



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 11:34 am BST

PDB ID : 1UIS  
Title : The 2.0 crystal structure of eqFP611, a far-red fluorescent protein from the sea anemone *Entacmaea quadricolor*  
Authors : Petersen, J.; Wilmann, P.G.; Beddoe, T.; Oakley, A.J.; Devenish, R.J.; Prescott, M.; Rossjohn, J.  
Deposited on : 2003-07-21  
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

<https://www.wwpdb.org/validation/2017/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  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

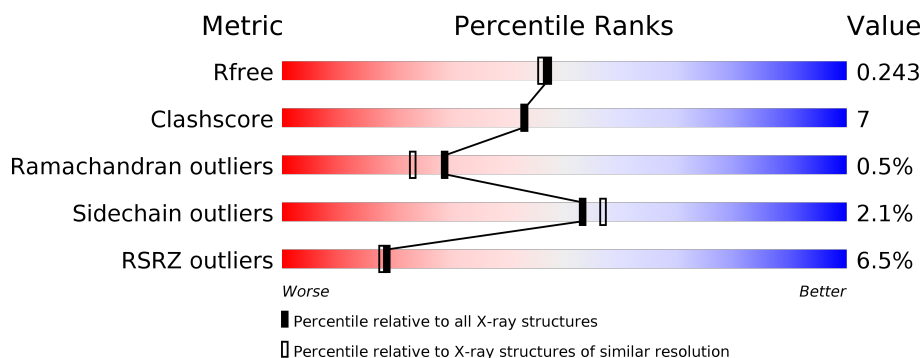
# 1 Overall quality at a glance

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))
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 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	231	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>• •</div> </div> </div>
1	B	231	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div>• •</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3887 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 red fluorescent protein FP611.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	0	0
			1796	1144	299	335	18			
1	B	225	Total	C	N	O	S	0	0	0
			1796	1144	299	335	18			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ILE	-	CLONING ARTIFACT	UNP Q8ISF8
A	0	HIS	-	CLONING ARTIFACT	UNP Q8ISF8
A	63	NRQ	MET	CHROMOPHORE	UNP Q8ISF8
A	63	NRQ	TYR	CHROMOPHORE	UNP Q8ISF8
A	63	NRQ	GLY	CHROMOPHORE	UNP Q8ISF8
B	-1	ILE	-	CLONING ARTIFACT	UNP Q8ISF8
B	0	HIS	-	CLONING ARTIFACT	UNP Q8ISF8
B	63	NRQ	MET	CHROMOPHORE	UNP Q8ISF8
B	63	NRQ	TYR	CHROMOPHORE	UNP Q8ISF8
B	63	NRQ	GLY	CHROMOPHORE	UNP Q8ISF8

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is ACETIC ACID (three-letter code: ACY) (formula: C<sub>2</sub>H<sub>4</sub>O<sub>2</sub>).



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

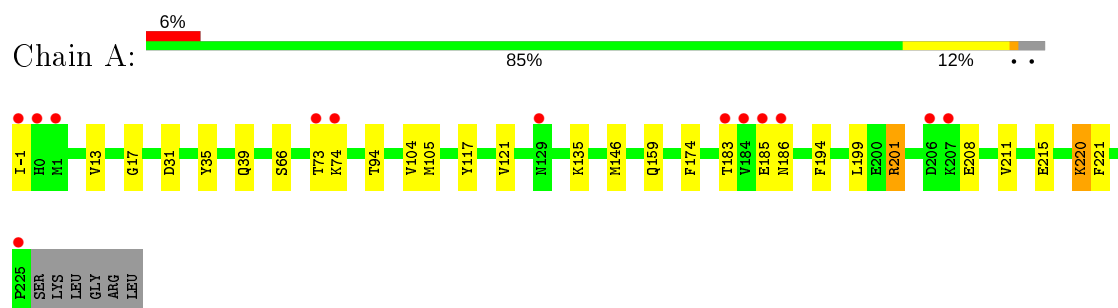
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	150	Total	O	0	0
			150	150		
4	B	135	Total	O	0	0
			135	135		

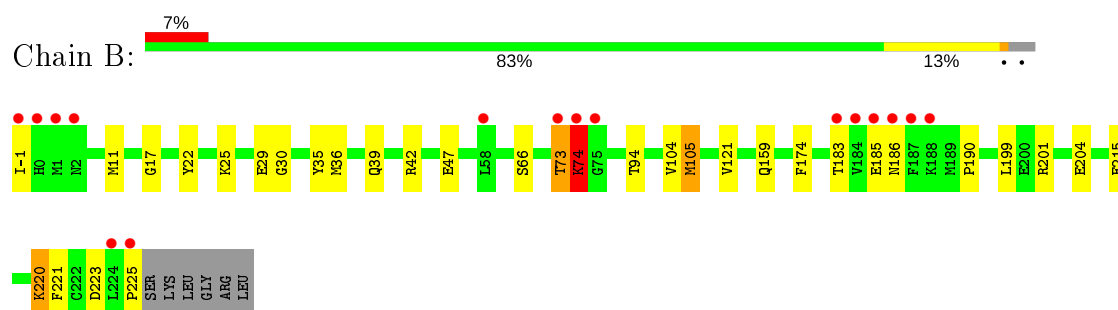
### 3 Residue-property plots [i](#)

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: red fluorescent protein FP611



- Molecule 1: red fluorescent protein FP611



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.94Å 77.94Å 327.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.00 19.72 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.00) 91.7 (19.72-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.69 (at 2.01Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.213 , 0.253 0.205 , 0.243	Depositor DCC
$R_{free}$ test set	1514 reflections (4.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.0	Xtriage
Anisotropy	0.470	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 43.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3887	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, NRQ, 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	0.39	0/1817	0.64	0/2444
1	B	0.38	0/1817	0.65	0/2444
All	All	0.38	0/3634	0.64	0/4888

There are no bond length outliers.

There are no bond angle outliers.

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	1796	0	1723	27	1
1	B	1796	0	1723	23	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
4	A	150	0	0	5	1
4	B	135	0	0	3	1
All	All	3887	0	3452	48	2

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 7.

All (48) 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:201:ARG:HD2	1:A:211:VAL:HG13	1.63	0.79
1:B:30:GLY:HA2	1:B:42:ARG:HH12	1.63	0.63
1:B:-1:ILE:HA	1:B:35:TYR:CZ	2.35	0.62
1:A:-1:ILE:HA	1:A:35:TYR:CZ	2.34	0.61
1:B:94:THR:HA	1:B:174:PHE:HB3	1.81	0.61
1:A:94:THR:HA	1:A:174:PHE:HB3	1.83	0.61
1:A:73:THR:HG22	1:A:74:LYS:H	1.69	0.58
1:A:183:THR:HB	1:A:186:ASN:HD22	1.71	0.55
1:A:199:LEU:HD22	1:A:215:GLU:HB2	1.88	0.54
1:A:13:VAL:HG13	1:A:117:TYR:HB2	1.90	0.54
1:B:223:ASP:O	1:B:225:PRO:HD3	2.08	0.54
1:A:39:GLN:HE22	1:A:66:SER:CB	2.22	0.53
1:B:29:GLU:OE1	1:B:42:ARG:HD2	2.08	0.53
1:B:74:LYS:HD2	1:B:74:LYS:H	1.75	0.52
1:B:39:GLN:HE22	1:B:66:SER:CB	2.22	0.52
1:A:183:THR:HB	1:A:186:ASN:ND2	2.25	0.51
1:A:183:THR:HG22	1:A:185:GLU:HG3	1.91	0.51
1:A:135:LYS:HE2	4:A:1113:HOH:O	2.12	0.50
1:A:201:ARG:HD2	1:A:211:VAL:CG1	2.37	0.49
1:B:201:ARG:HD3	1:B:204:GLU:HB3	1.95	0.48
1:B:199:LEU:HD22	1:B:215:GLU:HB2	1.94	0.48
1:B:183:THR:HB	1:B:186:ASN:HD22	1.77	0.48
1:A:159:GLN:NE2	4:A:1141:HOH:O	2.45	0.48
1:A:73:THR:HB	4:A:1145:HOH:O	2.13	0.48
1:B:73:THR:HB	1:B:74:LYS:HD2	1.96	0.47
1:A:104:VAL:HG22	1:A:121:VAL:HG12	1.96	0.47
1:B:22:TYR:HB2	4:B:1078:HOH:O	2.15	0.47
1:A:-1:ILE:HG23	1:A:35:TYR:CD2	2.51	0.46
1:A:159:GLN:OE1	1:B:159:GLN:HG3	2.17	0.45
1:A:220:LYS:HE2	1:A:221:PHE:O	2.17	0.45
1:B:183:THR:HB	1:B:186:ASN:ND2	2.31	0.45
1:A:17:GLY:HA3	1:A:121:VAL:O	2.17	0.44
1:B:220:LYS:HE2	1:B:221:PHE:O	2.17	0.43
1:A:159:GLN:HG3	1:B:159:GLN:OE1	2.19	0.43
1:B:220:LYS:NZ	4:B:1129:HOH:O	2.51	0.43
1:B:190:PRO:HG3	4:B:1079:HOH:O	2.19	0.42
1:A:201:ARG:HD3	4:A:1027:HOH:O	2.18	0.42
1:A:199:LEU:CD2	1:A:215:GLU:HB2	2.50	0.42
1:A:13:VAL:HG22	1:A:117:TYR:CD2	2.55	0.42
1:A:220:LYS:HB3	4:A:1145:HOH:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:VAL:HG22	1:B:121:VAL:HG12	2.02	0.41
1:A:-1:ILE:O	1:A:-1:ILE:HG22	2.20	0.41
1:A:146:MET:HE3	1:A:194:PHE:HE2	1.85	0.41
1:B:17:GLY:HA3	1:B:121:VAL:O	2.20	0.41
1:A:201:ARG:NH2	1:A:208:GLU:OE2	2.50	0.40
1:B:74:LYS:HD2	1:B:74:LYS:N	2.37	0.40
1:B:183:THR:HG22	1:B:185:GLU:HG3	2.02	0.40
1:B:25:LYS:HB3	1:B:47:GLU:HB2	2.03	0.40

All (2) 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:A:105:MET:CG	4:B:1138:HOH:O[12_566]	1.63	0.57
1:B:105:MET:CG	4:A:1152:HOH:O[12_566]	1.66	0.54

## 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	220/231 (95%)	215 (98%)	5 (2%)	0	100	100
1	B	220/231 (95%)	216 (98%)	2 (1%)	2 (1%)	17	11
All	All	440/462 (95%)	431 (98%)	7 (2%)	2 (0%)	29	23

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	73	THR
1	B	74	LYS

### 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	193/198 (98%)	190 (98%)	3 (2%)	62	67
1	B	193/198 (98%)	188 (97%)	5 (3%)	46	48
All	All	386/396 (98%)	378 (98%)	8 (2%)	53	57

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ASP
1	A	201	ARG
1	A	220	LYS
1	B	11	MET
1	B	36	MET
1	B	74	LYS
1	B	105	MET
1	B	220	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	39	GLN
1	A	159	GLN
1	A	186	ASN
1	B	0	HIS
1	B	8	ASN
1	B	23	GLN
1	B	39	GLN
1	B	118	HIS
1	B	159	GLN
1	B	186	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are 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	NRQ	B	63	1	23,24,25	3.91	11 (47%)	23,32,34	3.24	8 (34%)
1	NRQ	A	63	1	23,24,25	3.96	11 (47%)	23,32,34	3.24	9 (39%)

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	NRQ	B	63	1	-	4/9/31/32	0/2/2/2
1	NRQ	A	63	1	-	4/9/31/32	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	63	NRQ	CA3-C3	-8.25	1.21	1.49
1	A	63	NRQ	CA3-C3	-8.15	1.21	1.49
1	A	63	NRQ	CB2-CA2	8.12	1.41	1.35
1	B	63	NRQ	CB2-CA2	7.67	1.41	1.35
1	A	63	NRQ	CD2-CG2	5.95	1.51	1.39
1	B	63	NRQ	CE2-CZ	5.94	1.50	1.38
1	A	63	NRQ	CE1-CZ	5.88	1.50	1.38
1	B	63	NRQ	CE1-CZ	5.87	1.50	1.38
1	B	63	NRQ	CD2-CG2	5.83	1.50	1.39
1	A	63	NRQ	CE2-CZ	5.82	1.50	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	63	NRQ	CD1-CG2	5.57	1.50	1.39
1	B	63	NRQ	CD1-CG2	5.54	1.50	1.39
1	A	63	NRQ	CA3-N3	5.49	1.58	1.47
1	B	63	NRQ	CA3-N3	5.24	1.57	1.47
1	A	63	NRQ	OH-CZ	-4.86	1.25	1.37
1	B	63	NRQ	OH-CZ	-4.63	1.26	1.37
1	B	63	NRQ	C1-N3	4.47	1.45	1.38
1	A	63	NRQ	C1-N3	4.40	1.45	1.38
1	A	63	NRQ	O3-C3	2.29	1.32	1.19
1	B	63	NRQ	O3-C3	2.25	1.32	1.19
1	A	63	NRQ	CG2-CB2	-2.08	1.42	1.46
1	B	63	NRQ	CG2-CB2	-2.06	1.42	1.46

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	63	NRQ	CB2-CA2-C2	9.05	133.08	122.28
1	A	63	NRQ	CB2-CA2-C2	8.99	133.01	122.28
1	B	63	NRQ	CB2-CA2-N2	-8.35	117.25	128.83
1	A	63	NRQ	CB2-CA2-N2	-8.19	117.47	128.83
1	A	63	NRQ	CG2-CB2-CA2	6.72	138.18	129.94
1	B	63	NRQ	CG2-CB2-CA2	6.48	137.89	129.94
1	A	63	NRQ	CA2-N2-C1	2.68	109.23	104.33
1	B	63	NRQ	CA2-N2-C1	2.51	108.93	104.33
1	A	63	NRQ	O2-C2-CA2	-2.44	129.59	130.96
1	B	63	NRQ	CA2-C2-N3	2.27	104.44	103.37
1	B	63	NRQ	O3-C3-CA3	-2.26	119.58	126.39
1	A	63	NRQ	CE1-CZ-CE2	-2.21	116.04	119.77
1	B	63	NRQ	CE1-CZ-CE2	-2.21	116.05	119.77
1	A	63	NRQ	O3-C3-CA3	-2.16	119.86	126.39
1	A	63	NRQ	CD1-CE1-CZ	2.11	122.19	119.88
1	A	63	NRQ	CA2-C2-N3	2.05	104.34	103.37
1	B	63	NRQ	O2-C2-CA2	-2.03	129.82	130.96

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	63	NRQ	C2-CA2-CB2-CG2
1	A	63	NRQ	C2-CA2-CB2-CG2
1	B	63	NRQ	CB1-CG1-SD-CE
1	A	63	NRQ	CB1-CG1-SD-CE

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Mol	Chain	Res	Type	Atoms
1	B	63	NRQ	N2-CA2-CB2-CG2
1	A	63	NRQ	N2-CA2-CB2-CG2
1	B	63	NRQ	C3-CA3-N3-C1
1	A	63	NRQ	C3-CA3-N3-C1

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 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	ACY	B	1002	-	1,3,3	2.98	1 (100%)	0,3,3	0.00	-
3	ACY	A	1001	-	1,3,3	3.28	1 (100%)	0,3,3	0.00	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1001	ACY	CH3-C	3.28	1.52	1.48
3	B	1002	ACY	CH3-C	2.98	1.52	1.48

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	224/231 (96%)	0.08	13 (5%) 23 22	25, 36, 53, 78	0
1	B	224/231 (96%)	0.18	16 (7%) 16 15	26, 35, 61, 83	0
All	All	448/462 (96%)	0.13	29 (6%) 18 18	25, 35, 58, 83	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	7.3
1	B	-1	ILE	6.9
1	B	184	VAL	5.2
1	B	0	HIS	4.7
1	B	225	PRO	4.5
1	A	73	THR	4.4
1	A	185	GLU	3.8
1	B	73	THR	3.7
1	A	-1	ILE	3.4
1	A	129	ASN	3.3
1	B	2	ASN	3.2
1	B	74	LYS	3.1
1	A	184	VAL	3.1
1	A	225	PRO	3.1
1	A	74	LYS	3.1
1	B	185	GLU	3.1
1	A	183	THR	2.9
1	A	186	ASN	2.8
1	B	183	THR	2.7
1	B	58	LEU	2.7
1	A	206	ASP	2.6
1	B	75	GLY	2.6
1	A	1	MET	2.5
1	A	0	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	188	LYS	2.2
1	B	187	PHE	2.1
1	A	207	LYS	2.1
1	B	224	LEU	2.1
1	B	186	ASN	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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
1	NRQ	A	63	23/24	0.96	0.09	27,30,34,41	0
1	NRQ	B	63	23/24	0.97	0.09	27,30,37,40	0

## 6.3 Carbohydrates [i](#)

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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	ACY	A	1001	4/4	0.95	0.12	37,38,39,39	0
3	ACY	B	1002	4/4	0.98	0.13	39,40,41,44	0
2	CA	A	1004	1/1	0.98	0.05	44,44,44,44	0
2	CA	B	1003	1/1	0.98	0.24	53,53,53,53	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.