



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:56 PM UTC

PDB ID : 7F05 / pdb_00007f05
Title : Crystal structure of human cardiac calsequestrin bound with calcium
Authors : Fan, X.X.; Liu, X.X.; Su, X.D.; Wang, S.Q.
Deposited on : 2021-06-03
Resolution : 2.30 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

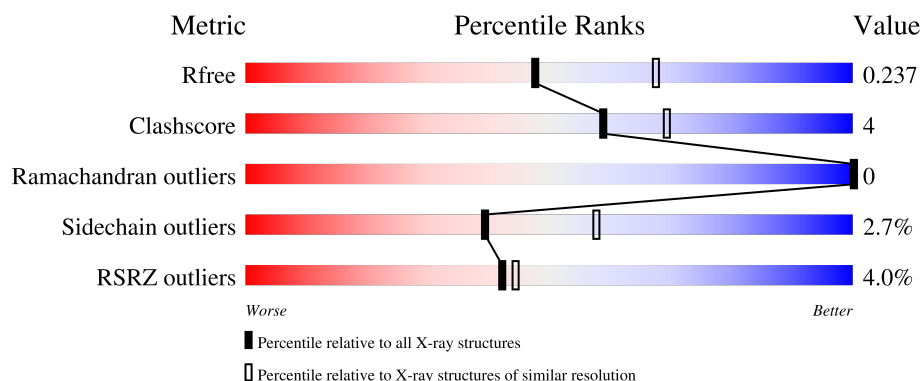
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	403	3% 78% 7% • 14%
1	B	403	4% 78% 8% 13%
1	C	403	2% 74% 13% • 12%
1	D	403	4% 75% 11% • 13%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12424 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 Calsequestrin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	1	0
			2850	1852	436	556	6			
1	C	355	Total	C	N	O	S	0	1	0
			2890	1872	440	572	6			
1	B	349	Total	C	N	O	S	0	0	0
			2825	1834	431	554	6			
1	D	349	Total	C	N	O	S	0	0	0
			2807	1820	425	556	6			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	MET	-	initiating methionine	UNP O14958
A	-2	GLY	-	expression tag	UNP O14958
A	-1	SER	-	expression tag	UNP O14958
A	0	SER	-	expression tag	UNP O14958
A	1	HIS	-	expression tag	UNP O14958
A	2	HIS	-	expression tag	UNP O14958
A	3	HIS	-	expression tag	UNP O14958
A	4	HIS	-	expression tag	UNP O14958
A	5	HIS	-	expression tag	UNP O14958
A	6	HIS	-	expression tag	UNP O14958
A	7	SER	-	expression tag	UNP O14958
A	8	SER	-	expression tag	UNP O14958
A	9	GLY	-	expression tag	UNP O14958
A	10	LEU	-	expression tag	UNP O14958
A	11	VAL	-	expression tag	UNP O14958
A	12	PRO	-	expression tag	UNP O14958
A	13	ARG	-	expression tag	UNP O14958
A	14	GLY	-	expression tag	UNP O14958
A	15	SER	-	expression tag	UNP O14958
A	16	HIS	-	expression tag	UNP O14958
A	17	MET	-	expression tag	UNP O14958

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Chain	Residue	Modelled	Actual	Comment	Reference
A	18	ALA	-	expression tag	UNP O14958
A	19	SER	-	expression tag	UNP O14958
A	385	GLU	SER	conflict	UNP O14958
A	393	GLU	SER	conflict	UNP O14958
C	-3	MET	-	initiating methionine	UNP O14958
C	-2	GLY	-	expression tag	UNP O14958
C	-1	SER	-	expression tag	UNP O14958
C	0	SER	-	expression tag	UNP O14958
C	1	HIS	-	expression tag	UNP O14958
C	2	HIS	-	expression tag	UNP O14958
C	3	HIS	-	expression tag	UNP O14958
C	4	HIS	-	expression tag	UNP O14958
C	5	HIS	-	expression tag	UNP O14958
C	6	HIS	-	expression tag	UNP O14958
C	7	SER	-	expression tag	UNP O14958
C	8	SER	-	expression tag	UNP O14958
C	9	GLY	-	expression tag	UNP O14958
C	10	LEU	-	expression tag	UNP O14958
C	11	VAL	-	expression tag	UNP O14958
C	12	PRO	-	expression tag	UNP O14958
C	13	ARG	-	expression tag	UNP O14958
C	14	GLY	-	expression tag	UNP O14958
C	15	SER	-	expression tag	UNP O14958
C	16	HIS	-	expression tag	UNP O14958
C	17	MET	-	expression tag	UNP O14958
C	18	ALA	-	expression tag	UNP O14958
C	19	SER	-	expression tag	UNP O14958
C	385	GLU	SER	conflict	UNP O14958
C	393	GLU	SER	conflict	UNP O14958
B	-3	MET	-	initiating methionine	UNP O14958
B	-2	GLY	-	expression tag	UNP O14958
B	-1	SER	-	expression tag	UNP O14958
B	0	SER	-	expression tag	UNP O14958
B	1	HIS	-	expression tag	UNP O14958
B	2	HIS	-	expression tag	UNP O14958
B	3	HIS	-	expression tag	UNP O14958
B	4	HIS	-	expression tag	UNP O14958
B	5	HIS	-	expression tag	UNP O14958
B	6	HIS	-	expression tag	UNP O14958
B	7	SER	-	expression tag	UNP O14958
B	8	SER	-	expression tag	UNP O14958
B	9	GLY	-	expression tag	UNP O14958

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Chain	Residue	Modelled	Actual	Comment	Reference
B	10	LEU	-	expression tag	UNP O14958
B	11	VAL	-	expression tag	UNP O14958
B	12	PRO	-	expression tag	UNP O14958
B	13	ARG	-	expression tag	UNP O14958
B	14	GLY	-	expression tag	UNP O14958
B	15	SER	-	expression tag	UNP O14958
B	16	HIS	-	expression tag	UNP O14958
B	17	MET	-	expression tag	UNP O14958
B	18	ALA	-	expression tag	UNP O14958
B	19	SER	-	expression tag	UNP O14958
B	385	GLU	SER	conflict	UNP O14958
B	393	GLU	SER	conflict	UNP O14958
D	-3	MET	-	initiating methionine	UNP O14958
D	-2	GLY	-	expression tag	UNP O14958
D	-1	SER	-	expression tag	UNP O14958
D	0	SER	-	expression tag	UNP O14958
D	1	HIS	-	expression tag	UNP O14958
D	2	HIS	-	expression tag	UNP O14958
D	3	HIS	-	expression tag	UNP O14958
D	4	HIS	-	expression tag	UNP O14958
D	5	HIS	-	expression tag	UNP O14958
D	6	HIS	-	expression tag	UNP O14958
D	7	SER	-	expression tag	UNP O14958
D	8	SER	-	expression tag	UNP O14958
D	9	GLY	-	expression tag	UNP O14958
D	10	LEU	-	expression tag	UNP O14958
D	11	VAL	-	expression tag	UNP O14958
D	12	PRO	-	expression tag	UNP O14958
D	13	ARG	-	expression tag	UNP O14958
D	14	GLY	-	expression tag	UNP O14958
D	15	SER	-	expression tag	UNP O14958
D	16	HIS	-	expression tag	UNP O14958
D	17	MET	-	expression tag	UNP O14958
D	18	ALA	-	expression tag	UNP O14958
D	19	SER	-	expression tag	UNP O14958
D	385	GLU	SER	conflict	UNP O14958
D	393	GLU	SER	conflict	UNP O14958

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	20	Total 20	Ca 20	0	0
2	C	22	Total 22	Ca 22	0	0
2	B	18	Total 18	Ca 18	0	0
2	D	14	Total 14	Ca 14	0	0

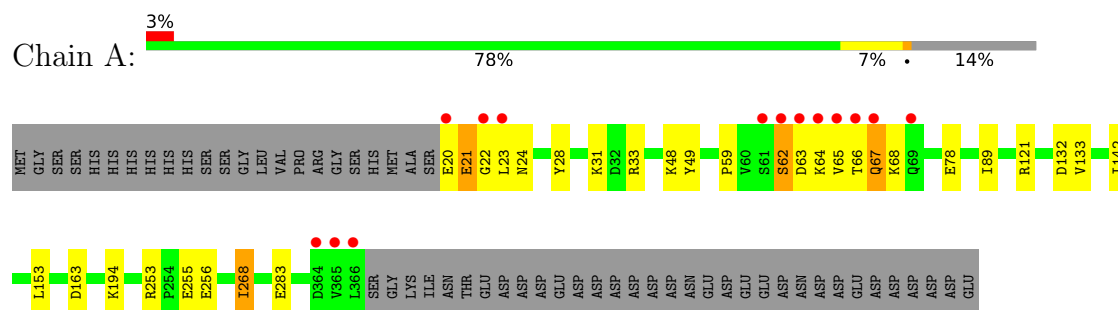
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	305	Total 305	O 305	0	0
3	C	243	Total 243	O 243	0	0
3	B	229	Total 229	O 229	0	0
3	D	201	Total 201	O 201	0	0

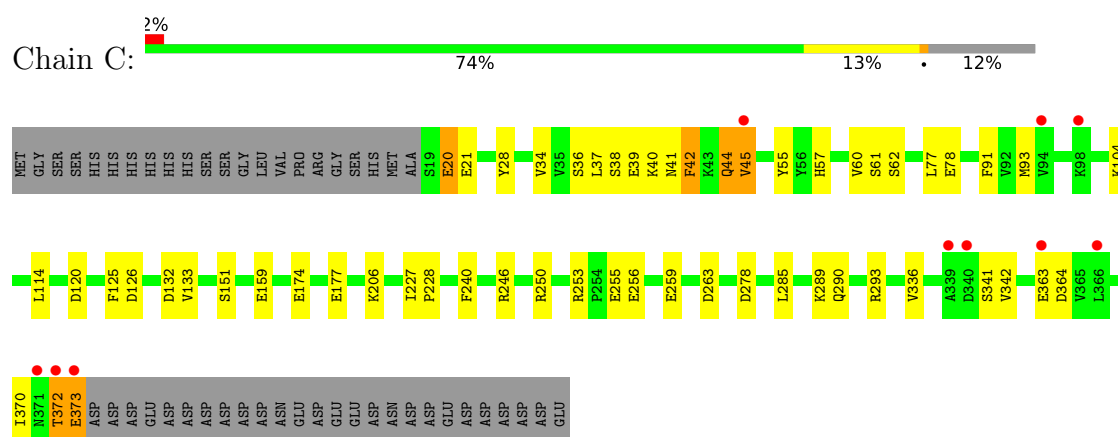
3 Residue-property plots [i](#)

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 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.

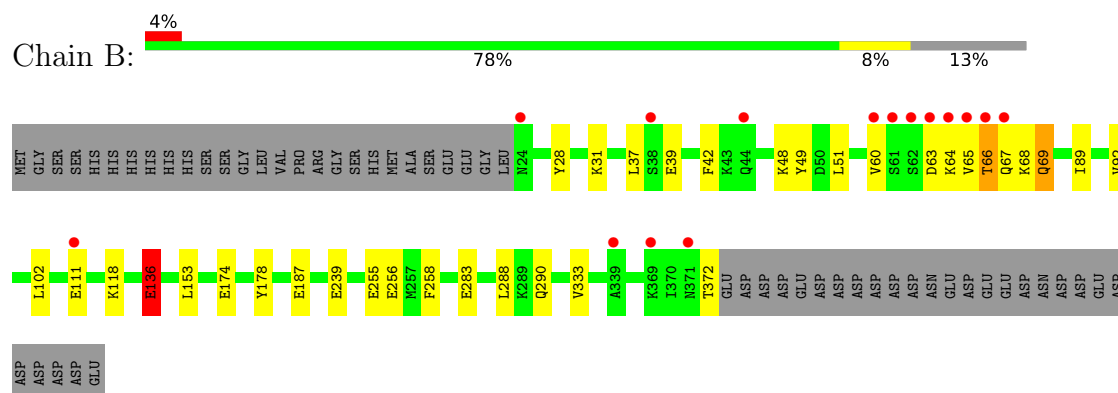
• Molecule 1: Calsequestrin-2



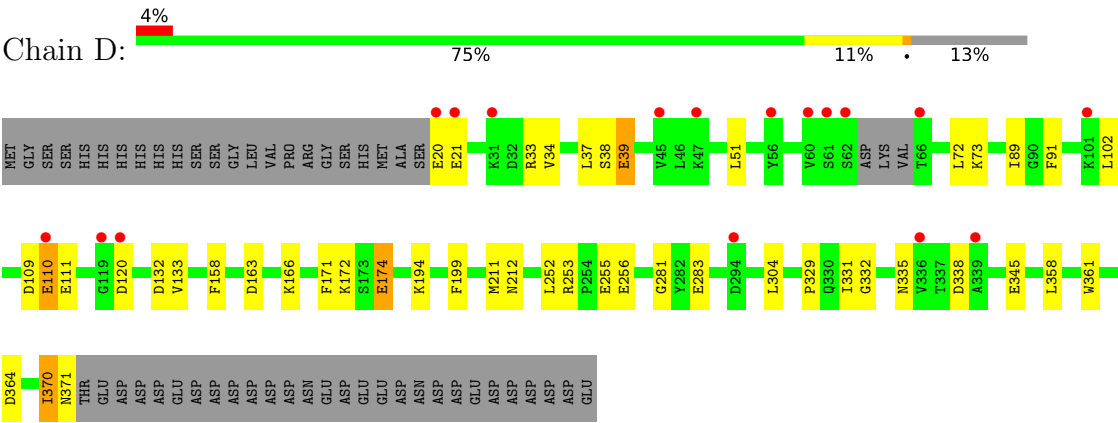
• Molecule 1: Calsequestrin-2



• Molecule 1: Calsequestrin-2



● Molecule 1: Calsequestrin-2



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	157.34Å 68.35Å 157.99Å 90.00° 91.20° 90.00°	Depositor
Resolution (Å)	19.74 – 2.30 19.74 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.74-2.30) 97.8 (19.74-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.24 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.208 , 0.234 0.213 , 0.237	Depositor DCC
R_{free} test set	3673 reflections (2.25%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.005 for l,k,-h 0.013 for l,-k,h 0.015 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12424	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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

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	0.45	4/2923 (0.1%)	0.85	5/3963 (0.1%)
1	B	0.46	2/2894 (0.1%)	0.74	1/3932 (0.0%)
1	C	0.59	4/2962 (0.1%)	0.84	6/4019 (0.1%)
1	D	0.50	0/2875	0.80	1/3911 (0.0%)
All	All	0.50	10/11654 (0.1%)	0.81	13/15825 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	290	GLN	C-O	-7.69	1.15	1.24
1	A	65	VAL	N-CA	-6.72	1.40	1.46
1	B	288	LEU	C-O	-6.19	1.16	1.24
1	A	121	ARG	C-O	-5.81	1.16	1.23
1	C	240	PHE	C-O	-5.78	1.17	1.24
1	C	36	SER	C-O	-5.50	1.17	1.24
1	B	136	GLU	C-O	-5.21	1.18	1.24
1	C	61	SER	CA-C	-5.12	1.47	1.53
1	A	59	PRO	C-O	-5.11	1.18	1.23
1	A	268	ILE	C-O	-5.05	1.18	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	VAL	N-CA-C	-15.63	97.68	113.47
1	C	44	GLN	N-CA-C	-12.96	99.85	114.62
1	A	62	SER	N-CA-C	-12.16	98.42	113.38
1	A	64	LYS	CA-C-N	-11.33	109.56	122.14
1	A	64	LYS	C-N-CA	-11.33	109.56	122.14
1	C	42	PHE	N-CA-C	-10.57	96.58	110.43
1	C	45	VAL	CA-C-N	8.34	131.46	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	45	VAL	C-N-CA	8.34	131.46	120.28
1	D	39	GLU	N-CA-C	6.08	120.43	112.89
1	C	372	THR	CB-CA-C	-5.78	109.37	117.23
1	A	21	GLU	N-CA-C	-5.69	105.44	112.90
1	B	372	THR	N-CA-C	5.50	126.39	111.00
1	C	341	SER	N-CA-C	5.20	116.98	109.07

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	2850	0	2734	19	0
1	B	2825	0	2683	24	1
1	C	2890	0	2751	28	1
1	D	2807	0	2622	36	0
2	A	20	0	0	0	0
2	B	18	0	0	1	0
2	C	22	0	0	0	0
2	D	14	0	0	0	0
3	A	305	0	0	1	0
3	B	229	0	0	1	0
3	C	243	0	0	3	0
3	D	201	0	0	6	0
All	All	12424	0	10790	98	1

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 4.

All (98) 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:D:120:ASP:OD2	3:D:501:HOH:O	1.65	1.14
1:A:194:LYS:HE3	3:A:756:HOH:O	1.46	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLU:N	1:B:69:GLN:HE22	1.67	0.92
1:B:64:LYS:C	1:B:66:THR:HA	1.97	0.88
1:B:65:VAL:N	1:B:66:THR:HA	1.90	0.87
1:A:20:GLU:HG3	1:A:21:GLU:H	1.38	0.87
1:D:110:GLU:OE1	3:D:502:HOH:O	1.94	0.86
1:D:20:GLU:O	1:D:21:GLU:HG2	1.77	0.84
1:D:172:LYS:N	1:D:211:MET:HE1	1.94	0.83
1:A:20:GLU:N	1:B:69:GLN:NE2	2.34	0.75
1:C:336:VAL:O	1:D:20:GLU:CB	2.37	0.73
1:C:77:LEU:HD12	1:C:93:MET:HE3	1.69	0.72
1:B:136:GLU:OE2	2:B:414:CA:CA	1.68	0.70
1:B:63:ASP:C	1:B:65:VAL:H	1.98	0.69
1:C:20:GLU:HG2	1:D:72:LEU:HD13	1.78	0.66
1:D:194:LYS:NZ	3:D:510:HOH:O	2.29	0.64
1:D:255:GLU:HG2	1:D:256:GLU:HG3	1.80	0.62
1:D:172:LYS:H	1:D:211:MET:HE1	1.64	0.60
1:D:335:ASN:OD1	1:D:338:ASP:N	2.34	0.60
1:B:28:TYR:OH	1:B:31:LYS:O	2.17	0.59
1:D:171:PHE:HA	1:D:211:MET:HE1	1.83	0.59
1:B:255:GLU:HG2	1:B:256:GLU:HG3	1.85	0.58
1:D:171:PHE:HA	1:D:211:MET:CE	2.33	0.58
1:D:33:ARG:NH1	3:D:516:HOH:O	2.36	0.57
1:C:336:VAL:HB	1:D:20:GLU:CB	2.34	0.56
1:C:253:ARG:NH1	3:C:512:HOH:O	2.38	0.56
1:D:211:MET:HE3	1:D:212:ASN:OD1	2.06	0.56
1:A:48:LYS:HD2	1:A:49:TYR:CZ	2.41	0.55
1:B:51:LEU:HB2	1:B:89:ILE:HD12	1.89	0.55
1:C:342:VAL:HG12	1:C:372:THR:HG22	1.89	0.55
1:A:255:GLU:HG2	1:A:256:GLU:HG3	1.89	0.54
1:C:255:GLU:HG2	1:C:256:GLU:HG3	1.89	0.53
1:D:171:PHE:C	1:D:211:MET:HE1	2.34	0.52
1:B:63:ASP:C	1:B:65:VAL:N	2.67	0.52
1:B:63:ASP:O	1:B:67:GLN:HG3	2.10	0.52
1:A:20:GLU:HG3	1:A:21:GLU:N	2.14	0.52
1:D:39:GLU:HG2	1:D:39:GLU:O	2.10	0.51
1:D:51:LEU:HB2	1:D:89:ILE:HD12	1.91	0.51
1:A:283:GLU:HG3	1:B:153:LEU:HB2	1.93	0.51
1:B:63:ASP:CB	1:B:65:VAL:H	2.24	0.50
1:D:20:GLU:C	1:D:21:GLU:HG2	2.37	0.50
1:C:38:SER:O	1:C:39:GLU:C	2.54	0.49
1:B:37:LEU:HD22	1:B:92:VAL:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:SER:O	1:C:40:LYS:N	2.46	0.48
1:B:48:LYS:HD2	1:B:49:TYR:CZ	2.49	0.48
1:B:65:VAL:N	1:B:66:THR:CA	2.72	0.48
1:C:41:ASN:C	1:C:42:PHE:O	2.52	0.48
1:D:158:PHE:O	1:D:166:LYS:NZ	2.47	0.47
1:A:23:LEU:O	1:A:24:ASN:HB2	2.13	0.47
1:C:293:ARG:O	3:C:502:HOH:O	2.20	0.47
1:C:372:THR:OG1	1:C:373:GLU:N	2.47	0.47
1:C:55:TYR:HD2	1:C:93:MET:HE2	1.79	0.47
1:D:370:ILE:HA	1:D:371:ASN:HA	1.45	0.47
1:D:171:PHE:CA	1:D:211:MET:HE1	2.45	0.47
1:C:278:ASP:OD2	3:C:501:HOH:O	2.20	0.47
1:B:187:GLU:OE1	3:B:501:HOH:O	2.21	0.47
1:C:126:ASP:HB3	1:C:259:GLU:HB2	1.98	0.46
1:C:364:ASP:HB3	1:C:370:ILE:HG12	1.98	0.46
1:C:159:GLU:OE2	1:C:206:LYS:NZ	2.48	0.46
1:C:20:GLU:H	1:C:20:GLU:HG3	1.40	0.46
1:D:332:GLY:HA2	1:D:361:TRP:CH2	2.52	0.45
1:D:73:LYS:NZ	3:D:514:HOH:O	2.43	0.45
1:C:132:ASP:OD1	1:C:133:VAL:N	2.50	0.45
1:A:67:GLN:HE21	1:A:67:GLN:HB2	1.58	0.45
1:D:252:LEU:HB2	1:D:304:LEU:HD11	1.99	0.45
1:B:37:LEU:HD21	1:B:42:PHE:HD1	1.82	0.44
1:B:39:GLU:HG3	1:B:102:LEU:HD13	1.97	0.44
1:D:364:ASP:HB3	1:D:370:ILE:HG13	1.99	0.44
1:B:51:LEU:HD21	1:B:118:LYS:HG3	1.99	0.44
1:C:28:TYR:CZ	1:C:78:GLU:HG2	2.52	0.44
1:C:34:VAL:HG22	1:C:91:PHE:HB2	1.99	0.44
1:A:20:GLU:CG	1:A:21:GLU:H	2.18	0.43
1:C:44:GLN:O	1:C:44:GLN:HG2	2.18	0.43
1:D:163:ASP:O	1:D:253:ARG:NH2	2.49	0.43
1:D:132:ASP:OD1	1:D:133:VAL:N	2.51	0.43
1:C:227:ILE:HA	1:C:228:PRO:HD3	1.83	0.43
1:D:39:GLU:HG3	1:D:102:LEU:HD13	1.99	0.43
1:D:331:ILE:HD13	1:D:358:LEU:HD22	2.01	0.43
1:D:281:GLY:HA2	1:D:329:PRO:HG3	2.01	0.43
1:C:114:LEU:HB3	1:C:125:PHE:HB3	2.01	0.43
1:C:285:LEU:HG	1:C:289:LYS:HE3	2.01	0.43
1:B:66:THR:C	1:B:68:LYS:N	2.73	0.43
1:D:174:GLU:CD	1:D:174:GLU:H	2.26	0.43
1:A:89:ILE:HD11	1:A:142:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:LEU:HB2	1:B:92:VAL:HB	2.02	0.42
1:D:34:VAL:HG22	1:D:91:PHE:HB2	2.02	0.42
1:D:174:GLU:HG3	1:D:199:PHE:CE2	2.55	0.42
1:A:132:ASP:OD1	1:A:133:VAL:N	2.52	0.42
1:A:163:ASP:O	1:A:253:ARG:NH2	2.50	0.42
1:A:153:LEU:HB2	1:B:283:GLU:HG3	2.02	0.42
1:D:194:LYS:NZ	3:D:535:HOH:O	2.52	0.41
1:A:31:LYS:HD3	1:A:33:ARG:CZ	2.51	0.41
1:C:151:SER:OG	1:D:283:GLU:OE2	2.32	0.41
1:C:246:ARG:HH21	1:C:263:ASP:CG	2.29	0.41
1:A:22:GLY:HA3	1:B:258:PHE:CZ	2.56	0.41
1:A:28:TYR:CE1	1:A:78:GLU:HB3	2.56	0.41
1:C:55:TYR:CE2	1:C:57:HIS:HB3	2.55	0.40
1:A:62:SER:O	1:A:63:ASP:C	2.63	0.40

All (1) 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)
1:C:177:GLU:OE1	1:B:178:TYR:OH[4_555]	2.08	0.12

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	346/403 (86%)	340 (98%)	6 (2%)	0	100	100
1	B	347/403 (86%)	338 (97%)	9 (3%)	0	100	100
1	C	354/403 (88%)	341 (96%)	13 (4%)	0	100	100
1	D	345/403 (86%)	337 (98%)	8 (2%)	0	100	100
All	All	1392/1612 (86%)	1356 (97%)	36 (3%)	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	308/366 (84%)	304 (99%)	4 (1%)	61	77
1	B	303/366 (83%)	294 (97%)	9 (3%)	36	53
1	C	312/366 (85%)	300 (96%)	12 (4%)	29	44
1	D	297/366 (81%)	289 (97%)	8 (3%)	39	58
All	All	1220/1464 (83%)	1187 (97%)	33 (3%)	39	58

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	THR
1	A	67	GLN
1	A	68	LYS
1	A	268	ILE
1	C	20	GLU
1	C	21	GLU
1	C	37	LEU
1	C	45	VAL
1	C	60	VAL
1	C	62	SER
1	C	104	LYS
1	C	120	ASP
1	C	174	GLU
1	C	250	ARG
1	C	363	GLU
1	C	373	GLU
1	B	60	VAL
1	B	66	THR
1	B	69	GLN
1	B	111	GLU
1	B	136	GLU
1	B	174	GLU

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Mol	Chain	Res	Type
1	B	239	GLU
1	B	290	GLN
1	B	333	VAL
1	D	37	LEU
1	D	38	SER
1	D	109	ASP
1	D	110	GLU
1	D	111	GLU
1	D	174	GLU
1	D	345	GLU
1	D	370	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	290	GLN
1	C	24	ASN
1	B	156	GLN
1	D	156	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 74 ligands modelled in this entry, 74 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

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	347/403 (86%)	0.10	14 (4%) 42 44	14, 27, 55, 124	1 (0%)
1	B	349/403 (86%)	0.36	15 (4%) 40 41	17, 33, 60, 106	0
1	C	355/403 (88%)	0.21	10 (2%) 55 57	14, 31, 65, 129	1 (0%)
1	D	349/403 (86%)	0.56	17 (4%) 35 36	16, 38, 76, 124	0
All	All	1400/1612 (86%)	0.31	56 (4%) 42 44	14, 32, 66, 129	2 (0%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	65	VAL	9.7
1	B	66	THR	5.5
1	A	67	GLN	5.4
1	C	339	ALA	5.3
1	A	65	VAL	4.9
1	A	64	LYS	4.6
1	D	66	THR	4.6
1	D	62	SER	4.6
1	A	62	SER	4.4
1	B	64	LYS	4.4
1	B	24	ASN	4.1
1	C	371	ASN	4.1
1	D	120	ASP	4.0
1	B	67	GLN	3.9
1	A	61	SER	3.8
1	D	20	GLU	3.8
1	A	63	ASP	3.7
1	B	111	GLU	3.7
1	A	66	THR	3.6
1	A	69	GLN	3.2
1	B	62	SER	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	22	GLY	3.2
1	B	38	SER	3.0
1	A	366	LEU	2.9
1	D	294	ASP	2.9
1	A	20	GLU	2.8
1	D	21	GLU	2.8
1	D	101	LYS	2.8
1	B	61	SER	2.8
1	D	110	GLU	2.7
1	D	339	ALA	2.6
1	B	44	GLN	2.6
1	D	56	TYR	2.6
1	D	60	VAL	2.6
1	D	61	SER	2.6
1	D	47	LYS	2.5
1	C	94	VAL	2.5
1	C	45	VAL	2.5
1	B	60	VAL	2.5
1	C	98	LYS	2.5
1	D	119	GLY	2.4
1	C	373	GLU	2.3
1	D	31	LYS	2.3
1	B	63	ASP	2.2
1	B	371	ASN	2.2
1	B	339	ALA	2.2
1	A	364	ASP	2.2
1	D	45	VAL	2.2
1	C	372	THR	2.2
1	A	365	VAL	2.1
1	A	23	LEU	2.1
1	C	366	LEU	2.1
1	C	340	ASP	2.1
1	C	363	GLU	2.1
1	B	369	LYS	2.0
1	D	336	VAL	2.0

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

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
2	CA	D	405	1/1	0.43	0.36	87,87,87,87	0
2	CA	B	413	1/1	0.51	0.20	114,114,114,114	0
2	CA	B	415	1/1	0.74	0.11	90,90,90,90	0
2	CA	C	419	1/1	0.74	0.13	84,84,84,84	0
2	CA	D	406	1/1	0.77	0.20	57,57,57,57	0
2	CA	B	417	1/1	0.78	0.11	86,86,86,86	0
2	CA	D	407	1/1	0.78	0.12	54,54,54,54	0
2	CA	C	422	1/1	0.81	0.09	80,80,80,80	0
2	CA	A	418	1/1	0.83	0.10	79,79,79,79	0
2	CA	A	420	1/1	0.83	0.09	90,90,90,90	0
2	CA	B	412	1/1	0.83	0.11	90,90,90,90	0
2	CA	B	414	1/1	0.84	0.14	71,71,71,71	0
2	CA	C	421	1/1	0.84	0.10	77,77,77,77	0
2	CA	C	420	1/1	0.84	0.12	81,81,81,81	0
2	CA	D	409	1/1	0.84	0.09	71,71,71,71	0
2	CA	C	416	1/1	0.85	0.10	58,58,58,58	0
2	CA	A	401	1/1	0.86	0.07	56,56,56,56	0
2	CA	C	418	1/1	0.86	0.12	66,66,66,66	0
2	CA	A	407	1/1	0.86	0.10	50,50,50,50	0
2	CA	B	404	1/1	0.87	0.09	47,47,47,47	0
2	CA	A	412	1/1	0.87	0.07	72,72,72,72	0
2	CA	D	414	1/1	0.87	0.11	69,69,69,69	0
2	CA	A	414	1/1	0.88	0.09	56,56,56,56	0
2	CA	C	410	1/1	0.88	0.11	65,65,65,65	0
2	CA	D	404	1/1	0.88	0.11	70,70,70,70	0
2	CA	D	413	1/1	0.88	0.09	59,59,59,59	0
2	CA	A	419	1/1	0.88	0.07	90,90,90,90	0
2	CA	B	401	1/1	0.89	0.07	44,44,44,44	0
2	CA	D	411	1/1	0.89	0.08	81,81,81,81	0
2	CA	A	417	1/1	0.89	0.07	60,60,60,60	0
2	CA	B	405	1/1	0.89	0.08	62,62,62,62	0
2	CA	D	412	1/1	0.91	0.13	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	B	402	1/1	0.91	0.06	38,38,38,38	0
2	CA	C	409	1/1	0.91	0.06	38,38,38,38	0
2	CA	C	403	1/1	0.92	0.07	46,46,46,46	0
2	CA	B	416	1/1	0.92	0.07	63,63,63,63	0
2	CA	A	403	1/1	0.92	0.06	30,30,30,30	0
2	CA	C	411	1/1	0.93	0.04	30,30,30,30	0
2	CA	D	408	1/1	0.94	0.09	70,70,70,70	0
2	CA	C	417	1/1	0.94	0.05	58,58,58,58	0
2	CA	A	404	1/1	0.95	0.04	23,23,23,23	0
2	CA	D	410	1/1	0.95	0.04	43,43,43,43	0
2	CA	A	410	1/1	0.95	0.08	45,45,45,45	0
2	CA	A	409	1/1	0.96	0.04	26,26,26,26	0
2	CA	C	407	1/1	0.96	0.06	37,37,37,37	0
2	CA	B	406	1/1	0.96	0.05	37,37,37,37	0
2	CA	C	412	1/1	0.96	0.03	25,25,25,25	0
2	CA	A	415	1/1	0.96	0.05	51,51,51,51	0
2	CA	A	406	1/1	0.97	0.05	20,20,20,20	0
2	CA	C	401	1/1	0.97	0.03	30,30,30,30	0
2	CA	B	407	1/1	0.97	0.04	26,26,26,26	0
2	CA	B	409	1/1	0.97	0.03	26,26,26,26	0
2	CA	B	411	1/1	0.97	0.03	43,43,43,43	0
2	CA	A	413	1/1	0.97	0.03	24,24,24,24	0
2	CA	C	405	1/1	0.97	0.05	26,26,26,26	0
2	CA	C	413	1/1	0.97	0.04	30,30,30,30	0
2	CA	C	414	1/1	0.97	0.04	33,33,33,33	0
2	CA	C	406	1/1	0.97	0.06	42,42,42,42	0
2	CA	A	416	1/1	0.97	0.05	45,45,45,45	0
2	CA	A	402	1/1	0.98	0.04	25,25,25,25	0
2	CA	B	408	1/1	0.98	0.03	24,24,24,24	0
2	CA	B	403	1/1	0.98	0.05	36,36,36,36	0
2	CA	B	410	1/1	0.98	0.03	24,24,24,24	0
2	CA	B	418	1/1	0.98	0.05	64,64,64,64	0
2	CA	C	408	1/1	0.98	0.04	26,26,26,26	0
2	CA	A	405	1/1	0.98	0.04	25,25,25,25	0
2	CA	C	402	1/1	0.98	0.05	34,34,34,34	0
2	CA	D	401	1/1	0.99	0.04	22,22,22,22	0
2	CA	D	402	1/1	0.99	0.02	22,22,22,22	0
2	CA	D	403	1/1	0.99	0.02	17,17,17,17	0
2	CA	C	404	1/1	0.99	0.02	25,25,25,25	0
2	CA	A	408	1/1	0.99	0.03	23,23,23,23	0
2	CA	C	415	1/1	0.99	0.03	30,30,30,30	0
2	CA	A	411	1/1	0.99	0.06	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.