



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 12:33 PM UTC

PDB ID : 2ARL / pdb_00002arl
Title : The 2.0 angstroms crystal structure of a pocilloporin at pH 3.5: the structural basis for the linkage between color transition and halide binding
Authors : Wilmann, P.G.; Battad, J.; Beddoe, T.; Olsen, S.; Smith, S.C.; Dove, S.; Devenish, R.J.; Rossjohn, J.; Prescott, M.
Deposited on : 2005-08-19
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

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

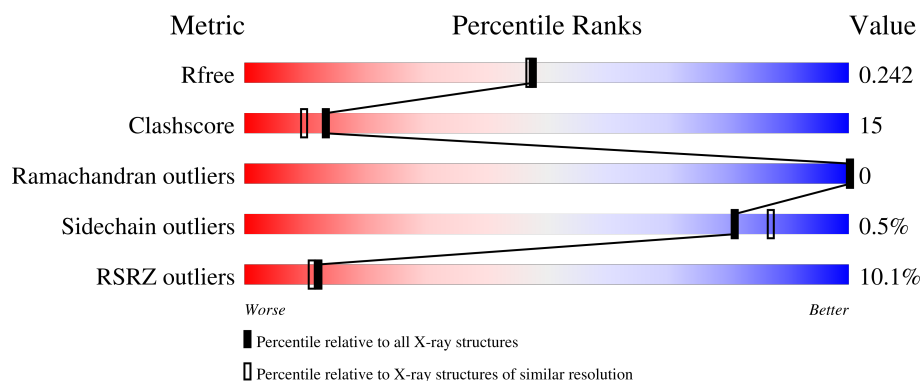
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 $\leq 5\%$. 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	219	10% 72% 27%

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
2	IOD	A	504	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACY	A	523	-	-	X	-
4	ACY	A	524	-	-	X	-
4	ACY	A	525	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 1922 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 GFP-like non-fluorescent chromoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1742	1116	287	327	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	GLY	-	expression tag	UNP P83690
A	36	LYS	ARG	SEE REMARK 999	UNP P83690
A	66	CRQ	GLN	chromophore	UNP P83690
A	66	CRQ	TYR	chromophore	UNP P83690
A	66	CRQ	GLY	chromophore	UNP P83690
A	146	SER	HIS	engineered mutation	UNP P83690

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	I	0	0
			4	4		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

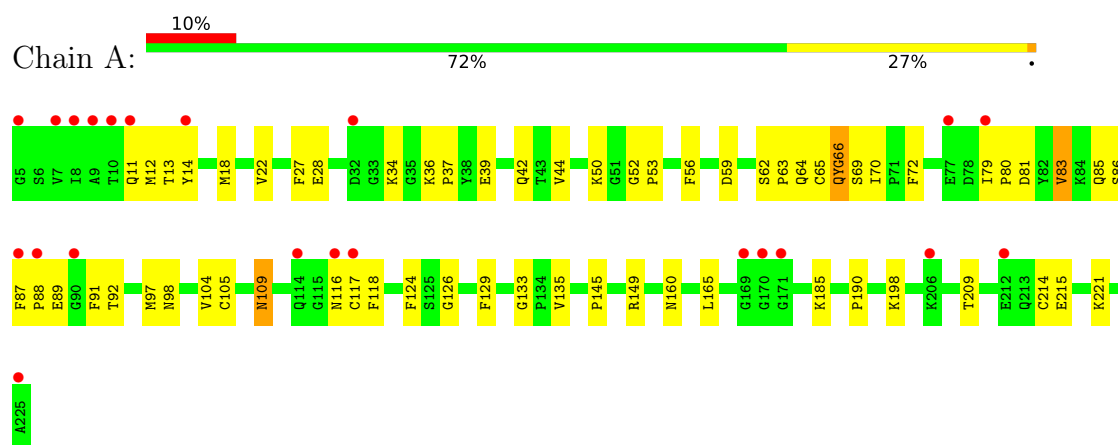
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	154	Total	O	0	0
			154	154		

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: GFP-like non-fluorescent chromoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 2 2	Depositor
Cell constants a, b, c, α , β , γ	92.77Å 92.77Å 75.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 50.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.00) 99.3 (50.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.237 , 0.253 0.223 , 0.242	Depositor DCC
R_{free} test set	1154 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtrriage
Anisotropy	(Not available)	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 42.3	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	1922	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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: CRQ, IOD, ACY, 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	0.39	0/1766	0.87	4/2387 (0.2%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	VAL	N-CA-C	5.67	115.86	110.42
1	A	209	THR	N-CA-C	-5.37	106.68	113.18
1	A	133	GLY	CA-C-N	5.04	124.82	119.28
1	A	133	GLY	C-N-CA	5.04	124.82	119.28

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	1742	0	1678	45	0
2	A	4	0	0	2	0
3	A	2	0	0	0	0
4	A	20	0	15	8	0
5	A	154	0	0	6	3
All	All	1922	0	1693	51	3

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 15.

All (51) 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:A:34:LYS:HD2	4:A:525:ACY:O	1.36	1.22
4:A:524:ACY:CH3	5:A:678:HOH:O	1.85	1.20
4:A:524:ACY:H3	5:A:678:HOH:O	1.41	1.15
1:A:65:CYS:SG	5:A:662:HOH:O	2.36	0.84
1:A:149:ARG:HH21	1:A:160:ASN:ND2	1.81	0.78
4:A:524:ACY:H2	5:A:678:HOH:O	1.64	0.77
1:A:62:SER:OG	1:A:63:PRO:HD3	1.92	0.70
1:A:36:LYS:HB2	1:A:39:GLU:HB2	1.78	0.64
1:A:12:MET:HA	1:A:116:ASN:HD21	1.63	0.62
4:A:523:ACY:OXT	4:A:525:ACY:OXT	2.19	0.61
1:A:18:MET:HE3	1:A:27:PHE:HE1	1.66	0.60
1:A:28:GLU:HB2	1:A:50:LYS:HB3	1.84	0.59
1:A:214:CYS:HB2	5:A:663:HOH:O	2.02	0.58
1:A:88:PRO:HG2	2:A:504:IOD:I	2.74	0.57
1:A:145:PRO:HB3	1:A:198:LYS:HD3	1.87	0.56
1:A:80:PRO:HG2	1:A:190:PRO:HA	1.89	0.55
1:A:149:ARG:HE	1:A:160:ASN:HD22	1.55	0.55
1:A:42:GLN:HE22	1:A:69:SER:CB	2.21	0.54
1:A:12:MET:HA	1:A:116:ASN:ND2	2.24	0.52
4:A:523:ACY:O	4:A:525:ACY:O	2.28	0.52
1:A:11:GLN:HE21	4:A:525:ACY:H1	1.75	0.51
1:A:37:PRO:HA	1:A:72:PHE:HA	1.91	0.51
1:A:12:MET:SD	1:A:116:ASN:ND2	2.81	0.50
1:A:83:VAL:HG13	1:A:91:PHE:CE1	2.47	0.49
1:A:88:PRO:HB2	2:A:504:IOD:I	2.83	0.49
1:A:66:CRQ:HB12	1:A:215:GLU:OE1	2.12	0.49
1:A:149:ARG:HD2	1:A:149:ARG:C	2.38	0.48
1:A:81:ASP:O	1:A:85:GLN:HG3	2.14	0.47
1:A:22:VAL:HA	1:A:126:GLY:O	2.15	0.47
1:A:79:ILE:HD11	1:A:221:LYS:HG2	1.97	0.47
1:A:52:GLY:HA2	1:A:53:PRO:C	2.40	0.46
1:A:87:PHE:HB3	1:A:88:PRO:HA	1.98	0.46
1:A:18:MET:HE3	1:A:27:PHE:CE1	2.50	0.46
1:A:59:ASP:HB3	1:A:165:LEU:HD21	1.97	0.46
1:A:14:TYR:OH	1:A:44:VAL:HB	2.17	0.45
4:A:524:ACY:O	5:A:631:HOH:O	2.21	0.45
1:A:98:ASN:HD22	1:A:104:VAL:HG22	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:TYR:HB3	1:A:118:PHE:HB2	1.98	0.44
1:A:86:SER:O	1:A:89:GLU:HB2	2.18	0.44
1:A:36:LYS:HA	1:A:37:PRO:HD2	1.87	0.44
1:A:89:GLU:OE1	1:A:185:LYS:HD3	2.17	0.44
1:A:109:ASN:ND2	1:A:109:ASN:C	2.76	0.43
1:A:13:THR:O	1:A:117:CYS:HA	2.20	0.42
1:A:64:GLN:HG3	1:A:124:PHE:CD2	2.54	0.42
1:A:70:ILE:CG2	1:A:83:VAL:HG21	2.50	0.41
1:A:80:PRO:HG2	1:A:190:PRO:CA	2.49	0.41
1:A:92:THR:HB	1:A:109:ASN:O	2.21	0.41
1:A:22:VAL:HG21	1:A:56:PHE:CZ	2.56	0.40
1:A:129:PHE:CE1	1:A:135:VAL:HG21	2.57	0.40
1:A:79:ILE:HA	1:A:80:PRO:HD2	1.96	0.40
1:A:97:MET:HB2	1:A:105:CYS:HB2	2.03	0.40

All (3) 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)
5:A:623:HOH:O	5:A:676:HOH:O[5_555]	1.47	0.73
5:A:637:HOH:O	5:A:670:HOH:O[5_555]	1.56	0.64
5:A:623:HOH:O	5:A:623:HOH:O[5_555]	2.12	0.08

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	214/219 (98%)	209 (98%)	5 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	A	188/188 (100%)	187 (100%)	1 (0%)	81	87

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	42	GLN
1	A	98	ASN
1	A	109	ASN
1	A	116	ASN
1	A	132	ASN
1	A	160	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](#)

1 non-standard protein/DNA/RNA residue is 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	CRQ	A	66	1	25,25,26	2.57	10 (40%)	28,34,36	2.42	5 (17%)

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	CRQ	A	66	1	-	5/10/32/33	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	CRQ	CA3-C3	-8.77	1.18	1.49
1	A	66	CRQ	CA3-N3	3.96	1.54	1.47
1	A	66	CRQ	C1-N3	3.78	1.44	1.38
1	A	66	CRQ	CB2-CA2	3.42	1.38	1.35
1	A	66	CRQ	C1-CA1	-2.44	1.44	1.48
1	A	66	CRQ	CG2-CB2	2.43	1.51	1.46
1	A	66	CRQ	CE1-CD1	2.31	1.42	1.38
1	A	66	CRQ	CE2-CD2	2.08	1.42	1.38
1	A	66	CRQ	O3-C3	2.03	1.31	1.20
1	A	66	CRQ	CE1-CZ	2.01	1.42	1.39

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CRQ	CG2-CB2-CA2	7.55	138.85	129.87
1	A	66	CRQ	CB2-CA2-N2	-6.75	119.60	128.76
1	A	66	CRQ	CB2-CA2-C2	6.07	129.71	122.36
1	A	66	CRQ	C3-CA3-N3	2.61	118.37	112.43
1	A	66	CRQ	C2-CA2-N2	2.42	110.69	108.95

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	66	CRQ	C1-CA1-CB1-CG1
1	A	66	CRQ	N2-CA2-CB2-CG2
1	A	66	CRQ	C2-CA2-CB2-CG2
1	A	66	CRQ	C3-CA3-N3-C1

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Mol	Chain	Res	Type	Atoms
1	A	66	CRQ	CA1-CB1-CG1-CD3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	66	CRQ	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 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	$\# Z > 2$
4	ACY	A	524	-	3,3,3	1.42	0	3,3,3	1.49	0
4	ACY	A	521	-	3,3,3	1.52	0	3,3,3	1.42	0
4	ACY	A	525	-	3,3,3	1.55	0	3,3,3	1.40	0
4	ACY	A	522	-	3,3,3	1.61	1 (33%)	3,3,3	1.30	0
4	ACY	A	523	-	3,3,3	1.58	0	3,3,3	1.34	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	522	ACY	OXT-C	-2.42	1.19	1.30

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	524	ACY	4	0
4	A	525	ACY	4	0
4	A	523	ACY	2	0

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	218/219 (99%)	0.69	22 (10%) 12 11	25, 37, 50, 56	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	114	GLN	4.2
1	A	88	PRO	4.0
1	A	169	GLY	3.6
1	A	170	GLY	3.6
1	A	77	GLU	3.4
1	A	116	ASN	3.3
1	A	32	ASP	3.0
1	A	87	PHE	2.9
1	A	212	GLU	2.8
1	A	9	ALA	2.7
1	A	225	ALA	2.6
1	A	8	ILE	2.5
1	A	10	THR	2.5
1	A	11	GLN	2.5
1	A	7	VAL	2.5
1	A	117	CYS	2.4
1	A	79	ILE	2.1
1	A	90	GLY	2.1
1	A	5	GLY	2.0
1	A	206	LYS	2.0
1	A	171	GLY	2.0
1	A	14	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CRQ	A	66	24/25	0.90	0.11	31,35,40,40	0

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	ACY	A	525	4/4	0.61	0.25	53,54,55,55	0
4	ACY	A	521	4/4	0.73	0.22	45,48,48,49	0
4	ACY	A	524	4/4	0.75	0.18	55,56,56,57	0
4	ACY	A	523	4/4	0.78	0.20	46,46,46,47	0
4	ACY	A	522	4/4	0.88	0.13	27,30,31,31	0
2	IOD	A	504	1/1	0.97	0.18	61,61,61,61	1
2	IOD	A	501	1/1	0.97	0.04	38,38,38,38	0
3	CL	A	512	1/1	0.98	0.13	38,38,38,38	0
2	IOD	A	503	1/1	0.98	0.15	56,56,56,56	0
3	CL	A	511	1/1	0.98	0.11	27,27,27,27	0
2	IOD	A	502	1/1	0.99	0.03	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.