



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 07:59 AM GMT

PDB ID : 3A8S
Title : Crystal structure analysis of the fluorescent protein KillerRed
Authors : Sakai, N.; Kitago, Y.; Takemoto, K.; Matsuda, T.; Nagai, T.
Deposited on : 2009-10-08
Resolution : 2.90 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

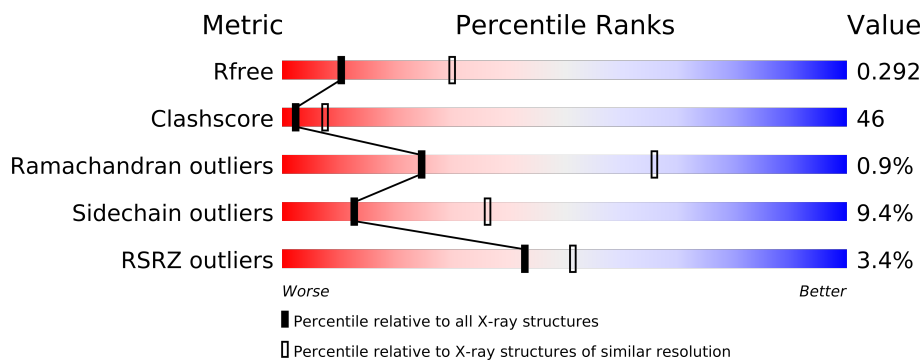
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance



The reported resolution of this entry is 2.90 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}	66092	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	238	
1	B	238	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3546 atoms, of which 0 are hydrogen and 0 are deuterium.

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 KillerRed.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1760	1113	305	327	15			
1	B	221	Total	C	N	O	S	0	0	0
			1758	1112	305	326	15			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	HIS	-	EXPRESSION TAG	UNP Q2TCH5
A	-1	MET	-	EXPRESSION TAG	UNP Q2TCH5
A	0	GLY	-	EXPRESSION TAG	UNP Q2TCH5
A	1	SER	-	EXPRESSION TAG	UNP Q2TCH5
A	65	CRQ	GLN	CHROMOPHORE	UNP Q2TCH5
A	65	CRQ	TYR	CHROMOPHORE	UNP Q2TCH5
A	65	CRQ	GLY	CHROMOPHORE	UNP Q2TCH5
B	-2	HIS	-	EXPRESSION TAG	UNP Q2TCH5
B	-1	MET	-	EXPRESSION TAG	UNP Q2TCH5
B	0	GLY	-	EXPRESSION TAG	UNP Q2TCH5
B	1	SER	-	EXPRESSION TAG	UNP Q2TCH5
B	65	CRQ	GLN	CHROMOPHORE	UNP Q2TCH5
B	65	CRQ	TYR	CHROMOPHORE	UNP Q2TCH5
B	65	CRQ	GLY	CHROMOPHORE	UNP Q2TCH5

- Molecule 2 is water.

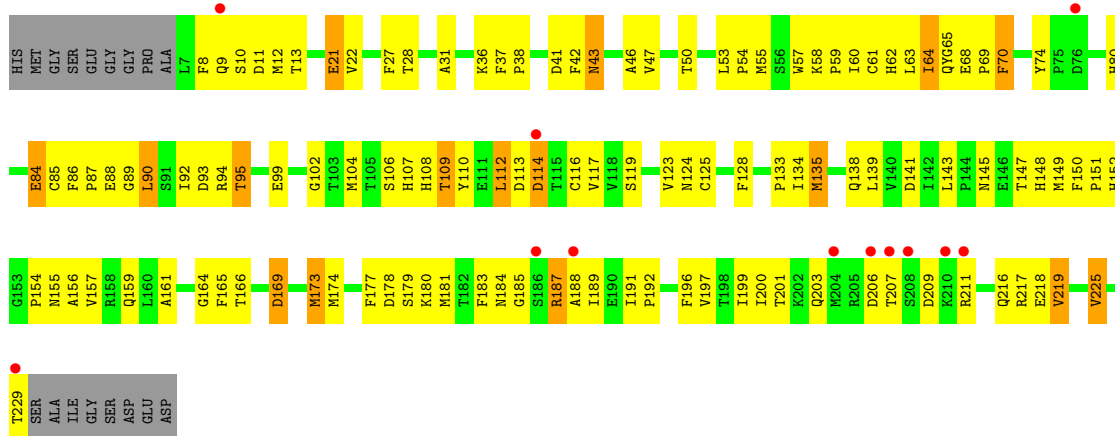
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	13	Total	O	0	0
			13	13		
2	B	15	Total	O	0	0
			15	15		

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 errors 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.

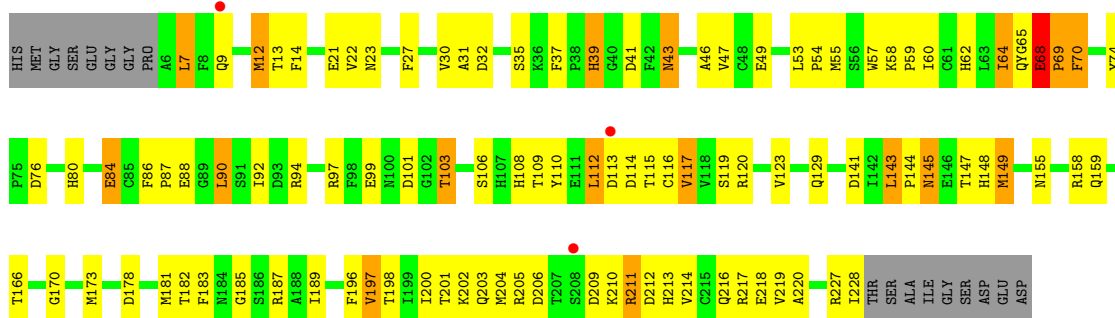
• Molecule 1: KillerRed

Chain A: 



• Molecule 1: KillerRed

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	123.46Å 123.46Å 110.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.88 – 2.90 19.88 – 2.90	Depositor EDS
% Data completeness (in resolution range)	94.9 (19.88-2.90) 94.8 (19.88-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	20.83 (at 2.88Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.245 , 0.298 0.244 , 0.292	Depositor DCC
R_{free} test set	1062 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	51.7	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 25.9	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 20685 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	3546	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CRQ

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	0.28	0/1783	0.44	0/2414
1	B	0.29	0/1781	0.47	0/2411
All	All	0.28	0/3564	0.45	0/4825

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	64	ILE	Mainchain
1	B	64	ILE	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1760	0	1670	180	0
1	B	1758	0	1668	139	0
2	A	13	0	0	2	0
2	B	15	0	0	4	0
All	All	3546	0	3338	315	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 46.

All (315) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:69:PRO:HG2	1:B:181:MET:CE	1.66	1.25
1:B:7:LEU:O	1:B:7:LEU:CD1	1.89	1.19
1:A:135:MET:HA	1:A:135:MET:CE	1.78	1.14
1:B:69:PRO:HG2	1:B:181:MET:HE3	1.12	1.09
1:A:141:ASP:HB2	1:A:166:THR:OG1	1.53	1.07
1:A:92:ILE:HB	1:A:108:HIS:HB2	1.33	1.05
1:B:7:LEU:O	1:B:7:LEU:HD12	1.51	1.04
1:A:57:TRP:CE3	1:A:60:ILE:HD11	1.93	1.04
1:B:69:PRO:CG	1:B:181:MET:HE3	1.88	1.03
1:A:58:LYS:HB2	1:A:59:PRO:HD3	1.41	1.02
1:B:84:GLU:HB3	1:B:189:ILE:HD11	1.41	1.02
1:A:135:MET:HA	1:A:135:MET:HE2	1.40	1.02
1:B:149:MET:HA	1:B:149:MET:HE2	1.42	1.00
1:B:7:LEU:O	1:B:7:LEU:HD13	1.63	0.98
1:B:68:GLU:H	1:B:69:PRO:HD3	1.29	0.96
1:B:7:LEU:C	1:B:7:LEU:HD12	1.85	0.96
1:A:147:THR:HG21	1:A:159:GLN:HE21	1.31	0.95
1:B:62:HIS:HD2	1:B:94:ARG:HH11	0.99	0.94
1:B:170:GLY:HA2	2:B:319:HOH:O	1.67	0.94
1:A:173:MET:HE3	1:A:173:MET:HA	1.51	0.92
1:A:88:GLU:OE2	1:A:187:ARG:NH2	2.03	0.91
1:B:149:MET:HA	1:B:149:MET:CE	2.01	0.90
1:B:205:ARG:NH2	1:B:212:ASP:OD1	2.04	0.90
1:A:135:MET:CA	1:A:135:MET:CE	2.51	0.88
1:A:147:THR:HG21	1:A:159:GLN:NE2	1.89	0.88
1:A:196:PHE:HE2	1:A:225:VAL:CG1	1.87	0.88
1:B:62:HIS:CD2	1:B:94:ARG:HH11	1.91	0.86
1:A:209:ASP:HB3	1:A:211:ARG:HG2	1.57	0.86
1:B:149:MET:HE1	1:B:159:GLN:HB2	1.59	0.84

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:135:MET:N	1:A:135:MET:HE3	1.92	0.84
1:A:196:PHE:HE2	1:A:225:VAL:HG12	1.40	0.84
1:B:92:ILE:HB	1:B:108:HIS:HB2	1.61	0.83
1:B:69:PRO:HG2	1:B:181:MET:HE1	1.59	0.83
1:A:64:ILE:O	1:A:65:CRQ:HG11	1.80	0.82
1:A:135:MET:HE3	1:A:135:MET:CA	2.10	0.81
1:A:84:GLU:HB3	1:A:189:ILE:HD11	1.59	0.81
1:A:112:LEU:C	1:A:112:LEU:HD13	2.01	0.81
1:A:173:MET:CE	1:A:173:MET:HA	2.11	0.80
1:B:22:VAL:HG23	1:B:27:PHE:HE2	1.46	0.80
1:A:65:CRQ:HB11	1:A:218:GLU:OE1	1.83	0.79
1:B:22:VAL:HG23	1:B:27:PHE:CE2	2.17	0.78
1:B:68:GLU:H	1:B:69:PRO:CD	1.96	0.77
1:B:147:THR:OG1	1:B:197:VAL:HG13	1.85	0.77
1:A:94:ARG:HH11	1:A:94:ARG:HG3	1.46	0.77
1:B:69:PRO:CG	1:B:181:MET:CE	2.51	0.77
1:B:62:HIS:HD2	1:B:94:ARG:NH1	1.79	0.77
1:A:88:GLU:N	1:A:88:GLU:OE1	2.18	0.76
1:A:147:THR:HB	1:A:197:VAL:CG1	2.15	0.76
1:A:183:PHE:CZ	1:A:185:GLY:HA3	2.21	0.76
1:B:149:MET:HE1	1:B:159:GLN:CB	2.16	0.75
1:A:31:ALA:HB2	1:A:46:ALA:CB	2.16	0.75
1:A:61:CYS:HA	1:A:216:GLN:HE22	1.48	0.75
1:A:68:GLU:HG3	1:A:70:PHE:CE1	2.21	0.75
1:A:57:TRP:CZ3	1:A:60:ILE:HD11	2.22	0.75
1:B:201:THR:HG23	2:B:303:HOH:O	1.86	0.75
1:A:209:ASP:HB3	1:A:211:ARG:CG	2.16	0.74
1:A:109:THR:O	1:A:119:SER:HA	1.88	0.74
1:B:65:CRQ:HD2	1:B:65:CRQ:N2	2.01	0.74
1:B:183:PHE:CZ	1:B:185:GLY:HA3	2.23	0.74
1:B:149:MET:HE3	1:B:159:GLN:HA	1.70	0.73
1:A:10:SER:O	1:A:12:MET:HG3	1.87	0.73
1:A:65:CRQ:HE2	1:A:201:THR:HG21	1.71	0.73
1:A:53:LEU:HD12	1:A:54:PRO:HD2	1.71	0.72
1:A:90:LEU:HD22	1:A:90:LEU:C	2.10	0.72
1:A:112:LEU:HD22	1:A:113:ASP:H	1.54	0.71
1:A:112:LEU:HD22	1:A:113:ASP:N	2.05	0.71
1:A:196:PHE:CE2	1:A:225:VAL:CG1	2.73	0.71
1:B:149:MET:CE	1:B:159:GLN:HA	2.21	0.70
1:B:203:GLN:O	1:B:204:MET:HG2	1.92	0.70
1:B:68:GLU:O	1:B:70:PHE:N	2.25	0.69
1:A:196:PHE:CE2	1:A:225:VAL:HG12	2.26	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:41:ASP:CG	1:B:217:ARG:HH21	1.96	0.68
1:B:86:PHE:HB3	1:B:87:PRO:HA	1.76	0.68
1:B:39:HIS:H	1:B:39:HIS:CD2	2.12	0.67
1:B:149:MET:HE3	1:B:159:GLN:HG3	1.75	0.67
1:A:58:LYS:HD3	1:A:203:GLN:NE2	2.10	0.67
1:A:147:THR:HB	1:A:197:VAL:HG13	1.75	0.67
1:A:219:VAL:HG12	1:B:228:ILE:HD11	1.75	0.67
1:A:90:LEU:N	1:A:184:ASN:ND2	2.43	0.67
1:A:61:CYS:CA	1:A:216:GLN:HE22	2.08	0.66
1:B:31:ALA:HB2	1:B:46:ALA:HB2	1.77	0.66
1:A:169:ASP:OD1	1:A:169:ASP:N	2.29	0.66
1:A:201:THR:HG22	1:A:218:GLU:HG3	1.76	0.66
1:A:31:ALA:HB2	1:A:46:ALA:HB2	1.76	0.66
1:A:65:CRQ:C	1:A:69:PRO:HD3	2.26	0.66
1:B:57:TRP:CD2	1:B:214:VAL:HG11	2.31	0.65
1:A:94:ARG:HG3	1:A:179:SER:HB2	1.77	0.65
1:B:65:CRQ:O	1:B:68:GLU:OE1	2.15	0.65
1:B:57:TRP:CE3	1:B:214:VAL:HG11	2.32	0.64
1:A:145:ASN:HA	1:B:196:PHE:HZ	1.62	0.64
1:B:84:GLU:HB3	1:B:189:ILE:CD1	2.23	0.64
1:B:149:MET:CE	1:B:159:GLN:HG3	2.27	0.64
1:A:62:HIS:HD2	1:A:94:ARG:HE	1.43	0.64
1:A:68:GLU:HG3	1:A:70:PHE:HE1	1.61	0.64
1:B:227:ARG:HB3	1:B:227:ARG:CZ	2.29	0.63
1:B:149:MET:CE	1:B:159:GLN:CB	2.76	0.63
1:A:62:HIS:HD2	1:A:94:ARG:NE	1.96	0.63
1:A:61:CYS:CB	1:A:216:GLN:HE22	2.12	0.63
1:A:135:MET:HE3	1:A:135:MET:HA	1.68	0.62
1:B:106:SER:OG	1:B:108:HIS:HE1	1.83	0.61
1:B:88:GLU:OE2	1:B:187:ARG:NH1	2.33	0.61
1:A:147:THR:HB	1:A:197:VAL:HG12	1.82	0.61
1:B:149:MET:CE	1:B:159:GLN:CA	2.79	0.60
1:B:58:LYS:HB2	1:B:59:PRO:HD3	1.83	0.60
1:A:58:LYS:HB2	1:A:59:PRO:CD	2.26	0.60
1:B:88:GLU:HG3	1:B:185:GLY:CA	2.31	0.60
1:A:58:LYS:HG3	1:A:139:LEU:HB2	1.84	0.60
1:A:99:GLU:OE1	1:A:174:MET:HB2	2.02	0.60
1:B:9:GLN:HA	1:B:9:GLN:NE2	2.17	0.59
1:B:206:ASP:HB3	1:B:213:HIS:ND1	2.16	0.59
1:B:86:PHE:CD1	1:B:112:LEU:HB2	2.35	0.59
1:A:206:ASP:O	1:A:207:THR:OG1	2.15	0.59
1:B:41:ASP:OD1	1:B:217:ARG:NH2	2.36	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:161:ALA:HB3	1:A:177:PHE:HB3	1.85	0.59
1:B:88:GLU:CD	1:B:187:ARG:HH11	2.04	0.59
1:A:57:TRP:HE3	1:A:60:ILE:HD11	1.65	0.59
1:A:36:LYS:O	1:A:38:PRO:O	2.21	0.59
1:A:94:ARG:NH1	1:A:177:PHE:CZ	2.71	0.58
1:B:57:TRP:CE3	1:B:60:ILE:HD11	2.39	0.58
1:A:55:MET:HB2	1:A:135:MET:HE1	1.85	0.58
1:A:173:MET:CA	1:A:173:MET:CE	2.82	0.58
1:B:68:GLU:N	1:B:69:PRO:CD	2.64	0.58
1:B:145:ASN:HD21	1:B:147:THR:HG22	1.68	0.58
1:A:90:LEU:O	1:A:90:LEU:HD22	2.04	0.58
1:B:201:THR:HG22	1:B:218:GLU:HB2	1.85	0.57
1:A:94:ARG:HG3	1:A:94:ARG:NH1	2.11	0.57
1:A:37:PHE:HA	1:A:38:PRO:C	2.23	0.57
1:A:93:ASP:OD1	1:A:107:HIS:HD2	1.87	0.57
1:A:109:THR:O	1:A:109:THR:HG22	2.04	0.57
1:A:65:CRQ:N2	1:A:65:CRQ:HD2	2.19	0.56
1:B:31:ALA:HB2	1:B:46:ALA:CB	2.34	0.56
1:A:53:LEU:HD12	1:A:54:PRO:CD	2.36	0.56
1:A:9:GLN:HA	1:A:37:PHE:CE1	2.41	0.56
1:B:14:PHE:HA	1:B:117:VAL:O	2.06	0.55
1:B:112:LEU:HD21	1:B:114:ASP:O	2.06	0.55
1:A:229:THR:O	1:B:202:LYS:NZ	2.40	0.55
1:A:147:THR:CB	1:A:197:VAL:HG13	2.37	0.55
1:A:86:PHE:HB3	1:A:87:PRO:HA	1.89	0.55
1:A:31:ALA:CB	1:A:46:ALA:HB2	2.36	0.55
1:A:145:ASN:HA	1:B:196:PHE:CZ	2.41	0.55
1:A:200:ILE:HB	1:A:219:VAL:HG13	1.90	0.54
1:A:149:MET:SD	1:A:159:GLN:HG3	2.48	0.54
1:B:106:SER:HB3	1:B:123:VAL:HG23	1.89	0.54
1:B:149:MET:HE2	1:B:149:MET:CA	2.23	0.54
1:A:88:GLU:O	1:A:184:ASN:HB2	2.08	0.53
1:B:149:MET:HE2	1:B:158:ARG:O	2.09	0.53
1:B:9:GLN:NE2	1:B:37:PHE:CE2	2.76	0.53
1:B:86:PHE:HD1	1:B:112:LEU:HB2	1.73	0.53
1:A:43:ASN:N	1:A:43:ASN:HD22	2.06	0.53
1:B:99:GLU:O	1:B:173:MET:HE1	2.08	0.53
1:A:173:MET:HE2	1:A:174:MET:H	1.75	0.53
1:B:88:GLU:HG3	1:B:185:GLY:HA3	1.89	0.53
1:B:58:LYS:CB	1:B:59:PRO:HD3	2.39	0.52
1:A:89:GLY:C	1:A:184:ASN:ND2	2.62	0.52
1:A:58:LYS:N	1:A:203:GLN:HE22	2.06	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:68:GLU:C	1:B:70:PHE:H	2.11	0.52
1:A:154:PRO:O	1:A:191:ILE:HD11	2.09	0.52
1:B:92:ILE:HD11	1:B:181:MET:CE	2.39	0.52
1:A:94:ARG:HG3	1:A:179:SER:CB	2.40	0.52
1:B:108:HIS:HB3	1:B:110:TYR:CE2	2.44	0.52
1:A:112:LEU:HD13	1:A:113:ASP:N	2.24	0.52
1:B:30:VAL:HG11	1:B:49:GLU:OE2	2.10	0.52
1:B:149:MET:HE1	1:B:159:GLN:CA	2.40	0.52
1:A:112:LEU:CD1	1:A:114:ASP:H	2.23	0.51
1:A:104:MET:HA	1:A:124:ASN:O	2.09	0.51
1:B:9:GLN:O	1:B:37:PHE:CD2	2.63	0.51
1:A:94:ARG:HD3	1:A:177:PHE:CE1	2.44	0.51
1:A:85:CYS:SG	1:A:183:PHE:HB2	2.50	0.51
1:A:187:ARG:HG2	1:A:188:ALA:N	2.25	0.51
1:A:86:PHE:CE1	1:A:89:GLY:HA2	2.46	0.51
1:A:84:GLU:HB3	1:A:189:ILE:CD1	2.35	0.51
1:A:53:LEU:HD22	1:A:57:TRP:CE2	2.46	0.51
1:A:94:ARG:NH1	1:A:177:PHE:HZ	2.09	0.51
1:A:74:TYR:CD1	1:A:80:HIS:CD2	2.98	0.51
1:B:65:CRQ:HG11	1:B:216:GLN:HE21	1.76	0.51
1:A:57:TRP:O	1:A:60:ILE:HG12	2.11	0.51
1:A:95:THR:HG22	1:A:178:ASP:HB3	1.92	0.51
1:B:68:GLU:C	1:B:70:PHE:N	2.63	0.50
1:B:53:LEU:HD22	1:B:57:TRP:CE2	2.46	0.50
1:A:133:PRO:HB2	1:A:139:LEU:CD2	2.41	0.50
1:B:227:ARG:HB3	1:B:227:ARG:NH1	2.26	0.50
1:A:11:ASP:OD1	1:A:36:LYS:HA	2.10	0.50
1:A:65:CRQ:CE2	1:A:199:ILE:HG21	2.42	0.50
1:A:62:HIS:CD2	1:A:94:ARG:NE	2.79	0.50
1:A:201:THR:HG22	1:A:218:GLU:CG	2.40	0.49
1:B:13:THR:OG1	1:B:116:CYS:HB2	2.12	0.49
1:B:30:VAL:CG1	1:B:49:GLU:OE2	2.60	0.49
1:A:55:MET:CB	1:A:135:MET:HE1	2.43	0.49
1:A:41:ASP:CG	1:A:217:ARG:HH21	2.16	0.49
1:B:123:VAL:O	1:B:123:VAL:HG13	2.12	0.49
1:A:134:ILE:C	1:A:135:MET:HE3	2.32	0.49
1:A:149:MET:CE	1:A:157:VAL:HG11	2.43	0.49
1:A:65:CRQ:CB1	1:A:218:GLU:OE1	2.59	0.49
1:A:85:CYS:O	1:A:183:PHE:HE2	1.95	0.49
1:B:60:ILE:HG13	1:B:60:ILE:O	2.12	0.49
1:B:99:GLU:O	1:B:173:MET:CE	2.61	0.49
1:A:134:ILE:HG22	1:A:135:MET:HE3	1.94	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:61:CYS:CB	1:A:216:GLN:NE2	2.76	0.48
1:B:64:ILE:O	1:B:65:CRQ:HG12	2.13	0.48
1:B:198:THR:O	1:B:220:ALA:HA	2.13	0.48
1:A:112:LEU:C	1:A:112:LEU:CD1	2.75	0.48
1:A:152:HIS:O	1:A:156:ALA:HB3	2.13	0.48
1:A:22:VAL:HG13	1:A:125:CYS:HB2	1.94	0.48
1:B:87:PRO:HD2	1:B:88:GLU:OE2	2.14	0.48
1:B:148:HIS:CD2	1:B:196:PHE:HE1	2.32	0.48
1:A:164:GLY:N	2:A:325:HOH:O	2.47	0.48
1:B:200:ILE:O	1:B:218:GLU:HB2	2.13	0.48
1:A:92:ILE:HG12	1:A:181:MET:HG2	1.95	0.48
1:B:149:MET:HE3	1:B:159:GLN:CA	2.43	0.48
1:B:97:ARG:HA	1:B:103:THR:HB	1.95	0.48
1:A:53:LEU:HD22	1:A:57:TRP:CD2	2.49	0.48
1:A:58:LYS:CB	1:A:59:PRO:HD3	2.25	0.47
1:A:112:LEU:HD11	1:A:114:ASP:H	1.79	0.47
1:B:112:LEU:HD23	1:B:113:ASP:N	2.29	0.47
1:A:183:PHE:CE1	1:A:185:GLY:HA3	2.48	0.47
1:A:8:PHE:HZ	1:A:86:PHE:CD2	2.33	0.47
1:B:43:ASN:N	1:B:43:ASN:ND2	2.62	0.47
1:B:57:TRP:O	1:B:60:ILE:HG12	2.15	0.47
1:B:149:MET:CE	1:B:159:GLN:CG	2.92	0.47
1:B:114:ASP:OD2	1:B:115:THR:HG23	2.15	0.47
1:A:13:THR:O	1:A:116:CYS:HA	2.15	0.47
1:A:92:ILE:HD12	1:A:110:TYR:CE2	2.50	0.47
1:B:43:ASN:HD22	1:B:43:ASN:N	2.13	0.47
1:B:149:MET:HE3	1:B:159:GLN:CG	2.44	0.47
1:A:155:ASN:C	1:A:155:ASN:OD1	2.53	0.46
1:B:92:ILE:HD11	1:B:181:MET:HE3	1.97	0.46
1:A:141:ASP:O	1:A:165:PHE:HA	2.15	0.46
1:A:22:VAL:CG2	1:A:27:PHE:CE2	2.98	0.46
1:A:102:GLY:HA3	1:A:128:PHE:CD2	2.50	0.46
1:A:134:ILE:HG22	1:A:135:MET:CE	2.44	0.46
1:A:65:CRQ:HE2	1:A:201:THR:CG2	2.44	0.46
1:B:35:SER:OG	1:B:39:HIS:HB3	2.16	0.46
1:A:108:HIS:HB3	1:A:110:TYR:CE2	2.51	0.46
1:A:209:ASP:C	1:A:211:ARG:H	2.18	0.46
1:B:112:LEU:CD2	1:B:112:LEU:C	2.84	0.46
1:B:12:MET:HG3	1:B:115:THR:O	2.15	0.46
1:B:65:CRQ:C	1:B:68:GLU:OE1	2.64	0.45
1:A:90:LEU:CA	1:A:184:ASN:ND2	2.79	0.45
1:A:61:CYS:HB2	1:A:216:GLN:NE2	2.31	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:227:ARG:NH1	1:B:227:ARG:CB	2.79	0.45
1:A:133:PRO:HB2	1:A:139:LEU:HD23	1.98	0.45
1:A:90:LEU:CD2	1:A:90:LEU:C	2.80	0.45
1:B:64:ILE:C	1:B:65:CRQ:HG12	2.38	0.45
1:B:22:VAL:O	1:B:23:ASN:C	2.53	0.45
1:B:21:GLU:C	1:B:21:GLU:CD	2.76	0.45
1:A:58:LYS:HD3	1:A:203:GLN:HE21	1.81	0.45
1:A:28:THR:H	1:A:50:THR:HG23	1.82	0.45
1:B:68:GLU:HG3	1:B:70:PHE:CE1	2.52	0.44
1:A:149:MET:HE3	1:A:157:VAL:HG11	1.99	0.44
1:A:22:VAL:HG23	1:A:27:PHE:CD2	2.52	0.44
1:B:14:PHE:O	1:B:32:ASP:HA	2.16	0.44
1:B:141:ASP:H	1:B:166:THR:HB	1.81	0.44
1:A:217:ARG:HG2	1:A:218:GLU:N	2.33	0.44
1:B:90:LEU:HD13	1:B:110:TYR:HB2	1.98	0.44
1:A:85:CYS:O	1:A:183:PHE:CE2	2.70	0.44
1:A:21:GLU:C	1:A:21:GLU:CD	2.75	0.44
1:B:13:THR:O	1:B:116:CYS:HA	2.18	0.44
1:A:42:PHE:N	1:A:42:PHE:CD2	2.86	0.44
1:A:173:MET:HE2	1:A:174:MET:N	2.32	0.44
1:A:70:PHE:CE2	1:A:110:TYR:CD1	3.06	0.43
1:A:133:PRO:HA	1:A:138:GLN:HB2	2.01	0.43
1:B:86:PHE:CB	1:B:87:PRO:HA	2.47	0.43
1:B:206:ASP:CB	1:B:213:HIS:ND1	2.82	0.43
1:B:158:ARG:HH21	1:B:178:ASP:CG	2.21	0.43
1:A:86:PHE:CD1	1:A:89:GLY:HA2	2.54	0.43
1:A:89:GLY:C	1:A:184:ASN:HD22	2.22	0.43
1:A:183:PHE:CE1	1:A:185:GLY:CA	3.02	0.43
1:B:90:LEU:HB2	1:B:182:THR:O	2.19	0.43
1:A:10:SER:O	1:A:11:ASP:C	2.57	0.43
1:A:147:THR:CG2	1:A:159:GLN:HG2	2.49	0.43
1:A:31:ALA:HB2	1:A:46:ALA:CA	2.49	0.43
1:A:53:LEU:CD1	1:A:54:PRO:HD2	2.47	0.43
1:B:203:GLN:C	1:B:204:MET:HG2	2.38	0.43
1:A:27:PHE:HA	1:A:50:THR:HG21	2.00	0.43
1:A:58:LYS:CB	1:A:59:PRO:CD	2.93	0.42
1:B:143:LEU:HA	1:B:144:PRO:HD3	1.95	0.42
1:A:92:ILE:HD12	1:A:110:TYR:HE2	1.84	0.42
1:A:62:HIS:CD2	1:A:94:ARG:HE	2.31	0.42
1:A:85:CYS:HA	1:A:183:PHE:CD2	2.54	0.42
1:A:143:LEU:HD11	1:A:166:THR:HG23	2.01	0.42
1:A:57:TRP:CE3	1:A:60:ILE:CD1	2.84	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:65:CRQ:CD2	1:A:199:ILE:HG21	2.49	0.42
1:B:22:VAL:HG21	1:B:55:MET:SD	2.59	0.42
1:B:53:LEU:HD12	1:B:54:PRO:HD2	2.01	0.42
1:A:63:LEU:HD11	1:A:104:MET:CE	2.49	0.42
1:A:106:SER:OG	1:A:108:HIS:HE1	2.03	0.42
1:A:209:ASP:C	1:A:211:ARG:N	2.72	0.42
1:A:209:ASP:HB3	1:A:211:ARG:HG3	1.98	0.42
1:A:22:VAL:HG22	1:A:27:PHE:CE2	2.55	0.42
1:B:155:ASN:CB	2:B:314:HOH:O	2.67	0.42
1:A:211:ARG:HB2	2:A:312:HOH:O	2.20	0.42
1:B:211:ARG:NH1	1:B:211:ARG:HB2	2.35	0.42
1:A:123:VAL:HG13	1:A:123:VAL:O	2.20	0.42
1:A:141:ASP:HB2	1:A:166:THR:HG1	1.75	0.42
1:B:109:THR:O	1:B:119:SER:HA	2.20	0.42
1:A:148:HIS:HB3	1:A:150:PHE:CE2	2.55	0.42
1:B:149:MET:CE	1:B:158:ARG:O	2.68	0.41
1:B:62:HIS:CD2	1:B:94:ARG:NH1	2.68	0.41
1:A:31:ALA:HB2	1:A:46:ALA:HA	2.01	0.41
1:A:197:VAL:HG13	1:A:197:VAL:O	2.20	0.41
1:B:209:ASP:OD2	1:B:211:ARG:NH1	2.52	0.41
1:B:101:ASP:HA	1:B:129:GLN:HG2	2.01	0.41
1:B:74:TYR:CD1	1:B:80:HIS:CD2	3.08	0.41
1:A:57:TRP:C	1:A:203:GLN:HE22	2.24	0.41
1:A:90:LEU:HD13	1:A:110:TYR:HB2	2.02	0.41
1:B:92:ILE:HB	1:B:108:HIS:CB	2.40	0.41
1:A:8:PHE:CZ	1:A:86:PHE:CD2	3.08	0.41
1:B:65:CRQ:HB12	1:B:218:GLU:OE1	2.20	0.41
1:B:68:GLU:N	1:B:69:PRO:HD3	2.11	0.41
1:A:41:ASP:OD2	1:A:217:ARG:NH2	2.49	0.41
1:A:151:PRO:HG3	1:A:192:PRO:O	2.20	0.41
1:B:155:ASN:HB2	2:B:314:HOH:O	2.21	0.40
1:B:149:MET:HA	1:B:149:MET:HE3	1.96	0.40
1:B:209:ASP:OD1	1:B:210:LYS:N	2.55	0.40
1:B:114:ASP:CG	1:B:115:THR:HG23	2.41	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	218/238 (92%)	198 (91%)	19 (9%)	1 (0%)	38	79
1	B	218/238 (92%)	202 (93%)	13 (6%)	3 (1%)	16	52
All	All	436/476 (92%)	400 (92%)	32 (7%)	4 (1%)	25	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	68	GLU
1	B	39	HIS
1	A	114	ASP
1	B	69	PRO

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	193/204 (95%)	176 (91%)	17 (9%)	14	40
1	B	192/204 (94%)	173 (90%)	19 (10%)	11	34
All	All	385/408 (94%)	349 (91%)	36 (9%)	13	36

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	GLU
1	A	43	ASN
1	A	47	VAL
1	A	70	PHE
1	A	84	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	90	LEU
1	A	95	THR
1	A	109	THR
1	A	112	LEU
1	A	117	VAL
1	A	135	MET
1	A	169	ASP
1	A	173	MET
1	A	180	LYS
1	A	187	ARG
1	A	219	VAL
1	A	225	VAL
1	B	7	LEU
1	B	12	MET
1	B	43	ASN
1	B	47	VAL
1	B	68	GLU
1	B	70	PHE
1	B	76	ASP
1	B	84	GLU
1	B	90	LEU
1	B	103	THR
1	B	112	LEU
1	B	117	VAL
1	B	120	ARG
1	B	143	LEU
1	B	145	ASN
1	B	149	MET
1	B	197	VAL
1	B	211	ARG
1	B	219	VAL

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

Mol	Chain	Res	Type
1	A	9	GLN
1	A	43	ASN
1	A	62	HIS
1	A	107	HIS
1	A	108	HIS
1	A	124	ASN
1	A	159	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	176	HIS
1	A	184	ASN
1	A	203	GLN
1	A	216	GLN
1	B	9	GLN
1	B	25	GLN
1	B	43	ASN
1	B	62	HIS
1	B	107	HIS
1	B	108	HIS
1	B	138	GLN
1	B	145	ASN
1	B	159	GLN
1	B	176	HIS
1	B	203	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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	CRQ	A	65	1	25,25,26	10.87	8 (32%)	32,34,36	5.24	14 (43%)
1	CRQ	B	65	1	25,25,26	10.73	8 (32%)	32,34,36	5.14	13 (40%)

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	65	1	-	0/10/32/33	0/2/2/2
1	CRQ	B	65	1	-	0/10/32/33	0/2/2/2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	65	CRQ	O-C	46.41	1.43	1.11
1	B	65	CRQ	O-C	45.73	1.43	1.11
1	A	65	CRQ	CB2-CA2	23.85	1.51	1.35
1	B	65	CRQ	CB2-CA2	23.53	1.51	1.35
1	B	65	CRQ	O2-C2	11.61	1.48	1.23
1	A	65	CRQ	O2-C2	11.39	1.48	1.23
1	B	65	CRQ	CA1-N	7.67	1.45	1.27
1	A	65	CRQ	CA1-N	7.54	1.45	1.27
1	A	65	CRQ	C2-N3	-3.51	1.32	1.39
1	B	65	CRQ	C2-N3	-3.42	1.32	1.39
1	A	65	CRQ	CA3-C	3.40	1.52	1.48
1	B	65	CRQ	CA3-C	3.15	1.52	1.48
1	A	65	CRQ	CG2-CB2	2.38	1.51	1.46
1	B	65	CRQ	CG2-CB2	2.28	1.51	1.46
1	A	65	CRQ	C1-N2	2.25	1.37	1.33
1	B	65	CRQ	C1-N2	2.16	1.37	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	65	CRQ	CG2-CB2-CA2	-15.65	112.12	130.10
1	A	65	CRQ	CA2-C2-N3	15.16	112.08	103.44
1	B	65	CRQ	CA2-C2-N3	14.15	111.50	103.44
1	A	65	CRQ	O2-C2-CA2	-13.35	123.24	130.96
1	A	65	CRQ	CG2-CB2-CA2	-12.77	115.44	130.10
1	B	65	CRQ	O2-C2-CA2	-10.92	124.64	130.96
1	A	65	CRQ	C1-CA1-N	-10.68	109.17	121.83
1	B	65	CRQ	C1-CA1-N	-10.42	109.49	121.83
1	A	65	CRQ	C2-CA2-N2	-9.07	102.08	108.91
1	B	65	CRQ	C2-CA2-N2	-8.74	102.33	108.91
1	A	65	CRQ	CA3-N3-C2	6.76	127.67	123.46
1	B	65	CRQ	CA3-N3-C2	5.32	126.77	123.46
1	B	65	CRQ	CB1-CA1-N	-3.35	110.03	125.36
1	A	65	CRQ	CB1-CA1-N	-3.28	110.32	125.36
1	B	65	CRQ	CB2-CA2-C2	3.03	127.12	122.46
1	B	65	CRQ	N3-C1-N2	-2.97	110.92	113.24
1	B	65	CRQ	CA2-N2-C1	2.75	109.45	104.22

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	CRQ	CA2-N2-C1	2.62	109.21	104.22
1	A	65	CRQ	CA1-C1-N2	2.62	126.78	122.17
1	A	65	CRQ	N3-C1-N2	-2.48	111.31	113.24
1	B	65	CRQ	CA1-C1-N2	2.47	126.53	122.17
1	A	65	CRQ	CB1-CA1-C1	-2.43	111.70	118.06
1	B	65	CRQ	CB1-CA1-C1	-2.26	112.14	118.06
1	B	65	CRQ	CA3-N3-C1	2.24	127.44	124.10
1	A	65	CRQ	C2-N3-C1	-2.16	105.33	109.01
1	A	65	CRQ	CB2-CA2-C2	2.10	125.69	122.46
1	A	65	CRQ	CA1-C1-N3	-2.04	121.92	124.78

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	221/238 (92%)	0.11	12 (5%) 25 30	37, 50, 64, 68	0
1	B	221/238 (92%)	-0.17	3 (1%) 72 80	32, 42, 52, 55	0
All	All	442/476 (92%)	-0.03	15 (3%) 43 51	32, 44, 59, 68	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	208	SER	5.1
1	A	207	THR	4.4
1	A	210	LYS	3.7
1	A	206	ASP	3.2
1	A	9	GLN	3.2
1	A	229	THR	3.1
1	A	114	ASP	3.1
1	B	9	GLN	3.0
1	A	76	ASP	2.9
1	A	188	ALA	2.8
1	A	186	SER	2.5
1	B	113	ASP	2.4
1	B	208	SER	2.4
1	A	204	MET	2.4
1	A	211	ARG	2.1

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

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, 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	CRQ	A	65	24/25	0.23	1.77	67,67,69,69	0
1	CRQ	B	65	24/25	0.20	1.06	52,53,57,57	0

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.