WORLDWIDE PDB

PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Nov 6, 2023 – 12:55 am GMT

PDB ID	:	2WHS
Title	:	Fluorescent Protein mKeima at pH 3.8
Authors	:	Violot, S.; Carpentier, P.; Blanchoin, L.; Bourgeois, D.
Deposited on	:	2009-05-06
Resolution	:	2.10 Å(reported)

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

|		
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
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)	:	2.36

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 2.10 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	Percentile Ranks	Value
Clashscore		8
Ramachandran outliers		0.2%
Sidechain outliers		3.8%
RSRZ outliers		2.3%

Worse   *Better*

▌ Percentile relative to all X-ray structures

▐ Percentile relative to X-ray structures of similar resolution
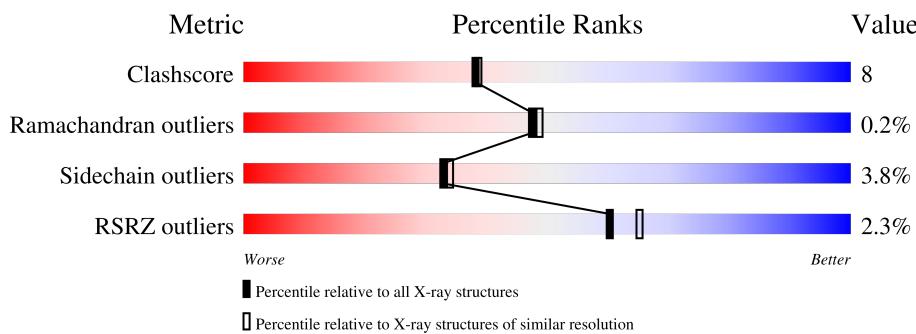
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 <=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	240	% 75% 13% · 10%
1	B	240	3% 81% 12% · 5%
1	C	240	78% 10% · 10%
1	D	240	3% 80% 13% · 5%

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	SO4	C	1221	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7919 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 LARGE STOKES SHIFT FLUORESCENT PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1740	1107	291	331	11			
1	B	227	Total	C	N	O	S	0	0	0
			1813	1153	306	342	12			
1	C	216	Total	C	N	O	S	0	0	0
			1733	1102	290	330	11			
1	D	227	Total	C	N	O	S	0	1	0
			1817	1155	306	344	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	62	CRQ	GLN	chromophore	UNP Q1JV70
A	62	CRQ	TYR	chromophore	UNP Q1JV70
A	62	CRQ	GLY	chromophore	UNP Q1JV70
B	62	CRQ	GLN	chromophore	UNP Q1JV70
B	62	CRQ	TYR	chromophore	UNP Q1JV70
B	62	CRQ	GLY	chromophore	UNP Q1JV70
C	62	CRQ	GLN	chromophore	UNP Q1JV70
C	62	CRQ	TYR	chromophore	UNP Q1JV70
C	62	CRQ	GLY	chromophore	UNP Q1JV70
D	62	CRQ	GLN	chromophore	UNP Q1JV70
D	62	CRQ	TYR	chromophore	UNP Q1JV70
D	62	CRQ	GLY	chromophore	UNP Q1JV70

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).
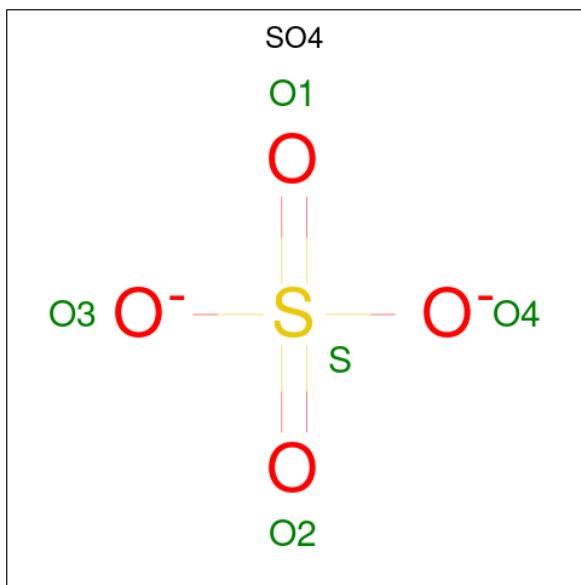
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 5	O 4	S 1	0	0
2	A	1	Total 5	O 4	S 1	0	0
2	A	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	159	Total 159	O 159	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	222	Total 222	O 222	0	0
3	C	166	Total 166	O 166	0	0
3	D	214	Total 214	O 214	0	0

3 Residue-property plots ⓘ

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: LARGE STOKES SHIFT FLUORESCENT PROTEIN
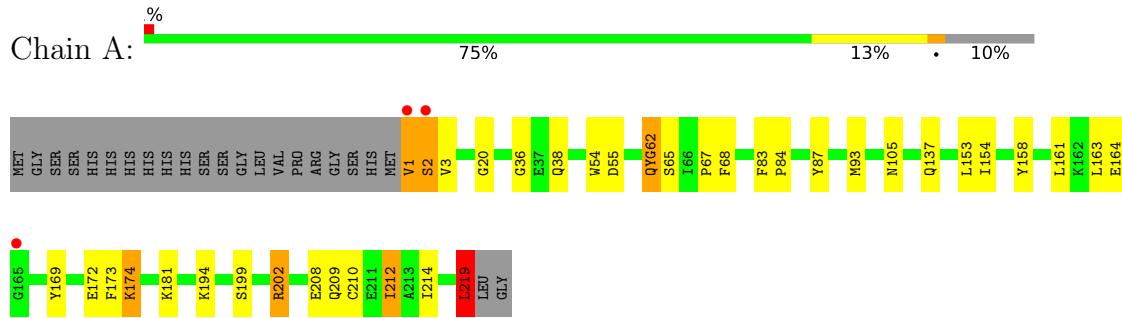
- Molecule 1: LARGE STOKES SHIFT FLUORESCENT PROTEIN
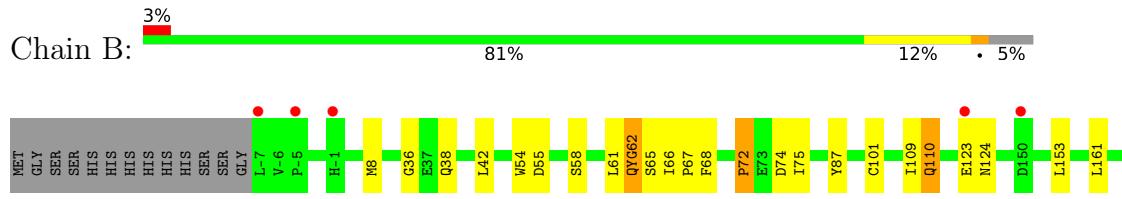
- Molecule 1: LARGE STOKES SHIFT FLUORESCENT PROTEIN
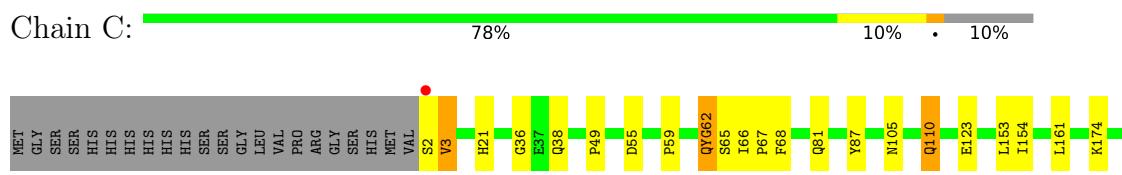
- Molecule 1: LARGE STOKES SHIFT FLUORESCENT PROTEIN
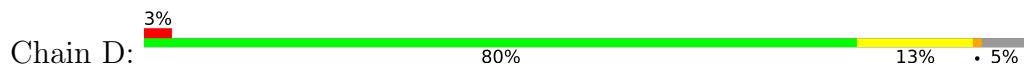
4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	92.57Å 98.10Å 123.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	76.70 – 2.10 46.28 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.2 (76.70-2.10) 99.2 (46.28-2.10)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.73 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.190 , 0.248 0.193 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	20.6	Xtriage
Anisotropy	1.084	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 49.7	EDS
L-test for twinning[2]	$< |L| > = 0.48, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7919	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.90 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.0715e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: SO4, 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	0.79	0/1761	0.78	1/2379 (0.0%)
1	B	0.79	1/1836 (0.1%)	0.80	1/2480 (0.0%)
1	C	0.79	0/1754	0.81	0/2369
1	D	0.78	0/1843	0.82	0/2488
All	All	0.79	1/7194 (0.0%)	0.81	2/9716 (0.0%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	101	CYS	CB-SG	-8.73	1.67	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	219	LEU	CA-CB-CG	5.89	128.85	115.30
1	B	193	ARG	NE-CZ-NH2	5.60	123.10	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	VAL	Peptide

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	1740	0	1673	45	0
1	B	1813	0	1751	26	0
1	C	1733	0	1661	23	0
1	D	1817	0	1756	23	0
2	A	15	0	0	0	0
2	B	15	0	0	0	0
2	C	10	0	0	3	0
2	D	15	0	0	0	0
3	A	159	0	0	2	0
3	B	222	0	0	4	0
3	C	166	0	0	2	0
3	D	214	0	0	6	0
All	All	7919	0	6841	107	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 8.

All (107) 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:212:ILE:HD11	1:A:214:ILE:CD1	1.65	1.25
1:A:212:ILE:HD11	1:A:214:ILE:HD12	1.03	1.02
1:A:212:ILE:CD1	1:A:214:ILE:HD12	1.95	0.95
1:C:38:GLN:HE22	1:C:65:SER:HB3	1.29	0.94
1:D:7:GLN:HB2	3:D:2017:HOH:O	1.70	0.92
1:A:1:VAL:HG12	1:A:1:VAL:O	1.74	0.85
1:B:36:GLY:HA2	1:B:68:PHE:O	1.78	0.83
1:B:123:GLU:O	1:B:124:ASN:OD1	1.96	0.83
1:A:202:ARG:H	1:A:202:ARG:CD	1.95	0.80
1:C:38:GLN:HE22	1:C:65:SER:CB	1.95	0.78
1:A:54:TRP:HE1	1:A:209:GLN:HE21	1.29	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:ILE:CD1	1:A:214:ILE:CD1	2.57	0.75
1:C:154:ILE:CG2	1:C:174:LYS:HE2	2.17	0.74
1:A:202:ARG:H	1:A:202:ARG:HD3	1.51	0.73
1:A:212:ILE:HD11	1:A:214:ILE:HD11	1.67	0.73
1:A:199:SER:HB3	1:A:208:GLU:HB2	1.69	0.73
1:A:38:GLN:HE22	1:A:65:SER:HB3	1.54	0.71
1:C:38:GLN:NE2	1:C:65:SER:HB3	2.03	0.70
1:D:54:TRP:HE1	1:D:209:GLN:HE21	1.41	0.69
1:B:8:MET:HE1	1:B:109:ILE:HD11	1.75	0.68
1:A:219:LEU:HD23	1:B:212:ILE:HG21	1.77	0.67
1:A:38:GLN:NE2	1:A:65:SER:HB3	2.10	0.67
1:B:62:CRQ:HD1	3:B:2172:HOH:O	1.93	0.66
1:B:38:GLN:HE22	1:B:65:SER:HB3	1.61	0.66
1:C:21:HIS:HB2	2:C:1221:SO4:O4	1.96	0.66
1:A:219:LEU:HD21	1:B:212:ILE:HB	1.78	0.65
1:C:81:GLN:HE22	1:C:184:ARG:H	1.45	0.65
1:D:54:TRP:HE1	1:D:209:GLN:NE2	1.93	0.64
1:A:54:TRP:HE1	1:A:209:GLN:NE2	1.95	0.64
1:A:202:ARG:HD3	1:A:202:ARG:N	2.12	0.64
1:B:38:GLN:HE22	1:B:65:SER:CB	2.12	0.62
1:A:219:LEU:C	1:A:219:LEU:HD22	2.20	0.62
1:C:154:ILE:HG23	1:C:174:LYS:HE2	1.82	0.61
1:C:38:GLN:NE2	1:C:65:SER:CB	2.64	0.61
1:C:2:SER:O	1:C:3:VAL:HB	2.01	0.60
1:A:154:ILE:HG23	1:A:174:LYS:HE2	1.84	0.59
1:A:1:VAL:O	1:A:1:VAL:CG1	2.47	0.59
1:A:38:GLN:HE22	1:A:65:SER:CB	2.15	0.59
1:A:154:ILE:CG2	1:A:174:LYS:HE2	2.34	0.58
1:A:55:ASP:HB3	1:A:161:LEU:HD21	1.85	0.58
1:B:74:ASP:OD1	1:B:217:HIS:HE1	1.87	0.58
1:B:72:PRO:HG2	1:B:75:ILE:HD12	1.85	0.57
1:D:74:ASP:OD1	1:D:217:HIS:HE1	1.87	0.57
1:A:36:GLY:HA2	1:A:68:PHE:O	2.04	0.57
1:D:7:GLN:CB	3:D:2017:HOH:O	2.42	0.57
1:C:62:CRQ:HE2	1:C:195:LEU:HB2	1.87	0.56
1:B:54:TRP:HE1	1:B:209:GLN:NE2	2.05	0.55
1:B:200:HIS:HD2	1:B:201:ASN:O	1.89	0.55
1:A:214:ILE:HD13	1:B:219:LEU:HB2	1.88	0.54
1:B:65:SER:N	3:B:2091:HOH:O	2.41	0.54
1:A:38:GLN:NE2	1:A:65:SER:CB	2.71	0.53
1:C:216:ARG:HD3	1:D:216:ARG:NH2	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:110:GLN:NE2	3:C:2088:HOH:O	2.41	0.52
1:B:110:GLN:NE2	3:B:2129:HOH:O	2.42	0.52
1:A:202:ARG:H	1:A:202:ARG:NE	2.07	0.52
1:C:21:HIS:HB2	2:C:1221:SO4:S	2.50	0.52
1:D:194:LYS:HE3	3:D:2189:HOH:O	2.10	0.52
1:C:188:ARG:HG3	3:C:2142:HOH:O	2.10	0.51
1:A:62:CRQ:HD1	3:A:2116:HOH:O	2.10	0.50
1:D:38:GLN:HE22	1:D:65:SER:HB3	1.77	0.50
1:A:93:MET:HE3	1:A:173:PHE:CZ	2.48	0.49
1:A:137:GLN:NE2	1:A:164:GLU:HG2	2.28	0.49
1:D:8:MET:HE1	1:D:109:ILE:HD11	1.95	0.49
1:A:67:PRO:HD3	1:A:87:TYR:OH	2.13	0.49
1:A:194:LYS:HB2	1:B:220:LEU:HG	1.95	0.48
1:B:123:GLU:C	1:B:124:ASN:OD1	2.52	0.48
1:A:219:LEU:CD2	1:B:212:ILE:HG21	2.42	0.48
1:C:81:GLN:HE22	1:C:184:ARG:HB2	1.77	0.48
1:B:67:PRO:HD3	1:B:87:TYR:OH	2.14	0.48
1:D:93:MET:HE3	1:D:173:PHE:CZ	2.50	0.47
1:A:219:LEU:HD23	1:B:212:ILE:CG2	2.43	0.47
1:A:83:PHE:HB3	1:A:84:PRO:HA	1.97	0.47
1:A:210:CYS:SG	3:A:2153:HOH:O	2.53	0.47
1:D:65:SER:N	3:D:2077:HOH:O	2.48	0.46
1:A:219:LEU:CD2	1:B:212:ILE:CG2	2.94	0.46
1:A:20:GLY:O	1:D:-3:GLY:HA2	2.15	0.46
1:B:55:ASP:HB3	1:B:161:LEU:HD21	1.98	0.46
1:A:212:ILE:CD1	1:A:214:ILE:HD11	2.38	0.45
1:A:202:ARG:CD	1:A:202:ARG:N	2.67	0.45
1:C:195:LEU:HD23	1:C:196:ASP:N	2.31	0.45
1:A:1:VAL:C	1:A:2:SER:HG	2.20	0.45
1:D:36:GLY:HA2	1:D:68:PHE:O	2.18	0.44
1:C:66:ILE:HG22	1:C:66:ILE:O	2.17	0.44
1:D:59:PRO:O	1:D:62:CRQ:C2	2.66	0.44
1:B:58:SER:CB	1:B:209:GLN:HE22	2.31	0.43
1:D:39:THR:HG23	3:D:2055:HOH:O	2.19	0.43
1:D:58:SER:OG	1:D:59:PRO:HD3	2.19	0.43
1:C:67:PRO:HD3	1:C:87:TYR:OH	2.19	0.43
1:D:55:ASP:HB3	1:D:161:LEU:HD21	2.01	0.43
1:A:2:SER:OG	1:A:3:VAL:N	2.50	0.42
1:A:1:VAL:O	1:A:2:SER:OG	2.33	0.42
1:A:158:TYR:CE1	1:A:172:GLU:HG3	2.54	0.42
1:C:36:GLY:HA2	1:C:68:PHE:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:PRO:HB3	1:C:203:ASP:HB2	2.02	0.42
1:D:4:ILE:HD12	1:D:80:LYS:HG2	2.01	0.42
1:B:174:LYS:HB3	3:B:2168:HOH:O	2.20	0.41
1:D:10:TYR:OH	1:D:40:VAL:HB	2.20	0.41
1:D:174:LYS:HD3	1:D:174:LYS:HA	1.85	0.41
1:D:58:SER:HA	1:D:209:GLN:HE22	1.85	0.41
1:D:62:CRQ:HD1	3:D:2159:HOH:O	2.21	0.41
1:C:21:HIS:HB2	2:C:1221:SO4:O2	2.20	0.41
1:D:93:MET:CE	1:D:173:PHE:CE1	3.03	0.41
1:A:163:LEU:HD21	1:A:169:TYR:HB2	2.02	0.41
1:B:42:LEU:HG	1:B:61:LEU:HD13	2.03	0.41
1:C:55:ASP:HB3	1:C:161:LEU:HD21	2.02	0.40
1:A:219:LEU:HD21	1:B:212:ILE:CB	2.48	0.40
1:C:62:CRQ:HD2	1:C:195:LEU:HD12	2.04	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	212/240 (88%)	208 (98%)	3 (1%)	1 (0%)	29	26
1	B	222/240 (92%)	219 (99%)	3 (1%)	0	100	100
1	C	211/240 (88%)	205 (97%)	5 (2%)	1 (0%)	29	26
1	D	223/240 (93%)	220 (99%)	3 (1%)	0	100	100
All	All	868/960 (90%)	852 (98%)	14 (2%)	2 (0%)	47	49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	SER

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Mol	Chain	Res	Type
1	C	3	VAL

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	188/207 (91%)	181 (96%)	7 (4%)	34	35
1	B	196/207 (95%)	191 (97%)	5 (3%)	46	50
1	C	187/207 (90%)	180 (96%)	7 (4%)	34	35
1	D	197/207 (95%)	187 (95%)	10 (5%)	24	22
All	All	768/828 (93%)	739 (96%)	29 (4%)	33	34

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	153	LEU
1	A	174	LYS
1	A	181	LYS
1	A	202	ARG
1	A	212	ILE
1	A	219	LEU
1	B	66	ILE
1	B	72	PRO
1	B	110	GLN
1	B	153	LEU
1	B	220	LEU
1	C	59	PRO
1	C	105	ASN
1	C	110	GLN
1	C	123	GLU
1	C	153	LEU
1	C	209	GLN
1	C	216	ARG
1	D	-7	LEU

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Mol	Chain	Res	Type
1	D	1	VAL
1	D	66	ILE
1	D	105	ASN
1	D	110	GLN
1	D	123	GLU
1	D	174	LYS
1	D	199	SER
1	D	202	ARG
1	D	209	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	38	GLN
1	A	112	ASN
1	A	124	ASN
1	A	200	HIS
1	A	209	GLN
1	B	38	GLN
1	B	110	GLN
1	B	112	ASN
1	B	128	ASN
1	B	156	ASN
1	B	200	HIS
1	B	209	GLN
1	B	217	HIS
1	C	38	GLN
1	C	81	GLN
1	C	110	GLN
1	C	112	ASN
1	C	124	ASN
1	C	200	HIS
1	D	38	GLN
1	D	110	GLN
1	D	112	ASN
1	D	156	ASN
1	D	200	HIS
1	D	209	GLN
1	D	217	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	C	62	1	24,25,26	4.01	6 (25%)	27,34,36	5.14	7 (25%)
1	CRQ	B	62	1	24,25,26	3.72	6 (25%)	27,34,36	3.91	7 (25%)
1	CRQ	A	62	1	24,25,26	3.76	6 (25%)	27,34,36	4.44	6 (22%)
1	CRQ	D	62	1	24,25,26	3.88	6 (25%)	27,34,36	4.44	9 (33%)

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	C	62	1	-	1/10/32/33	0/2/2/2
1	CRQ	B	62	1	-	1/10/32/33	0/2/2/2
1	CRQ	A	62	1	-	2/10/32/33	0/2/2/2
1	CRQ	D	62	1	-	1/10/32/33	0/2/2/2

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	62	CRQ	CB2-CA2	17.65	1.49	1.35
1	D	62	CRQ	CB2-CA2	17.05	1.49	1.35
1	A	62	CRQ	CB2-CA2	16.58	1.49	1.35
1	B	62	CRQ	CB2-CA2	15.94	1.48	1.35
1	B	62	CRQ	CA2-C2	-4.91	1.43	1.48
1	D	62	CRQ	CA2-C2	-4.16	1.44	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	62	CRQ	C1-N2	3.93	1.41	1.33
1	B	62	CRQ	OH-CZ	-3.91	1.27	1.37
1	D	62	CRQ	OH-CZ	-3.89	1.27	1.37
1	C	62	CRQ	C2-N3	-3.81	1.30	1.39
1	A	62	CRQ	C2-N3	-3.53	1.31	1.39
1	B	62	CRQ	C2-N3	-3.40	1.31	1.39
1	D	62	CRQ	C1-N2	3.35	1.40	1.33
1	C	62	CRQ	CA2-C2	-3.33	1.45	1.48
1	C	62	CRQ	OH-CZ	-3.21	1.29	1.37
1	A	62	CRQ	C1-N2	3.18	1.40	1.33
1	A	62	CRQ	CA2-C2	-3.08	1.45	1.48
1	B	62	CRQ	O2-C2	3.04	1.29	1.23
1	D	62	CRQ	C2-N3	-2.99	1.32	1.39
1	A	62	CRQ	OH-CZ	-2.95	1.30	1.37
1	B	62	CRQ	C1-N2	2.91	1.39	1.33
1	A	62	CRQ	O2-C2	2.77	1.29	1.23
1	D	62	CRQ	O2-C2	2.76	1.28	1.23
1	C	62	CRQ	O2-C2	2.01	1.27	1.23

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	62	CRQ	CA2-C2-N3	20.87	113.24	103.37
1	A	62	CRQ	CA2-C2-N3	16.84	111.33	103.37
1	D	62	CRQ	O2-C2-CA2	-16.63	121.62	130.96
1	D	62	CRQ	CA2-C2-N3	14.00	109.99	103.37
1	C	62	CRQ	O2-C2-CA2	-13.88	123.17	130.96
1	A	62	CRQ	O2-C2-CA2	-13.64	123.30	130.96
1	B	62	CRQ	O2-C2-CA2	-13.59	123.33	130.96
1	B	62	CRQ	CA2-C2-N3	13.02	109.53	103.37
1	C	62	CRQ	C2-CA2-N2	-5.59	105.02	108.93
1	A	62	CRQ	C2-CA2-N2	-4.59	105.72	108.93
1	C	62	CRQ	CA3-N3-C1	3.66	135.40	128.22
1	B	62	CRQ	O3-C3-CA3	-3.43	116.04	126.39
1	B	62	CRQ	CG1-CB1-CA1	-3.33	103.21	113.53
1	D	62	CRQ	O3-C3-CA3	-3.32	116.38	126.39
1	C	62	CRQ	O3-C3-CA3	-3.02	117.28	126.39
1	D	62	CRQ	CD2-CG2-CD1	2.89	121.91	117.64
1	A	62	CRQ	O3-C3-CA3	-2.77	118.02	126.39
1	C	62	CRQ	CG1-CB1-CA1	-2.70	105.15	113.53
1	A	62	CRQ	CG1-CB1-CA1	-2.66	105.28	113.53
1	C	62	CRQ	CB2-CA2-N2	2.64	132.49	128.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	62	CRQ	CD2-CG2-CD1	2.47	121.29	117.64
1	D	62	CRQ	CG1-CB1-CA1	-2.39	106.13	113.53
1	A	62	CRQ	CG2-CB2-CA2	2.38	132.86	129.94
1	B	62	CRQ	C2-CA2-N2	-2.32	107.31	108.93
1	D	62	CRQ	C2-CA2-N2	-2.31	107.32	108.93
1	D	62	CRQ	CB2-CA2-N2	2.27	131.98	128.83
1	D	62	CRQ	CE1-CD1-CG2	-2.21	118.36	121.25
1	B	62	CRQ	CE2-CD2-CG2	-2.10	118.51	121.25
1	D	62	CRQ	O2-C2-N3	2.03	128.38	124.35

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	62	CRQ	C1-CA1-CB1-CG1
1	D	62	CRQ	C1-CA1-CB1-CG1
1	A	62	CRQ	CA1-CB1-CG1-CD3
1	A	62	CRQ	C3-CA3-N3-C1
1	C	62	CRQ	C3-CA3-N3-C1

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	62	CRQ	2	0
1	B	62	CRQ	1	0
1	A	62	CRQ	1	0
1	D	62	CRQ	2	0

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

11 ligands 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	$\#|Z| > 2$
2	SO4	D	1222	-	4,4,4	0.32	0	6,6,6	0.72	0
2	SO4	B	1224	-	4,4,4	0.15	0	6,6,6	0.16	0
2	SO4	A	1222	-	4,4,4	0.18	0	6,6,6	0.34	0
2	SO4	C	1221	-	4,4,4	0.17	0	6,6,6	0.27	0
2	SO4	D	1224	-	4,4,4	0.17	0	6,6,6	0.16	0
2	SO4	A	1220	-	4,4,4	0.14	0	6,6,6	0.15	0
2	SO4	A	1221	-	4,4,4	0.11	0	6,6,6	0.22	0
2	SO4	D	1223	-	4,4,4	0.13	0	6,6,6	0.16	0
2	SO4	C	1220	-	4,4,4	0.20	0	6,6,6	0.09	0
2	SO4	B	1222	-	4,4,4	0.29	0	6,6,6	0.42	0
2	SO4	B	1223	-	4,4,4	0.18	0	6,6,6	0.21	0

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.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1221	SO4	3	0

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	216/240 (90%)	-0.08	3 (1%)	75	78	13, 21, 32, 39	1 (0%)
1	B	226/240 (94%)	-0.07	8 (3%)	44	50	12, 20, 35, 52	2 (0%)
1	C	215/240 (89%)	-0.17	1 (0%)	91	92	11, 20, 30, 44	1 (0%)
1	D	226/240 (94%)	-0.04	8 (3%)	44	50	14, 21, 37, 49	2 (0%)
All	All	883/960 (91%)	-0.09	20 (2%)	60	65	11, 20, 34, 52	6 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	VAL	5.0
1	D	165	GLY	4.8
1	D	-1	HIS	4.6
1	A	165	GLY	4.4
1	B	165	GLY	4.2
1	B	221	GLY	3.9
1	B	-1	HIS	3.6
1	D	166	GLY	3.4
1	B	202	ARG	3.1
1	C	2	SER	3.1
1	B	123	GLU	3.0
1	D	202	ARG	2.9
1	D	0	MET	2.8
1	A	2	SER	2.7
1	D	221	GLY	2.7
1	D	123	GLU	2.5
1	B	-7	LEU	2.3
1	B	150	ASP	2.2
1	D	-7	LEU	2.1
1	B	-5	PRO	2.1

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	C	62	24/25	0.90	0.16	21,25,32,34	0
1	CRQ	B	62	24/25	0.92	0.15	19,21,29,33	0
1	CRQ	A	62	24/25	0.92	0.15	18,22,29,31	0
1	CRQ	D	62	24/25	0.92	0.14	20,23,32,33	0

6.3 Carbohydrates (i)

There are no monosaccharides 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	SO4	C	1221	5/5	0.81	0.58	93,94,94,95	0
2	SO4	C	1220	5/5	0.82	0.19	88,88,88,89	0
2	SO4	B	1223	5/5	0.87	0.30	90,91,91,91	0
2	SO4	A	1221	5/5	0.88	0.19	86,86,87,87	0
2	SO4	D	1223	5/5	0.88	0.21	70,71,72,73	0
2	SO4	A	1222	5/5	0.89	0.23	80,80,80,81	0
2	SO4	D	1224	5/5	0.92	0.16	90,90,90,90	0
2	SO4	A	1220	5/5	0.94	0.20	82,82,83,83	0
2	SO4	B	1224	5/5	0.96	0.18	79,79,79,79	0
2	SO4	D	1222	5/5	0.97	0.15	36,36,38,41	0
2	SO4	B	1222	5/5	0.99	0.11	34,36,40,41	0

6.5 Other polymers (i)

There are no such residues in this entry.