



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 09:16 AM GMT

PDB ID : 3HOS
Title : Crystal structure of the mariner Mos1 paired end complex with Mg
Authors : Richardson, J.M.; Walkinshaw, M.D.
Deposited on : 2009-06-03
Resolution : 3.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

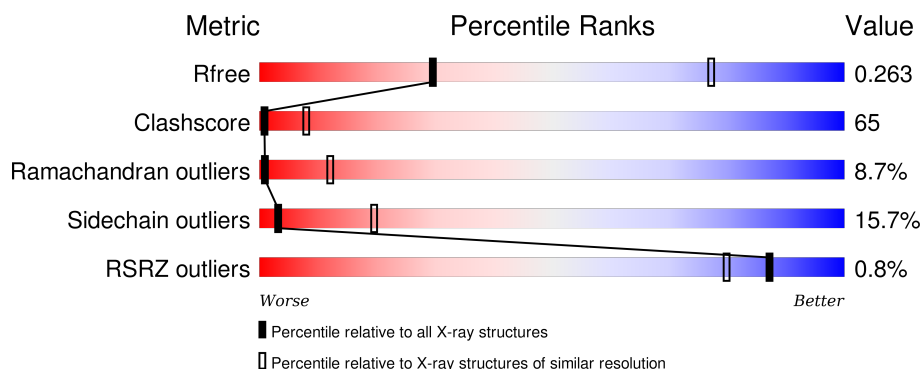
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	345	<div> <div>22%</div> <div>60%</div> <div>14%</div> <div>..</div> </div>
1	B	345	<div> <div>22%</div> <div>57%</div> <div>17%</div> <div>..</div> </div>
2	C	25	<div> <div>24%</div> <div>76%</div> </div>
2	E	25	<div> <div>32%</div> <div>64%</div> </div>
2	G	25	<div> <div>52%</div> <div>44%</div> </div>

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Mol	Chain	Length	Quality of chain
3	D	28	 <div> <div></div> <div>32%</div> <div>64%</div> </div>
3	F	28	 <div> <div></div> <div>29%</div> <div>68%</div> </div>
3	H	28	 <div> <div></div> <div>7%</div> <div>54%</div> <div>39%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8885 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 Transposable element mariner, complete cds.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	0	0
			2802	1778	507	507	10			
1	B	337	Total	C	N	O	S	0	0	0
			2812	1783	509	510	10			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	THR	LYS	SEE REMARK 999	UNP Q7JQ07
A	164	ASN	SER	SEE REMARK 999	UNP Q7JQ07
A	210	PRO	ARG	SEE REMARK 999	UNP Q7JQ07
A	216	ALA	THR	ENGINEERED	UNP Q7JQ07
A	344	PHE	LEU	SEE REMARK 999	UNP Q7JQ07
B	45	THR	LYS	SEE REMARK 999	UNP Q7JQ07
B	164	ASN	SER	SEE REMARK 999	UNP Q7JQ07
B	210	PRO	ARG	SEE REMARK 999	UNP Q7JQ07
B	216	ALA	THR	ENGINEERED	UNP Q7JQ07
B	344	PHE	LEU	SEE REMARK 999	UNP Q7JQ07

- Molecule 2 is a DNA chain called Mos1 NTS inverted repeat DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	25	Total	C	N	O	P	0	0	0
			520	248	94	153	25			
2	E	25	Total	C	N	O	P	0	0	0
			520	248	94	153	25			
2	G	25	Total	C	N	O	P	0	0	0
			520	248	94	153	25			

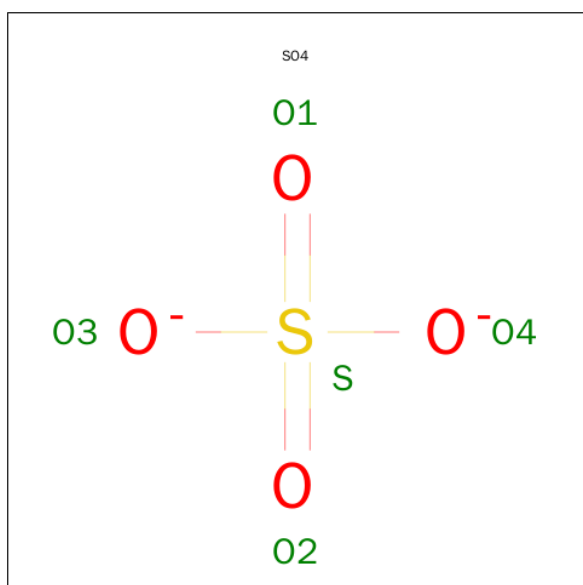
- Molecule 3 is a DNA chain called Mos1 TS inverted repeat DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	28	Total	C	N	O	P	0	0	0
			566	273	102	164	27			
3	F	28	Total	C	N	O	P	0	0	0
			566	273	102	164	27			
3	H	28	Total	C	N	O	P	0	0	0
			566	273	102	164	27			

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

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

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

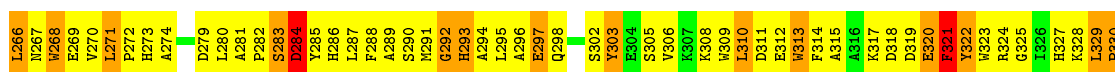
- Molecule 6 is water.

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	O	0	0
			1	1		



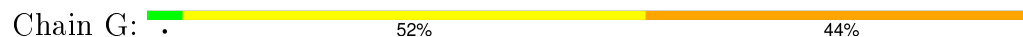
- Molecule 2: Mos1 NTS inverted repeat DNA



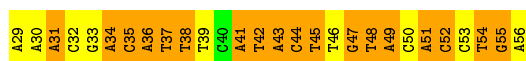
- Molecule 2: Mos1 NTS inverted repeat DNA



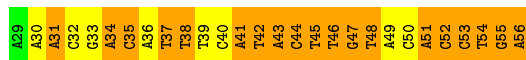
- Molecule 2: Mos1 NTS inverted repeat DNA



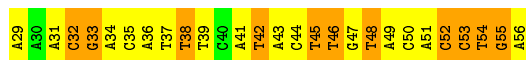
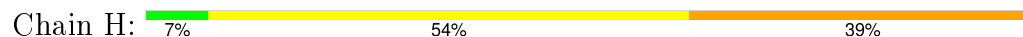
- Molecule 3: Mos1 TS inverted repeat DNA



- Molecule 3: Mos1 TS inverted repeat DNA



- Molecule 3: Mos1 TS inverted repeat DNA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	121.23Å 85.03Å 131.27Å 90.00° 98.93° 90.00°	Depositor
Resolution (Å)	29.87 – 3.50 29.87 – 3.50	Depositor EDS
% Data completeness (in resolution range)	96.7 (29.87-3.50) 96.7 (29.87-3.50)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 3.47Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.219 , 0.279 0.219 , 0.263	Depositor DCC
R_{free} test set	1647 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	89.1	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.22 , 116.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Outliers	0 of 32494 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	8885	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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, SO4

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	0.81	0/2875	0.83	0/3881
1	B	0.74	0/2885	0.82	0/3893
2	C	1.41	1/583 (0.2%)	2.39	52/898 (5.8%)
2	E	1.52	5/583 (0.9%)	2.51	44/898 (4.9%)
2	G	1.25	2/583 (0.3%)	2.20	41/898 (4.6%)
3	D	1.54	4/634 (0.6%)	2.48	48/975 (4.9%)
3	F	1.54	8/634 (1.3%)	2.46	67/975 (6.9%)
3	H	1.11	0/634	1.85	24/975 (2.5%)
All	All	1.06	20/9411 (0.2%)	1.63	276/13393 (2.1%)

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	56	DA	N9-C4	-9.95	1.31	1.37
2	G	4	DG	OP3-P	-9.83	1.49	1.61
2	C	4	DG	OP3-P	-9.48	1.49	1.61
2	E	4	DG	OP3-P	-7.87	1.51	1.61
2	G	28	DT	C4-C5	-7.11	1.38	1.45

The worst 5 of 276 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	4	DG	OP1-P-OP2	-24.98	82.14	119.60
2	E	4	DG	O5'-P-OP2	-16.21	91.11	105.70
2	G	28	DT	N3-C4-O4	12.95	127.67	119.90
2	G	4	DG	N1-C6-O6	12.74	127.55	119.90
3	F	48	DT	N3-C4-O4	12.30	127.28	119.90

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	2802	0	2739	398	0
1	B	2812	0	2747	415	0
2	C	520	0	285	54	0
2	E	520	0	285	33	0
2	G	520	0	285	52	0
3	D	566	0	318	76	0
3	F	566	0	317	48	0
3	H	566	0	318	61	0
4	A	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
All	All	8885	0	7294	1044	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 65.

The worst 5 of 1044 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:174:PRO:HA	1:B:113:ILE:HD11	1.26	1.17
3:H:36:DA:H2''	3:H:37:DT:H5'	1.23	1.15
3:H:45:DT:H2''	3:H:46:DT:C5'	1.75	1.14
3:D:43:DA:H2''	3:D:44:DC:H5''	1.27	1.14
3:H:33:DG:H2''	3:H:34:DA:H5''	1.31	1.11

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	332/345 (96%)	224 (68%)	86 (26%)	22 (7%)	1	19
1	B	333/345 (96%)	235 (71%)	62 (19%)	36 (11%)	0	8
All	All	665/690 (96%)	459 (69%)	148 (22%)	58 (9%)	1	12

5 of 58 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	62	ASP
1	A	126	GLU
1	A	137	GLU
1	A	150	HIS

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	298/307 (97%)	248 (83%)	50 (17%)	2	15
1	B	299/307 (97%)	255 (85%)	44 (15%)	4	22
All	All	597/614 (97%)	503 (84%)	94 (16%)	3	19

5 of 94 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	297	GLU
1	B	24	THR
1	B	271	LEU
1	A	304	GLU
1	A	340	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	293	HIS
1	B	95	GLN
1	B	254	HIS
1	B	10	GLN
1	B	100	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	$\# Z > 2$
5	SO4	A	346	-	4,4,4	0.21	0	6,6,6	0.24	0
5	SO4	B	346	-	4,4,4	0.19	0	6,6,6	0.31	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
5	SO4	A	346	-	-	0/0/0/0	0/0/0/0
5	SO4	B	346	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

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](#)

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	336/345 (97%)	-0.74	3 (0%) 85 78	14, 58, 131, 225	0
1	B	337/345 (97%)	-0.63	4 (1%) 81 72	18, 63, 137, 226	0
2	C	25/25 (100%)	-0.63	0 100 100	34, 53, 83, 122	0
2	E	25/25 (100%)	-0.55	0 100 100	35, 60, 80, 96	0
2	G	25/25 (100%)	-0.46	0 100 100	25, 94, 144, 156	0
3	D	28/28 (100%)	-0.74	0 100 100	36, 52, 76, 91	0
3	F	28/28 (100%)	-0.67	0 100 100	34, 62, 92, 105	0
3	H	28/28 (100%)	-0.43	0 100 100	28, 88, 150, 150	0
All	All	832/849 (97%)	-0.67	7 (0%) 87 80	14, 62, 136, 226	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	155	GLY	3.4
1	B	156	ASP	2.9
1	A	156	ASP	2.6
1	A	61	ASP	2.5
1	B	154	THR	2.1

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
4	MG	A	1001	1/1	0.97	0.16	-	2,2,2,2	0
5	SO4	B	346	5/5	0.93	0.17	-	103,112,125,135	0
5	SO4	A	346	5/5	0.95	0.32	-	117,126,130,139	0

6.5 Other polymers [i](#)

There are no such residues in this entry.