



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:46 PM UTC

PDB ID : 4YGB / pdb_00004ygb
Title : Crystal structure of ERGIC-53/MCFD2, monoclinic calcium-free form
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Deposited on : 2015-02-26
Resolution : 1.60 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

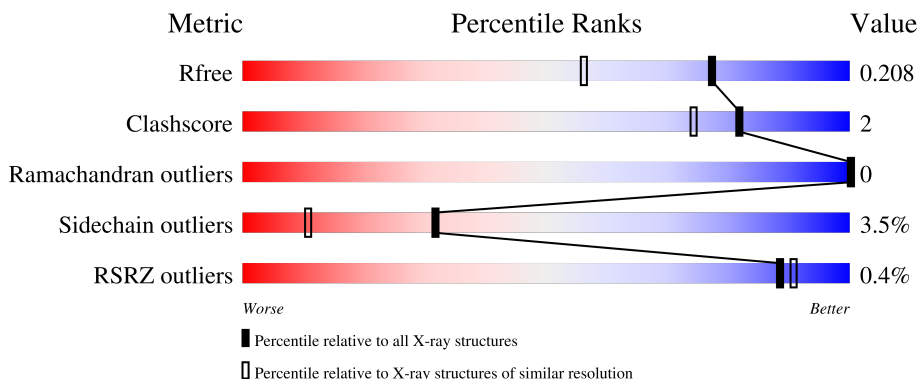
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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 $\leq 5\%$. 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	246	
1	C	246	
2	B	104	
2	D	104	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4333 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 Protein ERGIC-53.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	192	Total	C	N	O	S	0	1	0
			1529	982	268	275	4			
1	C	192	Total	C	N	O	S	0	1	0
			1529	982	268	275	4			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	expression tag	UNP P49257
A	25	ASN	-	expression tag	UNP P49257
A	26	HIS	-	expression tag	UNP P49257
A	27	LYS	-	expression tag	UNP P49257
A	28	VAL	-	expression tag	UNP P49257
A	29	HIS	-	expression tag	UNP P49257
A	30	MET	-	expression tag	UNP P49257
C	24	MET	-	expression tag	UNP P49257
C	25	ASN	-	expression tag	UNP P49257
C	26	HIS	-	expression tag	UNP P49257
C	27	LYS	-	expression tag	UNP P49257
C	28	VAL	-	expression tag	UNP P49257
C	29	HIS	-	expression tag	UNP P49257
C	30	MET	-	expression tag	UNP P49257

- Molecule 2 is a protein called Multiple coagulation factor deficiency protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	63	Total	C	N	O	S	0	0	0
			510	320	80	108	2			
2	D	63	Total	C	N	O	S	0	0	0
			510	320	80	108	2			

There are 48 discrepancies between the modelled and reference sequences:

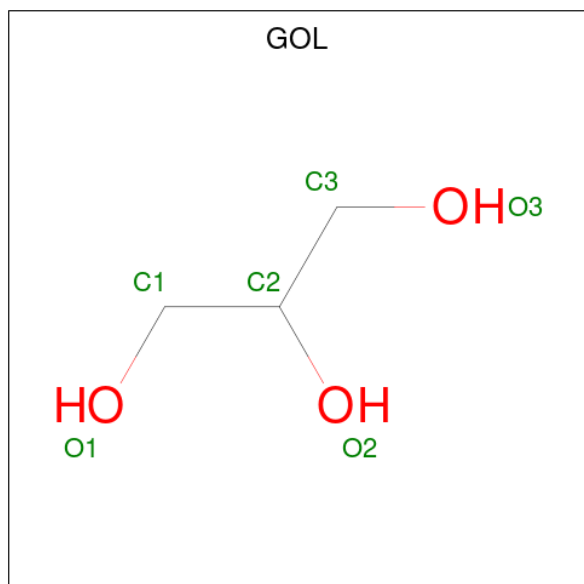
Chain	Residue	Modelled	Actual	Comment	Reference
B	43	MET	-	expression tag	UNP Q8NI22
B	44	GLY	-	expression tag	UNP Q8NI22
B	45	HIS	-	expression tag	UNP Q8NI22
B	46	HIS	-	expression tag	UNP Q8NI22
B	47	HIS	-	expression tag	UNP Q8NI22
B	48	HIS	-	expression tag	UNP Q8NI22
B	49	HIS	-	expression tag	UNP Q8NI22
B	50	HIS	-	expression tag	UNP Q8NI22
B	51	HIS	-	expression tag	UNP Q8NI22
B	52	HIS	-	expression tag	UNP Q8NI22
B	53	HIS	-	expression tag	UNP Q8NI22
B	54	HIS	-	expression tag	UNP Q8NI22
B	55	SER	-	expression tag	UNP Q8NI22
B	56	SER	-	expression tag	UNP Q8NI22
B	57	GLY	-	expression tag	UNP Q8NI22
B	58	HIS	-	expression tag	UNP Q8NI22
B	59	ILE	-	expression tag	UNP Q8NI22
B	60	GLU	-	expression tag	UNP Q8NI22
B	61	GLY	-	expression tag	UNP Q8NI22
B	62	ARG	-	expression tag	UNP Q8NI22
B	63	HIS	-	expression tag	UNP Q8NI22
B	64	MET	-	expression tag	UNP Q8NI22
B	65	LEU	-	expression tag	UNP Q8NI22
B	66	GLU	-	expression tag	UNP Q8NI22
D	43	MET	-	expression tag	UNP Q8NI22
D	44	GLY	-	expression tag	UNP Q8NI22
D	45	HIS	-	expression tag	UNP Q8NI22
D	46	HIS	-	expression tag	UNP Q8NI22
D	47	HIS	-	expression tag	UNP Q8NI22
D	48	HIS	-	expression tag	UNP Q8NI22
D	49	HIS	-	expression tag	UNP Q8NI22
D	50	HIS	-	expression tag	UNP Q8NI22
D	51	HIS	-	expression tag	UNP Q8NI22
D	52	HIS	-	expression tag	UNP Q8NI22
D	53	HIS	-	expression tag	UNP Q8NI22
D	54	HIS	-	expression tag	UNP Q8NI22
D	55	SER	-	expression tag	UNP Q8NI22
D	56	SER	-	expression tag	UNP Q8NI22
D	57	GLY	-	expression tag	UNP Q8NI22
D	58	HIS	-	expression tag	UNP Q8NI22
D	59	ILE	-	expression tag	UNP Q8NI22
D	60	GLU	-	expression tag	UNP Q8NI22
D	61	GLY	-	expression tag	UNP Q8NI22

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Chain	Residue	Modelled	Actual	Comment	Reference
D	62	ARG	-	expression tag	UNP Q8NI22
D	63	HIS	-	expression tag	UNP Q8NI22
D	64	MET	-	expression tag	UNP Q8NI22
D	65	LEU	-	expression tag	UNP Q8NI22
D	66	GLU	-	expression tag	UNP Q8NI22

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	2	Total Ca 2 2	0	0
4	D	2	Total Ca 2 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	95	Total 95	O 95	0	0
5	B	31	Total 31	O 31	0	0
5	C	81	Total 81	O 81	0	0
5	D	32	Total 32	O 32	0	0

- Molecule 1: Protein ERGIC-53

[illegible]

- Molecule 1: Protein ERGIC-53

ASN	M162	P163	M170	ASN	GLY	GLN	ILE	ILE	HIS	TYR	ASP	ASP	HIS	GLN	ASN	ASP	ASP	GLY	ALA	ALA	SER	Q185	L212	M218	K224	N225	M236	T268	GLU																			
MET	ASN	HIS	LYS	VAL	HIS	MET	ASP	GLY	VAL	GLY	ASP	PRO	ALA	VAL	ALA	LEU	P42	S60	W67	I74	V93	G116	L123	L129	ASN	GLN	GLY	LEU	GLU	GLY	PRO	VAL	PHE	GLY	SER	ALA	ASP	ALA	LEU	H144	V147	F151	F154	ASP	ASN	ASP	GLY	LYS

- Molecule 2: Multiple coagulation factor deficiency protein 2

MET	GLY	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	SER	SER	GLY	ILE	GLU	GLY	ARG	HIS	MET	MET	LEU	GLU	SER	GLU	GLN	ALA	PRO	L111	E116	A142	K143	SER	LEU	GLN
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- Molecule 2: Multiple coagulation factor deficiency protein 2

MET
 GLY
 HIS
 HIS
 HIS
 HIS
 HIS
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 HIS
 HIS
 HIS
 HIS
 SER
 SER
 SER
 GLY
 HIS
 ILE
 GLU
 GLY
 ARG
 HIS
 MET
 LEU
 GLU
 MET
 SER
 P69
 L72
 Q73
 T98
 HIS
 VAL
 HIS
 LYS
 GLU
 GLU
 GLY
 SER
 SER
 GLU
 GLN
 ALA
 ALA
 PRO
 L11
 E16
 K143
 SER
 LEU
 LUN

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.92Å 116.93Å 58.06Å 90.00° 120.13° 90.00°	Depositor
Resolution (Å)	20.00 – 1.60 20.00 – 1.60	Depositor EDS
% Data completeness (in resolution range)	98.8 (20.00-1.60) 98.7 (20.00-1.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.180 , 0.199 (Not available) , 0.208	Depositor DCC
R_{free} test set	4370 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.024 for -h-l,k,h 0.024 for l,k,-h-l 0.027 for l,-k,h 0.026 for h,-k,-h-l 0.468 for -h-l,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4333	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.90% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: GOL, CA

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	0.76	0/1574	0.80	0/2133
1	C	0.76	0/1574	0.81	0/2133
2	B	0.74	0/517	0.89	0/696
2	D	0.71	0/517	0.88	0/696
All	All	0.75	0/4182	0.83	0/5658

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	A	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	116	GLY	Peptide
1	C	116	GLY	Peptide

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	1529	0	1475	8	0
1	C	1529	0	1475	8	0
2	B	510	0	474	3	0
2	D	510	0	474	3	0
3	A	6	0	8	0	0
3	C	6	0	8	0	0
4	B	2	0	0	0	0
4	D	2	0	0	0	0
5	A	95	0	0	0	0
5	B	31	0	0	1	0
5	C	81	0	0	0	0
5	D	32	0	0	2	0
All	All	4333	0	3914	20	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 2.

The worst 5 of 20 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:C:123:LEU:HD23	1:C:151:PHE:HB2	1.69	0.72
1:A:60:SER:OG	2:B:111:LEU:HD21	1.92	0.70
1:C:60:SER:OG	2:D:111:LEU:HD21	1.94	0.67
1:C:162:ASN:HA	1:C:163:PRO:C	2.32	0.55
1:A:162:ASN:HA	1:A:163:PRO:C	2.32	0.54

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	185/246 (75%)	181 (98%)	4 (2%)	0	100	100
1	C	185/246 (75%)	181 (98%)	4 (2%)	0	100	100
2	B	59/104 (57%)	59 (100%)	0	0	100	100
2	D	59/104 (57%)	59 (100%)	0	0	100	100
All	All	488/700 (70%)	480 (98%)	8 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	159/199 (80%)	151 (95%)	8 (5%)	22	5
1	C	159/199 (80%)	155 (98%)	4 (2%)	42	18
2	B	56/92 (61%)	54 (96%)	2 (4%)	31	10
2	D	56/92 (61%)	55 (98%)	1 (2%)	51	28
All	All	430/582 (74%)	415 (96%)	15 (4%)	32	11

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	246	ILE
1	C	236	MET
2	B	70	GLN
2	D	72	LEU
1	C	123	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	225	ASN
1	C	235	ASN
2	D	75	HIS
1	A	225	ASN
1	A	235	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are 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	$\# Z > 2$
3	GOL	C	301	-	5,5,5	0.30	0	5,5,5	0.81	0
3	GOL	A	301	-	5,5,5	0.43	0	5,5,5	0.94	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
3	GOL	C	301	-	-	0/4/4/4	-
3	GOL	A	301	-	-	0/4/4/4	-

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(Å²)	Q<0.9	
1	A	192/246 (78%)	-0.61	0	100	100	15, 26, 47, 63	1 (0%)
1	C	192/246 (78%)	-0.51	0	100	100	15, 26, 47, 64	1 (0%)
2	B	63/104 (60%)	-0.47	0	100	100	20, 29, 50, 58	0
2	D	63/104 (60%)	-0.44	2 (3%)	50	53	20, 29, 49, 60	0
All	All	510/700 (72%)	-0.54	2 (0%)	88	91	15, 27, 49, 64	2 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	111	LEU	2.0
2	D	69	PRO	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	C	301	6/6	0.97	0.05	24,26,29,32	0
3	GOL	A	301	6/6	0.98	0.06	24,25,29,31	0
4	CA	B	501	1/1	0.99	0.03	20,20,20,20	0
4	CA	B	502	1/1	0.99	0.03	26,26,26,26	0
4	CA	D	502	1/1	0.99	0.03	26,26,26,26	0
4	CA	D	501	1/1	1.00	0.02	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.