



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:42 AM UTC

PDB ID : 4YGE / pdb_00004yge
Title : Crystal structure of ERGIC-53/MCFD2, trigonal calcium-bound form 2
Authors : Satoh, T.; Nishio, M.; Yagi-Utsumi, M.; Suzuki, K.; Anzai, T.; Mizushima, T.; Kamiya, Y.; Kato, K.
Deposited on : 2015-02-26
Resolution : 3.05 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

<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
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

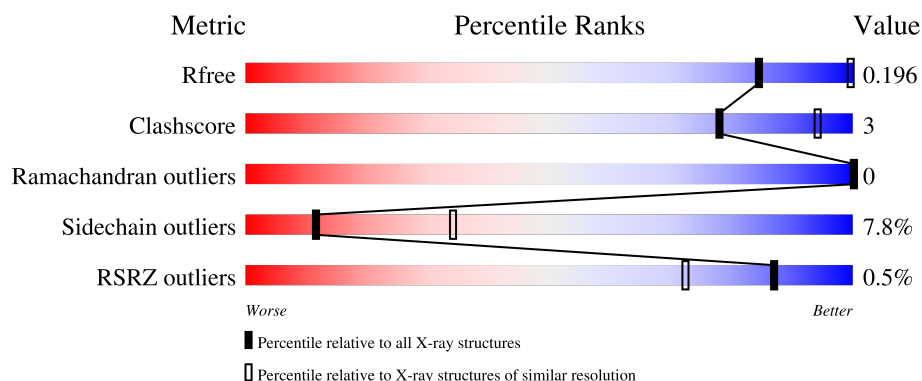
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	
1	E	246	
2	B	143	
2	D	143	

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Mol	Chain	Length	Quality of chain
2	F	143	39%7%53%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6911 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	227	Total	C	N	O	S	0	0	0
			1782	1131	314	333	4			
1	C	223	Total	C	N	O	S	0	0	0
			1753	1115	308	326	4			
1	E	223	Total	C	N	O	S	0	0	0
			1750	1115	308	323	4			

There are 21 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
E	24	MET	-	expression tag	UNP P49257
E	25	ASN	-	expression tag	UNP P49257
E	26	HIS	-	expression tag	UNP P49257
E	27	LYS	-	expression tag	UNP P49257
E	28	VAL	-	expression tag	UNP P49257
E	29	HIS	-	expression tag	UNP P49257
E	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	66	Total 532	C 331	N 85	O 113	S 3	0	0	0
2	D	66	Total 532	C 331	N 85	O 113	S 3	0	0	0
2	F	67	Total 540	C 337	N 86	O 114	S 3	0	0	0

There are 69 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	4	MET	-	expression tag	UNP Q8NI22
B	5	GLY	-	expression tag	UNP Q8NI22
B	6	HIS	-	expression tag	UNP Q8NI22
B	7	HIS	-	expression tag	UNP Q8NI22
B	8	HIS	-	expression tag	UNP Q8NI22
B	9	HIS	-	expression tag	UNP Q8NI22
B	10	HIS	-	expression tag	UNP Q8NI22
B	11	HIS	-	expression tag	UNP Q8NI22
B	12	HIS	-	expression tag	UNP Q8NI22
B	13	HIS	-	expression tag	UNP Q8NI22
B	14	HIS	-	expression tag	UNP Q8NI22
B	15	HIS	-	expression tag	UNP Q8NI22
B	16	SER	-	expression tag	UNP Q8NI22
B	17	SER	-	expression tag	UNP Q8NI22
B	18	GLY	-	expression tag	UNP Q8NI22
B	19	HIS	-	expression tag	UNP Q8NI22
B	20	ILE	-	expression tag	UNP Q8NI22
B	21	GLU	-	expression tag	UNP Q8NI22
B	22	GLY	-	expression tag	UNP Q8NI22
B	23	ARG	-	expression tag	UNP Q8NI22
B	24	HIS	-	expression tag	UNP Q8NI22
B	25	MET	-	expression tag	UNP Q8NI22
B	26	LEU	-	expression tag	UNP Q8NI22
D	4	MET	-	expression tag	UNP Q8NI22
D	5	GLY	-	expression tag	UNP Q8NI22
D	6	HIS	-	expression tag	UNP Q8NI22
D	7	HIS	-	expression tag	UNP Q8NI22
D	8	HIS	-	expression tag	UNP Q8NI22
D	9	HIS	-	expression tag	UNP Q8NI22
D	10	HIS	-	expression tag	UNP Q8NI22
D	11	HIS	-	expression tag	UNP Q8NI22
D	12	HIS	-	expression tag	UNP Q8NI22
D	13	HIS	-	expression tag	UNP Q8NI22

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Chain	Residue	Modelled	Actual	Comment	Reference
D	14	HIS	-	expression tag	UNP Q8NI22
D	15	HIS	-	expression tag	UNP Q8NI22
D	16	SER	-	expression tag	UNP Q8NI22
D	17	SER	-	expression tag	UNP Q8NI22
D	18	GLY	-	expression tag	UNP Q8NI22
D	19	HIS	-	expression tag	UNP Q8NI22
D	20	ILE	-	expression tag	UNP Q8NI22
D	21	GLU	-	expression tag	UNP Q8NI22
D	22	GLY	-	expression tag	UNP Q8NI22
D	23	ARG	-	expression tag	UNP Q8NI22
D	24	HIS	-	expression tag	UNP Q8NI22
D	25	MET	-	expression tag	UNP Q8NI22
D	26	LEU	-	expression tag	UNP Q8NI22
F	4	MET	-	expression tag	UNP Q8NI22
F	5	GLY	-	expression tag	UNP Q8NI22
F	6	HIS	-	expression tag	UNP Q8NI22
F	7	HIS	-	expression tag	UNP Q8NI22
F	8	HIS	-	expression tag	UNP Q8NI22
F	9	HIS	-	expression tag	UNP Q8NI22
F	10	HIS	-	expression tag	UNP Q8NI22
F	11	HIS	-	expression tag	UNP Q8NI22
F	12	HIS	-	expression tag	UNP Q8NI22
F	13	HIS	-	expression tag	UNP Q8NI22
F	14	HIS	-	expression tag	UNP Q8NI22
F	15	HIS	-	expression tag	UNP Q8NI22
F	16	SER	-	expression tag	UNP Q8NI22
F	17	SER	-	expression tag	UNP Q8NI22
F	18	GLY	-	expression tag	UNP Q8NI22
F	19	HIS	-	expression tag	UNP Q8NI22
F	20	ILE	-	expression tag	UNP Q8NI22
F	21	GLU	-	expression tag	UNP Q8NI22
F	22	GLY	-	expression tag	UNP Q8NI22
F	23	ARG	-	expression tag	UNP Q8NI22
F	24	HIS	-	expression tag	UNP Q8NI22
F	25	MET	-	expression tag	UNP Q8NI22
F	26	LEU	-	expression tag	UNP Q8NI22

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Ca 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total 2	Ca 2	0	0
3	C	1	Total 1	Ca 1	0	0
3	D	2	Total 2	Ca 2	0	0
3	F	2	Total 2	Ca 2	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Cl 1	0	0
4	C	1	Total 1	Cl 1	0	0
4	E	1	Total 1	Cl 1	0	0

- Molecule 5 is water.

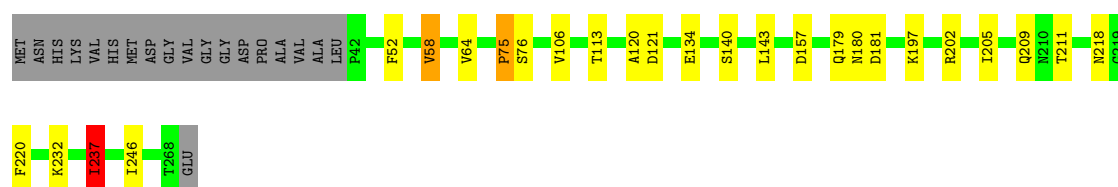
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	5	Total 5	O 5	0	0
5	C	2	Total 2	O 2	0	0
5	E	2	Total 2	O 2	0	0
5	F	1	Total 1	O 1	0	0

3 Residue-property plots

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 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.

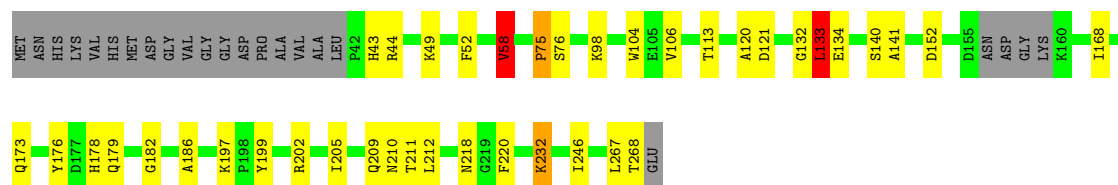
- Molecule 1: Protein ERGIC-53

Chain A: 



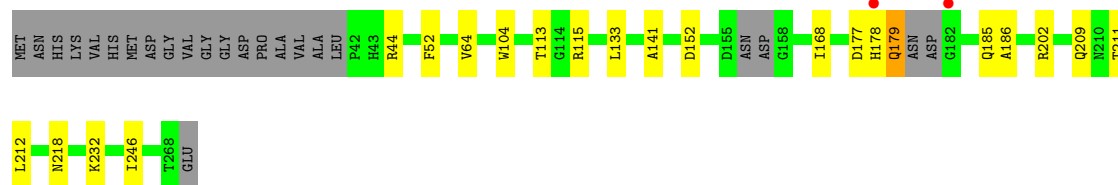
- Molecule 1: Protein ERGIC-53

Chain C: 



- Molecule 1: Protein ERGIC-53

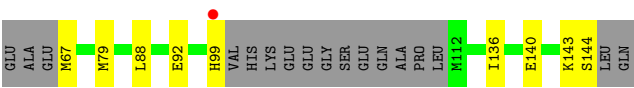
Chain E: 



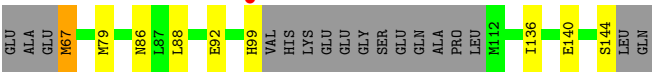
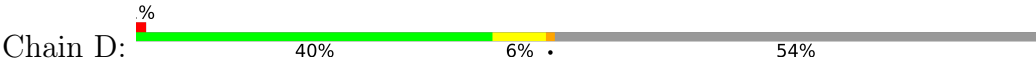
- Molecule 2: Multiple coagulation factor deficiency protein 2

Chain B: 

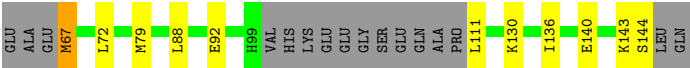
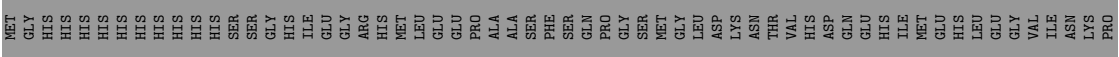
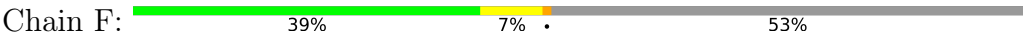




● Molecule 2: Multiple coagulation factor deficiency protein 2



● Molecule 2: Multiple coagulation factor deficiency protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	113.13Å 113.13Å 157.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.05 20.00 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-3.05) 99.5 (20.00-3.05)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.39 (at 3.04Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.173 , 0.211 (Not available) , 0.196	Depositor DCC
R_{free} test set	1156 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	39.2	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6911	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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: CA, CL

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	0.96	1/1832 (0.1%)	1.05	5/2487 (0.2%)
1	C	0.97	2/1802 (0.1%)	1.09	8/2446 (0.3%)
1	E	0.95	0/1798	1.03	5/2437 (0.2%)
2	B	0.77	0/540	0.93	0/727
2	D	0.76	0/540	0.97	2/727 (0.3%)
2	F	0.83	0/548	0.93	0/738
All	All	0.93	3/7060 (0.0%)	1.03	20/9562 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	75	PRO	C-O	-5.95	1.19	1.24
1	A	75	PRO	C-O	-5.55	1.19	1.24
1	C	199	TYR	N-CA	5.35	1.50	1.45

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	67	MET	CB-CG-SD	7.06	133.88	112.70
1	C	58	VAL	CB-CA-C	-7.01	100.60	111.26
1	A	58	VAL	CB-CA-C	-6.87	100.82	111.26
1	A	237	ILE	CB-CA-C	6.15	119.50	110.77
1	C	210	ASN	N-CA-C	6.05	120.25	112.86
1	C	141	ALA	CA-C-N	5.97	128.88	120.28
1	C	141	ALA	C-N-CA	5.97	128.88	120.28
2	D	86	ASN	N-CA-C	5.88	119.66	111.90
1	A	181	ASP	N-CA-C	-5.83	104.78	112.24
1	C	132	GLY	N-CA-C	5.72	124.82	115.73
1	E	104	TRP	N-CA-C	5.68	116.89	108.60
1	A	64	VAL	CA-C-N	-5.48	114.11	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64	VAL	C-N-CA	-5.48	114.11	119.76
1	E	141	ALA	CA-C-N	5.37	128.01	120.28
1	E	141	ALA	C-N-CA	5.37	128.01	120.28
1	E	64	VAL	CA-C-N	-5.25	114.36	119.76
1	E	64	VAL	C-N-CA	-5.25	114.36	119.76
1	C	133	LEU	CB-CA-C	-5.24	108.97	115.89
1	C	104	TRP	N-CA-C	5.21	116.20	108.60
1	C	182	GLY	N-CA-C	5.08	118.83	112.73

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	1782	0	1690	8	0
1	C	1753	0	1663	13	0
1	E	1750	0	1668	8	0
2	B	532	0	488	3	0
2	D	532	0	488	2	0
2	F	540	0	499	5	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	2	0	0	0	0
3	F	2	0	0	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
5	A	5	0	0	0	0
5	C	2	0	0	0	0
5	E	2	0	0	0	0
5	F	1	0	0	0	0
All	All	6911	0	6496	37	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 3.

All (37) 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)
2:F:67:MET:HE2	2:F:72:LEU:HD23	1.75	0.68
1:C:58:VAL:HG22	1:C:76:SER:O	1.98	0.63
1:A:58:VAL:HG22	1:A:76:SER:O	1.97	0.63
1:E:152:ASP:OD2	1:E:178:HIS:HE1	1.87	0.58
1:C:152:ASP:OD2	1:C:178:HIS:HE1	1.88	0.56
1:A:237:ILE:HG22	1:E:185:GLN:HB3	1.88	0.56
1:C:98:LYS:HE3	2:F:130:LYS:O	2.05	0.55
1:A:58:VAL:HG13	1:A:75:PRO:HB2	1.90	0.53
1:C:58:VAL:HG13	1:C:75:PRO:HB2	1.89	0.53
1:A:157:ASP:HB3	1:A:180:ASN:HA	1.91	0.52
1:E:177:ASP:OD2	1:E:179:GLN:HB3	2.10	0.51
1:E:179:GLN:O	1:E:179:GLN:HG2	2.11	0.50
1:C:152:ASP:OD2	1:C:178:HIS:CE1	2.68	0.47
1:E:152:ASP:OD2	1:E:178:HIS:CE1	2.69	0.45
2:F:88:LEU:HA	2:F:92:GLU:OE1	2.16	0.45
1:A:52:PHE:C	1:A:52:PHE:CD2	2.95	0.45
1:C:133:LEU:HB3	1:C:134:GLU:H	1.63	0.44
1:C:52:PHE:CD2	1:C:52:PHE:C	2.96	0.44
2:D:88:LEU:HA	2:D:92:GLU:OE1	2.19	0.43
1:A:106:VAL:CG2	1:A:205:ILE:HB	2.49	0.43
1:A:197:LYS:HE3	1:A:220:PHE:CZ	2.54	0.43
2:F:136:ILE:HA	2:F:140:GLU:OE1	2.19	0.43
2:B:88:LEU:HA	2:B:92:GLU:OE1	2.18	0.42
2:B:136:ILE:HA	2:B:140:GLU:OE1	2.19	0.42
2:D:136:ILE:HA	2:D:140:GLU:OE1	2.20	0.42
1:C:168:ILE:HD13	1:C:186:ALA:HA	2.01	0.42
1:C:197:LYS:HE3	1:C:220:PHE:CZ	2.55	0.42
1:E:168:ILE:HD13	1:E:186:ALA:HA	2.02	0.42
1:A:120:ALA:HB1	1:A:121:ASP:HA	2.02	0.42
1:E:52:PHE:CD2	1:E:52:PHE:C	2.98	0.41
2:B:136:ILE:HD13	2:B:136:ILE:HG21	1.87	0.41
1:C:120:ALA:HB1	1:C:121:ASP:HA	2.03	0.41
1:C:106:VAL:CG2	1:C:205:ILE:HB	2.51	0.41
2:F:136:ILE:HG21	2:F:136:ILE:HD13	1.87	0.41
1:E:212:LEU:O	1:E:232:LYS:HA	2.21	0.40
1:C:176:TYR:CE2	1:C:178:HIS:HA	2.57	0.40
1:C:212:LEU:O	1:C:232:LYS:HA	2.22	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	225/246 (92%)	216 (96%)	9 (4%)	0	100	100
1	C	219/246 (89%)	210 (96%)	9 (4%)	0	100	100
1	E	217/246 (88%)	206 (95%)	11 (5%)	0	100	100
2	B	62/143 (43%)	59 (95%)	3 (5%)	0	100	100
2	D	62/143 (43%)	59 (95%)	3 (5%)	0	100	100
2	F	63/143 (44%)	60 (95%)	3 (5%)	0	100	100
All	All	848/1167 (73%)	810 (96%)	38 (4%)	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	139/199 (70%)	127 (91%)	12 (9%)	10	31
1	C	182/199 (92%)	165 (91%)	17 (9%)	8	28
1	E	181/199 (91%)	171 (94%)	10 (6%)	19	46
2	B	59/125 (47%)	54 (92%)	5 (8%)	10	31
2	D	59/125 (47%)	55 (93%)	4 (7%)	14	39
2	F	60/125 (48%)	55 (92%)	5 (8%)	10	32
All	All	680/972 (70%)	627 (92%)	53 (8%)	11	34

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	THR
1	A	134	GLU
1	A	140	SER
1	A	143	LEU
1	A	179	GLN
1	A	202	ARG
1	A	209	GLN
1	A	211	THR
1	A	218	ASN
1	A	232	LYS
1	A	237	ILE
1	A	246	ILE
2	B	67	MET
2	B	79	MET
2	B	99	HIS
2	B	143	LYS
2	B	144	SER
1	C	43	HIS
1	C	44	ARG
1	C	49	LYS
1	C	58	VAL
1	C	113	THR
1	C	133	LEU
1	C	140	SER
1	C	173	GLN
1	C	179	GLN
1	C	202	ARG
1	C	209	GLN
1	C	211	THR
1	C	218	ASN
1	C	232	LYS
1	C	246	ILE
1	C	267	LEU
1	C	268	THR
2	D	67	MET
2	D	79	MET
2	D	99	HIS
2	D	144	SER
1	E	44	ARG
1	E	113	THR
1	E	115	ARG
1	E	133	LEU

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Mol	Chain	Res	Type
1	E	179	GLN
1	E	202	ARG
1	E	209	GLN
1	E	211	THR
1	E	218	ASN
1	E	246	ILE
2	F	67	MET
2	F	79	MET
2	F	111	LEU
2	F	143	LYS
2	F	144	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	130	ASN
1	A	178	HIS
1	A	191	GLN
1	A	218	ASN
1	A	225	ASN
1	A	257	HIS
2	B	119	ASN
1	C	178	HIS
1	C	191	GLN
1	C	209	GLN
1	C	218	ASN
1	C	225	ASN
1	C	257	HIS
2	D	119	ASN
1	E	173	GLN
1	E	178	HIS
1	E	185	GLN
1	E	191	GLN
1	E	225	ASN
1	E	235	ASN
1	E	257	HIS
2	F	119	ASN

5.3.3 RNA ⓘ

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 12 ligands modelled in this entry, 12 are monoatomic - leaving 0 for Mogul analysis.

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	227/246 (92%)	-0.58	0 100 100	17, 31, 56, 70	0
1	C	223/246 (90%)	-0.57	0 100 100	18, 32, 58, 85	0
1	E	223/246 (90%)	-0.55	2 (0%) 81 61	20, 32, 63, 97	0
2	B	66/143 (46%)	-0.18	1 (1%) 72 50	32, 53, 80, 100	0
2	D	66/143 (46%)	-0.13	1 (1%) 72 50	28, 50, 74, 102	0
2	F	67/143 (46%)	-0.43	0 100 100	23, 37, 60, 73	0
All	All	872/1167 (74%)	-0.49	4 (0%) 87 72	17, 34, 67, 102	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	99	HIS	2.3
2	B	99	HIS	2.2
1	E	178	HIS	2.1
1	E	182	GLY	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(Å ²)	Q<0.9
4	CL	C	502	1/1	0.90	0.15	51,51,51,51	0
4	CL	A	503	1/1	0.92	0.13	58,58,58,58	0
3	CA	C	501	1/1	0.93	0.07	56,56,56,56	0
4	CL	E	501	1/1	0.96	0.12	38,38,38,38	0
3	CA	D	502	1/1	0.97	0.04	45,45,45,45	0
3	CA	A	502	1/1	0.97	0.04	49,49,49,49	0
3	CA	B	501	1/1	0.98	0.03	59,59,59,59	0
3	CA	A	501	1/1	0.98	0.03	31,31,31,31	0
3	CA	F	502	1/1	0.99	0.02	32,32,32,32	0
3	CA	D	501	1/1	0.99	0.02	28,28,28,28	0
3	CA	B	502	1/1	0.99	0.02	38,38,38,38	0
3	CA	F	501	1/1	0.99	0.02	22,22,22,22	0

6.5 Other polymers ⓘ

There are no such residues in this entry.