



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:16 PM GMT

PDB ID : 4C01  
Title : Complete crystal structure of carboxylesterase Cest-2923 (lp\_2923) from *Lactobacillus plantarum* WCFS1  
Authors : Benavente, R.; Esteban-Torres, M.; Acebron, I.; de las Rivas, B.; Munoz, R.; Alvarez, Y.; Mancheno, J.M.  
Deposited on : 2013-07-30  
Resolution : 2.30 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

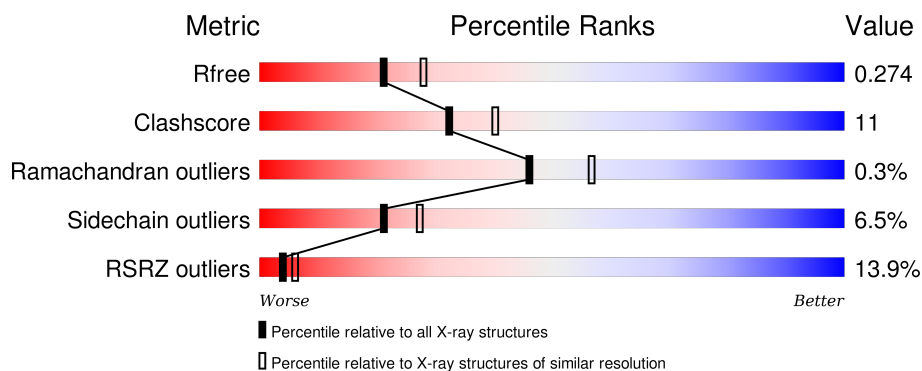
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	282	<div> <div>15%</div> <div>71%</div> <div>25%</div> <div>..</div> </div>
1	B	282	<div> <div>16%</div> <div>78%</div> <div>19%</div> <div>..</div> </div>
1	C	282	<div> <div>14%</div> <div>78%</div> <div>19%</div> <div>..</div> </div>
1	D	282	<div> <div>10%</div> <div>74%</div> <div>22%</div> <div>..</div> </div>
1	E	282	<div> <div>14%</div> <div>76%</div> <div>20%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	282	

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
3	SO4	A	1277	-	-	X	-
4	QY9	A	1278	-	-	-	X
4	QY9	B	1278	-	-	-	X
4	QY9	C	1279	-	-	-	X
4	QY9	E	1279	-	-	-	X
4	QY9	F	1277	-	-	-	X
5	CCN	B	1280	-	-	X	-
5	CCN	D	1279	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14061 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 CEST-2923.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	2	1
			2155	1373	381	393	8			
1	B	276	Total	C	N	O	S	0	1	1
			2144	1367	377	392	8			
1	C	276	Total	C	N	O	S	0	1	1
			2144	1367	377	392	8			
1	D	276	Total	C	N	O	S	0	1	1
			2144	1367	377	392	8			
1	E	276	Total	C	N	O	S	0	2	1
			2155	1373	381	393	8			
1	F	276	Total	C	N	O	S	0	2	1
			2155	1373	381	393	8			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	277	HIS	-	EXPRESSION TAG	UNP F9US10
A	278	HIS	-	EXPRESSION TAG	UNP F9US10
A	279	HIS	-	EXPRESSION TAG	UNP F9US10
A	280	HIS	-	EXPRESSION TAG	UNP F9US10
A	281	HIS	-	EXPRESSION TAG	UNP F9US10
A	282	HIS	-	EXPRESSION TAG	UNP F9US10
B	277	HIS	-	EXPRESSION TAG	UNP F9US10
B	278	HIS	-	EXPRESSION TAG	UNP F9US10
B	279	HIS	-	EXPRESSION TAG	UNP F9US10
B	280	HIS	-	EXPRESSION TAG	UNP F9US10
B	281	HIS	-	EXPRESSION TAG	UNP F9US10
B	282	HIS	-	EXPRESSION TAG	UNP F9US10
C	277	HIS	-	EXPRESSION TAG	UNP F9US10
C	278	HIS	-	EXPRESSION TAG	UNP F9US10
C	279	HIS	-	EXPRESSION TAG	UNP F9US10
C	280	HIS	-	EXPRESSION TAG	UNP F9US10
C	281	HIS	-	EXPRESSION TAG	UNP F9US10

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Chain	Residue	Modelled	Actual	Comment	Reference
C	282	HIS	-	EXPRESSION TAG	UNP F9US10
D	277	HIS	-	EXPRESSION TAG	UNP F9US10
D	278	HIS	-	EXPRESSION TAG	UNP F9US10
D	279	HIS	-	EXPRESSION TAG	UNP F9US10
D	280	HIS	-	EXPRESSION TAG	UNP F9US10
D	281	HIS	-	EXPRESSION TAG	UNP F9US10
D	282	HIS	-	EXPRESSION TAG	UNP F9US10
E	277	HIS	-	EXPRESSION TAG	UNP F9US10
E	278	HIS	-	EXPRESSION TAG	UNP F9US10
E	279	HIS	-	EXPRESSION TAG	UNP F9US10
E	280	HIS	-	EXPRESSION TAG	UNP F9US10
E	281	HIS	-	EXPRESSION TAG	UNP F9US10
E	282	HIS	-	EXPRESSION TAG	UNP F9US10
F	277	HIS	-	EXPRESSION TAG	UNP F9US10
F	278	HIS	-	EXPRESSION TAG	UNP F9US10
F	279	HIS	-	EXPRESSION TAG	UNP F9US10
F	280	HIS	-	EXPRESSION TAG	UNP F9US10
F	281	HIS	-	EXPRESSION TAG	UNP F9US10
F	282	HIS	-	EXPRESSION TAG	UNP F9US10

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C<sub>2</sub>H<sub>3</sub>O<sub>2</sub>).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



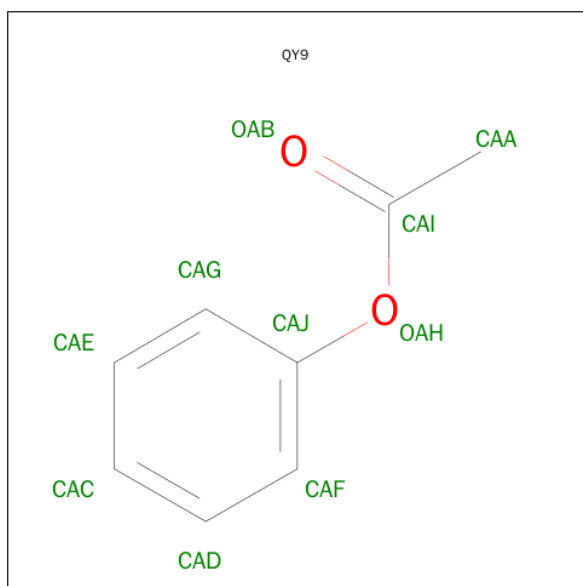
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		

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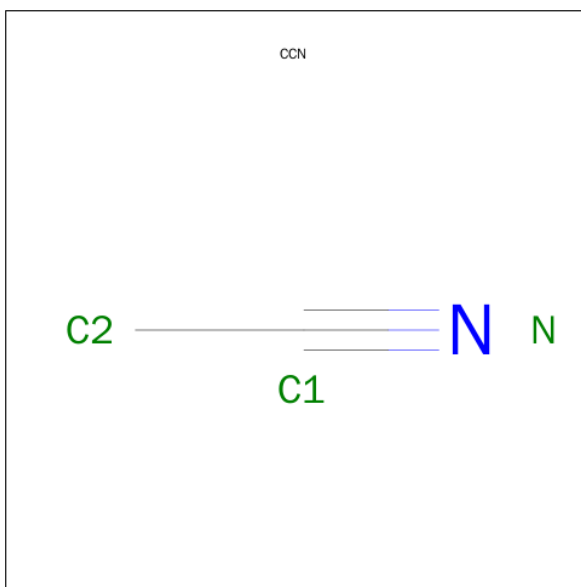
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is PHENYL ACETATE (three-letter code: QY9) (formula:  $C_8H_8O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	8	2		
4	B	1	Total	C	O	0	0
			10	8	2		
4	C	1	Total	C	O	0	0
			10	8	2		
4	E	1	Total	C	O	0	0
			10	8	2		
4	F	1	Total	C	O	0	0
			10	8	2		

- Molecule 5 is ACETONITRILE (three-letter code: CCN) (formula:  $C_2H_3N$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			3	2	1		
5	A	1	Total	C	N	0	0
			3	2	1		
5	A	1	Total	C	N	0	0
			3	2	1		
5	B	1	Total	C	N	0	0
			3	2	1		
5	B	1	Total	C	N	0	0
			3	2	1		
5	C	1	Total	C	N	0	0
			3	2	1		
5	D	1	Total	C	N	0	0
			3	2	1		
5	D	1	Total	C	N	0	0
			3	2	1		
5	E	1	Total	C	N	0	0
			3	2	1		
5	F	1	Total	C	N	0	0
			3	2	1		
5	F	1	Total	C	N	0	0
			3	2	1		
5	F	1	Total	C	N	0	0
			3	2	1		

- Molecule 6 is water.

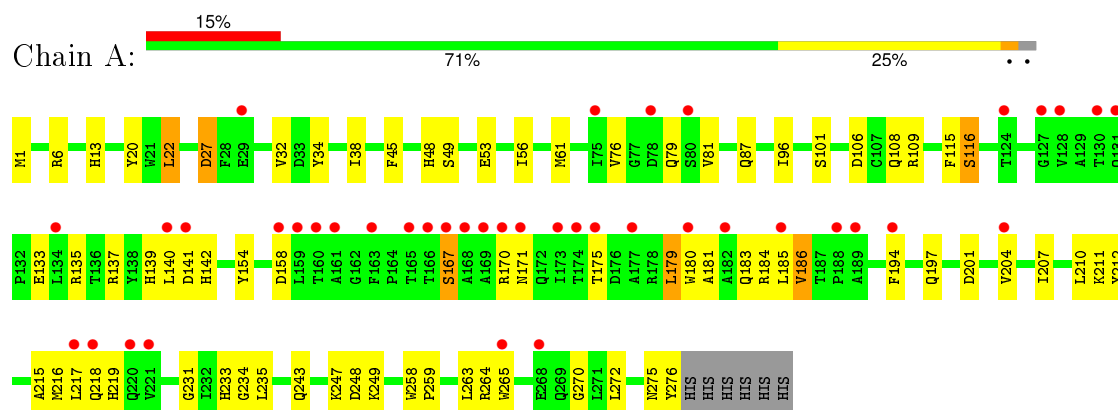


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	193	Total 193	O 193	0	0
6	B	182	Total 182	O 182	0	0
6	C	191	Total 191	O 191	0	0
6	D	137	Total 137	O 137	0	0
6	E	152	Total 152	O 152	0	0
6	F	159	Total 159	O 159	0	0

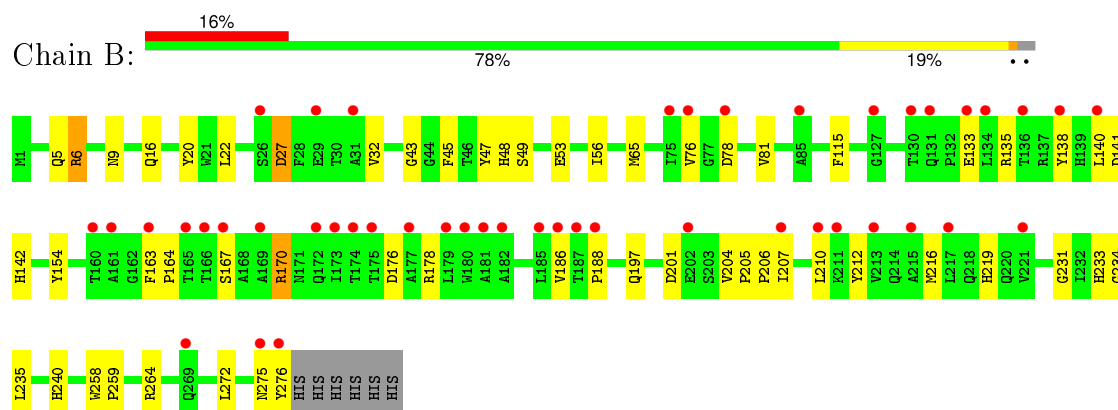
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

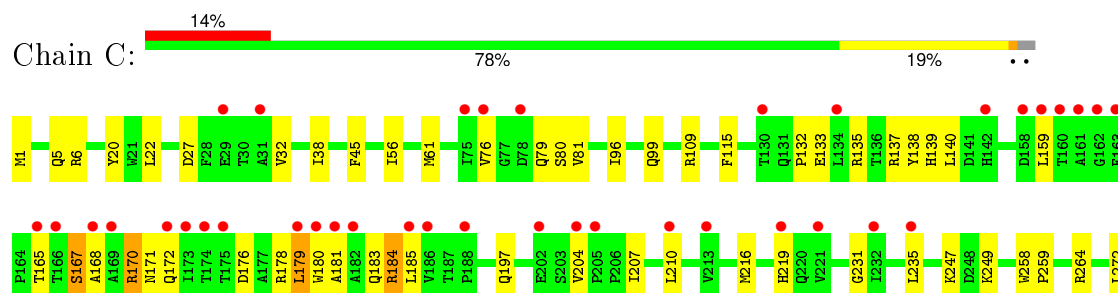
#### • Molecule 1: CEST-2923

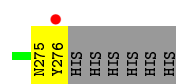


#### • Molecule 1: CEST-2923

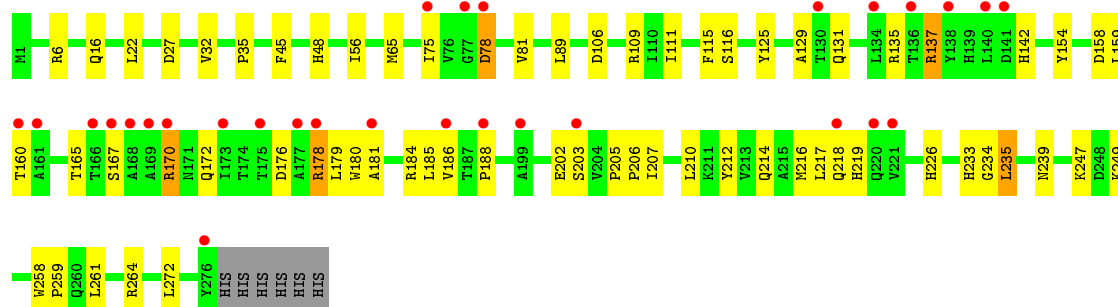
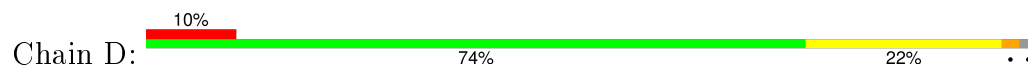


#### • Molecule 1: CEST-2923

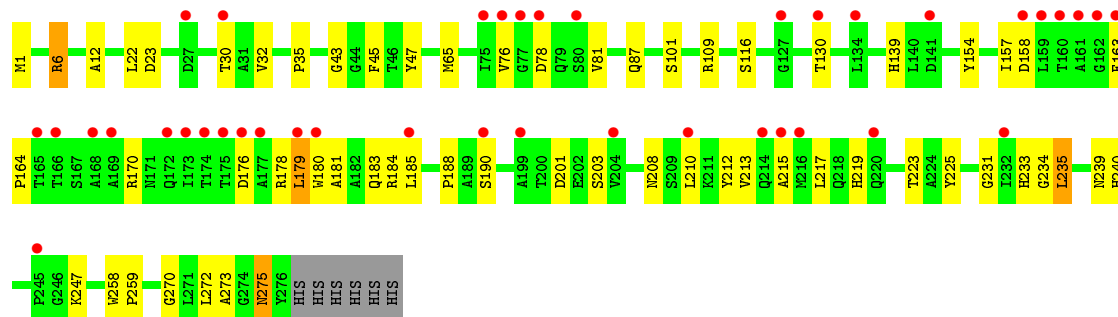
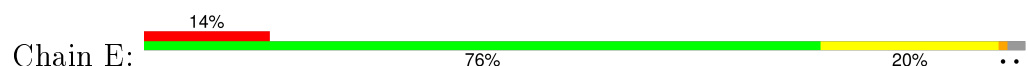




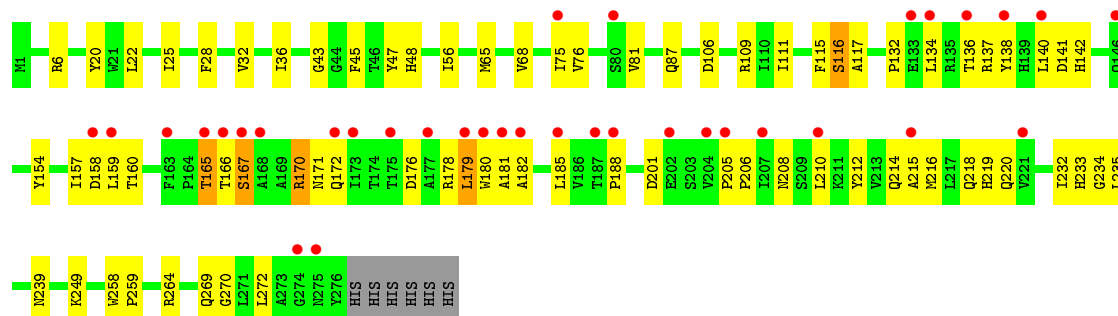
• Molecule 1: CEST-2923



• Molecule 1: CEST-2923



• Molecule 1: CEST-2923



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	244.89Å 141.38Å 82.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.20 – 2.30 49.20 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.20-2.30) 95.8 (49.20-2.30)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	43.02 (at 2.29Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.225 , 0.279 0.225 , 0.274	Depositor DCC
$R_{free}$ test set	6266 reflections (5.52%)	DCC
Wilson B-factor (Å <sup>2</sup> )	21.0	Xtriage
Anisotropy	0.378	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
Estimated twinning fraction	0.450 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.439 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.449 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.438 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.448 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	3 of 124507 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	14061	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CCN, QY9, SO4, ACT

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	0.43	0/2214	0.59	0/3030
1	B	0.43	0/2203	0.58	0/3016
1	C	0.42	0/2203	0.59	0/3016
1	D	0.43	0/2203	0.57	0/3016
1	E	0.43	0/2214	0.58	0/3030
1	F	0.43	0/2214	0.60	0/3030
All	All	0.43	0/13251	0.59	0/18138

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	A	2155	0	2103	51	0
1	B	2144	0	2091	39	0
1	C	2144	0	2091	38	0
1	D	2144	0	2091	43	0
1	E	2155	0	2103	49	0
1	F	2155	0	2103	61	0
2	A	4	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	4	0	3	0	0
2	C	4	0	3	1	0
2	D	4	0	3	0	0
2	E	4	0	3	0	0
2	F	4	0	3	0	0
3	A	5	0	0	3	0
3	B	5	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	2	0
3	E	10	0	0	1	0
4	A	10	0	8	0	0
4	B	10	0	8	3	0
4	C	10	0	8	0	0
4	E	10	0	8	0	0
4	F	10	0	8	0	0
5	A	9	0	9	1	0
5	B	6	0	6	2	0
5	C	3	0	3	0	0
5	D	6	0	6	3	0
5	E	3	0	3	0	0
5	F	9	0	9	2	0
6	A	193	0	0	11	0
6	B	182	0	0	10	0
6	C	191	0	0	13	0
6	D	137	0	0	9	0
6	E	152	0	0	11	0
6	F	159	0	0	18	0
All	All	14061	0	12676	271	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 11.

All (271) 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:170:ARG:NH1	1:A:171:ASN:OD1	2.03	0.90
1:D:178:ARG:HH12	1:D:185:LEU:HD21	1.39	0.88
1:C:167:SER:HA	1:C:170:ARG:HD3	1.57	0.86
1:F:208:ASN:ND2	6:F:2103:HOH:O	2.08	0.85
1:B:78:ASP:OD2	1:C:139:HIS:ND1	2.11	0.81
1:F:167:SER:HA	1:F:170:ARG:HD3	1.62	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:176:ASP:O	6:F:2112:HOH:O	1.99	0.78
1:C:20:TYR:HB2	1:C:61:MET:HE1	1.64	0.77
1:F:138:TYR:OH	1:F:178:ARG:NH2	2.18	0.77
1:C:184:ARG:NH1	6:C:2141:HOH:O	2.17	0.75
1:F:181:ALA:O	6:F:2116:HOH:O	2.05	0.74
1:E:213:VAL:O	6:E:2121:HOH:O	2.06	0.73
1:F:178:ARG:O	6:F:2114:HOH:O	2.08	0.71
1:B:9:ASN:H	5:F:1280:CCN:H22	1.55	0.70
1:E:208:ASN:ND2	6:E:2103:HOH:O	2.15	0.70
1:C:247:LYS:HD2	1:C:249:LYS:HE2	1.74	0.69
1:F:182:ALA:O	6:F:2084:HOH:O	2.11	0.69
1:A:247:LYS:HD2	1:A:249:LYS:HE2	1.76	0.68
1:E:188:PRO:O	6:E:2116:HOH:O	2.12	0.68
1:E:179:LEU:HB3	1:E:185:LEU:HD11	1.76	0.67
1:B:138:TYR:OH	1:B:178:ARG:NH2	2.27	0.67
1:A:247:LYS:NZ	3:A:1277:SO4:O3	2.28	0.66
1:A:183:GLN:HG3	1:A:211:LYS:HD2	1.77	0.66
1:B:167:SER:HA	1:B:170:ARG:HD3	1.76	0.66
1:B:186:VAL:HG12	6:B:2141:HOH:O	1.96	0.66
1:C:179:LEU:HB3	1:C:185:LEU:HD11	1.79	0.65
1:C:247:LYS:NZ	6:C:2167:HOH:O	2.29	0.65
1:F:219:HIS:HB2	6:F:2125:HOH:O	1.97	0.65
1:A:217:LEU:C	1:A:219:HIS:H	1.99	0.65
1:F:215:ALA:O	6:F:2125:HOH:O	2.14	0.65
1:C:170:ARG:NH1	1:C:171:ASN:OD1	2.30	0.64
1:A:276:TYR:N	6:A:2192:HOH:O	2.31	0.63
1:F:170:ARG:HB2	1:F:170:ARG:HH11	1.63	0.63
1:D:109:ARG:NH2	3:D:1278:SO4:O1	2.26	0.62
1:B:188:PRO:HD3	1:B:219:HIS:CD2	2.34	0.62
1:E:158:ASP:N	1:E:208:ASN:OD1	2.33	0.62
1:A:215:ALA:O	1:A:219:HIS:HB2	2.00	0.61
1:D:65:MET:HE1	1:D:111:ILE:HD12	1.82	0.61
1:D:188:PRO:HD3	1:D:219:HIS:CD2	2.35	0.61
1:A:247:LYS:NZ	3:A:1277:SO4:S	2.73	0.61
1:A:139:HIS:ND1	1:E:78:ASP:OD2	2.19	0.61
1:A:45:PHE:CD1	1:A:81:VAL:HG21	2.36	0.60
1:B:188:PRO:HD3	1:B:219:HIS:HD2	1.66	0.60
1:F:158:ASP:HA	1:F:181:ALA:HA	1.83	0.60
1:A:48:HIS:H	5:A:1281:CCN:H22	1.66	0.60
1:F:45:PHE:CD1	1:F:81:VAL:HG21	2.37	0.60
1:E:188:PRO:HD3	1:E:219:HIS:CD2	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:215:ALA:O	6:E:2122:HOH:O	2.17	0.60
1:A:87:GLN:NE2	1:A:139:HIS:HB2	2.17	0.59
1:F:185:LEU:N	6:F:2116:HOH:O	2.35	0.59
1:E:45:PHE:CD1	1:E:81:VAL:HG21	2.37	0.59
1:B:275:ASN:HD22	1:E:239:ASN:HB2	1.66	0.59
1:E:247:LYS:NZ	3:E:1278:SO4:S	2.75	0.59
1:F:176:ASP:HB3	6:F:2111:HOH:O	2.03	0.59
1:F:180:TRP:N	6:F:2112:HOH:O	2.36	0.58
1:F:109[A]:ARG:NH2	1:F:270:GLY:O	2.35	0.58
1:B:48:HIS:H	5:B:1280:CCN:H21	1.69	0.58
1:B:47:TYR:HB2	5:B:1280:CCN:H21	1.85	0.58
1:A:247:LYS:NZ	3:A:1277:SO4:O2	2.34	0.58
1:C:170:ARG:HB2	1:C:170:ARG:HH11	1.67	0.58
1:C:219:HIS:HB2	6:C:2144:HOH:O	2.04	0.58
1:F:214:GLN:HG2	1:F:218:GLN:OE1	2.05	0.57
1:B:275:ASN:HB3	1:E:240:HIS:H	1.71	0.56
1:D:186:VAL:HG13	6:D:2109:HOH:O	2.04	0.56
1:F:170:ARG:NH1	1:F:171:ASN:OD1	2.38	0.56
1:E:170:ARG:NH2	6:E:2108:HOH:O	2.30	0.56
1:A:186:VAL:HG12	6:A:2146:HOH:O	2.04	0.55
1:E:176:ASP:HB3	6:E:2110:HOH:O	2.06	0.55
1:C:137:ARG:HG2	1:C:138:TYR:CE2	2.41	0.55
1:E:130:THR:HG21	1:E:190:SER:HA	1.89	0.55
1:F:269:GLN:NE2	6:F:2154:HOH:O	2.40	0.55
1:F:176:ASP:OD2	1:F:178:ARG:NH2	2.39	0.55
1:B:45:PHE:CD1	1:B:81:VAL:HG21	2.42	0.55
1:F:158:ASP:N	1:F:208:ASN:OD1	2.40	0.54
1:A:116:SER:OG	1:A:233:HIS:NE2	2.40	0.54
1:D:247:LYS:NZ	3:D:1277:SO4:S	2.79	0.54
1:D:109:ARG:NE	6:D:2080:HOH:O	2.33	0.54
1:D:56:ILE:HD13	1:D:115:PHE:CZ	2.43	0.54
1:D:78:ASP:OD2	1:E:139:HIS:ND1	2.23	0.53
2:C:1276:ACT:H3	6:C:2058:HOH:O	2.07	0.53
6:A:2081:HOH:O	1:D:142:HIS:HB3	2.09	0.53
1:C:45:PHE:CD1	1:C:81:VAL:HG21	2.43	0.53
1:F:167:SER:O	1:F:170:ARG:NH1	2.39	0.52
1:B:5:GLN:NE2	6:B:2015:HOH:O	2.42	0.52
1:C:159:LEU:HD12	1:C:180:TRP:CD1	2.45	0.52
1:B:276:TYR:N	6:B:2181:HOH:O	2.42	0.52
1:A:264:ARG:HD3	6:A:2185:HOH:O	2.10	0.52
1:D:48:HIS:H	5:D:1279:CCN:C2	2.23	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:275:ASN:ND2	1:E:240:HIS:HD2	2.09	0.51
1:E:43:GLY:HA3	1:E:47:TYR:O	2.10	0.51
1:F:157:ILE:O	1:F:182:ALA:N	2.31	0.51
1:A:167:SER:HB2	1:A:170:ARG:NH2	2.26	0.51
1:D:212:TYR:O	1:D:216:MET:HG2	2.11	0.51
1:D:176:ASP:HB3	6:D:2101:HOH:O	2.10	0.50
1:C:197:GLN:NE2	1:C:204:VAL:O	2.44	0.50
1:C:216:MET:HA	6:C:2144:HOH:O	2.11	0.50
1:C:133:GLU:HG2	6:C:2119:HOH:O	2.12	0.50
1:C:132:PRO:HA	1:C:135:ARG:CB	2.42	0.50
1:C:176:ASP:OD1	1:C:178:ARG:HB3	2.12	0.50
1:D:159:LEU:N	6:D:2096:HOH:O	2.45	0.50
1:C:109:ARG:NE	6:C:2111:HOH:O	2.44	0.50
1:E:12:ALA:H	1:E:87:GLN:NE2	2.09	0.50
1:D:45:PHE:CD1	1:D:81:VAL:HG21	2.47	0.49
1:B:142:HIS:HB3	6:B:2122:HOH:O	2.11	0.49
1:D:75:ILE:HD13	1:D:172:GLN:HB3	1.93	0.49
1:F:170:ARG:HB2	1:F:170:ARG:NH1	2.27	0.49
1:C:20:TYR:CB	1:C:61:MET:HE1	2.39	0.49
1:E:217:LEU:HG	6:E:2121:HOH:O	2.11	0.49
1:A:197:GLN:NE2	1:A:204:VAL:O	2.46	0.49
1:E:188:PRO:HD3	1:E:219:HIS:HD2	1.76	0.49
1:F:56:ILE:HD13	1:F:115:PHE:CZ	2.48	0.49
1:B:56:ILE:HD13	1:B:115:PHE:CZ	2.47	0.49
1:F:48:HIS:H	5:F:1279:CCN:C2	2.25	0.49
1:D:249:LYS:NZ	6:D:2126:HOH:O	2.46	0.48
6:C:2060:HOH:O	1:D:27:ASP:OD1	2.20	0.48
1:A:154:TYR:CE2	1:A:234:GLY:HA2	2.48	0.48
1:F:43:GLY:HA3	1:F:47:TYR:O	2.13	0.48
1:E:213:VAL:HG21	1:E:225:TYR:CG	2.49	0.48
1:A:217:LEU:O	1:A:219:HIS:N	2.46	0.48
1:C:258:TRP:CG	1:C:259:PRO:HD3	2.49	0.48
1:F:188:PRO:HD3	1:F:219:HIS:CD2	2.49	0.48
1:F:136:THR:OG1	1:F:141:ASP:OD2	2.31	0.48
1:B:188:PRO:O	6:B:2142:HOH:O	2.20	0.47
1:E:154:TYR:CE2	1:E:234:GLY:HA2	2.49	0.47
1:D:35:PRO:HB3	1:D:109:ARG:HB3	1.96	0.47
1:F:116:SER:OG	1:F:233:HIS:NE2	2.44	0.47
1:B:275:ASN:ND2	1:E:239:ASN:HB2	2.29	0.47
1:D:205:PRO:HA	1:D:206:PRO:HD3	1.82	0.47
1:D:48:HIS:H	5:D:1279:CCN:H23	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1278:QY9:HAF	6:B:2030:HOH:O	2.14	0.47
1:F:75:ILE:HD13	1:F:172:GLN:HB3	1.96	0.47
1:B:197:GLN:NE2	1:B:204:VAL:O	2.47	0.47
1:B:27:ASP:N	1:B:27:ASP:OD1	2.46	0.47
1:F:157:ILE:HG23	1:F:182:ALA:HB3	1.96	0.47
1:F:106:ASP:OD2	1:F:109[A]:ARG:HD2	2.15	0.47
1:E:258:TRP:CG	1:E:259:PRO:HD3	2.50	0.47
1:E:35:PRO:HB3	1:E:109[A]:ARG:HB3	1.97	0.47
1:A:56:ILE:HD13	1:A:115:PHE:CZ	2.49	0.47
1:B:167:SER:O	1:B:170:ARG:NH1	2.46	0.46
1:F:116:SER:HB3	1:F:117:ALA:H	1.57	0.46
1:A:142:HIS:O	6:A:2125:HOH:O	2.21	0.46
1:F:159:LEU:N	1:F:180:TRP:O	2.43	0.46
4:B:1278:QY9:CAG	1:C:99:GLN:HE22	2.29	0.46
1:C:275:ASN:OD1	1:D:239:ASN:HB2	2.16	0.46
1:B:49:SER:O	1:B:53:GLU:HG3	2.15	0.46
1:A:1:MET:HE3	1:F:20:TYR:CD2	2.50	0.46
1:D:258:TRP:CG	1:D:259:PRO:HD3	2.50	0.46
1:B:176:ASP:OD2	1:B:178:ARG:NH2	2.49	0.46
1:C:168:ALA:O	1:C:172:GLN:HG3	2.15	0.46
1:E:1:MET:N	1:E:23:ASP:OD2	2.34	0.46
1:D:137:ARG:HH12	1:D:178:ARG:HD2	1.81	0.46
1:A:217:LEU:C	1:A:219:HIS:N	2.67	0.46
1:E:181:ALA:HB1	1:E:183:GLN:OE1	2.17	0.45
1:A:231:GLY:HA2	6:A:2162:HOH:O	2.16	0.45
1:D:158:ASP:HA	1:D:181:ALA:HA	1.99	0.45
6:C:2085:HOH:O	1:F:142:HIS:HB3	2.15	0.45
1:F:179:LEU:HA	6:F:2115:HOH:O	2.15	0.45
1:E:219:HIS:N	6:E:2122:HOH:O	2.48	0.45
1:E:163:PHE:HA	1:E:164:PRO:HA	1.79	0.45
1:B:133:GLU:OE2	6:B:2118:HOH:O	2.21	0.45
1:A:258:TRP:CG	1:A:259:PRO:HD3	2.51	0.45
1:D:217:LEU:C	1:D:219:HIS:H	2.20	0.45
1:B:163:PHE:HA	1:B:164:PRO:HA	1.81	0.45
1:C:167:SER:O	1:C:170:ARG:HB2	2.17	0.45
1:F:160:THR:N	6:F:2102:HOH:O	2.44	0.45
1:B:43:GLY:HA3	1:B:47:TYR:O	2.16	0.44
1:D:159:LEU:HD12	1:D:180:TRP:CD1	2.52	0.44
1:E:109[A]:ARG:NH2	1:E:270:GLY:O	2.50	0.44
1:F:165:THR:OG1	1:F:166:THR:N	2.50	0.44
1:E:176:ASP:OD2	1:E:178:ARG:NH2	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:201:ASP:OD1	1:E:233:HIS:HB2	2.17	0.44
1:A:106:ASP:OD2	1:A:109[A]:ARG:HD2	2.18	0.44
1:A:27:ASP:N	1:A:27:ASP:OD1	2.51	0.44
1:B:231:GLY:HA2	6:B:2154:HOH:O	2.16	0.44
1:E:87:GLN:OE1	1:E:139:HIS:HB2	2.17	0.44
1:C:135:ARG:HG3	1:C:140:LEU:HB2	2.00	0.44
1:A:212:TYR:O	1:A:216:MET:HG2	2.18	0.44
1:F:206:PRO:HD2	6:F:2059:HOH:O	2.16	0.44
6:B:2157:HOH:O	1:E:275:ASN:HB2	2.16	0.44
1:D:207:ILE:HA	1:D:210:LEU:HB2	2.00	0.44
1:D:170:ARG:NH2	6:D:2099:HOH:O	2.46	0.44
1:F:212:TYR:O	1:F:216:MET:HG2	2.18	0.44
1:C:276:TYR:N	6:C:2187:HOH:O	2.50	0.44
1:A:194:PHE:HB2	1:A:265:TRP:CE3	2.53	0.44
1:E:231:GLY:HA2	6:E:2127:HOH:O	2.17	0.44
1:C:79:GLN:HG3	6:C:2084:HOH:O	2.17	0.44
1:B:201:ASP:OD1	1:B:233:HIS:HB2	2.18	0.44
1:F:134:LEU:O	1:F:138:TYR:HD2	2.01	0.43
1:D:226:HIS:CE1	1:D:261:LEU:HD13	2.53	0.43
1:C:5:GLN:NE2	6:C:2013:HOH:O	2.51	0.43
1:C:132:PRO:HA	1:C:135:ARG:HB2	2.00	0.43
1:F:36:ILE:HD11	1:F:68:VAL:HG23	1.99	0.43
1:F:220:GLN:N	1:F:220:GLN:OE1	2.51	0.43
1:F:87:GLN:HA	1:F:140:LEU:CD2	2.48	0.43
4:B:1278:QY9:HAG	1:C:99:GLN:HE22	1.82	0.43
1:A:38:ILE:HD11	1:A:96:ILE:HD12	2.01	0.43
1:D:129:ALA:O	6:D:2085:HOH:O	2.21	0.43
1:C:132:PRO:HA	1:C:135:ARG:HB3	2.01	0.43
1:B:240:HIS:CD2	1:E:273:ALA:HB3	2.54	0.43
1:B:154:TYR:CE1	1:B:234:GLY:HA2	2.53	0.43
6:D:2012:HOH:O	1:E:6:ARG:NH1	2.36	0.43
1:E:223:THR:HG21	6:E:2121:HOH:O	2.17	0.43
1:F:154:TYR:CZ	1:F:234:GLY:HA2	2.53	0.43
1:F:132:PRO:HD3	6:F:2086:HOH:O	2.19	0.43
1:A:201:ASP:OD1	1:A:233:HIS:HB2	2.19	0.43
1:F:201:ASP:OD1	1:F:233:HIS:HB2	2.19	0.43
1:A:133:GLU:HG2	6:A:2116:HOH:O	2.18	0.43
1:B:212:TYR:O	1:B:216:MET:HG2	2.19	0.43
1:C:56:ILE:HD13	1:C:115:PHE:CZ	2.54	0.43
1:D:178:ARG:NH1	1:D:185:LEU:HD21	2.19	0.42
1:A:109[A]:ARG:NH2	1:A:270:GLY:O	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ARG:HG2	1:B:141:ASP:OD1	2.19	0.42
1:A:13:HIS:HE1	1:A:79:GLN:OE1	2.01	0.42
1:F:159:LEU:N	6:F:2102:HOH:O	2.52	0.42
1:E:158:ASP:HA	1:E:180:TRP:O	2.20	0.42
1:F:205:PRO:HA	1:F:206:PRO:HD3	1.75	0.42
1:C:38:ILE:HD11	1:C:96:ILE:HD12	2.01	0.42
1:A:135:ARG:HG2	1:A:141:ASP:OD1	2.20	0.42
1:D:154:TYR:CE2	1:D:234:GLY:HA2	2.54	0.42
1:A:175:THR:HG23	6:A:2086:HOH:O	2.19	0.42
1:F:178:ARG:HB2	1:F:178:ARG:HE	1.47	0.42
1:C:27:ASP:N	1:C:27:ASP:OD1	2.53	0.42
1:E:158:ASP:HA	1:E:181:ALA:HA	2.02	0.42
1:F:216:MET:HA	6:F:2125:HOH:O	2.19	0.42
1:E:12:ALA:H	1:E:87:GLN:HE21	1.67	0.42
1:D:116:SER:OG	1:D:233:HIS:NE2	2.42	0.42
1:A:20:TYR:CB	1:A:61:MET:HE1	2.49	0.42
1:A:137:ARG:NH2	6:A:2123:HOH:O	2.53	0.42
1:A:259:PRO:O	1:A:263:LEU:HG	2.20	0.41
1:F:154:TYR:CE2	1:F:234:GLY:HA2	2.55	0.41
1:E:116:SER:OG	1:E:233:HIS:NE2	2.47	0.41
1:B:258:TRP:CG	1:B:259:PRO:HD3	2.55	0.41
1:E:157:ILE:HG21	1:E:212:TYR:HB2	2.01	0.41
1:A:20:TYR:HB2	1:A:61:MET:HE1	2.02	0.41
1:D:131:GLN:O	1:D:135:ARG:N	2.49	0.41
1:E:235:LEU:HD12	1:E:235:LEU:HA	1.83	0.41
1:D:202:GLU:HB2	5:D:1280:CCN:C1	2.50	0.41
1:F:137:ARG:HD2	1:F:138:TYR:CE2	2.56	0.41
1:C:181:ALA:HB1	1:C:183:GLN:OE1	2.20	0.41
1:B:205:PRO:HA	1:B:206:PRO:HD3	1.88	0.41
1:A:181:ALA:HB1	1:A:183:GLN:OE1	2.20	0.41
1:D:188:PRO:HD3	1:D:219:HIS:HD2	1.82	0.41
1:D:158:ASP:OD1	1:D:160:THR:OG1	2.25	0.41
1:F:258:TRP:CG	1:F:259:PRO:HD3	2.55	0.41
1:E:219:HIS:HB2	6:E:2122:HOH:O	2.20	0.41
1:E:154:TYR:CZ	1:E:234:GLY:HA2	2.56	0.41
1:D:235:LEU:HD12	1:D:235:LEU:HA	1.96	0.41
1:D:184:ARG:NH2	6:D:2106:HOH:O	2.25	0.41
1:A:22:LEU:HD23	1:A:34:TYR:CG	2.56	0.41
1:B:135:ARG:HG3	1:B:140:LEU:HB2	2.02	0.41
1:A:108:GLN:NE2	6:A:2107:HOH:O	2.53	0.41
1:E:188:PRO:HB3	1:E:219:HIS:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:142:HIS:O	6:F:2094:HOH:O	2.22	0.41
1:A:135:ARG:HG3	1:A:140:LEU:HB2	2.03	0.41
1:A:158:ASP:HA	1:A:180:TRP:O	2.20	0.41
1:A:275:ASN:OD1	1:F:239:ASN:HB2	2.21	0.41
1:B:6:ARG:HG3	6:B:2010:HOH:O	2.20	0.41
1:C:167:SER:O	1:C:170:ARG:NH1	2.52	0.41
1:F:232:ILE:HD13	1:F:249:LYS:HG3	2.03	0.41
1:F:25:ILE:HD12	1:F:28:PHE:CZ	2.56	0.40
1:D:106:ASP:OD2	1:D:109:ARG:HD2	2.21	0.40
1:D:170:ARG:HG3	1:D:170:ARG:H	1.70	0.40
1:A:243:GLN:OE1	6:A:2171:HOH:O	2.22	0.40
1:A:179:LEU:HB3	1:A:185:LEU:HD11	2.02	0.40
1:B:20:TYR:CD2	1:E:1:MET:HE3	2.56	0.40
1:D:89:LEU:HD13	1:D:125:TYR:CG	2.57	0.40
1:B:178:ARG:HB2	1:B:178:ARG:HE	1.69	0.40
1:A:183:GLN:O	1:A:186:VAL:HG23	2.21	0.40
1:A:49:SER:O	1:A:53:GLU:HG3	2.22	0.40
1:F:65:MET:HE1	1:F:111:ILE:HD12	2.03	0.40
1:C:231:GLY:HA2	6:C:2157:HOH:O	2.21	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	276/282 (98%)	255 (92%)	20 (7%)	1 (0%)	39	48
1	B	275/282 (98%)	254 (92%)	21 (8%)	0	100	100
1	C	275/282 (98%)	251 (91%)	23 (8%)	1 (0%)	39	48
1	D	275/282 (98%)	254 (92%)	19 (7%)	2 (1%)	26	31
1	E	276/282 (98%)	256 (93%)	20 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	276/282 (98%)	255 (92%)	20 (7%)	1 (0%)	39	48
All	All	1653/1692 (98%)	1525 (92%)	123 (7%)	5 (0%)	46	57

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	218	GLN
1	D	165	THR
1	C	165	THR
1	D	218	GLN
1	F	165	THR

### 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	221/226 (98%)	205 (93%)	16 (7%)	18	22
1	B	220/226 (97%)	207 (94%)	13 (6%)	24	32
1	C	220/226 (97%)	204 (93%)	16 (7%)	17	22
1	D	220/226 (97%)	205 (93%)	15 (7%)	20	25
1	E	221/226 (98%)	207 (94%)	14 (6%)	22	29
1	F	221/226 (98%)	209 (95%)	12 (5%)	27	36
All	All	1323/1356 (98%)	1237 (94%)	86 (6%)	21	27

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	22	LEU
1	A	27	ASP
1	A	32	VAL
1	A	76	VAL
1	A	101	SER

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Mol	Chain	Res	Type
1	A	116	SER
1	A	167	SER
1	A	179	LEU
1	A	184	ARG
1	A	186	VAL
1	A	207	ILE
1	A	210	LEU
1	A	235	LEU
1	A	248	ASP
1	A	272	LEU
1	B	6	ARG
1	B	16	GLN
1	B	22	LEU
1	B	27	ASP
1	B	32	VAL
1	B	65	MET
1	B	76	VAL
1	B	170	ARG
1	B	207	ILE
1	B	210	LEU
1	B	235	LEU
1	B	264	ARG
1	B	272	LEU
1	C	1	MET
1	C	6	ARG
1	C	22	LEU
1	C	32	VAL
1	C	76	VAL
1	C	80[A]	SER
1	C	80[B]	SER
1	C	167	SER
1	C	170	ARG
1	C	179	LEU
1	C	184	ARG
1	C	207	ILE
1	C	210	LEU
1	C	235	LEU
1	C	264	ARG
1	C	272	LEU
1	D	6	ARG
1	D	16	GLN
1	D	22	LEU

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Mol	Chain	Res	Type
1	D	32	VAL
1	D	78	ASP
1	D	137	ARG
1	D	167	SER
1	D	170	ARG
1	D	178	ARG
1	D	179	LEU
1	D	203	SER
1	D	214	GLN
1	D	235	LEU
1	D	264	ARG
1	D	272	LEU
1	E	6	ARG
1	E	22	LEU
1	E	30	THR
1	E	32	VAL
1	E	65	MET
1	E	76	VAL
1	E	101	SER
1	E	179	LEU
1	E	184	ARG
1	E	203	SER
1	E	210	LEU
1	E	235	LEU
1	E	272	LEU
1	E	275	ASN
1	F	6	ARG
1	F	22	LEU
1	F	32	VAL
1	F	76	VAL
1	F	116	SER
1	F	167	SER
1	F	170	ARG
1	F	179	LEU
1	F	210	LEU
1	F	235	LEU
1	F	264	ARG
1	F	272	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:



Mol	Chain	Res	Type
1	C	99	GLN
1	E	226	HIS
1	E	240	HIS
1	F	275	ASN

### 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](#)

31 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	ACT	A	1276	-	1,3,3	1.14	0	0,3,3	0.00	-
3	SO4	A	1277	-	4,4,4	0.30	0	6,6,6	0.15	0
4	QY9	A	1278	-	10,10,10	0.70	0	12,12,12	1.51	1 (8%)
5	CCN	A	1279	-	2,2,2	0.64	0	1,1,1	0.61	0
5	CCN	A	1280	-	2,2,2	0.72	0	1,1,1	0.58	0
5	CCN	A	1281	-	2,2,2	0.67	0	1,1,1	0.62	0
2	ACT	B	1276	-	1,3,3	1.36	0	0,3,3	0.00	-
3	SO4	B	1277	-	4,4,4	0.29	0	6,6,6	0.09	0
4	QY9	B	1278	-	10,10,10	0.65	0	12,12,12	1.34	1 (8%)
5	CCN	B	1279	-	2,2,2	0.62	0	1,1,1	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	CCN	B	1280	-	2,2,2	0.65	0	1,1,1	0.60	0
2	ACT	C	1276	-	1,3,3	0.72	0	0,3,3	0.00	-
3	SO4	C	1277	-	4,4,4	0.25	0	6,6,6	0.24	0
3	SO4	C	1278	-	4,4,4	0.25	0	6,6,6	0.14	0
4	QY9	C	1279	-	10,10,10	0.66	0	12,12,12	1.35	1 (8%)
5	CCN	C	1280	-	2,2,2	0.69	0	1,1,1	0.59	0
2	ACT	D	1276	-	1,3,3	0.74	0	0,3,3	0.00	-
3	SO4	D	1277	-	4,4,4	0.27	0	6,6,6	0.20	0
3	SO4	D	1278	-	4,4,4	0.22	0	6,6,6	0.12	0
5	CCN	D	1279	-	2,2,2	0.67	0	1,1,1	0.60	0
5	CCN	D	1280	-	2,2,2	0.66	0	1,1,1	0.54	0
2	ACT	E	1276	-	1,3,3	1.20	0	0,3,3	0.00	-
3	SO4	E	1277	-	4,4,4	0.33	0	6,6,6	0.37	0
3	SO4	E	1278	-	4,4,4	0.28	0	6,6,6	0.11	0
4	QY9	E	1279	-	10,10,10	0.69	0	12,12,12	1.32	1 (8%)
5	CCN	E	1280	-	2,2,2	0.59	0	1,1,1	0.56	0
2	ACT	F	1276	-	1,3,3	1.14	0	0,3,3	0.00	-
4	QY9	F	1277	-	10,10,10	0.67	0	12,12,12	1.40	1 (8%)
5	CCN	F	1278	-	2,2,2	0.62	0	1,1,1	0.57	0
5	CCN	F	1279	-	2,2,2	0.72	0	1,1,1	0.57	0
5	CCN	F	1280	-	2,2,2	0.75	0	1,1,1	0.51	0

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	ACT	A	1276	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1277	-	-	0/0/0/0	0/0/0/0
4	QY9	A	1278	-	-	0/4/4/4	0/1/1/1
5	CCN	A	1279	-	-	0/0/0/0	0/0/0/0
5	CCN	A	1280	-	-	0/0/0/0	0/0/0/0
5	CCN	A	1281	-	-	0/0/0/0	0/0/0/0
2	ACT	B	1276	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1277	-	-	0/0/0/0	0/0/0/0
4	QY9	B	1278	-	-	0/4/4/4	0/1/1/1
5	CCN	B	1279	-	-	0/0/0/0	0/0/0/0
5	CCN	B	1280	-	-	0/0/0/0	0/0/0/0
2	ACT	C	1276	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1277	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1278	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	QY9	C	1279	-	-	0/4/4/4	0/1/1/1
5	CCN	C	1280	-	-	0/0/0/0	0/0/0/0
2	ACT	D	1276	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1277	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1278	-	-	0/0/0/0	0/0/0/0
5	CCN	D	1279	-	-	0/0/0/0	0/0/0/0
5	CCN	D	1280	-	-	0/0/0/0	0/0/0/0
2	ACT	E	1276	-	-	0/0/0/0	0/0/0/0
3	SO4	E	1277	-	-	0/0/0/0	0/0/0/0
3	SO4	E	1278	-	-	0/0/0/0	0/0/0/0
4	QY9	E	1279	-	-	0/4/4/4	0/1/1/1
5	CCN	E	1280	-	-	0/0/0/0	0/0/0/0
2	ACT	F	1276	-	-	0/0/0/0	0/0/0/0
4	QY9	F	1277	-	-	0/4/4/4	0/1/1/1
5	CCN	F	1278	-	-	0/0/0/0	0/0/0/0
5	CCN	F	1279	-	-	0/0/0/0	0/0/0/0
5	CCN	F	1280	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1278	QY9	OAH-CAI-CAA	3.59	118.74	110.90
4	E	1279	QY9	OAH-CAI-CAA	3.75	119.10	110.90
4	C	1279	QY9	OAH-CAI-CAA	4.05	119.74	110.90
4	F	1277	QY9	OAH-CAI-CAA	4.08	119.81	110.90
4	A	1278	QY9	OAH-CAI-CAA	4.42	120.54	110.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1277	SO4	3	0
5	A	1281	CCN	1	0
4	B	1278	QY9	3	0
5	B	1280	CCN	2	0
2	C	1276	ACT	1	0
3	D	1277	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1278	SO4	1	0
5	D	1279	CCN	2	0
5	D	1280	CCN	1	0
3	E	1278	SO4	1	0
5	F	1279	CCN	1	0
5	F	1280	CCN	1	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	276/282 (97%)	0.99	41 (14%) 3 5	14, 37, 72, 95	0
1	B	276/282 (97%)	1.04	46 (16%) 2 4	13, 37, 74, 104	0
1	C	276/282 (97%)	0.93	39 (14%) 4 6	13, 38, 73, 101	0
1	D	276/282 (97%)	0.89	29 (10%) 8 12	14, 36, 73, 105	0
1	E	276/282 (97%)	0.97	40 (14%) 3 5	14, 38, 75, 91	0
1	F	276/282 (97%)	0.91	35 (12%) 5 8	15, 37, 71, 87	0
All	All	1656/1692 (97%)	0.95	230 (13%) 4 6	13, 37, 73, 105	0

All (230) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	276	TYR	9.7
1	A	175	THR	6.4
1	B	161	ALA	5.9
1	F	168	ALA	5.6
1	C	159	LEU	5.1
1	E	172	GLN	5.1
1	B	160	THR	4.8
1	E	175	THR	4.8
1	B	140	LEU	4.7
1	F	159	LEU	4.7
1	C	31	ALA	4.6
1	A	188	PRO	4.3
1	E	168	ALA	4.3
1	A	204	VAL	4.2
1	F	134	LEU	4.2
1	C	202	GLU	4.0
1	C	180	TRP	3.9
1	F	204	VAL	3.9
1	D	173	ILE	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	166	THR	3.8
1	B	76	VAL	3.8
1	D	161	ALA	3.7
1	C	134	LEU	3.7
1	C	210	LEU	3.7
1	D	188	PRO	3.6
1	E	161	ALA	3.6
1	B	29	GLU	3.6
1	F	205	PRO	3.6
1	A	128	VAL	3.5
1	D	175	THR	3.5
1	A	80[A]	SER	3.5
1	C	175	THR	3.5
1	E	180	TRP	3.5
1	E	141	ASP	3.5
1	B	136	THR	3.4
1	D	140	LEU	3.4
1	E	30	THR	3.4
1	A	134	LEU	3.4
1	B	134	LEU	3.4
1	A	168	ALA	3.3
1	E	169	ALA	3.3
1	F	80[A]	SER	3.3
1	A	189	ALA	3.3
1	E	173	ILE	3.3
1	C	165	THR	3.3
1	B	78	ASP	3.3
1	C	204	VAL	3.2
1	F	75	ILE	3.2
1	E	159	LEU	3.2
1	E	78	ASP	3.2
1	D	78	ASP	3.2
1	E	166	THR	3.2
1	D	75	ILE	3.2
1	E	163	PHE	3.2
1	D	177	ALA	3.1
1	F	138	TYR	3.1
1	F	221	VAL	3.1
1	A	160	THR	3.1
1	B	166	THR	3.1
1	F	133	GLU	3.1
1	B	175	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	160	THR	3.1
1	C	172	GLN	3.1
1	A	161	ALA	3.1
1	D	170	ARG	3.1
1	B	180	TRP	3.1
1	D	220	GLN	3.0
1	B	221	VAL	3.0
1	C	168	ALA	3.0
1	D	136	THR	3.0
1	C	186	VAL	3.0
1	E	204	VAL	3.0
1	D	276	TYR	3.0
1	B	182	ALA	3.0
1	A	75	ILE	3.0
1	C	174	THR	3.0
1	F	180	TRP	3.0
1	D	186	VAL	3.0
1	D	168	ALA	2.9
1	E	130	THR	2.9
1	E	76	VAL	2.9
1	B	275	ASN	2.9
1	E	158	ASP	2.9
1	A	174	THR	2.9
1	D	218	GLN	2.9
1	E	75	ILE	2.9
1	F	163	PHE	2.9
1	D	221	VAL	2.9
1	B	130	THR	2.9
1	F	166	THR	2.9
1	D	160	THR	2.9
1	D	169	ALA	2.8
1	F	177	ALA	2.8
1	E	174	THR	2.8
1	A	158	ASP	2.8
1	B	215	ALA	2.8
1	C	169	ALA	2.8
1	C	76	VAL	2.8
1	A	177	ALA	2.8
1	C	162	GLY	2.8
1	D	77	GLY	2.8
1	F	188	PRO	2.8
1	F	140	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	169	ALA	2.8
1	D	203	SER	2.8
1	B	217	LEU	2.8
1	C	173	ILE	2.8
1	B	185	LEU	2.8
1	A	165	THR	2.8
1	F	172	GLN	2.7
1	C	185	LEU	2.7
1	C	205	PRO	2.7
1	A	29	GLU	2.7
1	D	130	THR	2.7
1	A	159	LEU	2.7
1	D	199	ALA	2.7
1	A	78	ASP	2.7
1	A	194	PHE	2.7
1	C	161	ALA	2.7
1	F	175	THR	2.7
1	F	185	LEU	2.6
1	A	166	THR	2.6
1	F	187	THR	2.6
1	B	213	VAL	2.6
1	B	165	THR	2.6
1	F	165	THR	2.6
1	B	131	GLN	2.6
1	E	190	SER	2.6
1	F	173	ILE	2.6
1	C	160	THR	2.6
1	B	75	ILE	2.6
1	B	188	PRO	2.5
1	C	158	ASP	2.5
1	F	167	SER	2.5
1	F	182	ALA	2.5
1	E	185	LEU	2.5
1	F	202	GLU	2.5
1	B	207	ILE	2.5
1	F	275	ASN	2.5
1	E	177	ALA	2.5
1	B	187	THR	2.5
1	A	180	TRP	2.5
1	C	130	THR	2.5
1	A	221	VAL	2.5
1	A	220	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	210	LEU	2.5
1	F	207	ILE	2.5
1	B	172	GLN	2.5
1	C	219	HIS	2.4
1	A	182	ALA	2.4
1	E	199	ALA	2.4
1	B	174	THR	2.4
1	D	141	ASP	2.4
1	C	142	HIS	2.4
1	E	214	GLN	2.4
1	B	163	PHE	2.4
1	B	169	ALA	2.4
1	E	245	PRO	2.4
1	B	173	ILE	2.4
1	E	232	ILE	2.4
1	C	29	GLU	2.4
1	F	136	THR	2.4
1	F	215	ALA	2.4
1	C	78	ASP	2.4
1	B	138	TYR	2.4
1	E	165	THR	2.4
1	A	140	LEU	2.3
1	A	217	LEU	2.3
1	B	133	GLU	2.3
1	F	179	LEU	2.3
1	D	138	TYR	2.3
1	A	170	ARG	2.3
1	E	220	GLN	2.3
1	F	181	ALA	2.3
1	A	130	THR	2.3
1	F	146	GLN	2.3
1	B	202	GLU	2.3
1	D	134	LEU	2.3
1	D	178	ARG	2.3
1	C	188	PRO	2.2
1	F	274	GLY	2.2
1	A	141	ASP	2.2
1	E	176	ASP	2.2
1	A	218	GLN	2.2
1	B	127	GLY	2.2
1	B	26	SER	2.2
1	C	276	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	163	PHE	2.2
1	A	268	GLU	2.2
1	B	210	LEU	2.2
1	C	232	ILE	2.2
1	B	186	VAL	2.2
1	E	77	GLY	2.2
1	E	127	GLY	2.2
1	E	162	GLY	2.2
1	E	134	LEU	2.2
1	F	210	LEU	2.2
1	A	163	PHE	2.1
1	D	167	SER	2.1
1	B	211	LYS	2.1
1	E	215	ALA	2.1
1	A	127	GLY	2.1
1	A	131	GLN	2.1
1	C	235	LEU	2.1
1	E	179	LEU	2.1
1	B	85	ALA	2.1
1	C	181	ALA	2.1
1	E	216	MET	2.1
1	D	181	ALA	2.1
1	F	158	ASP	2.1
1	B	31	ALA	2.1
1	B	177	ALA	2.1
1	C	221	VAL	2.1
1	A	171	ASN	2.1
1	A	167	SER	2.1
1	C	213	VAL	2.1
1	B	179	LEU	2.0
1	B	181	ALA	2.0
1	A	265	TRP	2.0
1	A	173	ILE	2.0
1	D	166	THR	2.0
1	C	179	LEU	2.0
1	C	182	ALA	2.0
1	E	80[A]	SER	2.0
1	C	75	ILE	2.0
1	B	269	GLN	2.0
1	E	27	ASP	2.0
1	A	124	THR	2.0
1	A	185	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	167	SER	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
4	QY9	B	1278	10/10	0.67	0.37	6.79	73,77,85,85	0
4	QY9	F	1277	10/10	0.68	0.30	5.19	70,74,81,84	0
4	QY9	A	1278	10/10	0.59	0.33	4.82	59,64,71,73	0
4	QY9	C	1279	10/10	0.70	0.33	4.76	75,78,87,90	0
4	QY9	E	1279	10/10	0.76	0.29	2.34	67,72,73,74	0
5	CCN	F	1280	3/3	0.73	0.21	1.76	19,19,30,34	0
2	ACT	D	1276	4/4	0.78	0.26	1.54	19,27,33,47	0
2	ACT	E	1276	4/4	0.88	0.20	1.42	28,31,37,50	0
5	CCN	E	1280	3/3	0.80	0.25	0.67	34,34,34,37	0
2	ACT	F	1276	4/4	0.82	0.19	0.60	32,35,40,51	0
5	CCN	A	1279	3/3	0.86	0.20	-0.03	35,35,37,40	0
5	CCN	F	1278	3/3	0.83	0.19	-0.12	34,34,36,38	0
2	ACT	C	1276	4/4	0.85	0.17	-0.21	31,39,41,55	0
5	CCN	C	1280	3/3	0.86	0.19	-0.46	38,38,39,42	0
5	CCN	B	1279	3/3	0.86	0.16	-0.65	34,34,35,39	0
3	SO4	C	1277	5/5	0.95	0.14	-0.77	39,40,46,59	0
2	ACT	B	1276	4/4	0.85	0.15	-0.80	38,39,45,53	0
2	ACT	A	1276	4/4	0.93	0.16	-0.98	31,32,40,54	0
5	CCN	D	1280	3/3	0.86	0.15	-1.04	34,34,37,38	0
3	SO4	E	1277	5/5	0.96	0.15	-1.05	37,44,47,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	CCN	B	1280	3/3	0.88	0.18	-	34,34,34,35	0
5	CCN	A	1280	3/3	0.94	0.10	-	32,32,36,36	0
3	SO4	B	1277	5/5	0.66	0.24	-	111,111,112,114	0
3	SO4	D	1278	5/5	0.80	0.15	-	108,109,110,110	0
5	CCN	D	1279	3/3	0.85	0.24	-	30,30,33,36	0
3	SO4	C	1278	5/5	0.60	0.27	-	111,111,113,114	0
5	CCN	F	1279	3/3	0.95	0.12	-	13,13,26,28	0
5	CCN	A	1281	3/3	0.84	0.20	-	26,26,27,28	0
3	SO4	D	1277	5/5	0.77	0.19	-	110,110,112,112	0
3	SO4	A	1277	5/5	0.77	0.22	-	105,106,107,108	0
3	SO4	E	1278	5/5	0.83	0.29	-	116,119,120,120	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.