wwPDB X-ray Structure Validation Summary Report (i)

Jun 15, 2024 – 11:45 AM EDT

PDB ID	:	2MCG
Title	:	THREE-DIMENSIONAL STRUCTURE OF A LIGHT CHAIN DIMER CRYSTALLIZED IN WATER. CONFORMATIONAL FLEXIBILITY OF A MOLECULE IN TWO CRYSTAL FORMS
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Deposited on	:	1989-05-09
Resolution	:	2.00 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

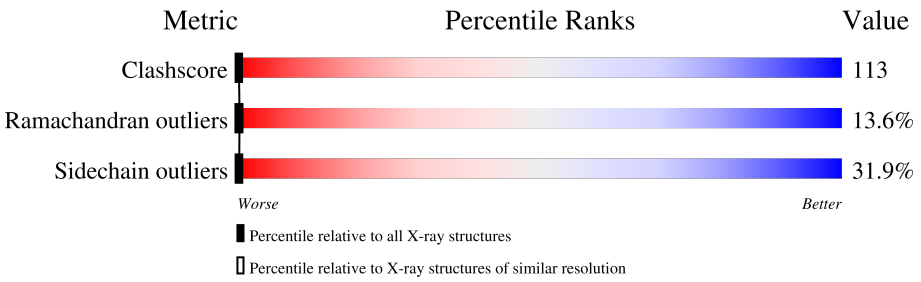
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 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))
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	1	216	17% 39% 34% 11%
1	2	216	14% 40% 31% 15%

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
1	PCA	1	1	-	-	X	-

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3530 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 IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	1	216	1606	1000	266	335	5	0	0	0
1	2	216	1606	1000	266	335	5	0	0	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	20	ILE	PHE	conflict	UNP P01709
1	23	THR	SER	conflict	UNP P01709
1	29	VAL	ILE	conflict	UNP P01709
1	31	GLY	ASN	conflict	UNP P01709
1	39	GLN	ARG	conflict	UNP P01709
1	42	ALA	PRO	conflict	UNP P01709
1	48	VAL	LEU	conflict	UNP P01709
1	49	ILE	MET	conflict	UNP P01709
1	54	ASN	THR	conflict	UNP P01709
1	62	ASP	ASN	conflict	UNP P01709
1	94	GLU	ALA	conflict	UNP P01709
1	97	ASP	ASN	conflict	UNP P01709
1	98	ASN	SER	conflict	UNP P01709
1	99	PHE	LEU	conflict	UNP P01709
1	100	VAL	ILE	conflict	UNP P01709
1	103	THR	GLY	conflict	UNP P01709
1	106	LYS	ARG	conflict	UNP P01709
1	107	VAL	LEU	conflict	UNP P01709
1	116	ASN	ALA	conflict	UNP P01709
1	118	THR	SER	conflict	UNP P01709
1	156	GLY	SER	conflict	UNP P01709
1	167	LYS	THR	conflict	UNP P01709
2	20	ILE	PHE	conflict	UNP P01709
2	23	THR	SER	conflict	UNP P01709

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Chain	Residue	Modelled	Actual	Comment	Reference
2	29	VAL	ILE	conflict	UNP P01709
2	31	GLY	ASN	conflict	UNP P01709
2	39	GLN	ARG	conflict	UNP P01709
2	42	ALA	PRO	conflict	UNP P01709
2	48	VAL	LEU	conflict	UNP P01709
2	49	ILE	MET	conflict	UNP P01709
2	54	ASN	THR	conflict	UNP P01709
2	62	ASP	ASN	conflict	UNP P01709
2	94	GLU	ALA	conflict	UNP P01709
2	97	ASP	ASN	conflict	UNP P01709
2	98	ASN	SER	conflict	UNP P01709
2	99	PHE	LEU	conflict	UNP P01709
2	100	VAL	ILE	conflict	UNP P01709
2	103	THR	GLY	conflict	UNP P01709
2	106	LYS	ARG	conflict	UNP P01709
2	107	VAL	LEU	conflict	UNP P01709
2	116	ASN	ALA	conflict	UNP P01709
2	118	THR	SER	conflict	UNP P01709
2	156	GLY	SER	conflict	UNP P01709
2	167	LYS	THR	conflict	UNP P01709

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	1	133	133	133	0	0
			Total	O		
2	2	185	185	185	0	0

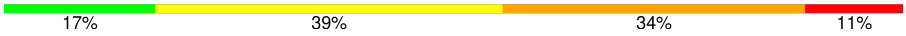
3 Residue-property plots (i)

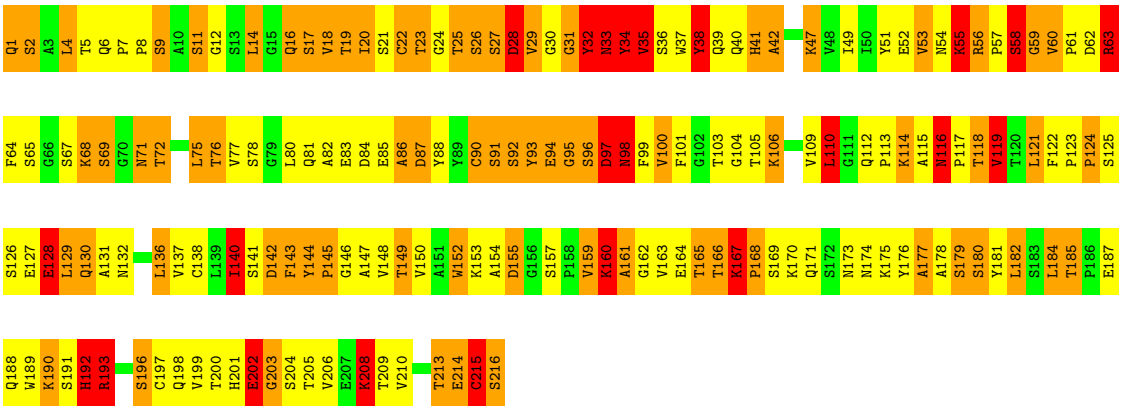
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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: IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN)

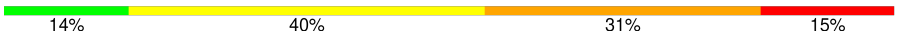
Chain 1:

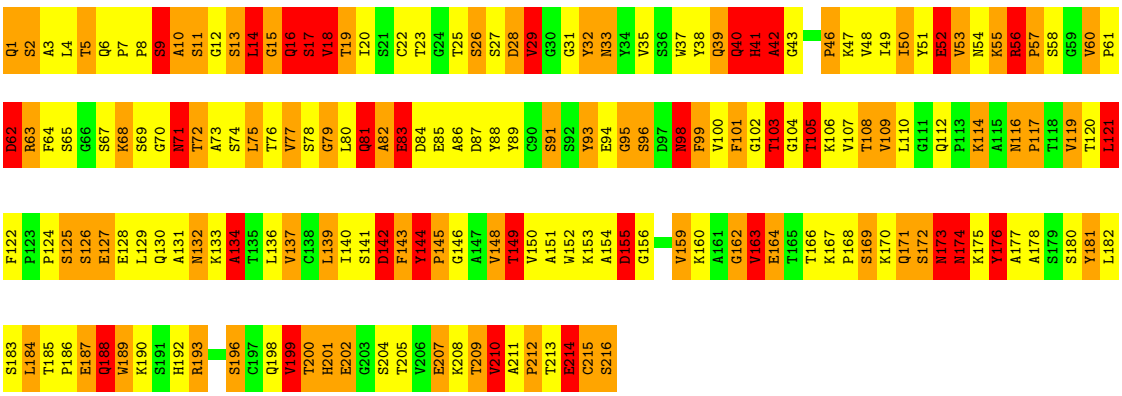




• Molecule 1: IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN)

Chain 2:





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	72.30Å 72.30Å 185.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	7.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (7.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.187 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3530	wwPDB-VP
Average B, all atoms (Å^2)	8.0	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: 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	1	1.42	7/1637 (0.4%)	2.51	82/2233 (3.7%)
1	2	1.40	1/1637 (0.1%)	2.44	94/2233 (4.2%)
All	All	1.41	8/3274 (0.2%)	2.48	176/4466 (3.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	1	0	2
1	2	0	1
All	All	0	3

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1	58	SER	C-O	6.03	1.34	1.23
1	1	4	LEU	CA-CB	-5.85	1.40	1.53
1	1	59	GLY	N-CA	5.62	1.54	1.46
1	1	196	SER	CB-OG	-5.54	1.35	1.42
1	1	157	SER	CB-OG	-5.25	1.35	1.42

The worst 5 of 176 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	1	63	ARG	CD-NE-CZ	23.28	156.19	123.60
1	1	4	LEU	CA-CB-CG	19.94	161.17	115.30
1	1	130	GLN	CB-CG-CD	17.04	155.90	111.60
1	1	55	LYS	CA-CB-CG	16.14	148.91	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	1	193	ARG	NE-CZ-NH2	-15.93	112.33	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	1	193	ARG	Sidechain
1	1	58	SER	Mainchain
1	2	163	VAL	Mainchain

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	1	1606	0	1538	368	4
1	2	1606	0	1536	356	1
2	1	133	0	0	35	2
2	2	185	0	0	41	1
All	All	3530	0	3074	712	5

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 113.

The worst 5 of 712 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)
1:1:177:ALA:HB3	2:1:347:HOH:O	1.32	1.26
1:1:19:THR:HG23	1:1:76:THR:CG2	1.66	1.26
2:1:314:HOH:O	1:2:141:SER:HB2	1.33	1.25
1:1:52:GLU:O	1:1:53:VAL:HG22	1.34	1.23
1:1:117:PRO:HB3	1:1:140:ILE:CD1	1.70	1.20

All (5) 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)
2:1:266:HOH:O	2:2:333:HOH:O[6_656]	1.54	0.66
1:1:81:GLN:NE2	2:1:242:HOH:O[4_646]	1.76	0.44
1:1:58:SER:O	1:1:215:CYS:CA[3_654]	2.11	0.09
1:1:32:TYR:OH	1:2:9:SER:CB[4_546]	2.14	0.06
1:1:59:GLY:O	1:1:216:SER:N[3_654]	2.19	0.01

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	1	214/216 (99%)	163 (76%)	28 (13%)	23 (11%)	0 0
1	2	214/216 (99%)	140 (65%)	39 (18%)	35 (16%)	0 0
All	All	428/432 (99%)	303 (71%)	67 (16%)	58 (14%)	0 0

5 of 58 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	2	SER
1	1	9	SER
1	1	14	LEU
1	1	26	SER
1	1	27	SER

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	1	180/180 (100%)	116 (64%)	64 (36%)	0	0
1	2	180/180 (100%)	129 (72%)	51 (28%)	0	0
All	All	360/360 (100%)	245 (68%)	115 (32%)	0	0

5 of 115 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	191	SER
1	2	200	THR
1	2	22	CYS
1	2	199	VAL
1	2	159	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	2	174	ASN
1	2	201	HIS
1	2	33	ASN
1	2	40	GLN
1	2	41	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are 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	$\#\|Z\| > 2$
1	PCA	2	1	1	7,8,9	3.31	3 (42%)	9,10,12	1.83	4 (44%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	PCA	1	1	1	7,8,9	3.79	2 (28%)	9,10,12	2.48	3 (33%)

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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1	1	PCA	CD-N	8.88	1.57	1.34
1	2	1	PCA	CD-N	7.47	1.53	1.34
1	2	1	PCA	CA-N	3.32	1.50	1.46
1	1	1	PCA	CA-N	3.25	1.50	1.46
1	2	1	PCA	CB-CG	2.55	1.58	1.53

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	1	1	PCA	OE-CD-N	4.19	134.03	124.96
1	1	1	PCA	OE-CD-CG	-3.80	119.94	126.72
1	1	1	PCA	CA-N-CD	-3.49	101.62	113.58
1	2	1	PCA	CB-CA-N	2.88	111.18	103.24
1	2	1	PCA	CA-N-CD	-2.47	105.11	113.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	2	1	PCA	1	0
1	1	1	PCA	8	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.