Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:47 PM GMT

PDB ID	:	1H9R
Title	:	TUNGSTATE BOUND COMPLEX DIMOP DOMAIN OF MODE FROM E.COLI
Authors	:	Gourley, D.G.; Hunter, W.N.
Deposited on	:	2001-03-19
Resolution	:	1.90 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 (i) symbol.

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

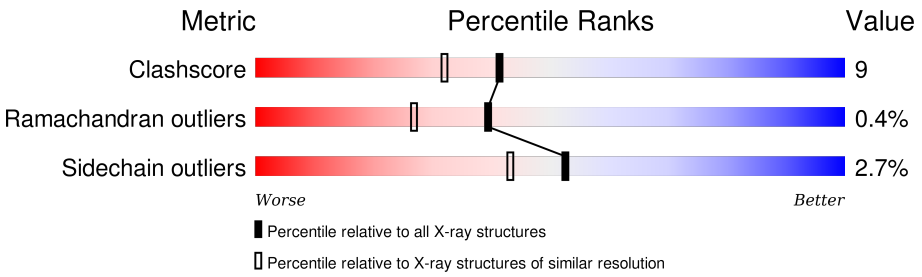
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	140	80% 18% ••
1	B	140	81% 16% ••

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2417 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.

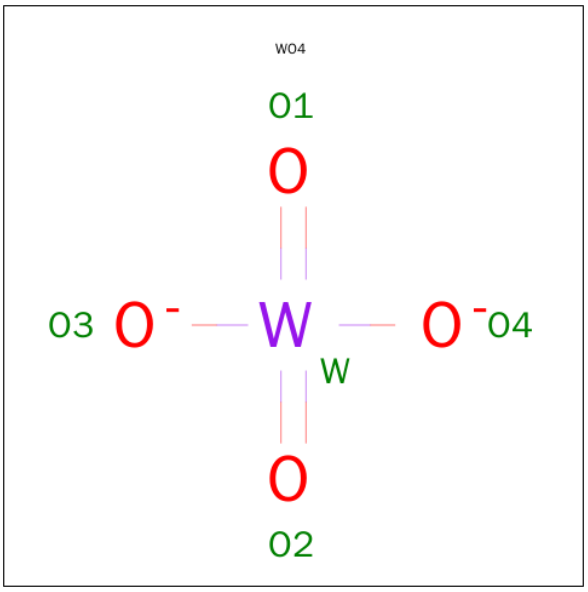
- Molecule 1 is a protein called MOLYBDENUM TRANSPORT PROTEIN MODE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	139	1047	649	185	209	4	17	0	0
1	B	139	1047	649	185	209	4	12	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	MET	LEU	ENGINEERED MUTATION	UNP P46930
B	123	MET	LEU	ENGINEERED MUTATION	UNP P46930

- Molecule 2 is TUNGSTATE(VI)ION (three-letter code: WO4) (formula: O_4W).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	W		
2	A	1	5	4	1	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	W		
2	B	1	5	4	1	0	0

- Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ni		
3	B	1	1	1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	165	165	165	0	0
			Total	O		
4	B	147	147	147	0	0

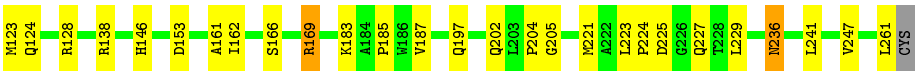
3 Residue-property plots (i)

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.

Note EDS was not executed.

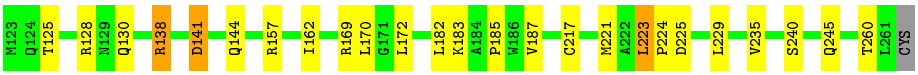
- Molecule 1: MOLYBDENUM TRANSPORT PROTEIN MODE

Chain A:



- Molecule 1: MOLYBDENUM TRANSPORT PROTEIN MODE

Chain B:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α, β, γ	72.23Å 73.23Å 49.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90	Depositor
% Data completeness (in resolution range)	97.9 (20.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.49	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.170 , 0.225	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2417	wwPDB-VP
Average B, all atoms ($Å^2$)	28.0	wwPDB-VP

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: NI, WO4

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	0.59	0/1060	1.22	6/1444 (0.4%)
1	B	0.73	1/1060 (0.1%)	1.31	10/1444 (0.7%)
All	All	0.66	1/2120 (0.0%)	1.26	16/2888 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	169	ARG	CG-CD	16.12	1.92	1.51

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	138	ARG	NE-CZ-NH2	-13.50	113.55	120.30
1	B	169	ARG	CG-CD-NE	13.03	139.17	111.80
1	A	169	ARG	CD-NE-CZ	7.50	134.10	123.60
1	B	169	ARG	CD-NE-CZ	7.44	134.02	123.60
1	B	157	ARG	CD-NE-CZ	7.15	133.62	123.60
1	B	128	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	B	157	ARG	NE-CZ-NH1	6.39	123.50	120.30
1	B	169	ARG	CB-CG-CD	-6.21	95.46	111.60
1	B	141	ASP	CB-CG-OD1	5.67	123.40	118.30
1	B	138	ARG	NH1-CZ-NH2	5.63	125.60	119.40
1	A	123	MET	CA-C-O	5.58	131.82	120.10
1	A	169	ARG	NE-CZ-NH1	5.47	123.04	120.30
1	A	123	MET	C-N-CA	5.47	135.37	121.70
1	B	225	ASP	CB-CG-OD1	5.35	123.12	118.30
1	A	128	ARG	NE-CZ-NH1	5.27	122.93	120.30
1	A	138	ARG	NE-CZ-NH2	-5.06	117.77	120.30

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	1047	0	1040	22	0
1	B	1047	0	1040	13	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	B	1	0	0	0	0
4	A	165	0	0	8	0
4	B	147	0	0	2	0
All	All	2417	0	2080	35	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 9.

All (35) 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:205:GLY:N	1:A:221:MET:HE2	1.97	0.79
1:A:153:ASP:HB3	4:A:2036:HOH:O	1.94	0.68
1:A:225:ASP:HA	4:A:2036:HOH:O	1.99	0.62
1:A:187:VAL:HG11	1:A:229:LEU:HD21	1.79	0.62
1:B:217:CYS:SG	1:B:235:VAL:HG12	2.44	0.57
1:B:138:ARG:HD2	1:B:172:LEU:O	2.05	0.56
1:A:225:ASP:HB3	1:A:227:GLN:H	1.69	0.56
1:A:225:ASP:HB3	1:A:227:GLN:HB2	1.91	0.53
1:A:236:ASN:H	1:A:236:ASN:HD22	1.55	0.52
1:A:225:ASP:CG	1:A:227:GLN:HE21	2.13	0.52
1:A:261:LEU:C	4:A:2163:HOH:O	2.48	0.51
1:A:223:LEU:HB3	4:A:2123:HOH:O	2.13	0.48
1:B:240:SER:HB2	4:B:2125:HOH:O	2.14	0.47
1:A:221:MET:CE	1:A:247:VAL:HG23	2.45	0.47
1:A:166:SER:HA	1:A:169:ARG:NH1	2.30	0.46
1:A:225:ASP:OD2	1:A:227:GLN:NE2	2.49	0.46
1:A:197:GLN:HG2	4:A:2082:HOH:O	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:HIS:HD2	4:A:2022:HOH:O	1.99	0.45
1:A:223:LEU:C	1:A:225:ASP:H	2.20	0.44
1:B:182:LEU:HD21	1:B:187:VAL:HG22	1.99	0.44
1:B:170:LEU:HD12	1:B:172:LEU:HD11	2.00	0.44
1:A:204:PRO:C	1:A:221:MET:HE2	2.37	0.44
1:B:187:VAL:HG11	1:B:229:LEU:HD21	2.00	0.44
1:A:241:LEU:HD11	4:A:2068:HOH:O	2.18	0.43
1:A:183:LYS:HG2	1:A:185:PRO:HD2	1.99	0.43
1:B:125:THR:OG1	1:B:130:GLN:NE2	2.51	0.43
1:B:141:ASP:HB2	1:B:144:GLN:O	2.19	0.43
1:B:245:GLN:HG2	4:B:2094:HOH:O	2.17	0.43
1:A:223:LEU:C	1:A:225:ASP:N	2.72	0.42
1:A:224:PRO:HD2	4:A:2123:HOH:O	2.18	0.42
1:B:183:LYS:HG2	1:B:185:PRO:HD2	2.02	0.42
1:B:170:LEU:HB2	1:B:172:LEU:HG	2.03	0.41
1:B:223:LEU:HB3	1:B:224:PRO:HD2	2.02	0.41
1:B:221:MET:CE	1:B:223:LEU:HD13	2.51	0.41
1:A:146:HIS:CE1	1:A:161:ALA:HB2	2.57	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	137/140 (98%)	134 (98%)	2 (2%)	1 (1%)	26 14
1	B	137/140 (98%)	136 (99%)	1 (1%)	0	100 100
All	All	274/280 (98%)	270 (98%)	3 (1%)	1 (0%)	39 27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	GLN

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	112/113 (99%)	109 (97%)	3 (3%)	52	43
1	B	112/113 (99%)	109 (97%)	3 (3%)	52	43
All	All	224/226 (99%)	218 (97%)	6 (3%)	52	43

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	ILE
1	A	202	GLN
1	A	236	ASN
1	B	162	ILE
1	B	223	LEU
1	B	260	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	HIS
1	A	227	GLN
1	A	236	ASN
1	B	130	GLN
1	B	146	HIS
1	B	198	ASN
1	B	243	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 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	WO4	A	1262	1	2,4,4	1.76	1 (50%)	0,6,6	0.00	-
2	WO4	B	1262	1	2,4,4	2.31	1 (50%)	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
2	WO4	A	1262	1	-	0/0/0/0	0/0/0/0
2	WO4	B	1262	1	-	0/0/0/0	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1262	WO4	W-O1	2.34	1.80	1.74
2	B	1262	WO4	W-O1	3.13	1.82	1.74

There are no bond angle outliers.

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)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates (i)

EDS was not executed - this section will therefore be empty.

6.4 Ligands (i)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers (i)

EDS was not executed - this section will therefore be empty.