



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2017 – 05:41 AM EDT

PDB ID : 2BE6  
Title : 2.0 Å crystal structure of the CaV1.2 IQ domain-Ca/CaM complex  
Authors : Van Petegem, F.; Chatelain, F.C.; Minor Jr., D.L.  
Deposited on : unknown  
Resolution : 2.00 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

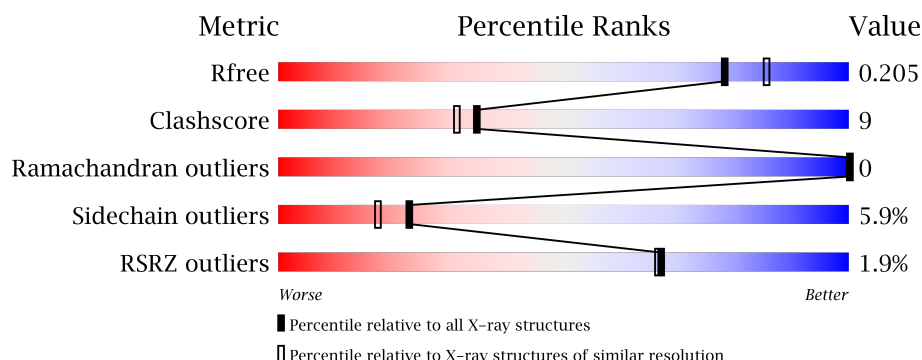
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	150	<div> <div>0.1%</div> <div>76%</div> <div>17%</div> <div>6%</div> </div>
1	B	150	<div> <div>3%</div> <div>74%</div> <div>18%</div> <div>7%</div> </div>
1	C	150	<div> <div>2%</div> <div>75%</div> <div>13%</div> <div>9%</div> </div>
2	D	37	<div> <div>3%</div> <div>62%</div> <div>16%</div> <div>22%</div> </div>
2	E	37	<div> <div>35%</div> <div>62%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	37	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: green (35%), yellow (19%), orange (5%), and grey (41%). The percentages are labeled below the bar.

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4120 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 Calmodulin 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	4	0
			1124	692	183	239	10			
1	B	140	Total	C	N	O	S	0	5	0
			1098	679	176	233	10			
1	C	137	Total	C	N	O	S	0	5	0
			1092	675	175	234	8			

- Molecule 2 is a protein called Voltage-dependent L-type calcium channel alpha-1C subunit.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	29	Total	C	N	O	0	0	0
			236	158	40	38			
2	E	14	Total	C	N	O	0	0	0
			106	74	15	17			
2	F	22	Total	C	N	O	0	0	0
			195	133	34	28			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1608	GLY	-	CLONING ARTIFACTS	UNP Q13933
D	1609	HIS	-	CLONING ARTIFACTS	UNP Q13933
D	1610	MET	-	CLONING ARTIFACTS	UNP Q13933
E	1608	GLY	-	CLONING ARTIFACTS	UNP Q13933
E	1609	HIS	-	CLONING ARTIFACTS	UNP Q13933
E	1610	MET	-	CLONING ARTIFACTS	UNP Q13933
F	1608	GLY	-	CLONING ARTIFACTS	UNP Q13933
F	1609	HIS	-	CLONING ARTIFACTS	UNP Q13933
F	1610	MET	-	CLONING ARTIFACTS	UNP Q13933

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	4	Total Ca 4 4	0	0
3	A	4	Total Ca 4 4	0	0
3	C	4	Total Ca 4 4	0	0

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ni 1 1	0	0
4	A	1	Total Ni 1 1	0	0
4	C	1	Total Ni 1 1	0	0

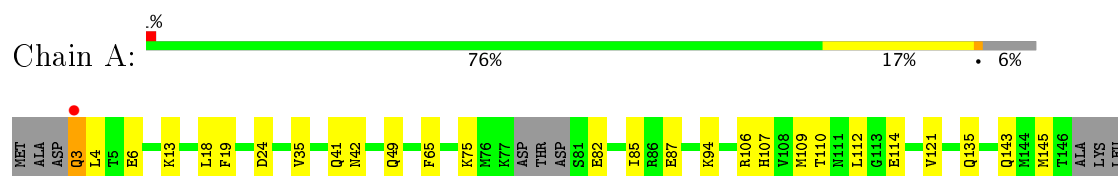
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	92	Total O 92 92	0	0
5	D	7	Total O 7 7	0	0
5	B	41	Total O 41 41	0	0
5	C	111	Total O 111 111	0	0
5	F	3	Total O 3 3	0	0

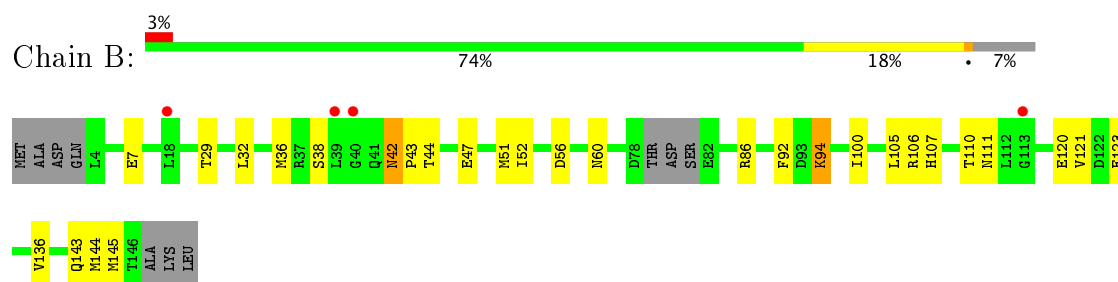
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

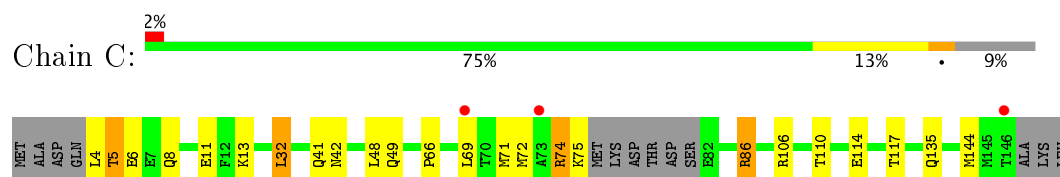
- Molecule 1: Calmodulin 2



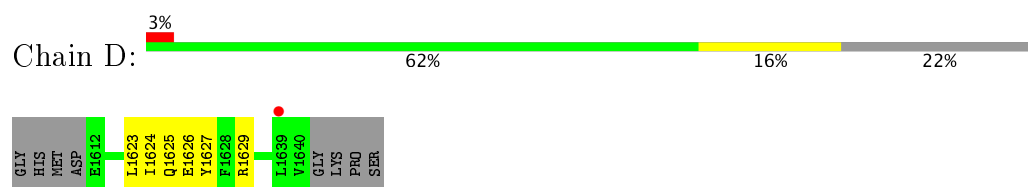
- Molecule 1: Calmodulin 2



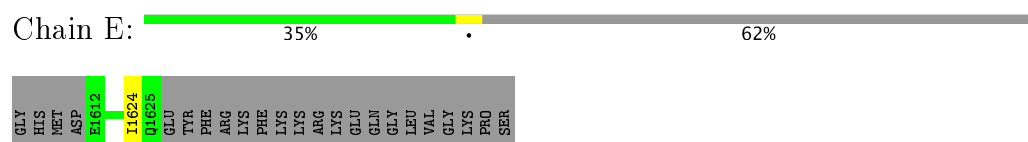
- Molecule 1: Calmodulin 2



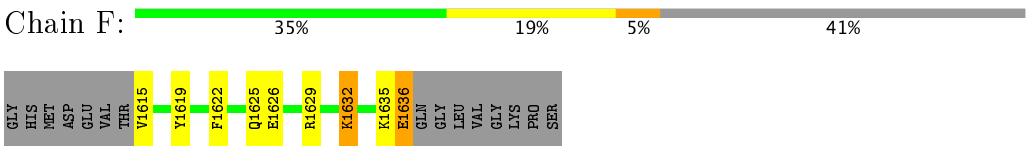
- Molecule 2: Voltage-dependent L-type calcium channel alpha-1C subunit



- Molecule 2: Voltage-dependent L-type calcium channel alpha-1C subunit



● Molecule 2: Voltage-dependent L-type calcium channel alpha-1C subunit



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.73Å 37.24Å 86.86Å 90.00° 97.77° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 40.59 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.4 (20.00-2.00) 96.9 (40.59-1.95)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.41 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.202 , 0.255 0.206 , 0.205	Depositor DCC
$R_{free}$ test set	1799 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.2	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 57.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.030 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4120	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.59% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: NI, 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	0.69	0/1144	0.75	0/1536
1	B	0.69	2/1126 (0.2%)	0.63	0/1514
1	C	0.68	0/1128	0.74	1/1517 (0.1%)
2	D	0.65	0/241	0.54	0/321
2	E	0.49	0/108	0.46	0/146
2	F	2.92	1/200 (0.5%)	0.68	0/262
All	All	0.93	3/3947 (0.1%)	0.69	1/5296 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1636	GLU	C-O	40.28	1.99	1.23
1	B	42	ASN	CG-ND2	15.11	1.70	1.32
1	B	42	ASN	CG-OD1	6.50	1.38	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	32	LEU	CB-CG-CD2	5.41	120.20	111.00

There are no chirality outliers.

There are no planarity outliers.

### 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	A	1124	0	1038	15	0
1	B	1098	0	1004	18	0
1	C	1092	0	1014	29	0
2	D	236	0	225	5	0
2	E	106	0	105	1	0
2	F	195	0	199	10	0
3	A	4	0	0	0	0
3	B	4	0	0	1	0
3	C	4	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	92	0	0	0	0
5	B	41	0	0	4	0
5	C	111	0	0	7	0
5	D	7	0	0	1	0
5	F	3	0	0	1	0
All	All	4120	0	3585	70	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 9.

All (70) 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:B:42:ASN:CG	1:B:42:ASN:ND2	1.70	1.44
1:C:71:MET:CE	1:C:72:MET:HE2	1.61	1.27
1:C:71:MET:HE2	1:C:72:MET:CE	1.83	1.09
1:C:71:MET:CE	1:C:72:MET:CE	2.33	1.06
1:C:135:GLN:OE1	5:C:581:HOH:O	1.71	1.04
1:B:106:ARG:O	1:B:110:THR:HG23	1.58	1.04
3:B:507:CA:CA	5:B:542:HOH:O	1.34	1.01
2:F:1636:GLU:C	2:F:1636:GLU:O	1.99	1.00
1:B:136:VAL:HG11	1:B:144:MET:CE	1.99	0.93
1:C:71:MET:HE2	1:C:72:MET:HE2	0.93	0.93
1:C:11:GLU:OE1	2:F:1629:ARG:NH2	2.07	0.87
1:A:94:LYS:HE2	1:A:107[B]:HIS:ND1	1.90	0.85
1:C:71:MET:HE1	1:C:72:MET:CE	2.05	0.85
1:A:94:LYS:HE2	1:A:107[B]:HIS:HD1	1.43	0.84
1:B:136:VAL:HG11	1:B:144:MET:HE3	1.57	0.83
1:A:112:LEU:HD11	2:D:1624:ILE:HG22	1.62	0.81
1:C:5:THR:HG22	1:C:8:GLN:H	1.45	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49[C]:GLN:OE1	5:C:528:HOH:O	2.01	0.78
1:C:49[B]:GLN:OE1	5:C:579:HOH:O	2.04	0.74
1:C:71:MET:HE1	1:C:72:MET:HE1	1.70	0.73
5:B:517:HOH:O	1:C:5:THR:HG21	1.87	0.73
1:B:105:LEU:HD23	1:B:121:VAL:HG13	1.70	0.72
1:C:106:ARG:O	1:C:110[B]:THR:HG23	1.91	0.70
1:B:60:ASN:OD1	5:B:542:HOH:O	2.09	0.69
1:C:48:LEU:C	1:C:49[C]:GLN:CA	2.62	0.68
2:F:1626:GLU:OE2	2:F:1629:ARG:NH1	2.26	0.68
1:C:144:MET:HE3	5:C:563:HOH:O	1.93	0.68
1:C:86:ARG:HG2	1:C:86:ARG:HH11	1.58	0.68
1:C:41:GLN:HE22	2:F:1615:VAL:CB	2.12	0.63
1:C:106:ARG:O	1:C:110[A]:THR:HG22	2.00	0.61
1:A:85:ILE:HG12	2:D:1623:LEU:HD13	1.84	0.60
1:C:114:GLU:OE2	1:C:114:GLU:HA	2.02	0.60
1:C:74:ARG:HB3	1:C:74:ARG:HH21	1.65	0.59
1:B:42:ASN:ND2	1:B:42:ASN:CB	2.64	0.57
1:C:117[B]:THR:HG21	5:C:571:HOH:O	2.06	0.55
1:B:43:PRO:HA	1:B:47:GLU:OE2	2.07	0.55
1:B:36[A]:MET:SD	1:B:51:MET:SD	3.05	0.55
1:C:106:ARG:HD3	5:C:623:HOH:O	2.07	0.55
1:A:94:LYS:HE2	1:A:107[B]:HIS:CE1	2.44	0.53
2:F:1625:GLN:NE2	5:F:31:HOH:O	2.41	0.52
1:B:92:PHE:HE1	2:E:1624:ILE:HG12	1.76	0.50
1:B:94:LYS:O	1:C:5:THR:HG23	2.11	0.50
1:C:4:LEU:HD22	1:C:69[B]:LEU:CD1	2.42	0.49
1:B:29:THR:HG22	1:B:52:ILE:HG13	1.93	0.49
1:A:94:LYS:CE	1:A:107[B]:HIS:HD1	2.19	0.48
1:A:145:MET:SD	2:D:1627:TYR:HD2	2.37	0.48
1:A:109:MET:HG3	1:A:114:GLU:HB2	1.96	0.48
1:A:109:MET:HE1	1:A:121:VAL:HG22	1.96	0.47
2:F:1622:PHE:CZ	2:F:1626:GLU:HG3	2.50	0.47
1:A:19:PHE:HA	1:A:35:VAL:HG21	1.96	0.47
2:D:1625:GLN:NE2	5:D:260:HOH:O	2.48	0.46
2:F:1622:PHE:CE1	2:F:1626:GLU:HG3	2.50	0.46
2:F:1626:GLU:OE1	2:F:1629:ARG:HD3	2.16	0.46
1:A:82:GLU:O	1:A:82:GLU:HG3	2.16	0.46
1:C:86:ARG:CG	1:C:86:ARG:HH11	2.27	0.45
1:B:44:THR:HG23	1:C:66:PRO:HG2	1.99	0.45
1:B:120:GLU:HA	1:B:123:GLU:HG2	1.99	0.45
2:D:1626:GLU:OE2	2:D:1629:ARG:NH2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107[B]:HIS:HE1	5:B:543:HOH:O	2.00	0.45
2:F:1632:LYS:O	2:F:1635:LYS:HB2	2.17	0.44
1:B:38:SER:O	1:B:111:ASN:HB3	2.18	0.44
1:C:42:ASN:ND2	5:C:538:HOH:O	2.50	0.43
1:A:3:GLN:CG	1:A:4:LEU:H	2.31	0.43
1:C:71:MET:HE3	2:F:1619:TYR:HD1	1.83	0.43
1:A:6:GLU:CD	1:A:6:GLU:H	2.22	0.43
1:A:13:LYS:HA	1:A:65:PHE:CE1	2.55	0.41
1:B:92:PHE:HB3	1:B:100:ILE:HD13	2.02	0.41
1:A:106:ARG:O	1:A:110:THR:HG23	2.21	0.41
1:C:74:ARG:HG3	1:C:75:LYS:N	2.36	0.40

There are no symmetry-related clashes.

## 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	A	141/150 (94%)	140 (99%)	1 (1%)	0	100	100
1	B	141/150 (94%)	140 (99%)	1 (1%)	0	100	100
1	C	139/150 (93%)	136 (98%)	3 (2%)	0	100	100
2	D	27/37 (73%)	27 (100%)	0	0	100	100
2	E	12/37 (32%)	12 (100%)	0	0	100	100
2	F	20/37 (54%)	20 (100%)	0	0	100	100
All	All	480/561 (86%)	475 (99%)	5 (1%)	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	120/128 (94%)	109 (91%)	11 (9%)	11	6
1	B	116/128 (91%)	110 (95%)	6 (5%)	27	22
1	C	119/128 (93%)	113 (95%)	6 (5%)	28	23
2	D	21/32 (66%)	21 (100%)	0	100	100
2	E	10/32 (31%)	10 (100%)	0	100	100
2	F	18/32 (56%)	17 (94%)	1 (6%)	25	19
All	All	404/480 (84%)	380 (94%)	24 (6%)	23	17

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	18	LEU
1	A	24	ASP
1	A	41	GLN
1	A	42	ASN
1	A	49	GLN
1	A	75	LYS
1	A	87	GLU
1	A	135[A]	GLN
1	A	135[B]	GLN
1	A	143	GLN
1	B	32	LEU
1	B	56	ASP
1	B	86	ARG
1	B	94	LYS
1	B	143	GLN
1	B	145	MET
1	C	5	THR
1	C	6	GLU
1	C	13	LYS
1	C	32	LEU
1	C	74	ARG

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Mol	Chain	Res	Type
1	C	86	ARG
2	F	1632	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	49	GLN
2	D	1625	GLN
1	B	42	ASN
1	B	49	GLN
1	B	111	ASN
1	B	135	GLN
1	C	41	GLN
1	C	42	ASN
1	C	111	ASN
1	C	135	GLN
2	F	1625	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 15 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	141/150 (94%)	0.12	1 (0%) 87 87	22, 30, 41, 49	0
1	B	140/150 (93%)	0.23	4 (2%) 52 52	24, 34, 45, 47	0
1	C	137/150 (91%)	-0.11	3 (2%) 62 61	21, 27, 39, 53	0
2	D	29/37 (78%)	0.30	1 (3%) 46 46	29, 33, 49, 53	0
2	E	14/37 (37%)	0.16	0 100 100	23, 26, 29, 30	0
2	F	22/37 (59%)	0.02	0 100 100	21, 28, 34, 36	0
All	All	483/561 (86%)	0.09	9 (1%) 67 66	21, 30, 44, 53	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	40	GLY	5.8
1	C	73	ALA	5.2
2	D	1639	LEU	4.2
1	B	18	LEU	2.9
1	C	146	THR	2.8
1	C	69[A]	LEU	2.6
1	A	3	GLN	2.4
1	B	113	GLY	2.4
1	B	39	LEU	2.2

### 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 ⓘ

There are no carbohydrates 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. 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, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	CA	C	514	1/1	0.96	0.04	-3.03	34,34,34,34	0
3	CA	C	511	1/1	0.97	0.05	-3.19	42,42,42,42	0
3	CA	B	509	1/1	0.98	0.04	-3.22	43,43,43,43	0
3	CA	A	502	1/1	0.96	0.04	-3.49	38,38,38,38	0
3	CA	A	504	1/1	1.00	0.04	-3.62	20,20,20,20	0
3	CA	B	508	1/1	0.98	0.03	-3.86	37,37,37,37	0
3	CA	B	506	1/1	0.98	0.05	-4.19	45,45,45,45	0
3	CA	C	512	1/1	0.98	0.05	-4.89	32,32,32,32	0
3	CA	C	513	1/1	0.99	0.02	-4.90	30,30,30,30	0
3	CA	B	507	1/1	0.97	0.07	-5.45	51,51,51,51	0
3	CA	A	501	1/1	0.98	0.02	-5.65	33,33,33,33	0
3	CA	A	503	1/1	1.00	0.02	-8.05	21,21,21,21	0
4	NI	C	515	1/1	0.99	0.04	-	44,44,44,44	0
4	NI	B	510	1/1	0.96	0.05	-	57,57,57,57	1
4	NI	A	505	1/1	0.97	0.04	-	38,38,38,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.