



Full wwPDB X-ray Structure Validation Report

Mar 31, 2014 – 03:52 PM BST

PDB ID : 4F3F
Title : Crystal Structure of Msln7-64 MORAb-009 FAB complex
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Deposited on : 2012-05-09
Resolution : 2.65 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

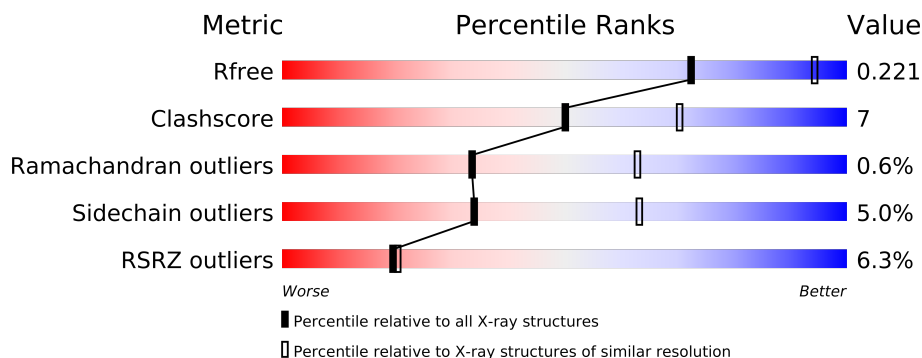
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable23004
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23004

1 Overall quality at a glance

The reported resolution of this entry is 2.65 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}	66092	2232 (2.70-2.62)
Clashscore	79885	2700 (2.70-2.62)
Ramachandran outliers	78287	2657 (2.70-2.62)
Sidechain outliers	78261	2657 (2.70-2.62)
RSRZ outliers	66119	2234 (2.70-2.62)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	213	
2	B	231	
3	C	69	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7476 atoms, of which 3627 are hydrogens and 0 are deuterium.

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 MORAb-009 Fab light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	211	Total	C	H	N	O	S	0	0	0
			3166	1006	1551	270	332	7			

- Molecule 2 is a protein called MORAb-009 Fab heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	220	Total	C	H	N	O	S	0	0	0
			3232	1033	1597	269	327	6			

- Molecule 3 is a protein called Mesothelin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	59	Total	C	H	N	O	S	0	0	0
			956	306	479	77	90	4			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	6	MET	-	EXPRESSION TAG	UNP Q13421
C	65	GLY	-	EXPRESSION TAG	UNP Q13421
C	66	SER	-	EXPRESSION TAG	UNP Q13421
C	67	LEU	-	EXPRESSION TAG	UNP Q13421
C	68	GLU	-	EXPRESSION TAG	UNP Q13421
C	69	HIS	-	EXPRESSION TAG	UNP Q13421
C	70	HIS	-	EXPRESSION TAG	UNP Q13421
C	71	HIS	-	EXPRESSION TAG	UNP Q13421
C	72	HIS	-	EXPRESSION TAG	UNP Q13421
C	73	HIS	-	EXPRESSION TAG	UNP Q13421
C	74	HIS	-	EXPRESSION TAG	UNP Q13421

- Molecule 4 is water.

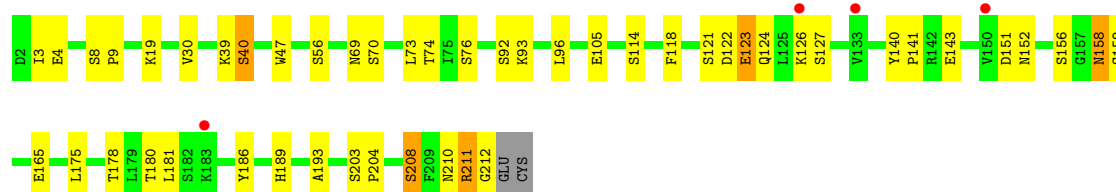
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	51	Total 51	O 51	0	0
4	B	45	Total 45	O 45	0	0
4	C	26	Total 26	O 26	0	0

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.

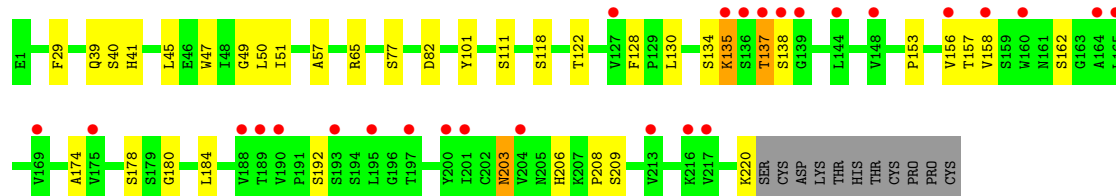
- Molecule 1: MORAb-009 Fab light chain

Chain A: 



- Molecule 2: MORAb-009 Fab heavy chain

Chain B: 



- Molecule 3: Mesothelin

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	146.16Å 146.16Å 80.87Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.08 – 2.65 49.84 – 2.61	Depositor EDS
% Data completeness (in resolution range)	98.9 (34.08-2.65) 97.7 (49.84-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_965)	Depositor
R, R_{free}	0.177 , 0.222 0.176 , 0.221	Depositor DCC
R_{free} test set	1484 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 22.7	EDS
Estimated twinning fraction	0.037 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 29240 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7476	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: PCA

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	0.82	1/1652 (0.1%)	0.79	0/2241
2	B	0.75	0/1669	0.74	0/2272
3	C	0.99	1/485 (0.2%)	0.85	0/653
All	All	0.81	2/3806 (0.1%)	0.78	0/5166

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	52	GLU	CB-CG	-5.72	1.41	1.52
1	A	143	GLU	CG-CD	5.32	1.59	1.51

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1615	1551	0	28	0
2	B	1635	1597	1	23	0
3	C	477	479	0	10	0
4	A	51	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	45	0	0	0	0
4	C	26	0	0	2	0
All	All	3849	3627	1	55	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (55) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:101:TYR:CE1	3:C:52:GLU:HG2	2.15	0.81
1:A:210:ASN:O	1:A:212:GLY:N	2.26	0.68
1:A:151:ASP:OD2	1:A:189:HIS:HB3	1.94	0.68
1:A:19:LYS:HG3	1:A:76:SER:HA	1.78	0.66
2:B:184:LEU:C	2:B:184:LEU:HD12	2.20	0.62
2:B:156:VAL:HG12	2:B:206:HIS:CD2	2.34	0.61
1:A:39:LYS:O	1:A:40:SER:C	2.39	0.61
2:B:101:TYR:CZ	3:C:52:GLU:HG2	2.36	0.60
1:A:121:SER:HG	2:B:128:PHE:HB3	1.67	0.60
1:A:40:SER:HB3	1:A:165:GLU:HG3	1.83	0.60
1:A:193:ALA:HB2	1:A:208:SER:HB3	1.83	0.59
1:A:123:GLU:N	1:A:123:GLU:OE1	2.35	0.59
3:C:45:ASN:HB3	4:C:121:HOH:O	2.03	0.58
2:B:174:ALA:HA	2:B:184:LEU:HB3	1.86	0.57
1:A:73:LEU:HD23	1:A:73:LEU:C	2.24	0.57
2:B:51:ILE:HG13	2:B:57:ALA:O	2.04	0.57
3:C:25:LYS:HD3	3:C:52:GLU:HB3	1.86	0.56
1:A:124:GLN:O	1:A:127:SER:HB2	2.05	0.56
1:A:40:SER:CB	1:A:165:GLU:HG3	2.37	0.54
2:B:50:LEU:C	2:B:50:LEU:HD12	2.30	0.52
2:B:206:HIS:ND1	2:B:209:SER:OG	2.43	0.52
2:B:135:LYS:HE3	2:B:135:LYS:HA	1.92	0.52
2:B:40:SER:O	2:B:41:HIS:C	2.48	0.52
1:A:8:SER:HA	1:A:9:PRO:C	2.31	0.51
2:B:101:TYR:CE1	3:C:52:GLU:CG	2.91	0.51
2:B:134:SER:HA	2:B:137:THR:OG1	2.12	0.49
1:A:73:LEU:HD23	1:A:74:THR:N	2.27	0.49
1:A:180:THR:O	1:A:181:LEU:HD23	2.11	0.49
1:A:151:ASP:O	1:A:152:ASN:HB2	2.13	0.48
3:C:8:PRO:O	3:C:9:SER:C	2.53	0.47
1:A:140:TYR:CG	1:A:141:PRO:HA	2.51	0.46
1:A:175:LEU:HD23	1:A:175:LEU:C	2.36	0.46
3:C:24:LYS:O	3:C:25:LYS:C	2.52	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:47:TRP:CZ2	2:B:49:GLY:HA2	2.51	0.46
1:A:30:VAL:HA	1:A:92:SER:HG	1.81	0.46
2:B:29:PHE:CD2	2:B:77:SER:HA	2.51	0.46
1:A:203:SER:HB2	1:A:204:PRO:CD	2.46	0.45
3:C:33:ASP:HB3	4:C:114:HOH:O	2.16	0.45
2:B:178:SER:C	2:B:180:GLY:N	2.69	0.45
1:A:122:ASP:N	4:A:343:HOH:O	2.49	0.45
1:A:122:ASP:O	1:A:126:LYS:HD2	2.18	0.44
1:A:159:SER:HA	1:A:178:THR:O	2.18	0.42
2:B:122:THR:CG2	2:B:209:SER:HB3	2.50	0.42
1:A:118:PHE:CB	2:B:130:LEU:HD22	2.50	0.42
2:B:39:GLN:HB2	2:B:45:LEU:HD23	2.02	0.41
1:A:158:ASN:OD1	1:A:158:ASN:N	2.53	0.41
1:A:3:ILE:HD12	1:A:93:LYS:HB2	2.03	0.41
2:B:156:VAL:HG12	2:B:206:HIS:HD2	1.81	0.40
3:C:37:LEU:HD23	3:C:61:LEU:CD2	2.52	0.40
2:B:137:THR:HG22	2:B:192:SER:HG	1.86	0.40
2:B:153:PRO:HD2	2:B:208:PRO:CB	2.51	0.40
1:A:96:LEU:HG	3:C:22:PHE:CZ	2.57	0.40
2:B:158:VAL:HA	2:B:203:ASN:O	2.22	0.40
1:A:186:TYR:CE2	1:A:211:ARG:HD2	2.57	0.40
1:A:69:ASN:OD1	1:A:69:ASN:N	2.54	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	209/213 (98%)	191 (91%)	15 (7%)	3 (1%)	16	37
2	B	218/231 (94%)	207 (95%)	11 (5%)	0	100	100
3	C	57/69 (83%)	53 (93%)	4 (7%)	0	100	100
All	All	484/513 (94%)	451 (93%)	30 (6%)	3 (1%)	33	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	211	ARG
1	A	40	SER
1	A	156	SER

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	185/187 (99%)	176 (95%)	9 (5%)	35	64
2	B	184/195 (94%)	173 (94%)	11 (6%)	27	53
3	C	52/61 (85%)	51 (98%)	1 (2%)	69	91
All	All	421/443 (95%)	400 (95%)	21 (5%)	34	63

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	47	TRP
1	A	56	SER
1	A	70	SER
1	A	105	GLU
1	A	114	SER
1	A	123	GLU
1	A	158	ASN
1	A	208	SER
2	B	65	ARG
2	B	82	ASP
2	B	111	SER
2	B	118	SER
2	B	135	LYS
2	B	137	THR
2	B	138	SER
2	B	157	THR
2	B	162	SER
2	B	203	ASN
2	B	220	LYS

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Mol	Chain	Res	Type
3	C	33	ASP

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

Mol	Chain	Res	Type
2	B	198	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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	PCA	B	1	2	8,8,9	6.03	2 (25%)	8,10,12	6.69	5 (62%)

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	PCA	B	1	2	-	0/0/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	PCA	O-C	16.49	1.22	1.11
2	B	1	PCA	CD-N	3.65	1.42	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	PCA	CA-N-CD	-17.77	100.48	114.37
2	B	1	PCA	OE-CD-N	-4.22	120.03	125.11
2	B	1	PCA	CB-CG-CD	-2.90	99.62	104.30
2	B	1	PCA	C-CA-N	-2.54	110.14	110.71
2	B	1	PCA	CG-CD-N	2.29	115.99	107.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	211/213 (99%)	-0.03	4 (1%) 64 68	16, 40, 88, 119	0
2	B	220/231 (95%)	0.40	27 (12%) 5 4	18, 48, 116, 162	0
3	C	59/69 (85%)	-0.19	0 100 100	19, 33, 58, 83	0
All	All	490/513 (95%)	0.14	31 (6%) 19 20	16, 41, 96, 162	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	136	SER	7.2
2	B	213	VAL	6.1
2	B	135	LYS	5.5
2	B	217	VAL	4.8
2	B	190	VAL	4.7
2	B	197	THR	3.9
2	B	201	ILE	3.7
2	B	138	SER	3.7
2	B	137	THR	3.6
2	B	144	LEU	3.4
1	A	126	LYS	3.3
2	B	156	VAL	3.1
2	B	127	VAL	3.0
2	B	139	GLY	2.9
2	B	188	VAL	2.8
2	B	165	LEU	2.7
2	B	169	VAL	2.6
1	A	150	VAL	2.6
2	B	216	LYS	2.4
2	B	200	TYR	2.4
2	B	193	SER	2.4
2	B	195	LEU	2.4
2	B	204	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	164	ALA	2.3
1	A	133	VAL	2.3
2	B	148	VAL	2.3
2	B	160	TRP	2.2
1	A	183	LYS	2.2
2	B	175	VAL	2.1
2	B	158	VAL	2.1
2	B	189	THR	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	PCA	B	1	8/9	0.18	0.58	33,40,47,51	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.