



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 01:52 PM UTC

PDB ID : 3CFH / pdb_00003cfh
Title : Photoswitchable red fluorescent protein psRFP, off-state
Authors : Kachalova, G.S.; Gundel, S.; Bartunik, H.D.; Wiedenmann, J.
Deposited on : 2008-03-03
Resolution : 1.75 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:

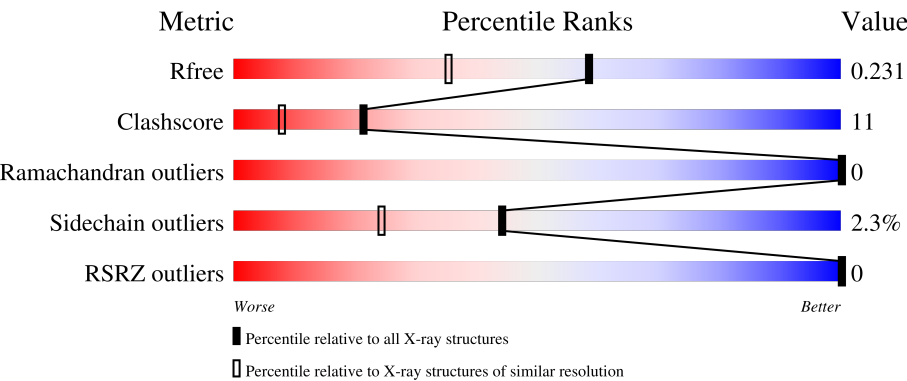
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	L	62	74%19%6%
1	M	62	73%19%6%
1	R	62	68%26%6%
1	S	62	84%10%6%
2	A	167	87%12%

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Mol	Chain	Length	Quality of chain
2	B	167	 87% 11% •
2	G	167	 89% 11% •
2	H	167	 90% 8% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8489 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 photoswitchable fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	58	Total	C	N	O	S	0	7	0
			481	304	77	91	9			
1	M	58	Total	C	N	O	S	0	10	0
			499	317	78	96	8			
1	R	58	Total	C	N	O	S	0	7	0
			495	313	80	93	9			
1	S	58	Total	C	N	O	S	0	5	0
			477	301	77	91	8			

- Molecule 2 is a protein called GFP-like photoswitchable fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	167	Total	C	N	O	S	0	11	0
			1413	906	234	257	16			
2	B	167	Total	C	N	O	S	0	9	0
			1408	900	235	257	16			
2	G	167	Total	C	N	O	S	0	9	0
			1406	899	235	256	16			
2	H	167	Total	C	N	O	S	0	7	0
			1402	894	236	256	16			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	GLY	SER	engineered mutation	PDB 3CFH
B	143	GLY	SER	engineered mutation	PDB 3CFH
G	143	GLY	SER	engineered mutation	PDB 3CFH
H	143	GLY	SER	engineered mutation	PDB 3CFH

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	51	Total 51	O 51	0	0
3	A	201	Total 201	O 201	0	0
3	M	41	Total 41	O 41	0	0
3	B	159	Total 159	O 159	0	0
3	R	59	Total 59	O 59	0	0
3	G	196	Total 196	O 196	0	0
3	S	38	Total 38	O 38	0	0
3	H	163	Total 163	O 163	0	1

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.

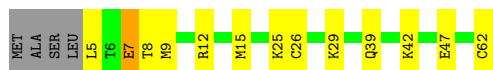
- Molecule 1: GFP-like photoswitchable fluorescent protein

Chain L: 



- Molecule 1: GFP-like photoswitchable fluorescent protein

Chain M: 



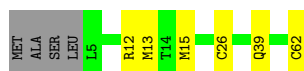
- Molecule 1: GFP-like photoswitchable fluorescent protein

Chain R: 



- Molecule 1: GFP-like photoswitchable fluorescent protein

Chain S: 



- Molecule 2: GFP-like photoswitchable fluorescent protein

Chain A: 

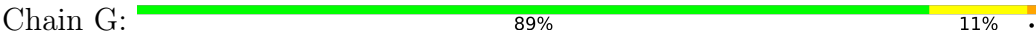


- Molecule 2: GFP-like photoswitchable fluorescent protein

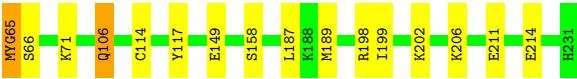
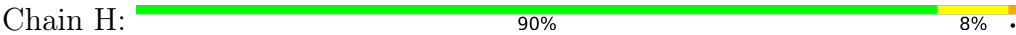
Chain B: 



- Molecule 2: GFP-like photoswitchable fluorescent protein



- Molecule 2: GFP-like photoswitchable fluorescent protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	97.82Å 97.85Å 241.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.27 – 1.75 12.27 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (12.27-1.75) 99.3 (12.27-1.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 1.74Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.202 , 0.230 0.203 , 0.231	Depositor DCC
R_{free} test set	5794 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	31.1	Xtriage
Anisotropy	0.494	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 59.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.469 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	8489	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: NRQ, CME

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	L	0.43	0/512	0.64	0/684
1	M	0.38	0/542	0.61	0/726
1	R	0.41	0/521	0.63	0/696
1	S	0.41	0/498	0.66	0/667
2	A	0.41	0/1412	0.67	0/1900
2	B	0.41	0/1398	0.67	0/1881
2	G	0.41	0/1393	0.66	0/1877
2	H	0.39	0/1382	0.65	0/1859
All	All	0.40	0/7658	0.66	0/10290

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 [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	L	481	0	475	11	0
1	M	499	0	490	26	0
1	R	495	0	493	17	0
1	S	477	0	456	19	0
2	A	1413	0	1377	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1408	0	1361	43	0
2	G	1406	0	1357	30	0
2	H	1402	0	1345	24	0
3	A	201	0	0	4	1
3	B	159	0	0	5	0
3	G	196	0	0	5	3
3	H	163	0	0	2	0
3	L	51	0	0	2	0
3	M	41	0	0	1	0
3	R	59	0	0	2	0
3	S	38	0	0	1	0
All	All	8489	0	7354	163	4

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 11.

All (163) 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:M:15[B]:MET:HG3	2:B:119[B]:VAL:CG2	1.44	1.45
1:S:13[B]:MET:CE	1:S:15[B]:MET:HB2	1.50	1.41
2:A:65[A]:NRQ:C3	2:A:66:SER:N	1.93	1.29
2:G:65[A]:NRQ:C3	2:G:66:SER:N	2.04	1.20
1:M:15[B]:MET:CG	2:B:119[B]:VAL:CG2	2.20	1.18
2:B:65[A]:NRQ:HE2	3:B:390:HOH:O	1.45	1.16
1:S:13[B]:MET:HE2	1:S:15[B]:MET:CB	1.79	1.11
1:S:13[B]:MET:CE	1:S:15[B]:MET:CB	2.30	1.09
2:B:149:GLU:HB2	2:B:189[A]:MET:HE2	1.39	1.03
1:S:13[B]:MET:HE2	1:S:15[B]:MET:HB2	1.04	1.02
1:M:15[B]:MET:HG3	2:B:119[B]:VAL:HG23	1.00	1.00
1:M:62:CYS:O	2:B:65[B]:NRQ:N1	1.97	0.97
2:H:65[A]:NRQ:C3	2:H:66:SER:N	2.32	0.92
1:M:62:CYS:O	2:B:65[A]:NRQ:N1	2.04	0.91
2:A:195[A]:GLU:OE2	3:A:327:HOH:O	1.89	0.90
1:M:7[A]:GLU:CD	1:M:8[A]:THR:HG23	1.95	0.90
2:B:65[A]:NRQ:C3	2:B:66:SER:N	2.33	0.90
1:M:15[B]:MET:SD	2:B:119[B]:VAL:HG21	2.13	0.89
1:M:15[B]:MET:CG	2:B:119[B]:VAL:HG21	2.00	0.89
1:S:13[B]:MET:HE3	1:S:15[B]:MET:HB2	1.55	0.85
1:L:33[B]:ASN:OD1	3:L:110:HOH:O	1.96	0.82
2:B:168:ARG:HH11	2:B:168:ARG:HG3	1.44	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:65[B]:NRQ:OH	2:A:158[B]:SER:OG	2.02	0.78
1:R:33[A]:ASN:OD1	3:R:102:HOH:O	2.01	0.78
1:R:62:CYS:O	2:G:65[B]:NRQ:N1	2.16	0.77
1:S:62:CYS:O	2:H:65[B]:NRQ:N1	2.18	0.77
2:G:65[B]:NRQ:OH	2:G:158[B]:SER:OG	2.03	0.75
2:G:65[A]:NRQ:HA31	2:G:65[A]:NRQ:N1	2.01	0.74
2:H:65[B]:NRQ:OH	2:H:158:SER:HB3	1.89	0.73
1:S:39:GLN:HE22	2:H:66:SER:HB3	1.54	0.72
2:A:65[B]:NRQ:OH	2:A:158[B]:SER:CB	2.37	0.72
1:R:39:GLN:HE22	2:G:66:SER:HB3	1.54	0.72
2:H:202:LYS:HE2	2:H:211:GLU:OE1	1.91	0.70
1:S:62:CYS:O	2:H:65[A]:NRQ:N1	2.24	0.70
2:G:67:LYS:HD2	3:G:321:HOH:O	1.92	0.70
1:M:7[A]:GLU:OE2	1:M:8[A]:THR:HG23	1.92	0.69
2:B:65[B]:NRQ:OH	2:B:158[B]:SER:CB	2.41	0.68
1:R:62:CYS:O	2:G:65[A]:NRQ:N1	2.28	0.67
2:B:65[A]:NRQ:CE2	3:B:390:HOH:O	2.21	0.67
2:G:65[B]:NRQ:HB12	2:G:214:GLU:OE1	1.95	0.65
2:A:120[B]:LYS:NZ	3:A:386:HOH:O	2.17	0.65
1:M:39:GLN:HE22	2:B:66:SER:HB3	1.60	0.65
2:G:65[B]:NRQ:OH	2:G:158[B]:SER:CB	2.44	0.65
2:B:65[B]:NRQ:N1	2:B:65[B]:NRQ:CA3	2.59	0.64
2:A:67:LYS:HD3	3:A:352:HOH:O	1.97	0.64
1:R:33[A]:ASN:ND2	1:R:36[A]:GLU:CD	2.56	0.64
1:M:5:LEU:HA	1:M:9:MET:HE2	1.80	0.63
1:M:15[B]:MET:CB	2:B:119[B]:VAL:CG2	2.77	0.63
2:G:197:HIS:HE1	3:G:321:HOH:O	1.80	0.63
1:M:15[B]:MET:CB	2:B:119[B]:VAL:HG22	2.28	0.63
2:H:65[B]:NRQ:N1	2:H:65[B]:NRQ:CA3	2.61	0.63
1:M:15[B]:MET:CG	2:B:119[B]:VAL:HG23	1.97	0.62
2:A:65[A]:NRQ:HA31	2:A:65[A]:NRQ:N1	2.15	0.62
1:M:15[B]:MET:HG3	2:B:119[B]:VAL:HG21	1.58	0.62
2:H:65[B]:NRQ:OH	2:H:158:SER:CB	2.48	0.62
2:B:65[B]:NRQ:OH	2:B:158[B]:SER:HB3	1.99	0.62
2:B:65[B]:NRQ:OH	2:B:158[B]:SER:OG	2.18	0.61
2:B:65[B]:NRQ:N1	2:B:65[B]:NRQ:HA31	2.16	0.60
2:B:65[B]:NRQ:N1	2:B:65[B]:NRQ:C3	2.65	0.60
2:B:168:ARG:HG3	2:B:168:ARG:NH1	2.11	0.60
2:A:197:HIS:HE1	3:A:352:HOH:O	1.84	0.60
1:R:25[B]:LYS:HB2	1:R:47:GLU:HB2	1.85	0.58
2:G:65[A]:NRQ:HE2	3:G:425:HOH:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:66:SER:OG	2:B:117:TYR:OH	2.22	0.58
2:H:66:SER:HG	2:H:117:TYR:HH	1.48	0.58
1:S:39:GLN:NE2	2:H:66:SER:HB3	2.18	0.57
1:M:12[A]:ARG:NH1	3:M:88:HOH:O	2.34	0.57
2:H:65[B]:NRQ:N1	2:H:65[B]:NRQ:HA31	2.19	0.57
1:M:7[A]:GLU:OE2	1:M:8[A]:THR:CG2	2.52	0.57
2:H:66:SER:N	3:H:358:HOH:O	2.36	0.57
1:R:39:GLN:NE2	2:G:66:SER:HB3	2.18	0.57
2:A:71[B]:LYS:HB3	2:A:217[B]:VAL:HG12	1.87	0.56
1:L:12[A]:ARG:HB3	2:A:114:CME:HE3	1.88	0.56
2:G:65[B]:NRQ:OH	2:G:158[B]:SER:HB3	2.06	0.56
1:R:25[A]:LYS:HB2	1:R:47:GLU:HB2	1.87	0.55
2:A:65[A]:NRQ:C3	2:A:66:SER:CA	2.83	0.55
2:H:65[B]:NRQ:N1	2:H:65[B]:NRQ:C3	2.70	0.55
1:S:12:ARG:NH1	3:S:82:HOH:O	2.31	0.55
2:G:66:SER:OG	2:G:117:TYR:OH	2.19	0.55
1:M:7[A]:GLU:OE1	1:M:8[A]:THR:HG23	2.06	0.55
2:B:65[A]:NRQ:N1	2:B:65[A]:NRQ:HA31	2.22	0.55
1:L:39:GLN:HE22	2:A:66:SER:HB3	1.73	0.55
2:G:65[A]:NRQ:HB12	2:G:214:GLU:OE1	2.06	0.55
2:G:181:LYS:NZ	3:G:278:HOH:O	2.07	0.54
1:M:39:GLN:NE2	2:B:66:SER:HB3	2.22	0.54
2:H:65[A]:NRQ:HE2	3:H:395:HOH:O	2.06	0.54
2:G:65[B]:NRQ:N1	2:G:65[B]:NRQ:HA31	2.22	0.54
1:M:15[B]:MET:HA	2:B:119[B]:VAL:HG22	1.90	0.53
2:B:66:SER:N	3:B:285:HOH:O	2.41	0.53
2:A:67:LYS:NZ	2:A:195[B]:GLU:OE2	2.42	0.52
2:A:67:LYS:HE3	2:A:178:TYR:CE1	2.44	0.52
1:R:5:LEU:HD22	1:R:9[A]:MET:HE3	1.91	0.52
2:G:65[A]:NRQ:N1	2:G:65[A]:NRQ:CA3	2.73	0.51
2:H:65[A]:NRQ:N1	2:H:65[A]:NRQ:HA31	2.26	0.51
2:G:65[B]:NRQ:N1	2:G:65[B]:NRQ:CA3	2.74	0.51
1:L:13[B]:MET:SD	1:L:62[B]:CYS:OXT	2.69	0.50
1:L:16:GLU:HG2	1:L:25:LYS:HG2	1.94	0.50
2:B:77:PRO:HG2	2:B:188:LYS:HD2	1.94	0.50
2:G:65[A]:NRQ:CA3	2:G:66:SER:N	2.60	0.50
2:G:197:HIS:ND1	2:G:214:GLU:OE2	2.39	0.50
1:M:15[B]:MET:CG	2:B:119[B]:VAL:HG22	2.30	0.49
1:L:49:GLY:HA2	1:L:50:PRO:C	2.37	0.49
2:B:82:GLN:HB3	2:B:187:LEU:HD12	1.95	0.49
1:M:25:LYS:HB2	1:M:47:GLU:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:R:74:HOH:O	2:G:209:HIS:HD2	1.96	0.49
2:A:65[B]:NRQ:N1	2:A:65[B]:NRQ:HA31	2.27	0.49
2:A:65[B]:NRQ:OH	2:A:158[B]:SER:HB3	2.12	0.49
2:G:71:LYS:HB3	2:G:217[B]:VAL:HG12	1.94	0.49
1:S:15[A]:MET:HB3	1:S:26:CYS:HB2	1.94	0.48
1:M:42:LYS:HE3	3:B:382:HOH:O	2.12	0.48
1:S:13[B]:MET:HE1	1:S:15[B]:MET:CB	2.36	0.48
1:M:15[B]:MET:CA	2:B:119[B]:VAL:HG22	2.44	0.48
2:G:67:LYS:NZ	2:G:195:GLU:OE1	2.35	0.48
2:B:150:VAL:HG13	2:B:155[A]:CYS:SG	2.53	0.48
1:S:13[B]:MET:CE	1:S:15[B]:MET:HB3	2.37	0.48
3:L:70:HOH:O	2:A:209:HIS:HD2	1.97	0.47
2:G:65[A]:NRQ:C3	2:G:66:SER:CA	2.90	0.47
2:A:65[A]:NRQ:N1	2:A:65[A]:NRQ:CA3	2.78	0.47
1:S:13[B]:MET:HE3	1:S:15[B]:MET:SD	2.55	0.47
1:R:15[A]:MET:HE3	1:R:26:CYS:SG	2.55	0.47
1:L:12[B]:ARG:HB3	2:A:114:CME:HE3	1.97	0.46
1:L:58:LEU:HD22	2:A:121[A]:ILE:HD13	1.97	0.46
1:S:13[B]:MET:CE	1:S:15[B]:MET:SD	3.03	0.46
1:M:15[A]:MET:HB3	1:M:26:CYS:HB2	1.97	0.46
1:M:39:GLN:OE1	2:B:66:SER:HB3	2.15	0.46
2:H:202:LYS:HG3	2:H:211:GLU:HB2	1.97	0.46
2:H:66:SER:OG	2:H:117:TYR:OH	2.24	0.45
1:R:58:LEU:HD22	2:G:121[B]:ILE:HD13	1.97	0.45
1:S:13[A]:MET:SD	1:S:15[A]:MET:HB2	2.56	0.45
2:B:197:HIS:CD2	3:B:377:HOH:O	2.68	0.45
2:B:77:PRO:CG	2:B:188:LYS:HD2	2.47	0.45
2:G:71:LYS:HE2	2:G:217[B]:VAL:HG12	1.99	0.45
2:A:228:LEU:HD21	2:B:202:LYS:HE3	1.99	0.45
2:H:198:ARG:O	2:H:214:GLU:HA	2.16	0.45
2:B:195[B]:GLU:OE2	2:B:197:HIS:NE2	2.44	0.45
1:R:45:VAL:HG21	1:R:51:LEU:HG	1.99	0.44
2:A:67:LYS:HE3	2:A:178:TYR:CZ	2.52	0.44
2:G:182:LYS:NZ	3:G:343:HOH:O	2.50	0.44
2:H:106:GLN:C	2:H:106:GLN:HE21	2.26	0.44
2:H:114:CME:HE2	2:H:114:CME:HB3	1.89	0.44
1:S:39:GLN:OE1	2:H:66:SER:CB	2.66	0.44
1:R:5:LEU:HD22	1:R:9[B]:MET:HE2	2.00	0.44
2:H:65[A]:NRQ:HD2	2:H:199:ILE:HD11	2.00	0.44
2:A:65[B]:NRQ:N1	2:A:65[B]:NRQ:CA3	2.81	0.43
1:R:33[A]:ASN:ND2	1:R:36[A]:GLU:OE1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:149:GLU:CB	2:B:189[A]:MET:HE2	2.27	0.43
1:L:25:LYS:HB2	1:L:47:GLU:HB2	2.01	0.43
2:A:149:GLU:HG3	2:A:189[A]:MET:HE2	2.01	0.43
1:R:49:GLY:HA2	1:R:50:PRO:C	2.43	0.42
1:R:25[C]:LYS:HB2	1:R:47:GLU:HB2	2.02	0.42
1:R:12[A]:ARG:HG2	2:G:114:CME:SG	2.60	0.41
2:B:65[A]:NRQ:N1	2:B:65[A]:NRQ:CA3	2.83	0.41
2:G:149:GLU:HG3	2:G:189[A]:MET:HE2	2.02	0.41
1:S:39:GLN:OE1	2:H:66:SER:HB3	2.21	0.41
1:L:39:GLN:NE2	2:A:66:SER:HB3	2.36	0.41
2:A:65[B]:NRQ:HB11	2:A:212:GLN:NE2	2.36	0.41
1:L:41:MET:SD	2:A:65[B]:NRQ:HE2A	2.61	0.41
2:B:65[A]:NRQ:C3	2:B:66:SER:CA	2.97	0.41
1:S:15[B]:MET:HB3	1:S:26:CYS:HB2	2.01	0.41
2:H:149:GLU:HG3	2:H:189[B]:MET:HE3	2.04	0.40
2:B:198:ARG:O	2:B:214:GLU:HA	2.21	0.40

All (4) 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)
3:G:426:HOH:O	3:G:426:HOH:O[3_555]	1.83	0.37
3:G:232:HOH:O	3:G:233:HOH:O[3_555]	1.98	0.22
3:G:233:HOH:O	3:G:426:HOH:O[3_555]	1.99	0.21
3:A:313:HOH:O	3:A:332:HOH:O[4_555]	2.15	0.05

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	L	62/62 (100%)	61 (98%)	1 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	M	66/62 (106%)	65 (98%)	1 (2%)	0	100	100
1	R	64/62 (103%)	63 (98%)	1 (2%)	0	100	100
1	S	61/62 (98%)	60 (98%)	1 (2%)	0	100	100
2	A	172/167 (103%)	171 (99%)	1 (1%)	0	100	100
2	B	170/167 (102%)	168 (99%)	2 (1%)	0	100	100
2	G	170/167 (102%)	167 (98%)	3 (2%)	0	100	100
2	H	168/167 (101%)	164 (98%)	4 (2%)	0	100	100
All	All	933/916 (102%)	919 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	56/52 (108%)	56 (100%)	0	100	100
1	M	59/52 (114%)	56 (95%)	3 (5%)	21	5
1	R	57/52 (110%)	57 (100%)	0	100	100
1	S	54/52 (104%)	54 (100%)	0	100	100
2	A	147/137 (107%)	143 (97%)	4 (3%)	39	19
2	B	145/137 (106%)	139 (96%)	6 (4%)	27	8
2	G	145/137 (106%)	143 (99%)	2 (1%)	59	44
2	H	143/137 (104%)	139 (97%)	4 (3%)	38	18
All	All	806/756 (107%)	787 (98%)	19 (2%)	44	23

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	106	GLN
2	A	119	VAL

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Mol	Chain	Res	Type
2	A	187	LEU
2	A	222	ASP
1	M	7[A]	GLU
1	M	7[B]	GLU
1	M	29	LYS
2	B	106	GLN
2	B	120[A]	LYS
2	B	120[B]	LYS
2	B	150	VAL
2	B	187	LEU
2	B	231	HIS
2	G	106	GLN
2	G	119	VAL
2	H	71	LYS
2	H	106	GLN
2	H	187	LEU
2	H	206	LYS

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

Mol	Chain	Res	Type
2	A	105	HIS
2	A	106	GLN
2	A	175	HIS
2	B	105	HIS
2	B	106	GLN
1	R	22	HIS
2	G	105	HIS
2	G	106	GLN
2	G	175	HIS
2	H	106	GLN
2	H	175	HIS
2	H	209	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

16 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
2	CME	A	114	2	8,9,10	0.90	0	6,9,11	0.54	0
2	CME	B	114	2	8,9,10	0.85	0	6,9,11	0.76	0
2	CME	G	221	2	8,9,10	0.83	0	6,9,11	0.69	0
2	NRQ	A	65[B]	2	24,24,25	2.58	7 (29%)	24,32,34	5.21	12 (50%)
2	CME	H	114	2	8,9,10	0.87	0	6,9,11	0.65	0
2	NRQ	B	65[B]	2	24,24,25	2.60	7 (29%)	24,32,34	4.98	11 (45%)
2	NRQ	G	65[A]	-	24,24,25	2.82	7 (29%)	24,32,34	4.44	9 (37%)
2	CME	G	114	2	8,9,10	0.89	0	6,9,11	0.52	0
2	NRQ	H	65[A]	-	24,24,25	2.83	7 (29%)	24,32,34	4.66	8 (33%)
2	CME	H	221	2	8,9,10	0.82	0	6,9,11	0.60	0
2	NRQ	B	65[A]	-	24,24,25	2.82	6 (25%)	24,32,34	4.76	9 (37%)
2	NRQ	A	65[A]	-	24,24,25	2.84	7 (29%)	24,32,34	4.57	8 (33%)
2	NRQ	G	65[B]	2	24,24,25	2.68	7 (29%)	24,32,34	4.94	12 (50%)
2	CME	A	221	2	8,9,10	0.81	0	6,9,11	0.67	0
2	CME	B	221	2	8,9,10	0.83	0	6,9,11	0.56	0
2	NRQ	H	65[B]	2	24,24,25	2.59	6 (25%)	24,32,34	4.84	11 (45%)

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
2	CME	A	114	2	-	2/5/8/10	-
2	CME	B	114	2	-	2/5/8/10	-
2	CME	G	221	2	-	1/5/8/10	-
2	NRQ	A	65[B]	2	-	4/9/31/32	0/2/2/2
2	CME	H	114	2	-	2/5/8/10	-
2	NRQ	B	65[B]	2	-	4/9/31/32	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NRQ	G	65[A]	-	-	1/9/31/32	0/2/2/2
2	CME	G	114	2	-	1/5/8/10	-
2	NRQ	H	65[A]	-	-	2/9/31/32	0/2/2/2
2	CME	H	221	2	-	1/5/8/10	-
2	NRQ	B	65[A]	-	-	1/9/31/32	0/2/2/2
2	NRQ	A	65[A]	-	-	1/9/31/32	0/2/2/2
2	NRQ	G	65[B]	2	-	3/9/31/32	0/2/2/2
2	CME	A	221	2	-	0/5/8/10	-
2	CME	B	221	2	-	1/5/8/10	-
2	NRQ	H	65[B]	2	-	4/9/31/32	0/2/2/2

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	65[A]	NRQ	CB2-CA2	9.43	1.44	1.35
2	G	65[A]	NRQ	CB2-CA2	9.34	1.44	1.35
2	B	65[A]	NRQ	CB2-CA2	9.26	1.44	1.35
2	A	65[A]	NRQ	CB2-CA2	9.16	1.44	1.35
2	G	65[B]	NRQ	CB2-CA2	8.69	1.43	1.35
2	B	65[B]	NRQ	CB2-CA2	7.99	1.42	1.35
2	H	65[B]	NRQ	CB2-CA2	7.72	1.42	1.35
2	A	65[B]	NRQ	CB2-CA2	7.54	1.42	1.35
2	A	65[A]	NRQ	CA2-C2	-5.83	1.42	1.48
2	B	65[A]	NRQ	CA2-C2	-5.61	1.42	1.48
2	H	65[A]	NRQ	CA2-C2	-5.44	1.42	1.48
2	A	65[B]	NRQ	OH-CZ	-5.36	1.24	1.37
2	G	65[A]	NRQ	CA2-C2	-5.31	1.42	1.48
2	B	65[B]	NRQ	OH-CZ	-5.28	1.25	1.37
2	G	65[B]	NRQ	OH-CZ	-5.28	1.25	1.37
2	H	65[B]	NRQ	OH-CZ	-5.14	1.25	1.37
2	A	65[A]	NRQ	OH-CZ	-5.12	1.25	1.37
2	H	65[B]	NRQ	CA2-C2	-5.07	1.43	1.48
2	B	65[A]	NRQ	OH-CZ	-5.00	1.25	1.37
2	H	65[A]	NRQ	OH-CZ	-4.99	1.25	1.37
2	G	65[A]	NRQ	OH-CZ	-4.98	1.25	1.37
2	A	65[B]	NRQ	CA2-C2	-4.87	1.43	1.48
2	B	65[B]	NRQ	CA2-C2	-4.55	1.43	1.48
2	H	65[B]	NRQ	C2-N3	-4.32	1.30	1.40
2	G	65[B]	NRQ	C2-N3	-4.22	1.30	1.40
2	B	65[B]	NRQ	C2-N3	-4.21	1.30	1.40
2	G	65[B]	NRQ	CA2-C2	-4.10	1.44	1.48
2	A	65[B]	NRQ	C2-N3	-3.96	1.30	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	65[A]	NRQ	C2-N3	-3.83	1.31	1.40
2	A	65[A]	NRQ	C2-N3	-3.76	1.31	1.40
2	B	65[A]	NRQ	C2-N3	-3.68	1.31	1.40
2	G	65[B]	NRQ	C1-CA1	-3.66	1.42	1.48
2	H	65[A]	NRQ	C2-N3	-3.65	1.31	1.40
2	A	65[B]	NRQ	C1-CA1	-3.56	1.42	1.48
2	B	65[A]	NRQ	C1-N2	3.43	1.40	1.33
2	H	65[A]	NRQ	C1-N2	3.43	1.40	1.33
2	G	65[A]	NRQ	C1-N2	3.39	1.40	1.33
2	A	65[A]	NRQ	C1-N2	3.38	1.40	1.33
2	B	65[B]	NRQ	C1-CA1	-3.33	1.43	1.48
2	H	65[B]	NRQ	C1-CA1	-3.06	1.43	1.48
2	A	65[B]	NRQ	CA2-N2	-2.76	1.32	1.38
2	G	65[B]	NRQ	CA2-N2	-2.69	1.32	1.38
2	A	65[A]	NRQ	O2-C2	2.51	1.28	1.23
2	B	65[A]	NRQ	O2-C2	2.47	1.28	1.23
2	G	65[A]	NRQ	O2-C2	2.43	1.28	1.23
2	H	65[A]	NRQ	O2-C2	2.37	1.27	1.23
2	B	65[B]	NRQ	CA2-N2	-2.28	1.33	1.38
2	H	65[B]	NRQ	CA2-N2	-2.27	1.33	1.38
2	A	65[A]	NRQ	C1-CA1	-2.21	1.44	1.48
2	A	65[B]	NRQ	O2-C2	2.19	1.27	1.23
2	G	65[A]	NRQ	C1-CA1	-2.09	1.45	1.48
2	B	65[B]	NRQ	O2-C2	2.06	1.27	1.23
2	H	65[A]	NRQ	CG2-CB2	2.03	1.50	1.46
2	G	65[B]	NRQ	O2-C2	2.03	1.27	1.23

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	65[B]	NRQ	CB2-CA2-C2	15.82	141.53	122.36
2	G	65[B]	NRQ	CB2-CA2-C2	15.32	140.92	122.36
2	B	65[B]	NRQ	CB2-CA2-C2	15.22	140.80	122.36
2	H	65[B]	NRQ	CB2-CA2-C2	14.56	140.01	122.36
2	B	65[A]	NRQ	O2-C2-CA2	-14.09	122.03	131.02
2	H	65[A]	NRQ	O2-C2-CA2	-13.62	122.33	131.02
2	A	65[A]	NRQ	O2-C2-CA2	-13.51	122.40	131.02
2	G	65[A]	NRQ	O2-C2-CA2	-12.88	122.80	131.02
2	A	65[B]	NRQ	CB2-CA2-N2	-11.69	112.88	128.76
2	H	65[A]	NRQ	CA2-C2-N3	11.61	113.25	103.50
2	B	65[A]	NRQ	CA2-C2-N3	11.53	113.19	103.50
2	A	65[A]	NRQ	CA2-C2-N3	11.00	112.73	103.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	65[B]	NRQ	CB2-CA2-N2	-10.96	113.88	128.76
2	G	65[A]	NRQ	CA2-C2-N3	10.95	112.69	103.50
2	B	65[B]	NRQ	CB2-CA2-N2	-10.93	113.92	128.76
2	A	65[A]	NRQ	CG2-CB2-CA2	10.21	142.01	129.87
2	H	65[B]	NRQ	CB2-CA2-N2	-10.20	114.92	128.76
2	B	65[A]	NRQ	CG2-CB2-CA2	10.19	141.98	129.87
2	H	65[A]	NRQ	CG2-CB2-CA2	9.81	141.52	129.87
2	G	65[A]	NRQ	CG2-CB2-CA2	9.70	141.40	129.87
2	H	65[B]	NRQ	CA2-C2-N3	8.45	110.59	103.50
2	G	65[B]	NRQ	CA2-C2-N3	8.41	110.56	103.50
2	A	65[B]	NRQ	O2-C2-CA2	-8.07	125.87	131.02
2	A	65[B]	NRQ	CA2-C2-N3	7.97	110.20	103.50
2	B	65[B]	NRQ	CA2-C2-N3	7.89	110.12	103.50
2	H	65[B]	NRQ	O2-C2-CA2	-7.50	126.23	131.02
2	B	65[B]	NRQ	O2-C2-CA2	-7.24	126.40	131.02
2	A	65[B]	NRQ	CG2-CB2-CA2	7.05	138.25	129.87
2	G	65[B]	NRQ	O2-C2-CA2	-6.88	126.63	131.02
2	H	65[A]	NRQ	C3-CA3-N3	6.38	126.94	112.43
2	B	65[A]	NRQ	C3-CA3-N3	6.36	126.91	112.43
2	A	65[A]	NRQ	C3-CA3-N3	6.17	126.47	112.43
2	G	65[A]	NRQ	C3-CA3-N3	6.06	126.23	112.43
2	H	65[A]	NRQ	C2-CA2-N2	-5.98	104.66	108.95
2	B	65[A]	NRQ	C2-CA2-N2	-5.70	104.87	108.95
2	G	65[B]	NRQ	CG2-CB2-CA2	5.44	136.33	129.87
2	H	65[B]	NRQ	C2-CA2-N2	-5.42	105.07	108.95
2	G	65[B]	NRQ	C2-CA2-N2	-5.27	105.18	108.95
2	G	65[A]	NRQ	C2-CA2-N2	-5.16	105.25	108.95
2	B	65[B]	NRQ	C2-CA2-N2	-5.13	105.27	108.95
2	A	65[A]	NRQ	C2-CA2-N2	-4.99	105.37	108.95
2	A	65[B]	NRQ	C2-CA2-N2	-4.74	105.55	108.95
2	B	65[B]	NRQ	CA3-N3-C2	4.48	133.52	123.67
2	B	65[B]	NRQ	CA3-N3-C1	-4.36	119.78	128.39
2	B	65[B]	NRQ	CG2-CB2-CA2	4.31	134.98	129.87
2	B	65[B]	NRQ	O3-C3-CA3	-4.27	106.27	125.47
2	H	65[B]	NRQ	O3-C3-CA3	-4.12	106.91	125.47
2	H	65[B]	NRQ	CA3-N3-C2	4.05	132.57	123.67
2	H	65[B]	NRQ	CA3-N3-C1	-3.82	120.85	128.39
2	A	65[B]	NRQ	CA3-N3-C2	3.73	131.87	123.67
2	H	65[B]	NRQ	CG2-CB2-CA2	3.69	134.26	129.87
2	G	65[B]	NRQ	CA3-N3-C2	3.50	131.37	123.67
2	A	65[B]	NRQ	CA3-N3-C1	-3.21	122.06	128.39
2	B	65[B]	NRQ	CD1-CG2-CD2	3.12	122.28	117.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	65[A]	NRQ	CB2-CA2-N2	3.05	132.90	128.76
2	A	65[B]	NRQ	O3-C3-CA3	-3.03	111.83	125.47
2	A	65[B]	NRQ	CD1-CG2-CD2	3.01	122.11	117.65
2	G	65[B]	NRQ	CA3-N3-C1	-2.87	122.73	128.39
2	G	65[B]	NRQ	O3-C3-CA3	-2.86	112.59	125.47
2	H	65[B]	NRQ	CD1-CG2-CD2	2.83	121.86	117.65
2	A	65[A]	NRQ	CB2-CA2-N2	2.83	132.60	128.76
2	H	65[A]	NRQ	CB2-CA2-N2	2.79	132.55	128.76
2	G	65[B]	NRQ	CD1-CG2-CD2	2.74	121.72	117.65
2	G	65[A]	NRQ	CB2-CA2-N2	2.42	132.05	128.76
2	A	65[A]	NRQ	CD1-CG2-CD2	2.31	121.08	117.65
2	A	65[B]	NRQ	CE2-CD2-CG2	-2.30	118.25	121.22
2	B	65[A]	NRQ	CD1-CG2-CD2	2.29	121.04	117.65
2	B	65[A]	NRQ	CE2-CD2-CG2	-2.21	118.37	121.22
2	G	65[A]	NRQ	CD1-CG2-CD2	2.19	120.91	117.65
2	H	65[B]	NRQ	N3-C1-N2	-2.13	110.09	112.62
2	B	65[B]	NRQ	CE2-CD2-CG2	-2.12	118.48	121.22
2	A	65[B]	NRQ	CA2-N2-C1	2.11	107.98	104.09
2	H	65[A]	NRQ	N3-C1-N2	-2.11	110.12	112.62
2	G	65[A]	NRQ	CE2-CD2-CG2	-2.09	118.52	121.22
2	B	65[A]	NRQ	N3-C1-N2	-2.08	110.15	112.62
2	A	65[A]	NRQ	N3-C1-N2	-2.07	110.16	112.62
2	G	65[B]	NRQ	C3-CA3-N3	2.05	117.09	112.43
2	G	65[B]	NRQ	CA2-N2-C1	2.03	107.83	104.09
2	H	65[A]	NRQ	CD1-CG2-CD2	2.01	120.64	117.65
2	G	65[A]	NRQ	CA3-N3-C2	2.01	128.09	123.67

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	65[A]	NRQ	C1-CA1-CB1-CG1
2	A	65[B]	NRQ	N2-CA2-CB2-CG2
2	A	65[B]	NRQ	C2-CA2-CB2-CG2
2	A	114	CME	SD-CE-CZ-OH
2	B	65[B]	NRQ	N2-CA2-CB2-CG2
2	B	65[B]	NRQ	C2-CA2-CB2-CG2
2	G	65[A]	NRQ	C1-CA1-CB1-CG1
2	G	65[B]	NRQ	C1-CA1-CB1-CG1
2	G	65[B]	NRQ	N2-CA2-CB2-CG2
2	G	65[B]	NRQ	C2-CA2-CB2-CG2
2	H	65[B]	NRQ	C2-CA2-CB2-CG2

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Mol	Chain	Res	Type	Atoms
2	H	114	CME	CE-SD-SG-CB
2	H	65[B]	NRQ	N2-CA2-CB2-CG2
2	A	114	CME	CE-SD-SG-CB
2	B	114	CME	CE-SD-SG-CB
2	H	65[B]	NRQ	CB1-CG1-SD-CE
2	B	65[A]	NRQ	CB1-CG1-SD-CE
2	H	65[A]	NRQ	CB1-CG1-SD-CE
2	A	65[B]	NRQ	CB1-CG1-SD-CE
2	B	65[B]	NRQ	CB1-CG1-SD-CE
2	B	114	CME	SD-CE-CZ-OH
2	G	114	CME	CE-SD-SG-CB
2	B	221	CME	CZ-CE-SD-SG
2	A	65[B]	NRQ	C1-CA1-CB1-CG1
2	H	65[A]	NRQ	C1-CA1-CB1-CG1
2	H	65[B]	NRQ	C1-CA1-CB1-CG1
2	B	65[B]	NRQ	C3-CA3-N3-C1
2	G	221	CME	CZ-CE-SD-SG
2	H	114	CME	CZ-CE-SD-SG
2	H	221	CME	CZ-CE-SD-SG

There are no ring outliers.

11 monomers are involved in 57 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	114	CME	2	0
2	A	65[B]	NRQ	8	0
2	H	114	CME	1	0
2	B	65[B]	NRQ	7	0
2	G	65[A]	NRQ	8	0
2	G	114	CME	1	0
2	H	65[A]	NRQ	5	0
2	B	65[A]	NRQ	7	0
2	A	65[A]	NRQ	5	0
2	G	65[B]	NRQ	7	0
2	H	65[B]	NRQ	6	0

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	B	1
2	H	1
2	G	1
2	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	65[A]:NRQ	C3	66:SER	N	2.33
1	H	65[A]:NRQ	C3	66:SER	N	2.32
1	G	65[A]:NRQ	C3	66:SER	N	2.04
1	A	65[A]:NRQ	C3	66:SER	N	1.93

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	L	58/62 (93%)	-1.11	0 100 100	18, 34, 39, 44	8 (13%)
1	M	58/62 (93%)	-0.96	0 100 100	23, 42, 55, 62	10 (17%)
1	R	58/62 (93%)	-1.05	0 100 100	14, 34, 44, 45	7 (12%)
1	S	58/62 (93%)	-0.92	0 100 100	20, 44, 55, 63	6 (10%)
2	A	164/167 (98%)	-1.21	0 100 100	15, 32, 50, 73	10 (6%)
2	B	164/167 (98%)	-1.09	0 100 100	20, 41, 53, 73	9 (5%)
2	G	164/167 (98%)	-1.18	0 100 100	13, 32, 51, 73	9 (5%)
2	H	164/167 (98%)	-1.07	0 100 100	22, 40, 52, 71	7 (4%)
All	All	888/916 (96%)	-1.11	0 100 100	13, 35, 53, 73	66 (7%)

There are no RSRZ outliers to report.

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
2	CME	B	114	10/11	0.97	0.07	45,46,47,47	4
2	CME	G	114	10/11	0.97	0.06	34,36,37,38	4
2	NRQ	B	65[A]	23/24	0.98	0.04	47,48,50,52	23
2	NRQ	G	65[A]	23/24	0.98	0.05	37,40,43,45	23
2	NRQ	G	65[B]	23/24	0.98	0.05	32,33,35,36	23
2	NRQ	B	65[B]	23/24	0.98	0.04	38,39,42,43	23
2	NRQ	H	65[A]	23/24	0.98	0.05	48,49,52,54	23
2	NRQ	H	65[B]	23/24	0.98	0.05	40,41,44,45	23

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CME	H	114	10/11	0.98	0.07	45,47,48,48	4
2	NRQ	A	65[A]	23/24	0.99	0.04	36,39,42,45	23
2	NRQ	A	65[B]	23/24	0.99	0.04	31,33,35,36	23
2	CME	G	221	10/11	0.99	0.04	41,44,45,45	0
2	CME	A	114	10/11	0.99	0.04	35,37,45,47	0
2	CME	B	221	10/11	0.99	0.05	45,45,47,48	0
2	CME	A	221	10/11	0.99	0.04	41,43,44,44	0
2	CME	H	221	10/11	0.99	0.04	45,46,47,47	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.