



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:27 AM GMT

PDB ID : 2CYD  
Title : Crystal structure of Lithium bound rotor ring of the V-ATPase from *Enterococcus hirae*  
Authors : Murata, T.; Yamato, I.; Kakinuma, Y.; Shirouzu, M.; Walker, J.E.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2005-07-06  
Resolution : 2.80 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  
<http://wwpdb.org/validation/2016/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	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

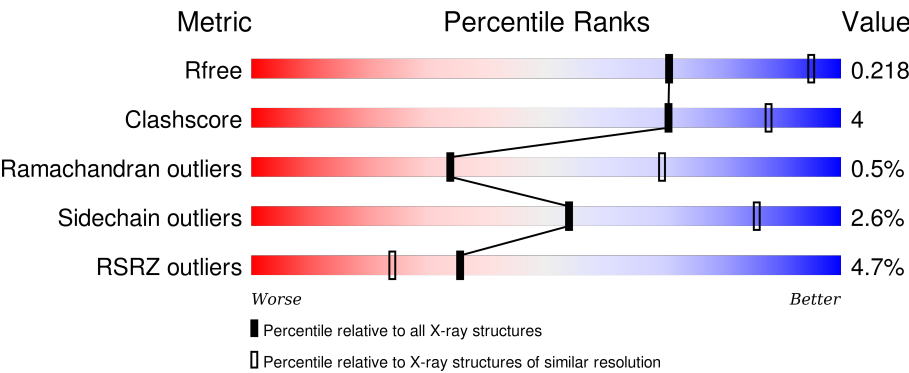
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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 <sub>free</sub>	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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. 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	156	<div><div>6%</div><div>93%</div><div>6% ..</div></div>
1	B	156	<div><div>4%</div><div>93%</div><div>6% ..</div></div>
1	C	156	<div><div>2%</div><div>89%</div><div>10% ..</div></div>
1	D	156	<div><div>13%</div><div>95%</div><div>..</div></div>
1	E	156	<div><div>8%</div><div>95%</div><div>..</div></div>

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Mol	Chain	Length	Quality of chain
1	F	156	
1	G	156	
1	H	156	
1	I	156	
1	J	156	

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	LI	A	157	-	-	-	X
2	LI	D	157	-	-	-	X
2	LI	G	157	-	-	-	X
2	LI	I	157	-	-	-	X
2	LI	J	157	-	-	-	X
3	LHG	A	558	-	-	-	X
3	LHG	A	559	-	-	-	X
3	LHG	B	1158	-	-	-	X
3	LHG	B	1159	-	-	-	X
3	LHG	C	2158	-	-	-	X
3	LHG	C	2159	-	-	-	X
3	LHG	D	3158	-	-	-	X
3	LHG	D	3159	-	-	-	X
3	LHG	E	4158	-	-	-	X
3	LHG	E	4159	-	-	-	X
3	LHG	F	5158	-	-	-	X
3	LHG	G	6158	-	-	-	X
3	LHG	G	6159	-	-	-	X
3	LHG	H	7158	-	-	-	X
3	LHG	H	7159	-	-	-	X
3	LHG	I	8158	-	-	-	X
3	LHG	I	8159	-	-	-	X
3	LHG	J	9158	-	-	-	X
3	LHG	J	9159	-	-	-	X
4	UMQ	C	1001	X	-	-	X
4	UMQ	F	1002	X	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 12831 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 V-type sodium ATP synthase subunit K.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	B	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	C	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	D	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	E	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	F	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	G	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	H	156	Total	C	N	O	S	0	9	0
			1155	760	179	208	8			
1	I	156	Total	C	N	O	S	0	7	0
			1148	757	178	205	8			
1	J	156	Total	C	N	O	S	0	8	0
			1153	759	180	207	7			

- Molecule 2 is LITHIUM ION (three-letter code: LI) (formula: Li).

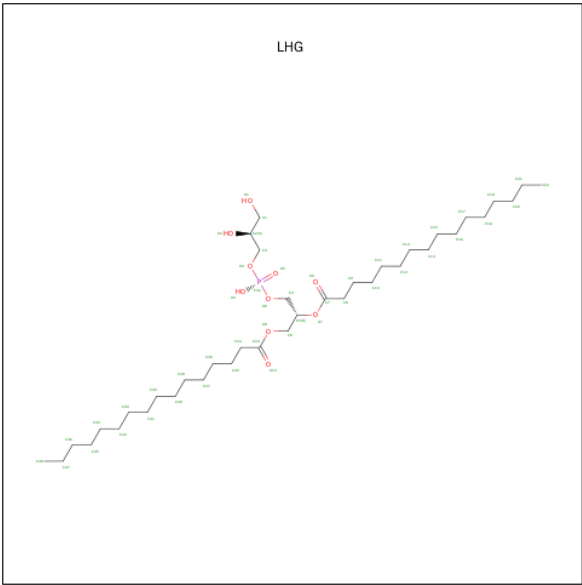
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	Li	0	0
			1	1		
2	J	1	Total	Li	0	0
			1	1		
2	D	1	Total	Li	0	0
			1	1		
2	E	1	Total	Li	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	H	1	Total	Li	0	0
			1	1		
2	B	1	Total	Li	0	0
			1	1		
2	I	1	Total	Li	0	0
			1	1		
2	C	1	Total	Li	0	0
			1	1		
2	A	1	Total	Li	0	0
			1	1		
2	F	1	Total	Li	0	0
			1	1		

- Molecule 3 is 1,2-DIPALMITOYL-PHOSPHATIDYL-GLYCEROLE (three-letter code: LHG) (formula: C<sub>38</sub>H<sub>75</sub>O<sub>10</sub>P).



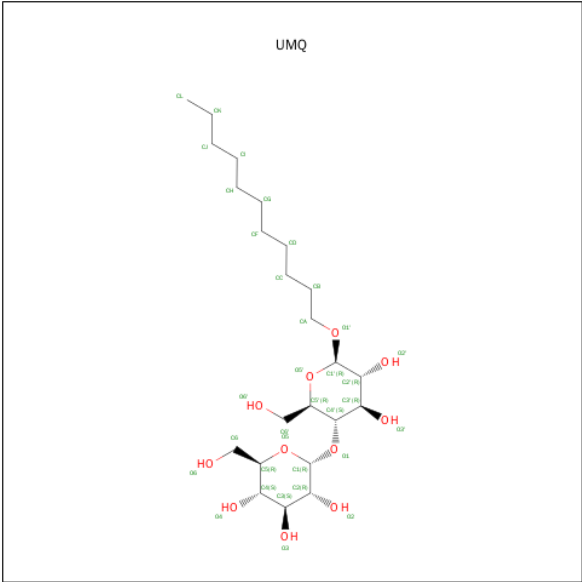
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			49	38	10	1		
3	A	1	Total	C	O		0	0
			40	35	5			
3	B	1	Total	C	O	P	0	0
			49	38	10	1		
3	B	1	Total	C	O		0	0
			40	35	5			
3	C	1	Total	C	O	P	0	0
			49	38	10	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total C O 40 35 5	0	0
3	D	1	Total C O P 49 38 10 1	0	0
3	D	1	Total C O 40 35 5	0	0
3	E	1	Total C O P 49 38 10 1	0	0
3	E	1	Total C O 40 35 5	0	0
3	F	1	Total C O P 49 38 10 1	0	0
3	F	1	Total C O 40 35 5	0	0
3	G	1	Total C O P 49 38 10 1	0	0
3	G	1	Total C O 40 35 5	0	0
3	H	1	Total C O P 49 38 10 1	0	0
3	H	1	Total C O 40 35 5	0	0
3	I	1	Total C O P 49 38 10 1	0	0
3	I	1	Total C O 40 35 5	0	0
3	J	1	Total C O P 49 38 10 1	0	0
3	J	1	Total C O 40 35 5	0	0

- Molecule 4 is UNDECYL-MALTOSIDE (three-letter code: UMQ) (formula: C<sub>23</sub>H<sub>44</sub>O<sub>11</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			34	23	11		
4	F	1	Total	C	O	0	0
			34	23	11		

- Molecule 5 is water.

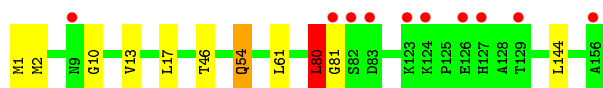
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	29	Total	O	0	0
			29	29		
5	B	29	Total	O	0	0
			29	29		
5	C	19	Total	O	0	0
			19	19		
5	D	24	Total	O	0	0
			24	24		
5	E	25	Total	O	0	0
			25	25		
5	F	32	Total	O	0	0
			32	32		
5	G	35	Total	O	0	0
			35	35		
5	H	37	Total	O	0	0
			37	37		
5	I	57	Total	O	0	0
			57	57		
5	J	35	Total	O	0	0
			35	35		

### 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 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: V-type sodium ATP synthase subunit K

Chain A: 




- Molecule 1: V-type sodium ATP synthase subunit K

Chain B: 



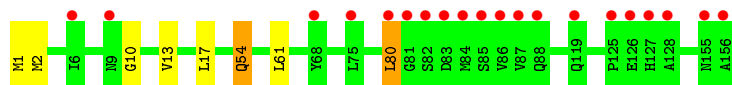
- Molecule 1: V-type sodium ATP synthase subunit K

Chain C: 



- Molecule 1: V-type sodium ATP synthase subunit K

Chain D: 




- Molecule 1: V-type sodium ATP synthase subunit K

Chain E: 



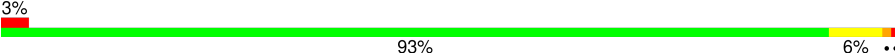
- Molecule 1: V-type sodium ATP synthase subunit K

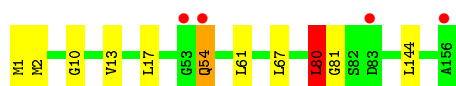


Chain F:  3% 89% 10% ..



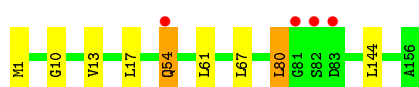
- Molecule 1: V-type sodium ATP synthase subunit K

Chain G:  3% 93% 6% ..




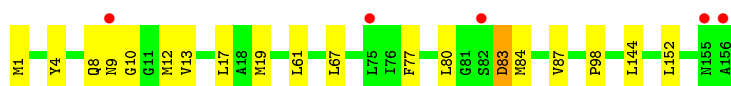
- Molecule 1: V-type sodium ATP synthase subunit K

Chain H:  3% 94% . .




- Molecule 1: V-type sodium ATP synthase subunit K

Chain I:  3% 88% 12% .



- Molecule 1: V-type sodium ATP synthase subunit K

Chain J:  % 91% 8% .



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.65Å 125.78Å 210.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	107.80 – 2.80 61.21 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.1 (107.80-2.80) 98.1 (61.21-2.80)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.93 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.208 , 0.218 0.209 , 0.218	Depositor DCC
$R_{free}$ test set	3890 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	62.0	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 62.3	EDS
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	18 of 77244 reflections (0.023%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	12831	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.54% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: LHG, UMQ, LI

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	0.47	0/1219	0.51	0/1649
1	B	0.48	0/1219	0.52	0/1649
1	C	0.48	0/1219	0.54	0/1649
1	D	0.44	0/1219	0.51	0/1649
1	E	0.46	0/1219	0.51	0/1649
1	F	0.48	0/1219	0.54	0/1649
1	G	0.47	0/1219	0.52	0/1649
1	H	0.51	0/1219	0.53	0/1649
1	I	0.54	0/1202	0.55	0/1626
1	J	0.53	0/1212	0.56	0/1640
All	All	0.49	0/12166	0.53	0/16458

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	1155	0	1218	11	0
1	B	1155	0	1218	11	0
1	C	1155	0	1218	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1155	0	1218	9	0
1	E	1155	0	1218	8	0
1	F	1155	0	1218	14	0
1	G	1155	0	1218	12	0
1	H	1155	0	1218	9	0
1	I	1148	0	1214	15	0
1	J	1153	0	1217	14	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
3	A	89	0	141	7	0
3	B	89	0	141	5	0
3	C	89	0	141	6	0
3	D	89	0	141	7	0
3	E	89	0	141	5	0
3	F	89	0	141	6	0
3	G	89	0	141	8	0
3	H	89	0	141	6	0
3	I	89	0	141	4	0
3	J	89	0	141	8	0
4	C	34	0	41	0	0
4	F	34	0	42	0	0
5	A	29	0	0	0	0
5	B	29	0	0	0	0
5	C	19	0	0	0	0
5	D	24	0	0	0	0
5	E	25	0	0	0	0
5	F	32	0	0	0	0
5	G	35	0	0	0	0
5	H	37	0	0	0	0
5	I	57	0	0	0	0
5	J	35	0	0	0	0
All	All	12831	0	13668	115	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 4.

All (115) 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)
3:A:559:LHG:HC62	3:J:9159:LHG:HC81	1.59	0.84
3:G:6159:LHG:HC81	3:H:7159:LHG:HC62	1.65	0.79
3:F:5159:LHG:HC81	3:G:6159:LHG:HC62	1.64	0.78
3:C:2159:LHG:HC81	3:D:3159:LHG:HC62	1.68	0.76
3:A:559:LHG:HC81	3:B:1159:LHG:HC62	1.71	0.71
3:E:4159:LHG:HC81	3:F:5159:LHG:HC62	1.73	0.69
3:I:8159:LHG:HC81	3:J:9159:LHG:HC62	1.75	0.68
3:D:3159:LHG:HC81	3:E:4159:LHG:HC62	1.77	0.67
3:B:1159:LHG:HC81	3:C:2159:LHG:HC62	1.78	0.65
1:I:10:GLY:O	1:I:13[A]:VAL:HG22	1.98	0.64
1:F:2:MET:HG2	1:G:1:MET:HB2	1.79	0.62
1:I:1:MET:HB3	3:I:8159:LHG:HC5	1.83	0.61
3:A:559:LHG:C6	3:J:9159:LHG:HC81	2.32	0.59
1:G:2:MET:HG2	1:H:1:MET:HB2	1.84	0.59
1:I:9:ASN:HD21	1:I:83:ASP:HA	1.68	0.58
1:A:1:MET:HB3	3:A:559:LHG:HC5	1.86	0.57
1:B:1:MET:HB3	3:B:1159:LHG:HC5	1.87	0.57
1:A:2:MET:HG2	1:B:1:MET:HB2	1.86	0.57
1:J:17[B]:LEU:HD21	3:J:9159:LHG:H192	1.87	0.56
3:H:7159:LHG:HC81	3:I:8159:LHG:HC62	1.87	0.56
1:B:17[B]:LEU:HD21	3:B:1159:LHG:H192	1.87	0.55
1:C:2:MET:HG2	1:D:1:MET:HB2	1.88	0.55
1:D:1:MET:HB3	3:D:3159:LHG:HC5	1.89	0.55
1:H:1:MET:HB3	3:H:7159:LHG:HC5	1.88	0.55
1:F:1:MET:HB3	3:F:5159:LHG:HC5	1.88	0.54
1:C:1:MET:HB3	3:C:2159:LHG:HC5	1.88	0.54
1:E:1:MET:HB3	3:E:4159:LHG:HC5	1.90	0.54
3:G:6159:LHG:HC81	3:H:7159:LHG:C6	2.36	0.53
1:F:17[B]:LEU:HD21	3:F:5159:LHG:H192	1.90	0.53
1:A:17[B]:LEU:HD21	3:A:559:LHG:H192	1.90	0.53
1:G:17[B]:LEU:HD21	3:G:6159:LHG:H192	1.91	0.53
1:E:17[B]:LEU:HD21	3:E:4159:LHG:H192	1.90	0.53
3:A:559:LHG:H341	3:J:9159:LHG:H212	1.91	0.52
1:G:1:MET:HB3	3:G:6159:LHG:HC5	1.89	0.52
1:F:50:GLU:HG2	1:F:51:LYS:HG3	1.90	0.52
1:H:17[B]:LEU:HD21	3:H:7159:LHG:H192	1.89	0.52
1:G:2:MET:CG	1:H:1:MET:HB2	2.40	0.52
1:F:2:MET:CG	1:G:1:MET:HB2	2.40	0.52
1:D:17[B]:LEU:HD21	3:D:3159:LHG:H192	1.92	0.51
1:C:17[B]:LEU:HD21	3:C:2159:LHG:H192	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:12[A]:MET:SD	1:I:80:LEU:HD22	2.52	0.50
3:F:5159:LHG:HC81	3:G:6159:LHG:C6	2.38	0.50
1:C:50:GLU:HG2	1:C:51:LYS:HG3	1.95	0.49
1:E:2:MET:HG2	1:F:1:MET:HB2	1.93	0.49
1:I:9:ASN:HD21	1:I:84:MET:H	1.61	0.48
1:C:2:MET:CG	1:D:1:MET:HB2	2.44	0.47
1:I:144:LEU:HD11	1:J:67:LEU:HD23	1.96	0.47
1:I:4:TYR:CD1	1:I:8:GLN:HG3	2.50	0.47
3:C:2159:LHG:HC81	3:D:3159:LHG:C6	2.40	0.47
1:A:2:MET:CG	1:B:1:MET:HB2	2.45	0.46
1:D:2:MET:HG2	1:E:1:MET:HB2	1.98	0.46
1:I:4:TYR:CE1	1:I:8:GLN:HG3	2.50	0.46
1:A:80:LEU:HB3	1:A:81:GLY:H	1.65	0.46
3:D:3159:LHG:HC81	3:E:4159:LHG:C6	2.46	0.45
1:D:10:GLY:O	1:D:13[A]:VAL:HG22	2.18	0.44
1:C:10:GLY:O	1:C:13[A]:VAL:HG22	2.17	0.44
1:F:10:GLY:O	1:F:13[A]:VAL:HG22	2.17	0.44
1:D:2:MET:CG	1:E:1:MET:HB2	2.48	0.43
1:I:17[B]:LEU:HD21	3:I:8159:LHG:H192	2.00	0.43
3:G:6159:LHG:H212	3:H:7159:LHG:H341	2.00	0.43
1:B:10:GLY:O	1:B:13[A]:VAL:HG22	2.18	0.43
1:E:10:GLY:O	1:E:13[A]:VAL:HG22	2.19	0.43
3:A:559:LHG:HC81	3:B:1159:LHG:C6	2.45	0.43
1:B:144:LEU:HD11	1:C:67:LEU:HD23	2.01	0.43
1:B:80:LEU:HB3	1:B:81:GLY:H	1.67	0.43
1:A:1:MET:HB2	1:J:2:MET:HG2	2.00	0.43
3:F:5159:LHG:H212	3:G:6159:LHG:H341	2.01	0.43
1:A:144:LEU:HD11	1:B:67:LEU:HD23	2.01	0.43
1:H:144:LEU:HD11	1:I:67:LEU:HD23	2.00	0.43
1:C:80:LEU:HB3	1:C:81:GLY:H	1.66	0.43
1:H:10:GLY:O	1:H:13[A]:VAL:HG22	2.19	0.43
1:F:144:LEU:HD11	1:G:67:LEU:HD23	2.01	0.42
1:J:2:MET:CE	3:J:9159:LHG:H121	2.49	0.42
1:A:10:GLY:O	1:A:13[A]:VAL:HG22	2.19	0.42
1:F:59:GLN:O	1:F:62:PRO:HD2	2.19	0.42
1:J:5:LEU:HD22	1:J:10:GLY:HA3	2.00	0.42
1:J:2:MET:HE1	3:J:9159:LHG:H121	2.01	0.42
1:B:2:MET:HG2	1:C:1:MET:HB2	2.00	0.42
1:F:15:ALA:O	1:F:18:ALA:HB3	2.19	0.42
1:I:87:VAL:HG23	1:J:4:TYR:CE2	2.55	0.42
1:A:46:THR:HG22	1:J:120:ILE:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:152:LEU:HD21	1:J:19:MET:SD	2.59	0.42
1:F:54[B]:GLN:NE2	1:F:128:ALA:HB3	2.34	0.42
1:G:144:LEU:HD11	1:H:67:LEU:HD23	2.02	0.42
1:J:61:LEU:HD23	1:J:135:ALA:HB3	2.02	0.41
1:G:80:LEU:HB3	1:G:81:GLY:H	1.66	0.41
1:C:15:ALA:O	1:C:18:ALA:HB3	2.20	0.41
1:I:9:ASN:ND2	1:I:84:MET:H	2.18	0.41
3:J:9158:LHG:H251	3:J:9158:LHG:HC82	2.03	0.41
3:C:2159:LHG:H212	3:D:3159:LHG:H341	2.03	0.41
1:I:98:PRO:HG3	1:J:20:ALA:HA	2.03	0.41
1:G:10:GLY:O	1:G:13[A]:VAL:HG22	2.21	0.41
1:I:19:MET:CE	1:I:77:PHE:HB2	2.51	0.41
1:C:54[B]:GLN:NE2	1:C:128:ALA:HB3	2.36	0.40
1:F:80:LEU:HB3	1:F:81:GLY:H	1.65	0.40
1:C:4:TYR:CE1	1:C:8:GLN:HG3	2.56	0.40

There are no symmetry-related clashes.

## 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	163/156 (104%)	162 (99%)	0	1 (1%)	30	65
1	B	163/156 (104%)	161 (99%)	1 (1%)	1 (1%)	30	65
1	C	163/156 (104%)	162 (99%)	0	1 (1%)	30	65
1	D	163/156 (104%)	161 (99%)	1 (1%)	1 (1%)	30	65
1	E	163/156 (104%)	161 (99%)	1 (1%)	1 (1%)	30	65
1	F	163/156 (104%)	162 (99%)	0	1 (1%)	30	65
1	G	163/156 (104%)	161 (99%)	1 (1%)	1 (1%)	30	65
1	H	163/156 (104%)	161 (99%)	1 (1%)	1 (1%)	30	65

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	I	161/156 (103%)	160 (99%)	1 (1%)	0	100	100
1	J	162/156 (104%)	160 (99%)	2 (1%)	0	100	100
All	All	1627/1560 (104%)	1611 (99%)	8 (0%)	8 (0%)	34	69

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	80	LEU
1	B	80	LEU
1	C	80	LEU
1	E	80	LEU
1	F	80	LEU
1	G	80	LEU
1	H	80	LEU
1	D	80	LEU

### 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	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	B	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	C	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	D	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	E	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	F	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	G	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	H	122/113 (108%)	118 (97%)	4 (3%)	45	79
1	I	120/113 (106%)	118 (98%)	2 (2%)	68	92
1	J	121/113 (107%)	117 (97%)	4 (3%)	45	79
All	All	1217/1130 (108%)	1179 (97%)	38 (3%)	54	81



All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54[A]	GLN
1	A	54[B]	GLN
1	A	61	LEU
1	A	80	LEU
1	B	54[A]	GLN
1	B	54[B]	GLN
1	B	61	LEU
1	B	80	LEU
1	C	54[A]	GLN
1	C	54[B]	GLN
1	C	61	LEU
1	C	80	LEU
1	D	54[A]	GLN
1	D	54[B]	GLN
1	D	61	LEU
1	D	80	LEU
1	E	54[A]	GLN
1	E	54[B]	GLN
1	E	61	LEU
1	E	80	LEU
1	F	54[A]	GLN
1	F	54[B]	GLN
1	F	61	LEU
1	F	80	LEU
1	G	54[A]	GLN
1	G	54[B]	GLN
1	G	61	LEU
1	G	80	LEU
1	H	54[A]	GLN
1	H	54[B]	GLN
1	H	61	LEU
1	H	80	LEU
1	I	61	LEU
1	I	83	ASP
1	J	7	THR
1	J	54[A]	GLN
1	J	54[B]	GLN
1	J	61	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	I	9	ASN
1	J	8	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 32 ligands modelled in this entry, 10 are monoatomic - leaving 22 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	LHG	A	558	-	48,48,48	0.96	2 (4%)	49,54,54	1.02	4 (8%)
3	LHG	A	559	-	39,39,48	1.03	2 (5%)	41,41,54	1.09	2 (4%)
3	LHG	B	1158	-	48,48,48	0.97	2 (4%)	49,54,54	0.99	2 (4%)
3	LHG	B	1159	-	39,39,48	1.03	2 (5%)	41,41,54	1.08	2 (4%)
4	UMQ	C	1001	-	35,35,35	1.19	3 (8%)	46,46,46	1.76	9 (19%)
3	LHG	C	2158	-	48,48,48	0.98	2 (4%)	49,54,54	0.97	2 (4%)
3	LHG	C	2159	-	39,39,48	1.02	2 (5%)	41,41,54	1.08	2 (4%)
3	LHG	D	3158	-	48,48,48	0.95	2 (4%)	49,54,54	1.01	3 (6%)
3	LHG	D	3159	-	39,39,48	1.03	2 (5%)	41,41,54	1.09	2 (4%)
3	LHG	E	4158	-	48,48,48	0.95	2 (4%)	49,54,54	1.02	3 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	LHG	E	4159	-	39,39,48	1.02	2 (5%)	41,41,54	1.08	2 (4%)
4	UMQ	F	1002	-	35,35,35	1.18	3 (8%)	46,46,46	1.44	6 (13%)
3	LHG	F	5158	-	48,48,48	0.97	2 (4%)	49,54,54	1.03	3 (6%)
3	LHG	F	5159	-	39,39,48	1.02	2 (5%)	41,41,54	1.09	2 (4%)
3	LHG	G	6158	-	48,48,48	0.97	2 (4%)	49,54,54	1.02	3 (6%)
3	LHG	G	6159	-	39,39,48	1.03	2 (5%)	41,41,54	1.07	2 (4%)
3	LHG	H	7158	-	48,48,48	0.96	2 (4%)	49,54,54	1.03	4 (8%)
3	LHG	H	7159	-	39,39,48	1.02	2 (5%)	41,41,54	1.08	2 (4%)
3	LHG	I	8158	-	48,48,48	0.96	2 (4%)	49,54,54	1.01	4 (8%)
3	LHG	I	8159	-	39,39,48	1.04	2 (5%)	41,41,54	1.11	3 (7%)
3	LHG	J	9158	-	48,48,48	0.94	2 (4%)	49,54,54	1.05	4 (8%)
3	LHG	J	9159	-	39,39,48	1.04	2 (5%)	41,41,54	1.01	2 (4%)

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
3	LHG	A	558	-	-	0/53/53/53	0/0/0/0
3	LHG	A	559	-	-	0/41/41/53	0/0/0/0
3	LHG	B	1158	-	-	0/53/53/53	0/0/0/0
3	LHG	B	1159	-	-	0/41/41/53	0/0/0/0
4	UMQ	C	1001	-	1/1/10/10	0/20/60/60	0/2/2/2
3	LHG	C	2158	-	-	0/53/53/53	0/0/0/0
3	LHG	C	2159	-	-	0/41/41/53	0/0/0/0
3	LHG	D	3158	-	-	0/53/53/53	0/0/0/0
3	LHG	D	3159	-	-	0/41/41/53	0/0/0/0
3	LHG	E	4158	-	-	0/53/53/53	0/0/0/0
3	LHG	E	4159	-	-	0/41/41/53	0/0/0/0
4	UMQ	F	1002	-	1/1/10/10	0/20/60/60	0/2/2/2
3	LHG	F	5158	-	-	0/53/53/53	0/0/0/0
3	LHG	F	5159	-	-	0/41/41/53	0/0/0/0
3	LHG	G	6158	-	-	0/53/53/53	0/0/0/0
3	LHG	G	6159	-	-	0/41/41/53	0/0/0/0
3	LHG	H	7158	-	-	0/53/53/53	0/0/0/0
3	LHG	H	7159	-	-	0/41/41/53	0/0/0/0
3	LHG	I	8158	-	-	0/53/53/53	0/0/0/0
3	LHG	I	8159	-	-	0/41/41/53	0/0/0/0
3	LHG	J	9158	-	-	0/53/53/53	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	LHG	J	9159	-	-	0/41/41/53	0/0/0/0

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	1002	UMQ	O3-C3	-4.44	1.32	1.43
4	F	1002	UMQ	O2-C2	-4.26	1.32	1.43
4	C	1001	UMQ	O3-C3	-4.25	1.32	1.43
4	C	1001	UMQ	O2-C2	-4.24	1.32	1.43
4	F	1002	UMQ	O1'-C1'	2.07	1.43	1.40
4	C	1001	UMQ	O1'-C1'	2.11	1.43	1.40
3	D	3158	LHG	O7-C7	3.93	1.46	1.34
3	I	8158	LHG	O7-C7	3.96	1.46	1.34
3	H	7158	LHG	O7-C7	3.98	1.46	1.34
3	E	4158	LHG	O7-C7	4.01	1.46	1.34
3	A	558	LHG	O7-C7	4.01	1.46	1.34
3	H	7159	LHG	O7-C7	4.02	1.46	1.34
3	J	9158	LHG	O7-C7	4.02	1.46	1.34
3	A	559	LHG	O7-C7	4.03	1.46	1.34
3	E	4159	LHG	O7-C7	4.04	1.46	1.34
3	F	5159	LHG	O7-C7	4.05	1.46	1.34
3	B	1158	LHG	O7-C7	4.06	1.46	1.34
3	G	6159	LHG	O7-C7	4.07	1.46	1.34
3	C	2159	LHG	O7-C7	4.08	1.46	1.34
3	B	1159	LHG	O7-C7	4.10	1.46	1.34
3	D	3159	LHG	O7-C7	4.10	1.46	1.34
3	F	5158	LHG	O7-C7	4.15	1.46	1.34
3	J	9158	LHG	O8-C23	4.16	1.45	1.33
3	J	9159	LHG	O8-C23	4.17	1.45	1.33
3	G	6158	LHG	O7-C7	4.18	1.46	1.34
3	A	558	LHG	O8-C23	4.24	1.46	1.33
3	F	5159	LHG	O8-C23	4.25	1.46	1.33
3	I	8159	LHG	O8-C23	4.26	1.46	1.33
3	C	2159	LHG	O8-C23	4.27	1.46	1.33
3	E	4158	LHG	O8-C23	4.27	1.46	1.33
3	I	8159	LHG	O7-C7	4.27	1.47	1.34
3	C	2158	LHG	O7-C7	4.27	1.47	1.34
3	F	5158	LHG	O8-C23	4.28	1.46	1.33
3	J	9159	LHG	O7-C7	4.29	1.47	1.34
3	I	8158	LHG	O8-C23	4.31	1.46	1.33
3	E	4159	LHG	O8-C23	4.32	1.46	1.33
3	C	2158	LHG	O8-C23	4.33	1.46	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1159	LHG	O8-C23	4.33	1.46	1.33
3	G	6159	LHG	O8-C23	4.34	1.46	1.33
3	G	6158	LHG	O8-C23	4.34	1.46	1.33
3	D	3159	LHG	O8-C23	4.35	1.46	1.33
3	H	7159	LHG	O8-C23	4.35	1.46	1.33
3	H	7158	LHG	O8-C23	4.35	1.46	1.33
3	B	1158	LHG	O8-C23	4.36	1.46	1.33
3	D	3158	LHG	O8-C23	4.36	1.46	1.33
3	A	559	LHG	O8-C23	4.39	1.46	1.33

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	9158	LHG	O7-C7-O9	-2.24	117.67	123.67
3	F	5158	LHG	O8-C23-O10	-2.23	117.72	123.49
3	I	8158	LHG	O8-C23-O10	-2.21	117.80	123.49
3	A	558	LHG	O8-C23-O10	-2.20	117.81	123.49
3	I	8158	LHG	O7-C7-O9	-2.19	117.80	123.67
4	C	1001	UMQ	C3'-C4'-C5'	-2.13	106.02	110.84
3	H	7158	LHG	O7-C7-O9	-2.12	117.97	123.67
3	J	9158	LHG	O8-C23-O10	-2.12	118.02	123.49
3	A	558	LHG	O7-C7-O9	-2.08	118.09	123.67
3	G	6158	LHG	O8-C23-O10	-2.07	118.16	123.49
3	D	3158	LHG	O7-C7-O9	-2.05	118.16	123.67
3	H	7158	LHG	O8-C23-O10	-2.04	118.23	123.49
3	E	4158	LHG	O8-C23-O10	-2.04	118.23	123.49
4	C	1001	UMQ	O5-C5-C6	2.12	111.72	106.36
4	F	1002	UMQ	O3-C3-C2	2.14	115.16	110.34
3	I	8159	LHG	C6-O8-C23	2.25	123.14	116.85
4	F	1002	UMQ	O2-C2-C3	2.58	116.16	110.34
3	J	9159	LHG	O8-C23-C24	2.60	119.83	111.90
3	I	8159	LHG	O8-C23-C24	2.69	120.10	111.90
3	B	1159	LHG	O8-C23-C24	2.75	120.28	111.90
3	G	6159	LHG	O8-C23-C24	2.75	120.29	111.90
3	D	3159	LHG	O8-C23-C24	2.75	120.29	111.90
4	C	1001	UMQ	O2-C2-C3	2.78	116.60	110.34
3	F	5159	LHG	O8-C23-C24	2.79	120.39	111.90
3	E	4159	LHG	O8-C23-C24	2.80	120.42	111.90
3	A	559	LHG	O8-C23-C24	2.82	120.48	111.90
3	H	7159	LHG	O8-C23-C24	2.82	120.49	111.90
3	C	2159	LHG	O8-C23-C24	2.82	120.50	111.90
3	C	2158	LHG	O8-C23-C24	2.82	120.50	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1158	LHG	O8-C23-C24	2.83	120.53	111.90
3	J	9158	LHG	O8-C23-C24	2.84	120.56	111.90
3	G	6158	LHG	O8-C23-C24	2.87	120.65	111.90
3	D	3158	LHG	O8-C23-C24	2.89	120.70	111.90
3	A	558	LHG	O8-C23-C24	2.90	120.73	111.90
3	E	4158	LHG	O8-C23-C24	2.97	120.96	111.90
3	I	8158	LHG	O8-C23-C24	3.03	121.12	111.90
3	H	7158	LHG	O8-C23-C24	3.07	121.25	111.90
3	F	5158	LHG	O8-C23-C24	3.20	121.64	111.90
4	F	1002	UMQ	O3-C3-C4	3.46	118.12	110.34
4	F	1002	UMQ	O2-C2-C1	3.57	117.84	110.02
4	F	1002	UMQ	C4-C3-C2	3.61	117.53	110.79
4	F	1002	UMQ	C1-C2-C3	3.65	117.16	109.97
4	C	1001	UMQ	O1'-C1'-C2'	3.71	112.73	108.04
3	J	9159	LHG	O7-C7-C8	3.92	120.05	111.53
4	C	1001	UMQ	O3-C3-C2	3.93	119.19	110.34
3	A	559	LHG	O7-C7-C8	4.01	120.24	111.53
3	C	2159	LHG	O7-C7-C8	4.01	120.25	111.53
3	E	4159	LHG	O7-C7-C8	4.02	120.27	111.53
3	H	7159	LHG	O7-C7-C8	4.04	120.30	111.53
3	F	5159	LHG	O7-C7-C8	4.06	120.36	111.53
3	G	6159	LHG	O7-C7-C8	4.07	120.37	111.53
3	C	2158	LHG	O7-C7-C8	4.10	120.44	111.53
4	C	1001	UMQ	O2-C2-C1	4.12	119.04	110.02
3	D	3159	LHG	O7-C7-C8	4.16	120.57	111.53
3	B	1159	LHG	O7-C7-C8	4.17	120.59	111.53
3	I	8159	LHG	O7-C7-C8	4.18	120.61	111.53
3	I	8158	LHG	O7-C7-C8	4.21	120.69	111.53
3	B	1158	LHG	O7-C7-C8	4.29	120.85	111.53
3	F	5158	LHG	O7-C7-C8	4.34	120.97	111.53
3	H	7158	LHG	O7-C7-C8	4.35	120.97	111.53
3	G	6158	LHG	O7-C7-C8	4.36	121.00	111.53
3	A	558	LHG	O7-C7-C8	4.36	121.01	111.53
3	E	4158	LHG	O7-C7-C8	4.37	121.02	111.53
3	D	3158	LHG	O7-C7-C8	4.38	121.04	111.53
4	C	1001	UMQ	O3-C3-C4	4.58	120.64	110.34
3	J	9158	LHG	O7-C7-C8	4.64	121.62	111.53
4	C	1001	UMQ	C1-C2-C3	4.70	119.23	109.97
4	C	1001	UMQ	C4-C3-C2	4.96	120.05	110.79

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	F	1002	UMQ	C2
4	C	1001	UMQ	C3

There are no torsion outliers.

There are no ring outliers.

11 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	559	LHG	7	0
3	B	1159	LHG	5	0
3	C	2159	LHG	6	0
3	D	3159	LHG	7	0
3	E	4159	LHG	5	0
3	F	5159	LHG	6	0
3	G	6159	LHG	8	0
3	H	7159	LHG	6	0
3	I	8159	LHG	4	0
3	J	9158	LHG	1	0
3	J	9159	LHG	7	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 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	156/156 (100%)	0.44	10 (6%) 23 14	44, 49, 69, 86	0
1	B	156/156 (100%)	0.24	7 (4%) 37 26	44, 49, 69, 86	0
1	C	156/156 (100%)	0.17	3 (1%) 70 59	44, 49, 69, 86	0
1	D	156/156 (100%)	0.75	20 (12%) 5 2	44, 49, 69, 86	0
1	E	156/156 (100%)	0.46	13 (8%) 14 7	44, 49, 69, 86	0
1	F	156/156 (100%)	0.32	5 (3%) 51 39	44, 49, 69, 86	0
1	G	156/156 (100%)	0.24	4 (2%) 59 47	44, 49, 69, 86	0
1	H	156/156 (100%)	0.21	4 (2%) 59 47	44, 49, 69, 86	0
1	I	156/156 (100%)	0.41	5 (3%) 51 39	44, 49, 66, 86	0
1	J	156/156 (100%)	-0.04	2 (1%) 79 71	44, 49, 59, 72	0
All	All	1560/1560 (100%)	0.32	73 (4%) 35 24	44, 49, 69, 86	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	82	SER	8.8
1	E	81	GLY	8.5
1	D	81	GLY	6.9
1	H	82	SER	6.0
1	D	82	SER	5.6
1	E	83[A]	ASP	5.4
1	H	81	GLY	5.4
1	F	81	GLY	5.1
1	F	83[A]	ASP	5.0
1	E	4	TYR	4.9
1	H	83[A]	ASP	4.8
1	A	83[A]	ASP	4.6
1	D	156	ALA	4.6

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Mol	Chain	Res	Type	RSRZ
1	C	82	SER	4.4
1	B	82	SER	4.4
1	E	5	LEU	4.3
1	F	9	ASN	4.3
1	D	126	GLU	4.3
1	E	9	ASN	4.1
1	E	8	GLN	4.0
1	A	81	GLY	3.9
1	E	7	THR	3.8
1	D	127	HIS	3.8
1	F	82	SER	3.7
1	A	126	GLU	3.7
1	D	83[A]	ASP	3.6
1	A	9	ASN	3.5
1	D	85	SER	3.5
1	E	46	THR	3.4
1	B	54[A]	GLN	3.4
1	D	88	GLN	3.4
1	E	6	ILE	3.3
1	D	125	PRO	3.3
1	D	9	ASN	3.3
1	A	82	SER	3.3
1	B	53	GLY	3.2
1	D	6	ILE	3.2
1	D	84	MET	3.2
1	F	10	GLY	3.2
1	D	68	TYR	3.1
1	A	156	ALA	3.0
1	H	54[A]	GLN	3.0
1	D	155[A]	ASN	3.0
1	C	83[A]	ASP	2.9
1	D	128	ALA	2.9
1	B	57	ILE	2.8
1	B	83[A]	ASP	2.8
1	B	126	GLU	2.7
1	A	123	LYS	2.7
1	E	10	GLY	2.7
1	I	75[A]	LEU	2.5
1	I	9	ASN	2.5
1	I	156	ALA	2.5
1	J	81	GLY	2.5
1	D	86	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	129	THR	2.4
1	E	1	MET	2.4
1	D	75[A]	LEU	2.4
1	E	49	PRO	2.4
1	B	52	PHE	2.3
1	D	80	LEU	2.3
1	G	54[A]	GLN	2.3
1	A	127	HIS	2.2
1	I	155[A]	ASN	2.2
1	D	87	VAL	2.2
1	A	124	LYS	2.2
1	G	156	ALA	2.2
1	G	53	GLY	2.1
1	I	82	SER	2.1
1	D	119[A]	GLN	2.1
1	J	83	ASP	2.1
1	C	85	SER	2.0
1	G	83[A]	ASP	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 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. 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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	LHG	A	558	49/49	0.73	0.43	14.73	87,106,123,124	0
3	LHG	C	2158	49/49	0.73	0.42	11.94	87,106,123,124	0
3	LHG	J	9158	49/49	0.76	0.40	10.68	87,106,123,124	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	LHG	H	7158	49/49	0.66	0.43	10.57	87,106,123,124	0
3	LHG	E	4158	49/49	0.74	0.38	10.09	87,106,123,124	0
3	LHG	I	8158	49/49	0.73	0.42	9.65	87,106,122,124	0
3	LHG	G	6158	49/49	0.75	0.39	9.59	87,106,123,124	0
3	LHG	B	1158	49/49	0.78	0.40	9.35	87,106,123,124	0
3	LHG	D	3158	49/49	0.71	0.37	8.98	87,106,123,124	0
3	LHG	D	3159	40/49	0.62	0.49	7.95	116,128,142,143	0
3	LHG	B	1159	40/49	0.65	0.43	7.04	116,128,142,143	0
3	LHG	F	5158	49/49	0.74	0.40	6.17	87,106,122,124	0
3	LHG	H	7159	40/49	0.79	0.38	6.04	116,128,142,143	0
2	LI	I	157	1/1	0.91	0.34	5.20	47,47,47,47	0
3	LHG	I	8159	40/49	0.70	0.39	4.49	116,128,142,143	0
2	LI	G	157	1/1	0.91	0.34	4.44	47,47,47,47	0
3	LHG	C	2159	40/49	0.76	0.34	4.19	116,128,142,143	0
3	LHG	J	9159	40/49	0.70	0.41	4.17	116,128,142,143	0
2	LI	D	157	1/1	0.88	0.40	4.12	47,47,47,47	0
2	LI	J	157	1/1	0.92	0.27	3.83	47,47,47,47	0
3	LHG	A	559	40/49	0.67	0.41	3.79	116,128,142,143	0
2	LI	A	157	1/1	0.95	0.32	3.22	47,47,47,47	0
3	LHG	G	6159	40/49	0.75	0.35	3.07	116,128,142,143	0
4	UMQ	F	1002	34/34	0.86	0.41	2.27	77,88,90,90	0
4	UMQ	C	1001	34/34	0.86	0.37	2.01	91,97,98,99	0
3	LHG	E	4159	40/49	0.59	0.45	1.88	116,128,142,143	0
3	LHG	F	5159	40/49	0.74	0.32	1.80	116,128,142,143	0
2	LI	E	157	1/1	0.89	0.20	1.08	47,47,47,47	0
2	LI	C	157	1/1	0.96	0.18	-0.10	47,47,47,47	0
2	LI	B	157	1/1	0.90	0.17	-0.91	47,47,47,47	0
2	LI	H	157	1/1	0.91	0.12	-4.08	47,47,47,47	0
2	LI	F	157	1/1	0.82	0.11	-10.23	47,47,47,47	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.