
1:B:132:ASP:OD1	1:B:172:TYR:OH	2.30	0.43
1:A:705:PHE:HB3	1:A:757:PHE:HB3	2.00	0.43
1:B:10:PRO:HD3	1:B:47:GLN:O	2.19	0.43
1:B:188:LEU:HD12	1:B:188:LEU:HA	1.81	0.43
1:B:479:LYS:HE3	1:B:479:LYS:HB3	1.90	0.42
1:A:135:PHE:HA	1:A:452:ILE:HD11	2.02	0.42
1:B:484:PHE:HZ	1:B:697:VAL:HG21	1.84	0.42
1:A:732:ARG:HD2	6:A:1148:HOH:O	2.19	0.42
1:B:370:GLN:HG2	1:B:374:TYR:CE1	2.55	0.41
1:B:714:LEU:N	6:B:1083:HOH:O	2.43	0.41
1:B:622:GLU:O	6:B:1060:HOH:O	2.22	0.41
1:B:653:LEU:HA	1:B:653:LEU:HD23	1.94	0.41
1:B:18:VAL:HG11	1:B:34:LEU:HD22	2.02	0.41
1:B:330:ASP:OD2	1:B:333:GLU:HG3	2.20	0.41
1:B:243:ARG:HB3	1:B:249:TYR:CZ	2.55	0.41
1:B:243:ARG:HB3	1:B:249:TYR:CE1	2.56	0.40
1:A:567:ILE:HG12	6:A:1144:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:GLU:CD	1:B:216:GLU:H	2.24	0.40
1:A:232:LYS:HE3	1:A:429:SER:CB	2.51	0.40
1:B:766:PHE:O	1:B:770:MET:HG2	2.22	0.40
1:B:311:LEU:HD23	1:B:311:LEU:HA	1.86	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	754/798 (94%)	733 (97%)	19 (2%)	2 (0%)	46	57
1	B	736/798 (92%)	706 (96%)	27 (4%)	3 (0%)	39	48
All	All	1490/1596 (93%)	1439 (97%)	46 (3%)	5 (0%)	46	57

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	408	LYS
1	A	26	PRO
1	B	720	LYS
1	B	712	HIS
1	B	140	VAL

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	651/705 (92%)	639 (98%)	12 (2%)	66	82
1	B	617/705 (88%)	612 (99%)	5 (1%)	86	94
All	All	1268/1410 (90%)	1251 (99%)	17 (1%)	76	87

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	34	LEU
1	A	57	LYS
1	A	158	THR
1	A	179	ASP
1	A	276	ARG
1	A	480	LEU
1	A	603	MET
1	A	708	ARG
1	A	733	LEU
1	A	736	LYS
1	A	783	ARG
1	B	450	TYR
1	B	606	GLU
1	B	638	LEU
1	B	707	SER
1	B	708	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	20	ASN
1	B	290	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 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
2	ADP	A	801	3,4	22,29,29	1.06	2 (9%)	27,45,45	1.74	2 (7%)
3	VO4	A	802	2,4	1,4,4	4.80	1 (100%)	0,6,6	0.00	-
5	GOL	A	804	-	5,5,5	0.50	0	5,5,5	0.60	0
5	GOL	A	805	-	5,5,5	0.39	0	5,5,5	0.73	0
5	GOL	A	806	-	5,5,5	0.25	0	5,5,5	0.44	0
5	GOL	A	807	-	5,5,5	0.35	0	5,5,5	0.61	0
2	ADP	B	801	3,4	22,29,29	1.01	1 (4%)	27,45,45	1.90	4 (14%)
3	VO4	B	802	2,4	1,4,4	4.78	1 (100%)	0,6,6	0.00	-
5	GOL	B	804	-	5,5,5	0.30	0	5,5,5	0.35	0
5	GOL	B	805	-	5,5,5	0.35	0	5,5,5	0.24	0

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	801	3,4	-	0/12/32/32	0/3/3/3
3	VO4	A	802	2,4	-	0/0/0/0	0/0/0/0
5	GOL	A	804	-	-	0/4/4/4	0/0/0/0
5	GOL	A	805	-	-	0/4/4/4	0/0/0/0
5	GOL	A	806	-	-	0/4/4/4	0/0/0/0
5	GOL	A	807	-	-	0/4/4/4	0/0/0/0
2	ADP	B	801	3,4	-	0/12/32/32	0/3/3/3
3	VO4	B	802	2,4	-	0/0/0/0	0/0/0/0
5	GOL	B	804	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	B	805	-	-	0/4/4/4	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	ADP	O4'-C1'	2.03	1.43	1.41
2	B	801	ADP	C5-C4	2.49	1.46	1.40
2	A	801	ADP	C5-C4	2.69	1.46	1.40
3	B	802	VO4	O1-V	4.78	1.94	1.63
3	A	802	VO4	O1-V	4.80	1.94	1.63

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	ADP	N3-C2-N1	-7.32	123.29	128.89
2	A	801	ADP	N3-C2-N1	-6.72	123.75	128.89
2	A	801	ADP	C4-C5-N7	-3.41	106.34	109.48
2	B	801	ADP	C4-C5-N7	-2.79	106.91	109.48
2	B	801	ADP	C2-N1-C6	2.00	122.34	118.77
2	B	801	ADP	O2B-PB-O1B	3.14	120.70	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	804	GOL	4	0
5	A	805	GOL	1	0
5	A	806	GOL	1	0
5	B	805	GOL	3	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	759/798 (95%)	-0.21	19 (2%) 61 70	23, 39, 61, 90	0
1	B	745/798 (93%)	-0.02	29 (3%) 43 52	24, 43, 80, 114	0
All	All	1504/1596 (94%)	-0.12	48 (3%) 51 60	23, 41, 71, 114	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	638	LEU	5.7
1	B	638	LEU	5.5
1	A	35	ASN	5.4
1	B	786	HIS	4.7
1	A	406	VAL	4.3
1	B	410	PRO	4.2
1	B	35	ASN	3.9
1	A	37	LYS	3.7
1	B	785	ASN	3.2
1	B	361	GLY	3.2
1	B	366	ASN	3.1
1	B	409	VAL	3.1
1	B	38	GLY	3.0
1	A	407	ILE	3.0
1	A	747	ILE	3.0
1	B	731	PRO	2.9
1	B	772	SER	2.9
1	B	774	PRO	2.8
1	B	787	TRP	2.8
1	B	497	GLN	2.8
1	B	394	VAL	2.8
1	B	776	HIS	2.7
1	A	395	MET	2.7
1	B	502	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	411	LEU	2.6
1	B	712	HIS	2.6
1	A	161	THR	2.6
1	B	753	LEU	2.6
1	B	754	THR	2.6
1	B	749	TYR	2.5
1	A	503	VAL	2.5
1	B	447	THR	2.5
1	A	38	GLY	2.4
1	A	753	LEU	2.4
1	A	603	MET	2.4
1	A	396	LEU	2.3
1	B	248	PHE	2.3
1	B	727	ALA	2.3
1	A	341	VAL	2.2
1	A	303	ALA	2.2
1	A	721	TYR	2.1
1	B	721	TYR	2.1
1	A	177	ASP	2.1
1	B	639	SER	2.1
1	A	537	GLN	2.1
1	A	344	VAL	2.1
1	B	93	ILE	2.1
1	B	344	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
5	GOL	A	806	6/6	0.93	0.22	7.16	41,49,55,57	0
5	GOL	B	805	6/6	0.84	0.36	6.46	42,53,55,59	0
4	MG	A	803	1/1	0.97	0.38	5.34	32,32,32,32	0
5	GOL	B	804	6/6	0.97	0.20	5.32	49,51,52,55	0
4	MG	B	803	1/1	0.95	0.33	5.18	32,32,32,32	0
5	GOL	A	807	6/6	0.90	0.17	1.80	32,48,54,61	0
3	VO4	B	802	5/5	0.99	0.15	0.43	22,23,28,30	0
3	VO4	A	802	5/5	0.99	0.16	0.17	22,23,26,27	0
2	ADP	B	801	27/27	0.99	0.12	0.13	22,25,32,34	0
5	GOL	A	804	6/6	0.91	0.17	0.10	42,42,43,48	0
5	GOL	A	805	6/6	0.91	0.13	-0.00	36,40,42,53	0
2	ADP	A	801	27/27	0.99	0.14	-0.08	20,24,28,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.