



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 07:20 pm GMT

PDB ID : 4U3F  
Title : Cytochrome bc1 complex from chicken with designed inhibitor bound  
Authors : Huang, L.-S.; Zhu, X.-L.; Yang, G.F.; Berry, E.A.  
Deposited on : 2014-07-21  
Resolution : 3.23 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

<https://www.wwpdb.org/validation/2017/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 : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : 1.13  
EDS : trunk30967  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30967

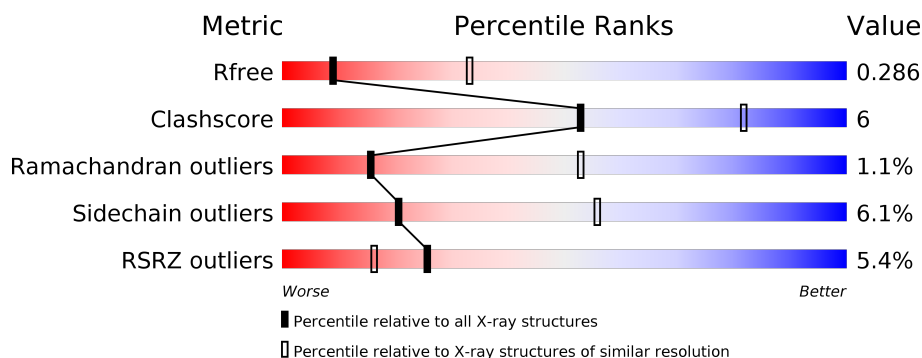
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.23 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))
$R_{free}$	111664	1379 (3.28-3.20)
Clashscore	122126	1510 (3.28-3.20)
Ramachandran outliers	120053	1485 (3.28-3.20)
Sidechain outliers	120020	1484 (3.28-3.20)
RSRZ outliers	108989	1331 (3.28-3.20)

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	446	<div> <div>2%</div> <div>81%16%..</div> </div>
1	N	446	<div> <div>2%</div> <div>81%16%..</div> </div>
2	B	441	<div> <div>2%</div> <div>72%21%5%</div> </div>
2	O	441	<div> <div>2%</div> <div>73%20%..</div> </div>
3	C	380	<div> <div>%</div> <div>84%14%. .</div> </div>
3	P	380	<div> <div>%</div> <div>83%15%. .</div> </div>

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Mol	Chain	Length	Quality of chain
4	D	241	
4	Q	241	
5	E	196	
5	R	196	
6	F	110	
6	S	110	
7	G	81	
7	T	81	
8	H	77	
8	U	77	
9	I	76	
9	V	76	
10	J	61	
10	W	61	

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
11	PEE	P	411	-	-	-	X
3	FME	C	1	-	-	-	X

## 2 Entry composition

There are 21 unique types of molecules in this entry. The entry contains 32790 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 Mitochondrial ubiquinol-cytochrome-c reductase complex core protein i.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	0	0
			3442	2157	606	658	21			
1	N	442	Total	C	N	O	S	0	0	0
			3437	2154	605	657	21			

- Molecule 2 is a protein called Mitochondrial ubiquinol-cytochrome-c reductase complex core protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	421	Total	C	N	O	S	0	0	0
			3137	1971	544	613	9			
2	O	422	Total	C	N	O	S	0	0	0
			3143	1974	545	614	10			

- Molecule 3 is a protein called Cytochrome b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	380	Total	C	N	O	S	0	0	0
			3021	2025	478	505	13			
3	P	379	Total	C	N	O	S	0	1	0
			3022	2025	480	505	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	FME	-	expression tag	UNP P18946
P	1	FME	-	expression tag	UNP P18946

- Molecule 4 is a protein called Mitochondrial cytochrome c1, heme protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	241	Total	C	N	O	S	0	0	0
			1898	1212	327	347	12			
4	Q	241	Total	C	N	O	S	0	0	0
			1898	1212	327	347	12			

- Molecule 5 is a protein called Cytochrome b-c1 complex subunit Rieske, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	196	Total	C	N	O	S	0	0	0
			1513	952	263	292	6			
5	R	196	Total	C	N	O	S	0	0	0
			1509	950	263	290	6			

- Molecule 6 is a protein called Cytochrome b-c1 complex subunit 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	101	Total	C	N	O	S	0	0	0
			891	570	159	159	3			
6	S	101	Total	C	N	O	S	0	0	0
			891	570	159	159	3			

- Molecule 7 is a protein called Mitochondrial ubiquinol-cytochrome c reductase ubiquinone-binding protein qp-c.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	G	80	Total	C	N	O	0	0	0
			672	437	119	116			
7	T	79	Total	C	N	O	0	0	0
			662	432	117	113			

- Molecule 8 is a protein called Cytochrome b-c1 complex subunit 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	68	Total	C	N	O	S	0	0	0
			562	343	104	110	5			
8	U	68	Total	C	N	O	S	0	0	0
			558	341	104	108	5			

- Molecule 9 is a protein called Cytochrome b-c1 complex subunit Rieske, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	44	Total	C	N	O	S	0	0	1
			266	157	54	53	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	V	42	Total	C	N	O	S	0	0	1
			265	157	55	51	2			

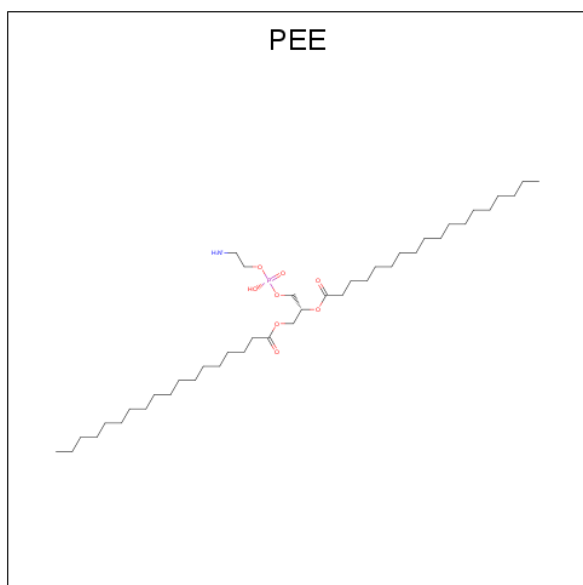
There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	1	AME	-	expression tag	UNP Q5ZLR5
I	29	UNK	LEU	see remark 999	UNP Q5ZLR5
I	30	UNK	ALA	see remark 999	UNP Q5ZLR5
I	31	UNK	PRO	see remark 999	UNP Q5ZLR5
I	32	UNK	ALA	see remark 999	UNP Q5ZLR5
I	33	UNK	ALA	see remark 999	UNP Q5ZLR5
I	37	UNK	LEU	see remark 999	UNP Q5ZLR5
I	38	UNK	ARG	see remark 999	UNP Q5ZLR5
I	39	UNK	ALA	see remark 999	UNP Q5ZLR5
I	40	UNK	GLU	see remark 999	UNP Q5ZLR5
I	41	UNK	LYS	see remark 999	UNP Q5ZLR5
I	42	UNK	VAL	see remark 999	UNP Q5ZLR5
I	43	UNK	VAL	see remark 999	UNP Q5ZLR5
I	44	UNK	LEU	see remark 999	UNP Q5ZLR5
I	45	UNK	ASP	see remark 999	UNP Q5ZLR5
V	-1	AME	-	expression tag	UNP Q5ZLR5
V	27	UNK	LEU	see remark 999	UNP Q5ZLR5
V	28	UNK	ALA	see remark 999	UNP Q5ZLR5
V	29	UNK	PRO	see remark 999	UNP Q5ZLR5
V	30	UNK	ALA	see remark 999	UNP Q5ZLR5
V	31	UNK	ALA	see remark 999	UNP Q5ZLR5
V	32	UNK	LEU	see remark 999	UNP Q5ZLR5
V	33	UNK	ARG	see remark 999	UNP Q5ZLR5
V	37	UNK	ALA	see remark 999	UNP Q5ZLR5
V	38	UNK	GLU	see remark 999	UNP Q5ZLR5
V	39	UNK	LYS	see remark 999	UNP Q5ZLR5
V	40	UNK	VAL	see remark 999	UNP Q5ZLR5
V	41	UNK	VAL	see remark 999	UNP Q5ZLR5
V	42	UNK	LEU	see remark 999	UNP Q5ZLR5
V	43	UNK	ASP	see remark 999	UNP Q5ZLR5

- Molecule 10 is a protein called Mitochondrial ubiquinol-cytochrome c reductase 7.2 kda protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
10	J	61	Total	C	N	O	0	0	0
			497	321	87	89			
10	W	60	Total	C	N	O	0	0	1
			479	311	86	82			

- Molecule 11 is 1,2-Dioleoyl-sn-glycero-3-phosphoethanolamine (three-letter code: PEE) (formula: C<sub>41</sub>H<sub>83</sub>NO<sub>8</sub>P).



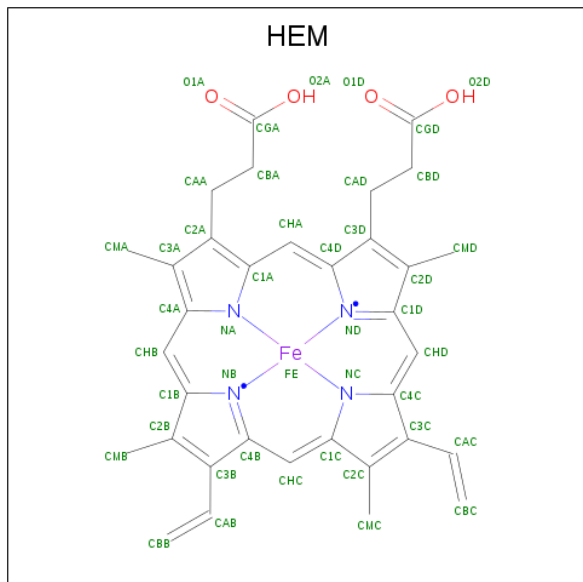
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	A	1	Total	C	N	O	P	0	0
			26	16	1	8	1		
11	C	1	Total	C	N	O	P	0	0
			49	39	1	8	1		
11	C	1	Total	C	N	O	P	0	0
			11	5	1	4	1		
11	C	1	Total	C	N	O	P	0	0
			47	37	1	8	1		
11	C	1	Total	C	O			0	0
			15	13	2				
11	E	1	Total	C	N	O	P	0	0
			48	38	1	8	1		
11	E	1	Total	C				0	0
			15	15					
11	N	1	Total	C	N	O	P	0	0
			8	2	1	4	1		
11	P	1	Total	C	O			0	0
			15	13	2				

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	P	1	Total	C	N	O	P	0	0
			49	39	1	8	1		
11	P	1	Total	C	N	O	P	0	0
			41	31	1	8	1		
11	P	1	Total	C	O			0	0
			12	10	2				
11	P	1	Total	C	N	O	P	0	0
			25	15	1	8	1		
11	R	1	Total	C	N	O	P	0	0
			49	39	1	8	1		

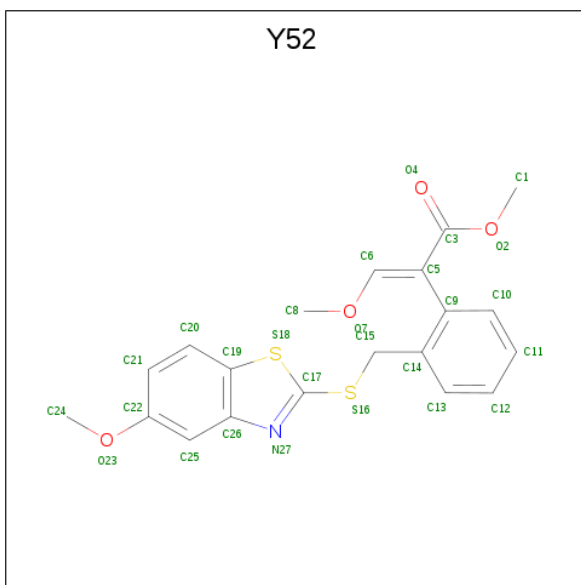
- Molecule 12 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
12	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
12	P	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
12	P	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

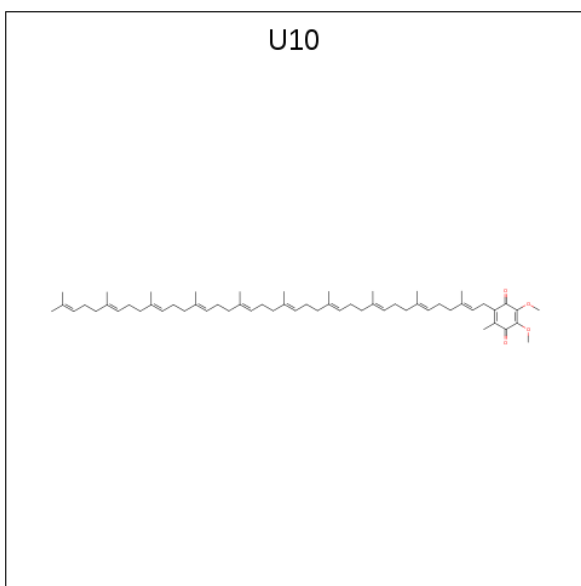
- Molecule 13 is methyl (2E)-3-methoxy-2-(2-[(5-methoxy-1,3-benzothiazol-2-yl)sulfanyl]methyl}phenyl)prop-2-enoate (three-letter code: Y52) (formula:  $C_{20}H_{19}NO_4S_2$ ).





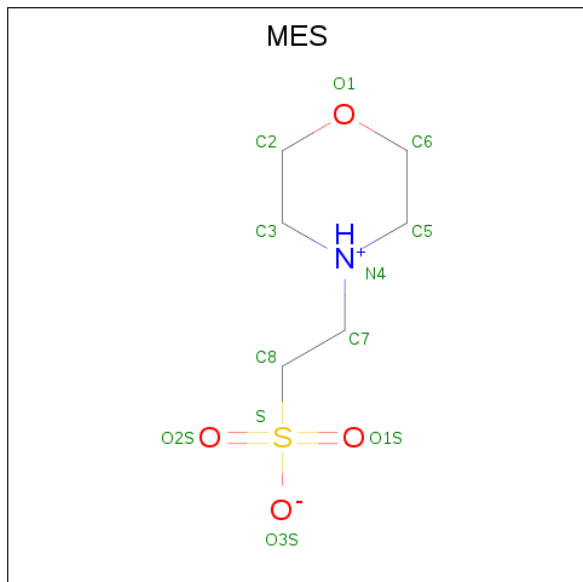
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
13	C	1	Total 27	C 20	N 1	O 4	S 2	0	0
13	P	1	Total 27	C 20	N 1	O 4	S 2	0	0

- Molecule 14 is UBIQUINONE-10 (three-letter code: U10) (formula:  $C_{59}H_{90}O_4$ ).



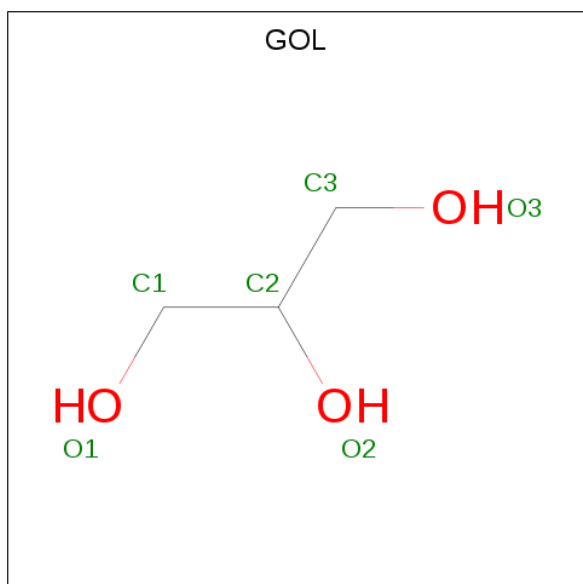
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
14	C	1	Total	C	O	0	0
			19	15	4		
14	P	1	Total	C	O	0	0
			19	15	4		

- Molecule 15 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula:  $C_6H_{13}NO_4S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
15	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 16 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



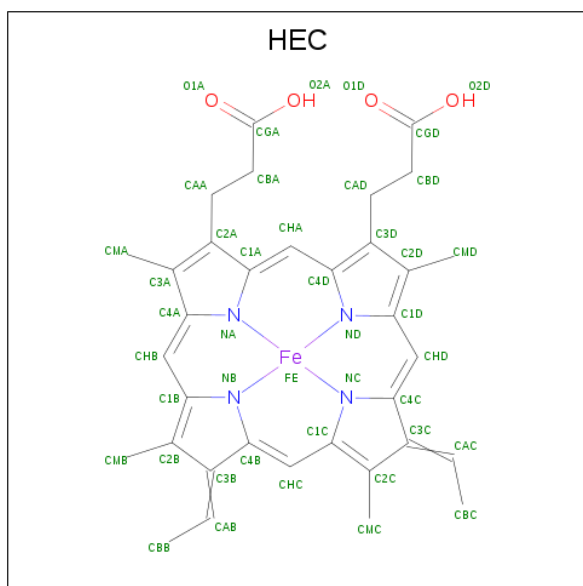
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
16	C	1	Total	C	O	0	0
			6	3	3		

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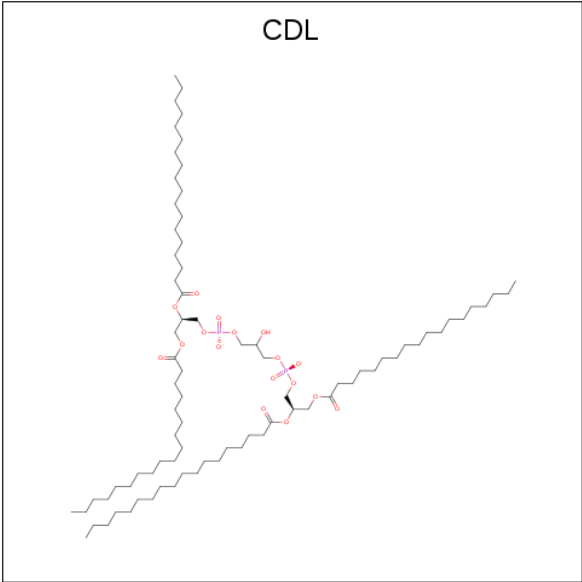
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
16	P	1	Total	C	O	0	0
			6	3	3		

- Molecule 17 is HEME C (three-letter code: HEC) (formula:  $\text{C}_{34}\text{H}_{34}\text{FeN}_4\text{O}_4$ ).



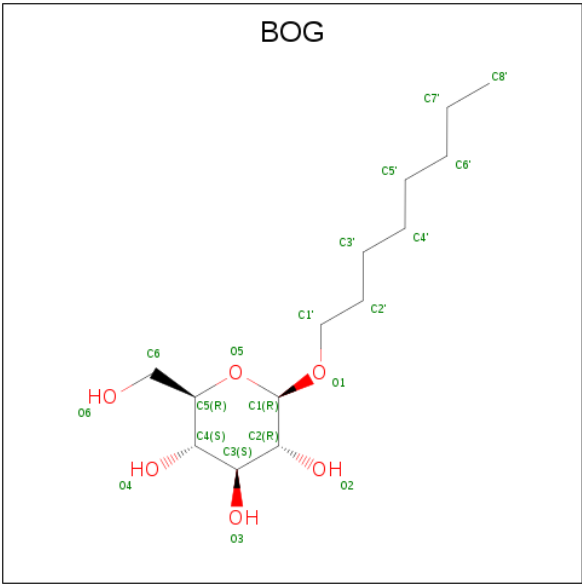
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
17	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
17	Q	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 18 is CARDIOLIPIN (three-letter code: CDL) (formula:  $\text{C}_{81}\text{H}_{156}\text{O}_{17}\text{P}_2$ ).



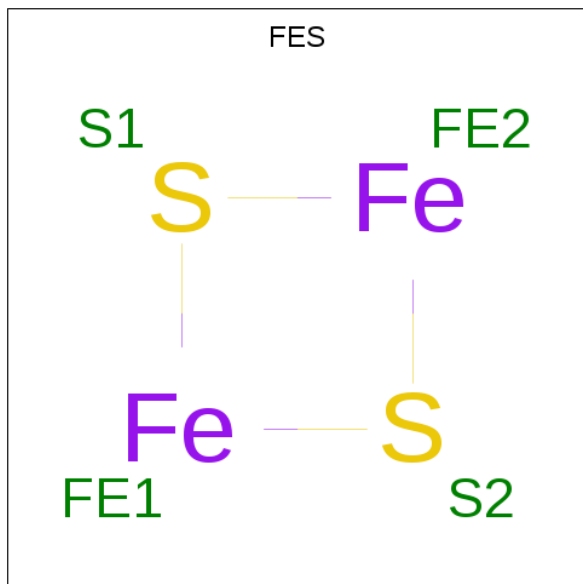
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
18	D	1	Total	C	O	P	0	0
			42	23	17	2		
18	G	1	Total	C	O	P	0	0
			40	21	17	2		
18	Q	1	Total	C	O	P	0	0
			42	23	17	2		
18	T	1	Total	C	O	P	0	0
			40	21	17	2		

- Molecule 19 is B-OCTYLGLUCOSIDE (three-letter code: BOG) (formula: C<sub>14</sub>H<sub>28</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
19	D	1	Total	C	O	0	0
			20	14	6		
19	P	1	Total	C	O	0	0
			20	14	6		
19	Q	1	Total	C	O	0	0
			20	14	6		

- Molecule 20 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe<sub>2</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
20	E	1	Total	Fe	S	0	0
			4	2	2		
20	R	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 21 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
21	A	2	Total	O	0	0
			2	2		
21	B	1	Total	O	0	0
			1	1		
21	C	2	Total	O	0	0
			2	2		
21	F	1	Total	O	0	0
			1	1		
21	N	1	Total	O	0	0
			1	1		

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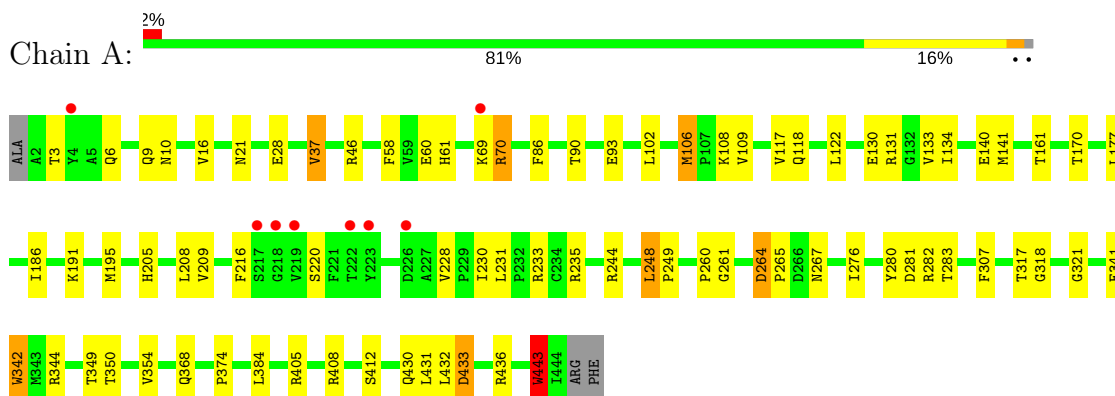
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
21	O	2	Total	O	0	0
			2	2		
21	P	2	Total	O	0	0
			2	2		

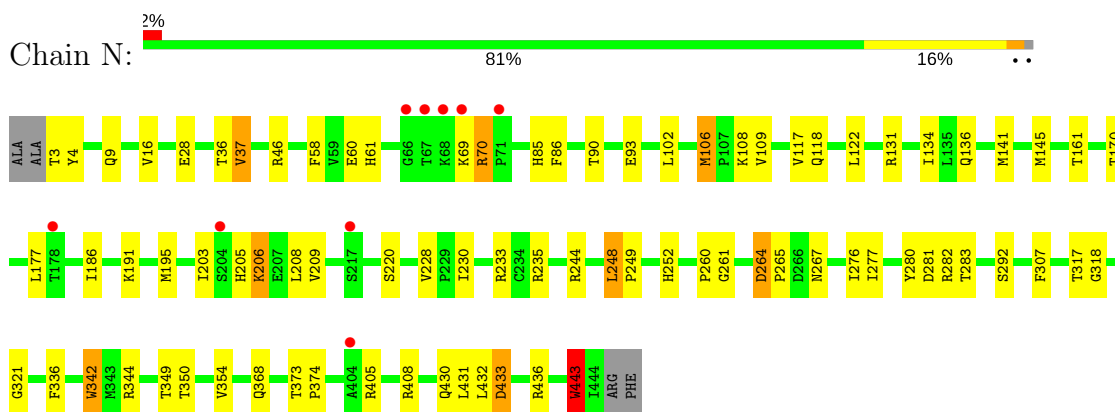
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

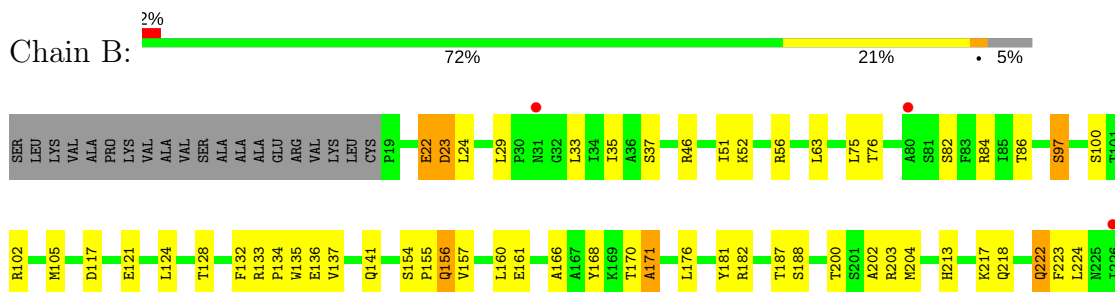
- Molecule 1: Mitochondrial ubiquinol-cytochrome-c reductase complex core protein i

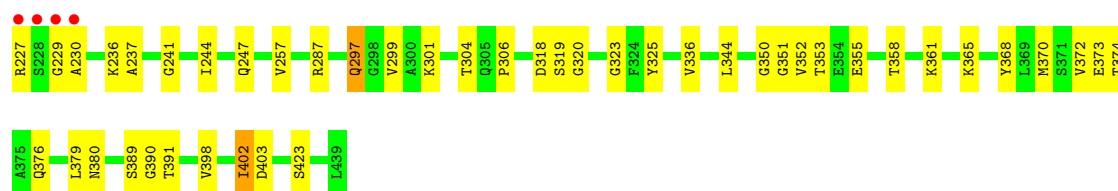


- Molecule 1: Mitochondrial ubiquinol-cytochrome-c reductase complex core protein i

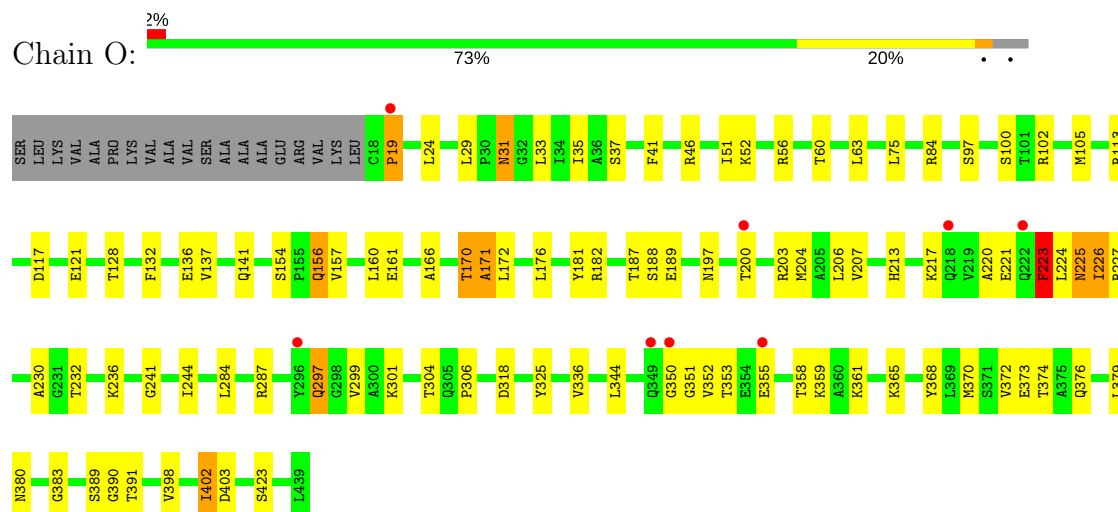


- Molecule 2: Mitochondrial ubiquinol-cytochrome-c reductase complex core protein 2

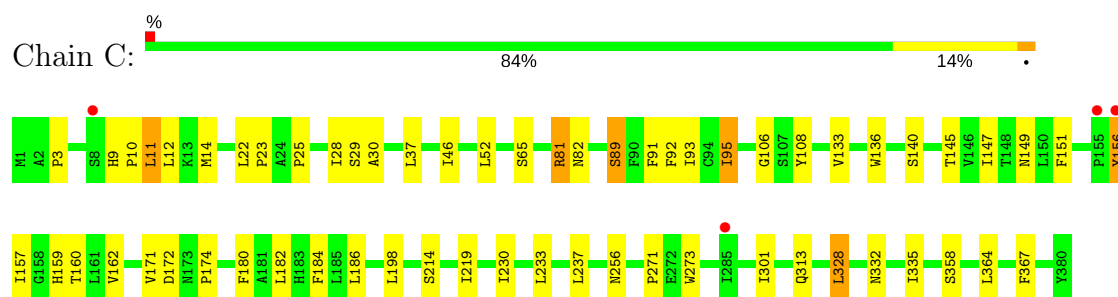




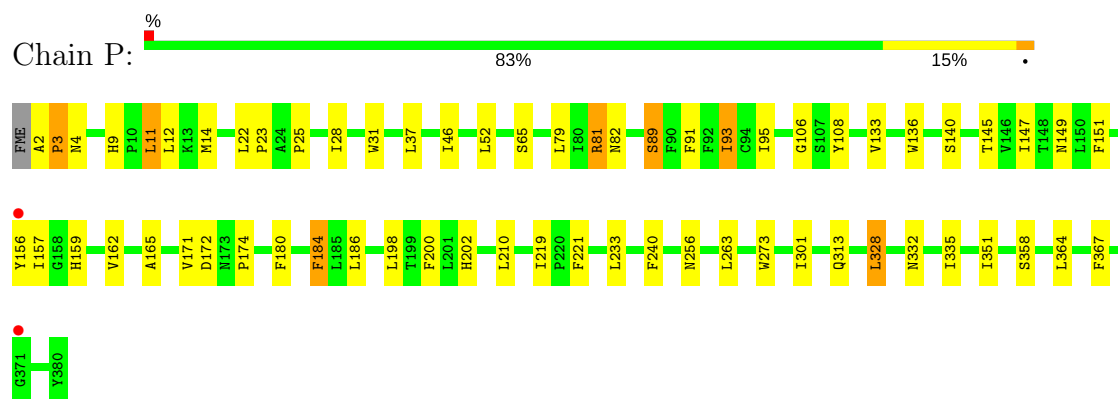
- Molecule 2: Mitochondrial ubiquinol-cytochrome-c reductase complex core protein 2



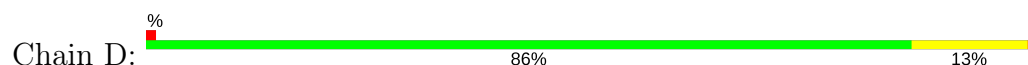
- Molecule 3: Cytochrome b



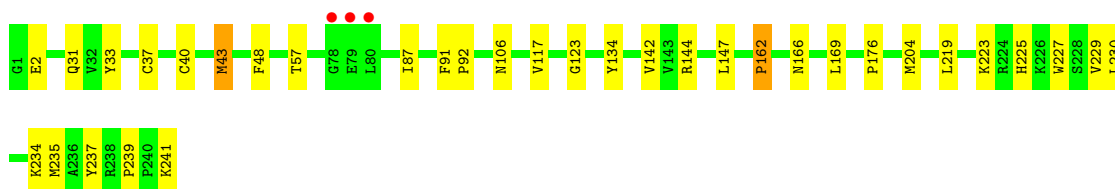
- Molecule 3: Cytochrome b



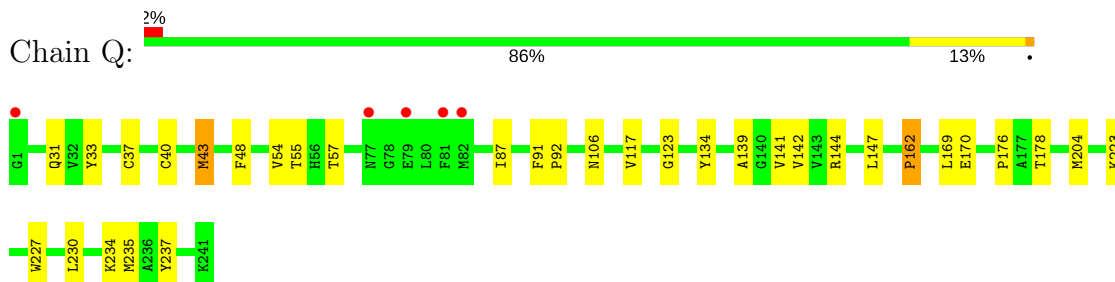
- Molecule 4: Mitochondrial cytochrome c1, heme protein



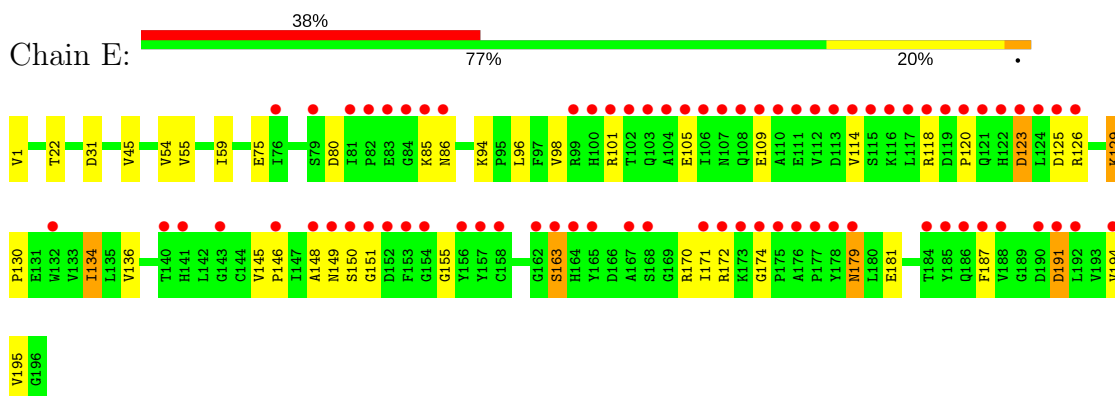




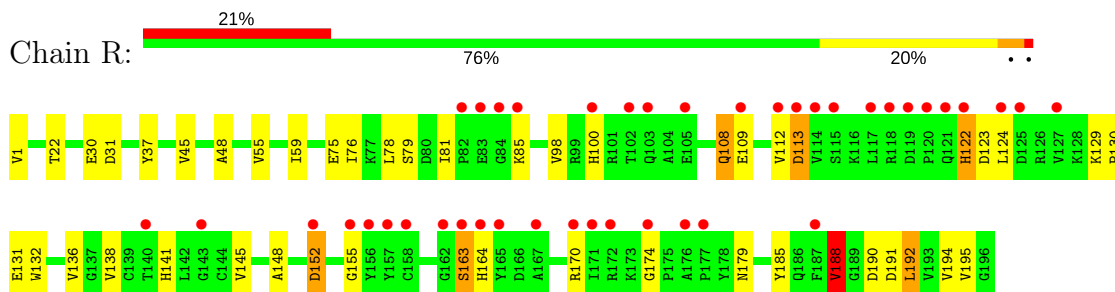
- Molecule 4: Mitochondrial cytochrome c1, heme protein



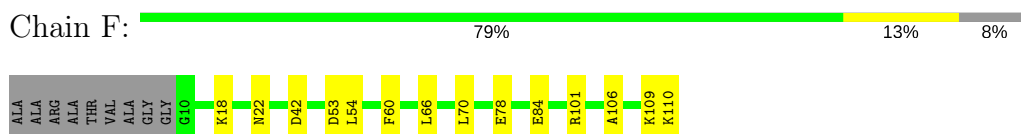
- Molecule 5: Cytochrome b-c1 complex subunit Rieske, mitochondrial



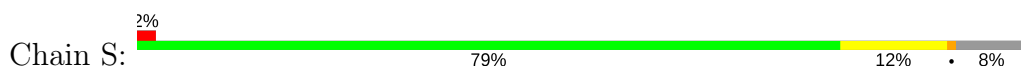
- Molecule 5: Cytochrome b-c1 complex subunit Rieske, mitochondrial

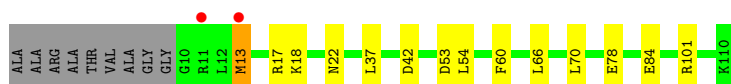


- Molecule 6: Cytochrome b-c1 complex subunit 7

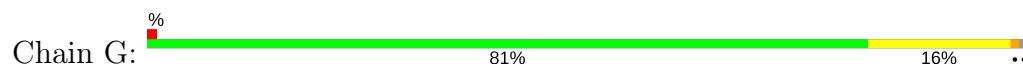


- Molecule 6: Cytochrome b-c1 complex subunit 7

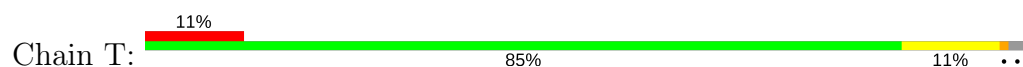




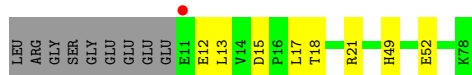
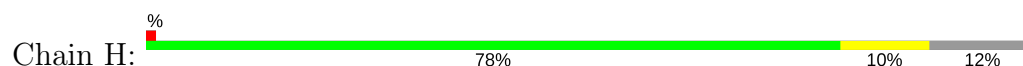
- Molecule 7: Mitochondrial ubiquinol-cytochrome c reductase ubiquinone-binding protein qp-c



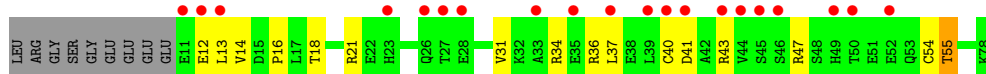
- Molecule 7: Mitochondrial ubiquinol-cytochrome c reductase ubiquinone-binding protein qp-c



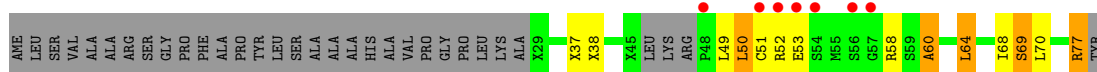
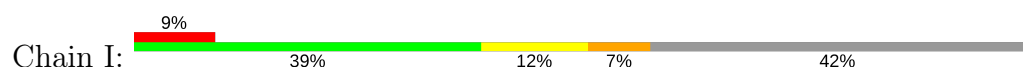
- Molecule 8: Cytochrome b-c1 complex subunit 6



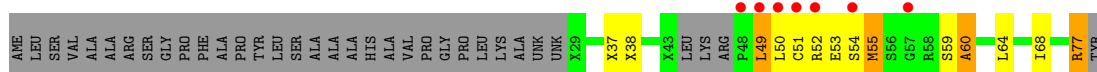
- Molecule 8: Cytochrome b-c1 complex subunit 6



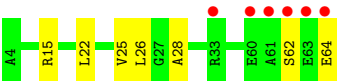
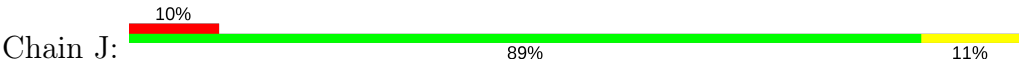
- Molecule 9: Cytochrome b-c1 complex subunit Rieske, mitochondrial



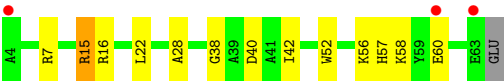
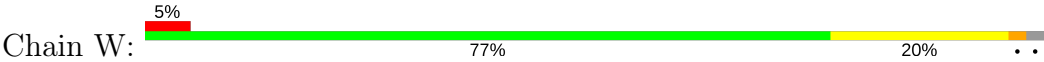
- Molecule 9: Cytochrome b-c1 complex subunit Rieske, mitochondrial



- Molecule 10: Mitochondrial ubiquinol-cytochrome c reductase 7.2 kda protein



● Molecule 10: Mitochondrial ubiquinol-cytochrome c reductase 7.2 kda protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	167.25Å 181.48Å 239.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.23 54.30 – 3.23	Depositor EDS
% Data completeness (in resolution range)	98.3 (15.00-3.23) 82.2 (54.30-3.23)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.30	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.05 (at 3.26Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1745)	Depositor
R, $R_{free}$	0.236 , 0.284 0.243 , 0.286	Depositor DCC
$R_{free}$ test set	2286 reflections (1.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	49.5	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 55.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	32790	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.07% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CDL, Y52, U10, FES, MES, HEC, PEE, FME, HEM, BOG

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	0.21	0/3513	0.40	0/4760
1	N	0.21	0/3508	0.39	0/4753
2	B	0.21	0/3192	0.40	0/4330
2	O	0.22	0/3198	0.40	0/4339
3	C	0.22	0/3114	0.39	0/4263
3	P	0.22	0/3102	0.39	0/4245
4	D	0.20	0/1956	0.37	0/2658
4	Q	0.20	0/1956	0.37	0/2658
5	E	0.20	0/1547	0.38	0/2103
5	R	0.20	0/1543	0.38	0/2098
6	F	0.21	0/911	0.35	0/1219
6	S	0.21	0/911	0.35	0/1219
7	G	0.22	0/694	0.39	0/941
7	T	0.22	0/684	0.39	0/929
8	H	0.21	0/570	0.36	0/763
8	U	0.21	0/566	0.40	0/758
9	I	0.20	0/208	0.48	0/279
9	V	0.21	0/215	0.56	0/288
10	J	0.21	0/508	0.34	0/682
10	W	0.21	0/490	0.35	0/660
All	All	0.21	0/32386	0.39	0/43945

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	216	PHE	Peptide

## 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	3442	0	3354	41	0
1	N	3437	0	3349	40	0
2	B	3137	0	3131	55	0
2	O	3143	0	3135	56	0
3	C	3021	0	3068	35	0
3	P	3022	0	3064	39	0
4	D	1898	0	1846	17	0
4	Q	1898	0	1846	16	0
5	E	1513	0	1480	21	0
5	R	1509	0	1476	22	0
6	F	891	0	893	6	0
6	S	891	0	893	8	0
7	G	672	0	653	11	0
7	T	662	0	645	10	0
8	H	562	0	545	3	0
8	U	558	0	541	8	0
9	I	266	0	223	14	0
9	V	265	0	239	11	0
10	J	497	0	490	3	0
10	W	479	0	478	6	0
11	A	26	0	26	1	0
11	C	122	0	174	2	0
11	E	63	0	99	5	0
11	N	8	0	7	0	0
11	P	142	0	193	7	0
11	R	49	0	75	5	0
12	C	86	0	60	7	0
12	P	86	0	60	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	C	27	0	19	2	0
13	P	27	0	19	1	0
14	C	19	0	17	1	0
14	P	19	0	17	3	0
15	C	12	0	13	0	0
16	C	6	0	8	1	0
16	P	6	0	8	1	0
17	D	43	0	30	0	0
17	Q	43	0	30	0	0
18	D	42	0	28	1	0
18	G	40	0	24	1	0
18	Q	42	0	28	0	0
18	T	40	0	24	3	0
19	D	20	0	28	0	0
19	P	20	0	28	3	0
19	Q	20	0	28	0	0
20	E	4	0	0	0	0
20	R	4	0	0	0	0
21	A	2	0	0	0	0
21	B	1	0	0	1	0
21	C	2	0	0	0	0
21	F	1	0	0	0	0
21	N	1	0	0	0	0
21	O	2	0	0	1	0
21	P	2	0	0	0	0
All	All	32790	0	32392	385	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 6.

All (385) 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)
9:V:37:UNK:HG3	9:V:38:UNK:H	1.37	0.90
9:V:49:LEU:HB2	9:V:55:MET:HB3	1.57	0.85
9:I:37:UNK:HG3	9:I:38:UNK:N	1.94	0.81
9:I:37:UNK:HG3	9:I:38:UNK:H	1.47	0.78
2:B:361:LYS:HD2	2:B:403:ASP:HA	1.67	0.76
2:O:361:LYS:HD2	2:O:403:ASP:HA	1.66	0.76
9:V:37:UNK:CG	9:V:38:UNK:H	2.00	0.75
2:O:389:SER:O	2:O:391:THR:N	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:328:LEU:HD12	7:G:51:PRO:HB3	1.71	0.72
2:O:226:ILE:HG22	2:O:227:ARG:H	1.55	0.72
2:B:389:SER:O	2:B:391:THR:N	2.22	0.71
2:B:306:PRO:HB3	9:I:52:ARG:HB2	1.73	0.70
4:D:166:ASN:HD21	8:H:15:ASP:HB2	1.56	0.70
8:U:40:CYS:SG	8:U:43:ARG:NH2	2.64	0.70
1:N:248:LEU:HD13	1:N:249:PRO:HD2	1.73	0.70
2:O:46:ARG:HG2	2:O:379:LEU:HD22	1.73	0.69
1:A:248:LEU:HD13	1:A:249:PRO:HD2	1.73	0.69
2:B:52:LYS:O	2:B:203:ARG:NH2	2.26	0.69
2:B:236:LYS:HE3	2:B:318:ASP:HB2	1.75	0.69
9:V:37:UNK:HG3	9:V:38:UNK:N	2.01	0.68
5:R:188:VAL:HG23	5:R:192:LEU:HB3	1.76	0.68
3:C:89:SER:HG	3:C:273:TRP:HE1	1.42	0.68
9:I:37:UNK:CG	9:I:38:UNK:H	2.07	0.67
2:O:46:ARG:NH2	2:O:376:GLN:OE1	2.28	0.67
2:O:52:LYS:O	2:O:203:ARG:NH2	2.26	0.67
3:C:81:ARG:HH12	16:C:508:GOL:H2	1.60	0.66
1:A:261:GLY:O	1:A:267:ASN:ND2	2.28	0.66
1:A:244:ARG:HE	7:G:10:VAL:HB	1.59	0.66
2:B:46:ARG:HG2	2:B:379:LEU:HD22	1.78	0.66
11:E:502:PEE:H62	10:J:25:VAL:HG22	1.78	0.66
1:N:261:GLY:O	1:N:267:ASN:ND2	2.29	0.66
1:A:106:MET:HE3	1:A:208:LEU:HA	1.80	0.64
3:C:23:PRO:HG2	7:G:3:HIS:HB3	1.78	0.64
3:P:328:LEU:HD12	7:T:51:PRO:HB3	1.79	0.64
2:B:86:THR:HA	9:I:70:LEU:HD11	1.81	0.63
4:D:239:PRO:HG2	4:D:241:LYS:HB2	1.79	0.63
1:A:131:ARG:NH2	1:A:177:LEU:O	2.31	0.63
1:N:131:ARG:NH2	1:N:177:LEU:O	2.32	0.63
9:I:60:ALA:HA	9:I:77:ARG:HH22	1.64	0.62
3:P:89:SER:HG	3:P:273:TRP:HE1	1.44	0.62
2:B:46:ARG:NH2	2:B:376:GLN:OE1	2.33	0.61
6:F:42:ASP:OD1	6:F:101:ARG:NH1	2.34	0.61
12:C:502:HEM:HMC2	12:C:502:HEM:HBC2	1.84	0.60
5:E:101:ARG:NH2	5:E:130:PRO:O	2.34	0.60
6:F:106:ALA:HA	6:F:109:LYS:HE2	1.82	0.60
12:C:501:HEM:HBC2	12:C:501:HEM:HMC1	1.84	0.60
7:T:40:ARG:NH1	18:T:101:CDL:OA4	2.35	0.60
6:S:42:ASP:OD1	6:S:101:ARG:NH1	2.34	0.59
1:A:412:SER:O	10:J:15:ARG:NH2	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:P:403:HEM:HBC2	12:P:403:HEM:HMC2	1.84	0.59
19:P:406:BOG:H4	11:R:502:PEE:H24	1.84	0.59
4:Q:178:THR:HG21	8:U:16:PRO:HD3	1.84	0.59
11:E:502:PEE:H42	11:E:503:PEE:H32	1.84	0.58
2:O:225:ASN:HD22	2:O:226:ILE:H	1.50	0.58
12:P:402:HEM:HBC2	12:P:402:HEM:HMC1	1.84	0.58
11:A:501:PEE:H57	11:E:502:PEE:H57	1.86	0.58
4:D:144:ARG:HB3	4:D:147:LEU:HD12	1.85	0.58
4:Q:144:ARG:HB3	4:Q:147:LEU:HD12	1.85	0.58
3:C:151:PHE:HB2	3:C:162:VAL:HG22	1.86	0.58
7:G:36:ASN:OD1	7:G:39:ARG:NH1	2.37	0.58
1:N:60:GLU:OE1	2:O:287:ARG:NH2	2.37	0.57
7:T:36:ASN:OD1	7:T:39:ARG:NH1	2.37	0.57
1:A:321:GLY:HA2	1:A:342:TRP:HZ2	1.70	0.57
5:E:118:ARG:NH1	5:E:172:ARG:O	2.37	0.57
3:P:202:HIS:NE2	14:P:405:U10:O2	2.38	0.57
2:B:56:ARG:HG2	2:B:171:ALA:HB1	1.87	0.57
3:P:23:PRO:HG2	7:T:3:HIS:HB3	1.87	0.57
1:N:106:MET:HE3	1:N:208:LEU:HA	1.87	0.57
3:P:301:ILE:HD11	3:P:364:LEU:HD21	1.86	0.57
3:P:335:ILE:HD13	7:T:58:LEU:HD23	1.87	0.57
1:N:321:GLY:HA2	1:N:342:TRP:HZ2	1.69	0.56
3:C:301:ILE:HD11	3:C:364:LEU:HD21	1.86	0.56
12:P:402:HEM:HMB1	12:P:402:HEM:HBB2	1.88	0.56
12:C:501:HEM:HMB1	12:C:501:HEM:HBB2	1.87	0.56
5:R:113:ASP:OD2	5:R:113:ASP:N	2.34	0.56
12:P:403:HEM:HMB1	12:P:403:HEM:HBB2	1.87	0.56
2:O:287:ARG:HA	9:V:53:GLU:HB3	1.86	0.56
2:B:29:LEU:HD12	2:B:33:LEU:HD23	1.88	0.56
2:B:157:VAL:HG23	9:I:64:LEU:HD21	1.87	0.56
2:B:156:GLN:NE2	21:B:501:HOH:O	2.38	0.55
1:N:244:ARG:HE	7:T:10:VAL:HB	1.70	0.55
14:P:405:U10:H1M1	14:P:405:U10:C8	2.37	0.55
2:O:117:ASP:N	2:O:117:ASP:OD1	2.39	0.55
2:O:56:ARG:HG2	2:O:171:ALA:HB1	1.87	0.55
8:U:18:THR:HA	8:U:21:ARG:HB2	1.88	0.55
5:E:163:SER:HA	5:E:174:GLY:HA3	1.88	0.55
2:O:128:THR:HG21	2:O:224:LEU:HG	1.89	0.55
3:P:151:PHE:HB2	3:P:162:VAL:HG22	1.87	0.55
5:R:76:ILE:HD13	5:R:98:VAL:HG21	1.89	0.55
3:C:198:LEU:HD21	12:C:502:HEM:HMA3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:THR:O	1:A:408:ARG:NH1	2.40	0.54
2:B:63:LEU:HB2	2:B:182:ARG:HD3	1.89	0.54
12:C:502:HEM:HBB2	12:C:502:HEM:HMB1	1.88	0.54
1:A:161:THR:HG21	1:A:235:ARG:H	1.71	0.54
3:P:198:LEU:HD21	12:P:403:HEM:HMA3	1.88	0.54
3:P:93:ILE:HD12	3:P:240:PHE:HZ	1.72	0.54
5:E:136:VAL:HB	5:E:181:GLU:HB3	1.89	0.54
2:O:350:GLY:O	2:O:352:VAL:N	2.39	0.54
3:C:335:ILE:HD13	7:G:58:LEU:HD23	1.89	0.54
1:N:349:THR:O	1:N:408:ARG:NH1	2.40	0.54
3:C:180:PHE:HE1	3:P:180:PHE:HE1	1.56	0.53
1:A:6:GLN:HA	1:A:9:GLN:HE21	1.73	0.53
8:U:13:LEU:HB2	8:U:14:VAL:HA	1.90	0.53
2:B:353:THR:HG22	2:B:355:GLU:H	1.73	0.53
3:C:25:PRO:HB2	3:C:28:ILE:HG23	1.91	0.53
2:O:63:LEU:HB2	2:O:182:ARG:HD3	1.90	0.53
3:C:147:ILE:HG13	13:C:503:Y52:H12	1.91	0.52
1:N:4:TYR:HB2	2:O:113:ARG:HB2	1.92	0.52
2:O:353:THR:HG22	2:O:355:GLU:H	1.73	0.52
2:O:75:LEU:HD22	2:O:136:GLU:HB3	1.92	0.52
3:P:25:PRO:HB2	3:P:28:ILE:HG23	1.91	0.52
2:O:306:PRO:HB3	9:V:52:ARG:N	2.25	0.51
2:O:157:VAL:HG23	9:V:64:LEU:HD21	1.91	0.51
1:A:117:VAL:HG23	1:A:118:GLN:HG3	1.93	0.51
3:P:81:ARG:HH12	16:P:408:GOL:H2	1.74	0.51
5:R:79:SER:OG	5:R:191:ASP:OD2	2.29	0.51
2:O:189:GLU:OE2	2:O:189:GLU:N	2.43	0.51
3:P:172:ASP:HB3	3:P:174:PRO:HD2	1.93	0.51
14:P:405:U10:H1M1	14:P:405:U10:H8	1.93	0.51
2:B:299:VAL:HG11	2:B:336:VAL:HG13	1.92	0.51
3:C:182:LEU:HD21	11:C:509:PEE:H3	1.92	0.51
3:P:147:ILE:HG13	13:P:404:Y52:H12	1.91	0.51
2:B:75:LEU:HD22	2:B:136:GLU:HB3	1.92	0.51
5:E:45:VAL:HG13	10:J:28:ALA:HA	1.92	0.51
2:O:37:SER:HB3	2:O:213:HIS:ND1	2.26	0.51
1:N:145:MET:HB3	1:N:252:HIS:CE1	2.46	0.51
1:A:244:ARG:NE	7:G:10:VAL:HB	2.26	0.51
2:O:157:VAL:HG23	9:V:64:LEU:HD11	1.93	0.51
1:N:117:VAL:HG23	1:N:118:GLN:HG3	1.93	0.51
2:B:117:ASP:OD1	2:B:117:ASP:N	2.39	0.50
2:B:154:SER:O	2:B:157:VAL:HG12	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:299:VAL:HG11	2:O:336:VAL:HG13	1.92	0.50
2:B:350:GLY:O	2:B:352:VAL:N	2.39	0.50
2:O:84:ARG:NH2	2:O:121:GLU:OE2	2.45	0.50
3:P:52:LEU:HD11	3:P:81:ARG:HA	1.93	0.50
2:O:154:SER:O	2:O:157:VAL:HG12	2.11	0.50
4:Q:237:TYR:HB2	6:S:60:PHE:CD1	2.47	0.50
5:E:171:ILE:HG22	5:E:179:ASN:OD1	2.12	0.49
2:B:132:PHE:O	2:B:188:SER:OG	2.27	0.49
2:B:84:ARG:NH2	2:B:121:GLU:OE2	2.45	0.49
3:C:29:SER:HB2	18:G:101:CDL:HB21	1.94	0.49
2:O:31:ASN:N	2:O:31:ASN:OD1	2.43	0.49
5:R:122:HIS:CG	5:R:123:ASP:H	2.31	0.49
3:C:172:ASP:HB3	3:C:174:PRO:HD2	1.94	0.49
1:N:161:THR:HG21	1:N:235:ARG:H	1.77	0.49
1:N:443:TRP:HA	1:N:443:TRP:CE3	2.48	0.49
2:O:161:GLU:OE1	2:O:176:LEU:N	2.45	0.49
2:O:365:LYS:HG3	2:O:402:ILE:HD12	1.95	0.49
1:A:443:TRP:HA	1:A:443:TRP:CE3	2.48	0.49
2:B:128:THR:HG21	2:B:224:LEU:HD22	1.95	0.49
2:O:181:TYR:CZ	2:O:182:ARG:HG3	2.48	0.49
1:A:140:GLU:HG3	9:I:51:CYS:SG	2.52	0.49
2:B:137:VAL:O	2:B:141:GLN:HG2	2.13	0.48
9:I:49:LEU:HD21	9:I:58:ARG:NE	2.28	0.48
19:P:406:BOG:H3'2	11:R:502:PEE:H46	1.95	0.48
3:C:332:ASN:ND2	3:C:358:SER:OG	2.45	0.48
2:B:181:TYR:CZ	2:B:182:ARG:HG3	2.48	0.48
2:B:365:LYS:HG3	2:B:402:ILE:HD12	1.95	0.48
5:E:125:ASP:HB3	5:E:126:ARG:HD2	1.95	0.48
2:B:247:GLN:NE2	2:O:60:THR:OG1	2.46	0.48
2:B:161:GLU:OE1	2:B:176:LEU:N	2.46	0.48
2:B:157:VAL:HG23	9:I:64:LEU:HD11	1.96	0.48
3:C:219:ILE:HG21	4:D:230:LEU:HD11	1.94	0.48
3:C:52:LEU:HD11	3:C:81:ARG:HA	1.93	0.48
2:O:132:PHE:O	2:O:188:SER:OG	2.27	0.48
10:W:57:HIS:CE1	10:W:58:LYS:HG3	2.49	0.48
1:A:191:LYS:O	1:A:195:MET:HG3	2.14	0.48
2:B:35:ILE:HD13	2:B:217:LYS:HA	1.95	0.48
4:D:235:MET:HB3	7:G:15:THR:HG22	1.96	0.48
1:N:102:LEU:H	1:N:102:LEU:HD12	1.79	0.48
2:B:203:ARG:HD2	2:B:230:ALA:O	2.14	0.48
5:E:129:LYS:HD2	5:E:130:PRO:HD2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:191:LYS:O	1:N:195:MET:HG3	2.14	0.48
2:B:37:SER:HB3	2:B:213:HIS:ND1	2.29	0.47
4:D:33:TYR:CD1	4:D:37:CYS:HB2	2.49	0.47
3:P:9:HIS:HB3	3:P:12:LEU:HB2	1.96	0.47
5:R:81:ILE:HG22	5:R:100:HIS:HB2	1.95	0.47
2:O:137:VAL:O	2:O:141:GLN:HG2	2.14	0.47
4:Q:33:TYR:CD1	4:Q:37:CYS:HB2	2.49	0.47
5:E:55:VAL:O	5:E:59:ILE:HG12	2.14	0.47
1:N:368:GLN:O	1:N:374:PRO:HB2	2.14	0.47
2:O:35:ILE:HD13	2:O:217:LYS:HA	1.96	0.47
3:C:52:LEU:HD13	12:C:501:HEM:HBD1	1.96	0.47
5:R:55:VAL:O	5:R:59:ILE:HG12	2.15	0.47
6:F:53:ASP:OD1	6:F:54:LEU:N	2.48	0.47
3:C:145:THR:O	3:C:149:ASN:HB2	2.15	0.47
5:E:75:GLU:HB3	5:E:194:VAL:HG22	1.96	0.47
1:N:205:HIS:HB3	1:N:206:LYS:HE3	1.97	0.47
1:N:46:ARG:NH1	1:N:93:GLU:OE2	2.44	0.47
2:O:217:LYS:O	2:O:221:GLU:HG2	2.14	0.47
2:B:241:GLY:HA2	2:B:423:SER:HB3	1.97	0.46
2:O:203:ARG:HD2	2:O:230:ALA:O	2.15	0.46
2:B:22:GLU:H	2:B:22:GLU:HG3	1.54	0.46
2:B:56:ARG:NH1	2:B:318:ASP:OD2	2.47	0.46
3:C:9:HIS:HB3	3:C:12:LEU:HB2	1.97	0.46
5:E:191:ASP:N	5:E:191:ASP:OD2	2.47	0.46
3:P:52:LEU:HD13	12:P:402:HEM:HBD1	1.96	0.46
3:P:351:ILE:HD11	7:T:61:TRP:HZ3	1.79	0.46
3:P:332:ASN:ND2	3:P:358:SER:OG	2.45	0.46
5:R:30:GLU:OE1	10:W:7:ARG:NH2	2.49	0.46
3:C:10:PRO:HG2	3:P:200:PHE:CE1	2.50	0.46
5:R:152:ASP:HB2	5:R:164:HIS:CG	2.50	0.46
5:R:30:GLU:HB2	10:W:7:ARG:HG2	1.97	0.46
1:A:368:GLN:O	1:A:374:PRO:HB2	2.14	0.46
1:A:102:LEU:H	1:A:102:LEU:HD12	1.79	0.46
3:P:145:THR:O	3:P:149:ASN:HB2	2.15	0.46
5:E:145:VAL:HA	5:E:146:PRO:HD3	1.79	0.46
5:E:94:LYS:HG2	3:P:263:LEU:HD21	1.98	0.46
2:O:56:ARG:NH1	2:O:318:ASP:OD2	2.49	0.46
1:N:37:VAL:HG13	1:N:109:VAL:HG11	1.98	0.46
5:R:136:VAL:HG12	5:R:138:VAL:HG23	1.98	0.45
6:S:53:ASP:OD1	6:S:54:LEU:N	2.48	0.45
5:R:131:GLU:HG2	5:R:132:TRP:CD1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:GLU:OE1	1:A:344:ARG:NH1	2.45	0.45
1:A:46:ARG:NH1	1:A:93:GLU:OE2	2.44	0.45
2:B:23:ASP:N	2:B:23:ASP:OD1	2.34	0.45
7:G:72:LYS:HZ2	8:H:52:GLU:CG	2.30	0.45
5:R:108:GLN:HB3	5:R:108:GLN:HE21	1.55	0.45
1:A:37:VAL:HG13	1:A:109:VAL:HG11	1.98	0.45
3:C:186:LEU:HD21	11:C:509:PEE:H2	1.99	0.45
5:E:150:SER:OG	5:E:151:GLY:N	2.49	0.45
5:R:37:TYR:CE2	11:R:502:PEE:H10	2.52	0.45
3:P:79:LEU:HD22	4:Q:204:MET:HE1	1.97	0.45
3:C:46:ILE:HA	12:C:501:HEM:HMC2	1.98	0.45
1:A:141:MET:HB3	1:A:141:MET:HE2	1.79	0.45
1:A:317:THR:OG1	1:A:318:GLY:N	2.51	0.45
1:N:69:LYS:HE3	1:N:70:ARG:HH21	1.81	0.45
1:A:69:LYS:HE3	1:A:70:ARG:HH21	1.81	0.44
1:A:321:GLY:HA2	1:A:342:TRP:CZ2	2.52	0.44
1:N:280:TYR:HB3	1:N:307:PHE:CE2	2.52	0.44
1:A:191:LYS:HA	1:A:191:LYS:HE2	2.00	0.44
1:A:281:ASP:O	1:A:283:THR:N	2.50	0.44
3:C:133:VAL:HA	3:C:140:SER:HB3	1.99	0.44
2:O:206:LEU:HD23	2:O:220:ALA:HB2	2.00	0.44
3:P:133:VAL:HA	3:P:140:SER:HB3	2.00	0.44
3:P:89:SER:OG	3:P:273:TRP:NE1	2.30	0.44
4:Q:48:PHE:HB2	4:Q:87:ILE:HA	1.99	0.44
6:S:13:MET:O	6:S:17:ARG:HG2	2.18	0.44
3:P:165:ALA:HB1	11:P:409:PEE:H60	1.99	0.44
4:Q:134:TYR:CG	4:Q:162:PRO:HG3	2.53	0.44
4:Q:237:TYR:HB2	6:S:60:PHE:CG	2.53	0.44
5:R:48:ALA:HB1	11:R:502:PEE:H66	1.99	0.44
2:B:168:TYR:CD2	2:B:237:ALA:HB1	2.53	0.44
4:D:48:PHE:HB2	4:D:87:ILE:HA	2.00	0.44
3:C:89:SER:OG	3:C:273:TRP:NE1	2.31	0.44
1:A:61:HIS:CE1	1:A:134:ILE:HG12	2.53	0.44
1:N:191:LYS:HE2	1:N:191:LYS:HA	2.00	0.44
6:S:18:LYS:O	6:S:22:ASN:ND2	2.50	0.44
1:A:280:TYR:HB3	1:A:307:PHE:CE2	2.52	0.44
2:B:370:MET:HA	2:B:373:GLU:HG3	2.00	0.44
2:B:287:ARG:HG2	9:I:53:GLU:HB3	2.00	0.44
1:N:281:ASP:O	1:N:283:THR:N	2.50	0.44
1:N:433:ASP:OD1	1:N:436:ARG:HG2	2.18	0.44
2:O:370:MET:HA	2:O:373:GLU:HG3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:207:VAL:HG21	2:O:383:GLY:HA2	2.00	0.44
3:P:219:ILE:HG21	4:Q:230:LEU:HD11	2.00	0.44
1:A:433:ASP:OD1	1:A:436:ARG:HG2	2.18	0.43
2:O:197:ASN:HB3	2:O:232:THR:HB	2.00	0.43
5:E:98:VAL:HG22	5:E:134:ILE:HG22	1.99	0.43
5:E:123:ASP:OD1	5:E:170:ARG:NH2	2.51	0.43
5:R:75:GLU:HG2	5:R:194:VAL:HG22	1.99	0.43
2:B:166:ALA:HB2	2:B:244:ILE:HG13	2.00	0.43
5:E:118:ARG:HD3	5:E:171:ILE:HG23	2.00	0.43
1:N:61:HIS:CE1	1:N:134:ILE:HG12	2.53	0.43
11:P:409:PEE:H64	11:P:409:PEE:H10	2.00	0.43
1:N:136:GLN:HE21	9:V:50:LEU:HB2	1.83	0.43
1:A:264:ASP:HA	1:A:265:PRO:HD3	1.90	0.43
4:D:225:HIS:O	7:G:20:PRO:HB3	2.18	0.43
5:R:185:TYR:HB3	5:R:195:VAL:HG13	2.01	0.43
7:T:40:ARG:HD2	18:T:101:CDL:OA4	2.17	0.43
1:A:46:ARG:HD3	1:A:231:LEU:HD13	2.01	0.43
1:A:3:THR:OG1	1:A:6:GLN:OE1	2.23	0.43
14:C:504:U10:C8	14:C:504:U10:H1M1	2.49	0.43
4:D:204:MET:HB3	4:D:204:MET:HE2	1.84	0.43
3:P:210:LEU:HD12	6:S:66:LEU:HD23	2.01	0.43
3:P:31:TRP:NE1	11:P:407:PEE:O4	2.51	0.43
3:P:46:ILE:HA	12:P:402:HEM:HMC2	1.99	0.43
4:D:134:TYR:CG	4:D:162:PRO:HG3	2.53	0.43
6:F:18:LYS:O	6:F:22:ASN:ND2	2.50	0.43
1:N:36:THR:HG21	1:N:373:THR:HA	2.01	0.43
8:U:31:VAL:HA	8:U:34:ARG:HB3	1.99	0.43
1:N:264:ASP:HA	1:N:265:PRO:HD3	1.90	0.43
3:P:106:GLY:HA2	3:P:108:TYR:CE2	2.53	0.43
1:N:317:THR:OG1	1:N:318:GLY:N	2.50	0.43
5:E:86:ASN:HD22	5:E:148:ALA:HB1	1.84	0.42
2:O:29:LEU:HD12	2:O:33:LEU:HD23	2.01	0.42
4:D:223:LYS:HE2	4:D:227:TRP:CD1	2.54	0.42
1:N:122:LEU:HD11	1:N:186:ILE:HD12	2.02	0.42
1:N:244:ARG:NE	7:T:10:VAL:HB	2.34	0.42
2:B:368:TYR:O	2:B:372:VAL:HG23	2.20	0.42
3:C:106:GLY:HA2	3:C:108:TYR:CE2	2.53	0.42
2:O:166:ALA:HB2	2:O:244:ILE:HG13	2.01	0.42
2:O:306:PRO:HB3	9:V:52:ARG:H	1.85	0.42
2:O:241:GLY:HA2	2:O:423:SER:HB3	2.01	0.42
4:D:91:PHE:HA	4:D:92:PRO:HD3	1.74	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:V:60:ALA:HA	9:V:77:ARG:HH22	1.85	0.42
3:P:9:HIS:CE1	3:P:11:LEU:HB2	2.55	0.42
2:B:319:SER:OG	2:B:320:GLY:N	2.53	0.42
3:C:136:TRP:HH2	3:C:171:VAL:HG12	1.85	0.42
3:C:230:ILE:HG22	4:D:219:LEU:HD13	2.01	0.42
5:E:96:LEU:HD21	5:E:195:VAL:HG21	2.01	0.42
19:P:406:BOG:H1'1	11:R:502:PEE:H33	2.02	0.42
1:A:60:GLU:OE1	2:B:287:ARG:NH2	2.49	0.42
4:D:33:TYR:CZ	4:D:43:MET:HG3	2.55	0.42
1:N:60:GLU:OE2	1:N:90:THR:HG22	2.20	0.42
3:P:221:PHE:HE2	11:P:401:PEE:H26	1.84	0.42
4:Q:139:ALA:HB3	8:U:54:CYS:SG	2.60	0.42
4:Q:223:LYS:HE2	4:Q:227:TRP:CD1	2.54	0.42
3:C:37:LEU:HD21	3:C:233:LEU:HA	2.02	0.42
1:N:321:GLY:HA2	1:N:342:TRP:CZ2	2.52	0.42
1:A:205:HIS:O	1:A:209:VAL:HG13	2.20	0.41
1:A:21:ASN:OD1	1:A:21:ASN:N	2.50	0.41
1:A:60:GLU:OE2	1:A:90:THR:HG22	2.20	0.41
3:C:9:HIS:CE1	3:C:11:LEU:HB2	2.55	0.41
1:A:140:GLU:CD	9:I:50:LEU:H	2.20	0.41
1:A:122:LEU:HD11	1:A:186:ILE:HD12	2.02	0.41
3:P:186:LEU:HD21	11:P:409:PEE:H1	2.02	0.41
5:R:163:SER:HA	5:R:174:GLY:HA3	2.01	0.41
3:C:271:PRO:HB3	13:C:503:Y52:C14	2.51	0.41
1:N:336:PHE:CE2	3:P:4:ASN:HB3	2.55	0.41
5:R:78:LEU:HB3	5:R:132:TRP:CZ2	2.54	0.41
3:C:92:PHE:HA	3:C:95:ILE:HG22	2.02	0.41
3:C:30:ALA:HB2	18:D:502:CDL:HA31	2.03	0.41
2:O:156:GLN:NE2	21:O:502:HOH:O	2.46	0.41
4:Q:91:PHE:HA	4:Q:92:PRO:HD3	1.74	0.41
7:T:40:ARG:HB3	18:T:101:CDL:HA31	2.02	0.41
2:B:257:VAL:O	2:B:323:GLY:HA3	2.21	0.41
1:N:277:ILE:O	1:N:292:SER:OG	2.28	0.41
2:O:368:TYR:O	2:O:372:VAL:HG23	2.20	0.41
5:R:170:ARG:HA	5:R:179:ASN:HB3	2.02	0.41
2:B:124:LEU:HD21	2:B:223:PHE:HB3	2.02	0.41
1:N:141:MET:HB3	1:N:141:MET:HE2	1.77	0.41
1:N:205:HIS:O	1:N:209:VAL:HG13	2.20	0.41
1:N:276:ILE:HD11	1:N:354:VAL:HG12	2.02	0.41
11:P:411:PEE:H13	11:P:411:PEE:H50	2.03	0.41
4:D:117:VAL:O	4:D:123:GLY:HA2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:223:PHE:HA	2:O:223:PHE:HD2	1.77	0.41
2:O:297:GLN:O	2:O:301:LYS:HG3	2.20	0.41
3:P:136:TRP:HH2	3:P:171:VAL:HG12	1.85	0.41
3:P:37:LEU:HD21	3:P:233:LEU:HA	2.03	0.41
4:Q:54:VAL:HG12	4:Q:55:THR:HG23	2.03	0.41
4:Q:141:VAL:HG21	8:U:55:THR:HG23	2.03	0.41
2:B:202:ALA:HB3	2:B:229:GLY:O	2.21	0.41
2:B:76:THR:HG22	2:B:82:SER:H	1.86	0.41
3:C:237:LEU:HD21	11:E:503:PEE:H29	2.02	0.41
1:N:106:MET:HG3	1:N:203:ILE:HD13	2.03	0.41
2:O:51:ILE:HG12	2:O:204:MET:HG2	2.03	0.41
11:P:407:PEE:H25	11:P:407:PEE:H32	1.92	0.41
4:Q:117:VAL:O	4:Q:123:GLY:HA2	2.20	0.41
5:R:45:VAL:HG13	10:W:28:ALA:HA	2.03	0.41
2:B:133:ARG:HA	2:B:134:PRO:HD3	1.87	0.41
2:B:218:GLN:O	2:B:222:GLN:HB2	2.20	0.41
3:C:214:SER:HA	6:F:66:LEU:HD11	2.03	0.41
8:H:17:LEU:HD21	8:H:21:ARG:HH21	1.86	0.41
2:O:200:THR:O	2:O:204:MET:HG3	2.20	0.41
3:P:2:ALA:HA	3:P:3:PRO:HD2	1.72	0.41
4:Q:33:TYR:CZ	4:Q:43:MET:HG3	2.55	0.41
1:A:276:ILE:HD11	1:A:354:VAL:HG12	2.03	0.41
4:D:237:TYR:HB2	6:F:60:PHE:CD1	2.56	0.41
5:E:54:VAL:HG11	11:E:502:PEE:H43	2.03	0.41
4:D:229:VAL:HG23	7:G:20:PRO:HG3	2.02	0.41
7:G:24:ARG:NH2	7:G:28:ASN:H	2.18	0.41
9:I:51:CYS:HB3	9:I:52:ARG:H	1.70	0.41
5:R:129:LYS:HA	5:R:130:PRO:HD3	1.85	0.41
2:B:100:SER:HB2	2:B:105:MET:HG2	2.02	0.41
2:B:154:SER:HA	2:B:155:PRO:HD3	1.94	0.41
2:O:100:SER:HB2	2:O:105:MET:HG2	2.03	0.41
3:P:184:PHE:CZ	12:P:402:HEM:HBC1	2.56	0.41
2:B:133:ARG:HD3	2:B:135:TRP:CZ2	2.57	0.40
2:O:19:PRO:HB3	2:O:41:PHE:CD1	2.56	0.40
1:N:85:HIS:CD2	2:O:284:LEU:HD22	2.55	0.40
10:W:38:GLY:O	10:W:42:ILE:HG13	2.21	0.40
2:B:200:THR:O	2:B:204:MET:HG3	2.21	0.40
2:B:51:ILE:HG12	2:B:204:MET:HG2	2.03	0.40
2:O:170:THR:O	2:O:172:LEU:N	2.54	0.40
2:O:359:LYS:HB2	2:O:359:LYS:HE3	1.91	0.40
10:W:52:TRP:O	10:W:56:LYS:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:LEU:HD12	1:A:384:LEU:HA	1.95	0.40
2:B:297:GLN:O	2:B:301:LYS:HG3	2.20	0.40
2:B:97:SER:HA	9:I:69:SER:HA	2.03	0.40
2:O:56:ARG:NH2	2:O:236:LYS:HA	2.37	0.40
8:U:37:LEU:O	8:U:41:ASP:N	2.49	0.40
5:E:105:GLU:O	5:E:109:GLU:HB2	2.21	0.40
1:A:130:GLU:HA	1:A:133:VAL:HB	2.04	0.40
6:S:37:LEU:HA	6:S:37:LEU:HD23	1.93	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	441/446 (99%)	421 (96%)	16 (4%)	4 (1%)	19	57
1	N	440/446 (99%)	419 (95%)	17 (4%)	4 (1%)	19	57
2	B	419/441 (95%)	397 (95%)	18 (4%)	4 (1%)	17	55
2	O	420/441 (95%)	397 (94%)	17 (4%)	6 (1%)	12	47
3	C	378/380 (100%)	365 (97%)	11 (3%)	2 (0%)	31	68
3	P	378/380 (100%)	364 (96%)	12 (3%)	2 (0%)	31	68
4	D	239/241 (99%)	226 (95%)	11 (5%)	2 (1%)	21	59
4	Q	239/241 (99%)	225 (94%)	12 (5%)	2 (1%)	21	59
5	E	194/196 (99%)	168 (87%)	20 (10%)	6 (3%)	4	27
5	R	194/196 (99%)	175 (90%)	14 (7%)	5 (3%)	6	32
6	F	99/110 (90%)	97 (98%)	2 (2%)	0	100	100
6	S	99/110 (90%)	98 (99%)	1 (1%)	0	100	100
7	G	78/81 (96%)	74 (95%)	4 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
7	T	77/81 (95%)	72 (94%)	5 (6%)	0	100	100
8	H	66/77 (86%)	64 (97%)	2 (3%)	0	100	100
8	U	66/77 (86%)	56 (85%)	9 (14%)	1 (2%)	11	46
9	I	28/76 (37%)	17 (61%)	9 (32%)	2 (7%)	1	8
9	V	28/76 (37%)	19 (68%)	7 (25%)	2 (7%)	1	8
10	J	59/61 (97%)	55 (93%)	3 (5%)	1 (2%)	10	43
10	W	58/61 (95%)	54 (93%)	3 (5%)	1 (2%)	10	43
All	All	4000/4218 (95%)	3763 (94%)	193 (5%)	44 (1%)	16	53

All (44) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	390	GLY
2	O	19	PRO
2	O	226	ILE
2	O	390	GLY
9	V	59	SER
2	B	171	ALA
2	B	351	GLY
5	E	149	ASN
2	O	171	ALA
2	O	351	GLY
5	R	148	ALA
5	R	188	VAL
1	A	282	ARG
1	A	433	ASP
3	C	3	PRO
3	C	156	TYR
4	D	176	PRO
10	J	62	SER
1	N	282	ARG
1	N	433	ASP
2	O	223	PHE
3	P	3	PRO
4	Q	176	PRO
1	A	443	TRP
5	E	114	VAL
5	E	179	ASN
9	I	60	ALA
1	N	443	TRP

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Mol	Chain	Res	Type
8	U	12	GLU
2	B	227	ARG
5	E	120	PRO
5	E	155	GLY
5	E	163	SER
9	I	64	LEU
3	P	156	TYR
5	R	155	GLY
5	R	163	SER
9	V	60	ALA
10	W	15	ARG
5	R	141	HIS
4	D	162	PRO
4	Q	162	PRO
1	A	260	PRO
1	N	260	PRO

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	365/368 (99%)	342 (94%)	23 (6%)	20	54
1	N	365/368 (99%)	339 (93%)	26 (7%)	16	49
2	B	331/347 (95%)	312 (94%)	19 (6%)	23	58
2	O	332/347 (96%)	313 (94%)	19 (6%)	23	58
3	C	328/328 (100%)	309 (94%)	19 (6%)	22	57
3	P	327/328 (100%)	310 (95%)	17 (5%)	25	61
4	D	200/200 (100%)	191 (96%)	9 (4%)	30	65
4	Q	200/200 (100%)	190 (95%)	10 (5%)	27	62
5	E	166/166 (100%)	156 (94%)	10 (6%)	21	56
5	R	165/166 (99%)	150 (91%)	15 (9%)	10	35
6	F	93/96 (97%)	89 (96%)	4 (4%)	32	66

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
6	S	93/96 (97%)	89 (96%)	4 (4%)	32	66
7	G	71/71 (100%)	67 (94%)	4 (6%)	23	59
7	T	70/71 (99%)	67 (96%)	3 (4%)	32	66
8	H	64/71 (90%)	60 (94%)	4 (6%)	20	54
8	U	63/71 (89%)	60 (95%)	3 (5%)	28	63
9	I	22/45 (49%)	18 (82%)	4 (18%)	2	9
9	V	23/45 (51%)	17 (74%)	6 (26%)	0	1
10	J	49/49 (100%)	46 (94%)	3 (6%)	20	56
10	W	47/49 (96%)	42 (89%)	5 (11%)	7	29
All	All	3374/3482 (97%)	3167 (94%)	207 (6%)	20	56

All (207) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	16	VAL
1	A	28	GLU
1	A	37	VAL
1	A	58	PHE
1	A	70	ARG
1	A	86	PHE
1	A	106	MET
1	A	108	LYS
1	A	170	THR
1	A	220	SER
1	A	228	VAL
1	A	230	ILE
1	A	233	ARG
1	A	248	LEU
1	A	264	ASP
1	A	342	TRP
1	A	350	THR
1	A	405	ARG
1	A	430	GLN
1	A	431	LEU
1	A	432	LEU
1	A	443	TRP
2	B	22	GLU
2	B	23	ASP

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Mol	Chain	Res	Type
2	B	24	LEU
2	B	97	SER
2	B	102	ARG
2	B	156	GLN
2	B	160	LEU
2	B	170	THR
2	B	187	THR
2	B	222	GLN
2	B	297	GLN
2	B	304	THR
2	B	325	TYR
2	B	344	LEU
2	B	358	THR
2	B	374	THR
2	B	380	ASN
2	B	398	VAL
2	B	402	ILE
3	C	11	LEU
3	C	14	MET
3	C	22	LEU
3	C	65	SER
3	C	81	ARG
3	C	82	ASN
3	C	89	SER
3	C	91	PHE
3	C	93	ILE
3	C	95	ILE
3	C	156	TYR
3	C	157	ILE
3	C	159	HIS
3	C	160	THR
3	C	184	PHE
3	C	256	ASN
3	C	313	GLN
3	C	328	LEU
3	C	367	PHE
4	D	2	GLU
4	D	31	GLN
4	D	40	CYS
4	D	43	MET
4	D	57	THR
4	D	106	ASN

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Mol	Chain	Res	Type
4	D	142	VAL
4	D	169	LEU
4	D	234	LYS
5	E	1	VAL
5	E	22	THR
5	E	31	ASP
5	E	80	ASP
5	E	85	LYS
5	E	123	ASP
5	E	129	LYS
5	E	134	ILE
5	E	187	PHE
5	E	191	ASP
6	F	70	LEU
6	F	78	GLU
6	F	84	GLU
6	F	110	LYS
7	G	2	ILE
7	G	3	HIS
7	G	4	PHE
7	G	6	ASN
8	H	12	GLU
8	H	13	LEU
8	H	18	THR
8	H	49	HIS
9	I	50	LEU
9	I	68	ILE
9	I	69	SER
9	I	77	ARG
10	J	22	LEU
10	J	26	LEU
10	J	64	GLU
1	N	3	THR
1	N	9	GLN
1	N	16	VAL
1	N	28	GLU
1	N	37	VAL
1	N	58	PHE
1	N	70	ARG
1	N	86	PHE
1	N	106	MET
1	N	108	LYS

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Mol	Chain	Res	Type
1	N	170	THR
1	N	206	LYS
1	N	220	SER
1	N	228	VAL
1	N	230	ILE
1	N	233	ARG
1	N	248	LEU
1	N	264	ASP
1	N	342	TRP
1	N	344	ARG
1	N	350	THR
1	N	405	ARG
1	N	430	GLN
1	N	431	LEU
1	N	432	LEU
1	N	443	TRP
2	O	24	LEU
2	O	31	ASN
2	O	97	SER
2	O	102	ARG
2	O	156	GLN
2	O	160	LEU
2	O	170	THR
2	O	187	THR
2	O	223	PHE
2	O	225	ASN
2	O	297	GLN
2	O	304	THR
2	O	325	TYR
2	O	344	LEU
2	O	358	THR
2	O	374	THR
2	O	380	ASN
2	O	398	VAL
2	O	402	ILE
3	P	11	LEU
3	P	14	MET
3	P	22	LEU
3	P	65	SER
3	P	81	ARG
3	P	82	ASN
3	P	89	SER

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Mol	Chain	Res	Type
3	P	91	PHE
3	P	93	ILE
3	P	95	ILE
3	P	157	ILE
3	P	159	HIS
3	P	184	PHE
3	P	256	ASN
3	P	313	GLN
3	P	328	LEU
3	P	367	PHE
4	Q	31	GLN
4	Q	40	CYS
4	Q	43	MET
4	Q	57	THR
4	Q	106	ASN
4	Q	142	VAL
4	Q	169	LEU
4	Q	170	GLU
4	Q	234	LYS
4	Q	235	MET
5	R	1	VAL
5	R	22	THR
5	R	31	ASP
5	R	85	LYS
5	R	108	GLN
5	R	109	GLU
5	R	112	VAL
5	R	113	ASP
5	R	122	HIS
5	R	124	LEU
5	R	145	VAL
5	R	152	ASP
5	R	188	VAL
5	R	190	ASP
5	R	192	LEU
6	S	13	MET
6	S	70	LEU
6	S	78	GLU
6	S	84	GLU
7	T	3	HIS
7	T	4	PHE
7	T	6	ASN

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Mol	Chain	Res	Type
8	U	36	ARG
8	U	47	ARG
8	U	55	THR
9	V	49	LEU
9	V	51	CYS
9	V	54	SER
9	V	55	MET
9	V	68	ILE
9	V	77	ARG
10	W	15	ARG
10	W	16	ARG
10	W	22	LEU
10	W	40	ASP
10	W	60	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	118	GLN
1	A	173	ASN
1	A	271	HIS
2	B	392	HIS
3	C	313	GLN
3	C	332	ASN
4	D	105	ASN
4	D	166	ASN
5	E	122	HIS
7	G	23	GLN
1	N	118	GLN
1	N	136	GLN
1	N	173	ASN
1	N	271	HIS
2	O	225	ASN
2	O	363	GLN
3	P	313	GLN
3	P	332	ASN
4	Q	105	ASN
5	R	86	ASN
5	R	108	GLN
7	T	23	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z  > 2$
3	FME	C	1	3	8,8,10	1.94	2 (25%)	7,8,11	1.26	1 (14%)

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	FME	C	1	3	-	0/4/8/11	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	FME	CN-N	-4.89	1.33	1.46
3	C	1	FME	CA-C	2.20	1.53	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	FME	CN-N-CA	2.54	121.87	113.60

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

36 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
11	PEE	A	501	-	25,25,50	1.35	2 (8%)	28,30,55	1.62	5 (17%)
12	HEM	C	501	3	27,50,50	1.83	4 (14%)	17,82,82	1.55	4 (23%)
12	HEM	C	502	3	27,50,50	1.81	4 (14%)	17,82,82	1.54	1 (5%)
13	Y52	C	503	-	27,29,29	0.67	0	29,39,39	1.10	3 (10%)
14	U10	C	504	-	19,19,63	1.87	2 (10%)	21,26,79	1.53	2 (9%)
11	PEE	C	505	-	48,48,50	0.93	2 (4%)	51,53,55	1.15	4 (7%)
15	MES	C	506	-	12,12,12	2.26	1 (8%)	14,16,16	1.24	3 (21%)
11	PEE	C	507	-	10,10,50	0.51	0	11,12,55	0.54	0
16	GOL	C	508	-	5,5,5	0.35	0	5,5,5	0.34	0
11	PEE	C	509	-	46,46,50	1.02	2 (4%)	49,51,55	1.12	4 (8%)
11	PEE	C	510	-	11,14,50	0.25	0	10,14,55	0.56	0
17	HEC	D	501	4	26,50,50	1.98	2 (7%)	18,82,82	1.55	3 (16%)
18	CDL	D	502	-	41,41,99	1.30	4 (9%)	47,53,111	1.33	5 (10%)
19	BOG	D	503	-	20,20,20	0.54	0	25,25,25	0.63	0
20	FES	E	501	5	0,4,4	0.00	-	0,4,4	0.00	-
11	PEE	E	502	-	47,47,50	0.97	2 (4%)	50,52,55	1.08	4 (8%)
11	PEE	E	503	-	14,14,50	0.22	0	13,13,55	0.63	0
18	CDL	G	101	-	39,39,99	1.33	4 (10%)	45,51,111	1.37	5 (11%)
11	PEE	N	501	-	7,7,50	1.34	1 (14%)	9,9,55	1.22	1 (11%)
11	PEE	P	401	-	11,14,50	0.23	0	10,14,55	0.58	0
12	HEM	P	402	3	27,50,50	1.83	4 (14%)	17,82,82	1.54	4 (23%)
12	HEM	P	403	3	27,50,50	1.82	4 (14%)	17,82,82	1.52	2 (11%)
13	Y52	P	404	-	27,29,29	0.66	0	29,39,39	1.05	2 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
14	U10	P	405	-	19,19,63	1.92	2 (10%)	21,26,79	1.62	4 (19%)
19	BOG	P	406	-	20,20,20	0.51	0	25,25,25	0.61	0
11	PEE	P	407	-	48,48,50	0.96	2 (4%)	51,53,55	1.01	3 (5%)
16	GOL	P	408	-	5,5,5	0.37	0	5,5,5	0.28	0
11	PEE	P	409	-	40,40,50	1.03	2 (5%)	43,45,55	1.17	3 (6%)
11	PEE	P	410	-	8,11,50	0.22	0	7,11,55	0.77	0
11	PEE	P	411	-	24,24,50	1.29	2 (8%)	27,29,55	1.43	4 (14%)
17	HEC	Q	501	4	26,50,50	1.99	2 (7%)	18,82,82	1.56	3 (16%)
18	CDL	Q	502	-	41,41,99	1.29	4 (9%)	47,53,111	1.31	6 (12%)
19	BOG	Q	503	-	20,20,20	0.54	1 (5%)	25,25,25	0.59	0
20	FES	R	501	5	0,4,4	0.00	-	0,4,4	0.00	-
11	PEE	R	502	-	48,48,50	0.98	2 (4%)	51,53,55	1.11	6 (11%)
18	CDL	T	101	-	39,39,99	1.34	4 (10%)	45,51,111	1.39	5 (11%)

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
11	PEE	A	501	-	-	0/29/29/54	0/0/0/0
12	HEM	C	501	3	-	0/6/54/54	0/0/8/8
12	HEM	C	502	3	-	0/6/54/54	0/0/8/8
13	Y52	C	503	-	-	0/18/20/20	0/3/3/3
14	U10	C	504	-	-	0/11/35/87	0/1/1/1
11	PEE	C	505	-	-	0/52/52/54	0/0/0/0
15	MES	C	506	-	-	0/6/14/14	0/1/1/1
11	PEE	C	507	-	-	0/10/10/54	0/0/0/0
16	GOL	C	508	-	-	0/4/4/4	0/0/0/0
11	PEE	C	509	-	-	0/50/50/54	0/0/0/0
11	PEE	C	510	-	-	0/10/12/54	0/0/0/0
17	HEC	D	501	4	-	0/6/54/54	0/0/8/8
18	CDL	D	502	-	-	0/51/51/110	0/0/0/0
19	BOG	D	503	-	-	0/11/31/31	0/1/1/1
20	FES	E	501	5	-	0/0/4/4	0/1/1/1
11	PEE	E	502	-	-	0/51/51/54	0/0/0/0
11	PEE	E	503	-	-	0/12/12/54	0/0/0/0
18	CDL	G	101	-	-	0/49/49/110	0/0/0/0
11	PEE	N	501	-	-	0/5/5/54	0/0/0/0
11	PEE	P	401	-	-	0/10/12/54	0/0/0/0
12	HEM	P	402	3	-	0/6/54/54	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	HEM	P	403	3	-	0/6/54/54	0/0/8/8
13	Y52	P	404	-	-	0/18/20/20	0/3/3/3
14	U10	P	405	-	-	0/11/35/87	0/1/1/1
19	BOG	P	406	-	-	0/11/31/31	0/1/1/1
11	PEE	P	407	-	-	0/52/52/54	0/0/0/0
16	GOL	P	408	-	-	0/4/4/4	0/0/0/0
11	PEE	P	409	-	-	0/44/44/54	0/0/0/0
11	PEE	P	410	-	-	0/7/9/54	0/0/0/0
11	PEE	P	411	-	-	0/28/28/54	0/0/0/0
17	HEC	Q	501	4	-	0/6/54/54	0/0/8/8
18	CDL	Q	502	-	-	0/51/51/110	0/0/0/0
19	BOG	Q	503	-	-	0/11/31/31	0/1/1/1
20	FES	R	501	5	-	0/0/4/4	0/1/1/1
11	PEE	R	502	-	-	0/52/52/54	0/0/0/0
18	CDL	T	101	-	-	0/49/49/110	0/0/0/0

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	C	506	MES	C8-S	-7.58	1.66	1.77
17	Q	501	HEC	C3B-C2B	-5.90	1.34	1.40
17	D	501	HEC	C3B-C2B	-5.87	1.34	1.40
17	Q	501	HEC	C3C-C2C	-5.74	1.34	1.40
17	D	501	HEC	C3C-C2C	-5.70	1.34	1.40
12	P	402	HEM	C3C-C2C	-3.78	1.35	1.40
12	C	501	HEM	C3C-C2C	-3.71	1.35	1.40
12	C	501	HEM	C3B-C2B	-3.64	1.35	1.40
12	P	403	HEM	C3B-C2B	-3.63	1.35	1.40
12	P	402	HEM	C3B-C2B	-3.61	1.35	1.40
12	C	502	HEM	C3C-C2C	-3.57	1.35	1.40
12	P	403	HEM	C3C-C2C	-3.54	1.35	1.40
12	C	502	HEM	C3B-C2B	-3.52	1.35	1.40
19	Q	503	BOG	O1-C1	2.00	1.43	1.40
18	D	502	CDL	OA8-CA7	2.43	1.45	1.33
18	T	101	CDL	OA8-CA7	2.49	1.45	1.33
18	G	101	CDL	OA8-CA7	2.50	1.45	1.33
18	Q	502	CDL	OA8-CA7	2.51	1.45	1.33
14	P	405	U10	C4-C3	3.18	1.48	1.36
11	N	501	PEE	P-O1P	3.20	1.61	1.50
14	C	504	U10	C4-C3	3.26	1.49	1.36
18	G	101	CDL	OA6-CA5	3.93	1.45	1.34
18	T	101	CDL	OB6-CB5	3.96	1.45	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	P	402	HEM	C3B-CAB	3.97	1.55	1.47
18	Q	502	CDL	OB6-CB5	3.98	1.45	1.34
18	G	101	CDL	OB6-CB5	3.98	1.45	1.34
18	Q	502	CDL	OA6-CA5	3.98	1.45	1.34
18	D	502	CDL	OB6-CB5	3.98	1.45	1.34
12	P	402	HEM	C3C-CAC	3.98	1.55	1.47
12	C	502	HEM	C3B-CAB	3.99	1.55	1.47
12	C	501	HEM	C3C-CAC	4.00	1.55	1.47
12	C	501	HEM	C3B-CAB	4.01	1.55	1.47
12	P	403	HEM	C3B-CAB	4.01	1.55	1.47
11	A	501	PEE	O2-C10	4.04	1.46	1.34
18	T	101	CDL	OA6-CA5	4.04	1.46	1.34
18	D	502	CDL	OA6-CA5	4.05	1.46	1.34
11	C	505	PEE	O2-C10	4.05	1.46	1.34
11	P	411	PEE	O2-C10	4.06	1.46	1.34
12	P	403	HEM	C3C-CAC	4.09	1.55	1.47
12	C	502	HEM	C3C-CAC	4.09	1.55	1.47
11	P	409	PEE	O2-C10	4.13	1.46	1.34
11	P	407	PEE	O2-C10	4.18	1.46	1.34
11	C	505	PEE	O3-C30	4.19	1.45	1.33
18	Q	502	CDL	OB8-CB7	4.24	1.45	1.33
11	P	407	PEE	O3-C30	4.24	1.45	1.33
11	P	411	PEE	O3-C30	4.25	1.45	1.33
11	E	502	PEE	O3-C30	4.28	1.45	1.33
11	E	502	PEE	O2-C10	4.29	1.46	1.34
18	D	502	CDL	OB8-CB7	4.31	1.45	1.33
11	P	409	PEE	O3-C30	4.32	1.45	1.33
18	G	101	CDL	OB8-CB7	4.32	1.45	1.33
11	R	502	PEE	O3-C30	4.35	1.45	1.33
18	T	101	CDL	OB8-CB7	4.38	1.46	1.33
11	R	502	PEE	O2-C10	4.43	1.47	1.34
11	C	509	PEE	O3-C30	4.48	1.46	1.33
11	C	509	PEE	O2-C10	4.53	1.47	1.34
11	A	501	PEE	O3-C30	4.70	1.46	1.33
14	C	504	U10	C6-C1	7.04	1.50	1.35
14	P	405	U10	C6-C1	7.29	1.51	1.35

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	C	504	U10	C7-C8-C9	-4.37	119.38	126.79
14	P	405	U10	C7-C8-C9	-3.90	120.19	126.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	T	101	CDL	CB4-OB6-CB5	-3.56	109.48	117.88
18	G	101	CDL	CB4-OB6-CB5	-3.31	110.05	117.88
12	C	502	HEM	CBD-CAD-C3D	-2.94	106.86	112.47
12	P	403	HEM	CBD-CAD-C3D	-2.92	106.89	112.47
18	D	502	CDL	CB4-OB6-CB5	-2.81	111.24	117.88
11	A	501	PEE	O2-C10-O4	-2.77	116.85	123.69
17	Q	501	HEC	CMB-C2B-C1B	-2.73	124.27	128.46
18	Q	502	CDL	CB4-OB6-CB5	-2.73	111.44	117.88
14	P	405	U10	C7-C6-C5	-2.69	115.12	118.48
17	D	501	HEC	CMB-C2B-C1B	-2.69	124.33	128.46
11	C	505	PEE	O2-C10-O4	-2.51	117.50	123.69
11	P	411	PEE	O2-C10-O4	-2.38	117.82	123.69
13	C	503	Y52	C15-S16-C17	-2.35	98.97	101.10
11	C	505	PEE	O3-C30-O5	-2.34	117.86	123.58
11	A	501	PEE	O3-C30-O5	-2.31	117.95	123.58
11	E	502	PEE	O3-C30-O5	-2.29	118.00	123.58
17	D	501	HEC	CMC-C2C-C1C	-2.25	125.01	128.46
18	Q	502	CDL	CA4-OA6-CA5	-2.24	112.58	117.88
11	E	502	PEE	C3-C2-C1	-2.19	106.93	111.86
17	Q	501	HEC	CMC-C2C-C1C	-2.17	125.14	128.46
13	C	503	Y52	C25-C26-N27	-2.15	124.51	130.75
18	D	502	CDL	CA4-OA6-CA5	-2.12	112.87	117.88
18	Q	502	CDL	OB6-CB5-OB7	-2.12	118.45	123.69
11	P	411	PEE	O3-C30-O5	-2.11	118.44	123.58
11	R	502	PEE	O3-C30-O5	-2.11	118.44	123.58
12	P	402	HEM	CBD-CAD-C3D	-2.10	108.46	112.47
18	G	101	CDL	CA4-OA6-CA5	-2.10	112.92	117.88
12	C	501	HEM	CBD-CAD-C3D	-2.09	108.47	112.47
13	P	404	Y52	C25-C26-N27	-2.09	124.69	130.75
11	P	409	PEE	O3-C30-O5	-2.09	118.47	123.58
11	P	407	PEE	C2-O2-C10	-2.09	112.94	117.88
18	T	101	CDL	OB6-CB5-OB7	-2.03	118.69	123.69
12	C	501	HEM	CMC-C2C-C3C	2.04	128.60	124.88
12	P	403	HEM	CMC-C2C-C3C	2.05	128.61	124.88
17	D	501	HEC	C1D-C2D-C3D	2.06	108.43	107.00
15	C	506	MES	O1S-S-C8	2.08	109.42	106.92
12	P	402	HEM	CMC-C2C-C3C	2.09	128.69	124.88
12	P	402	HEM	C4A-C3A-C2A	2.10	108.46	107.00
17	Q	501	HEC	C1D-C2D-C3D	2.15	108.49	107.00
12	P	402	HEM	CMB-C2B-C3B	2.16	128.82	124.88
15	C	506	MES	O2S-S-C8	2.17	109.52	106.92
12	C	501	HEM	C4A-C3A-C2A	2.19	108.52	107.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	C	501	HEM	CMB-C2B-C3B	2.22	128.93	124.88
15	C	506	MES	O3S-S-C8	2.27	109.44	105.77
11	R	502	PEE	O3-C3-C2	2.30	114.38	108.64
11	R	502	PEE	C2-O2-C10	2.42	123.60	117.88
14	P	405	U10	C8-C7-C6	2.48	118.78	111.95
18	Q	502	CDL	OB8-CB7-C71	2.53	119.23	111.92
11	R	502	PEE	O2-C2-C1	2.54	117.63	108.43
11	C	509	PEE	C2-O2-C10	2.54	123.89	117.88
11	R	502	PEE	O3-C30-C31	2.58	119.39	111.92
11	C	509	PEE	O2-C2-C1	2.59	117.83	108.43
11	P	411	PEE	O3-C30-C31	2.70	119.73	111.92
18	G	101	CDL	OA6-CA5-C11	2.71	118.56	110.74
11	P	407	PEE	O3-C30-C31	2.83	120.11	111.92
18	T	101	CDL	OA6-CA5-C11	2.87	119.01	110.74
18	D	502	CDL	OB8-CB7-C71	2.97	120.51	111.92
11	C	509	PEE	O3-C30-C31	2.97	120.52	111.92
13	C	503	Y52	C19-C26-N27	2.98	114.91	108.04
13	P	404	Y52	C19-C26-N27	3.02	115.00	108.04
18	G	101	CDL	OB8-CB7-C71	3.14	119.61	111.38
18	Q	502	CDL	OA6-CA5-C11	3.14	119.81	110.74
11	N	501	PEE	O3P-P-O2P	3.17	120.12	107.59
11	E	502	PEE	O3-C30-C31	3.27	121.38	111.92
14	P	405	U10	C10-C9-C11	3.30	119.87	115.96
18	D	502	CDL	OA6-CA5-C11	3.43	120.63	110.74
11	C	505	PEE	O3-C30-C31	3.44	121.86	111.92
11	P	407	PEE	O2-C10-C11	3.46	118.84	111.55
18	T	101	CDL	OB8-CB7-C71	3.53	120.64	111.38
14	C	504	U10	C10-C9-C11	3.58	120.21	115.96
11	P	409	PEE	O2-C10-C11	3.65	119.25	111.55
11	A	501	PEE	O3-C3-C2	3.69	117.83	108.64
11	C	509	PEE	O2-C10-C11	3.74	119.43	111.55
11	A	501	PEE	O3-C30-C31	3.81	122.93	111.92
11	R	502	PEE	O2-C10-C11	3.91	119.78	111.55
11	P	409	PEE	O3-C30-C31	4.01	123.50	111.92
11	E	502	PEE	O2-C10-C11	4.11	120.21	111.55
18	D	502	CDL	OB6-CB5-C51	4.24	120.49	111.55
11	A	501	PEE	O2-C10-C11	4.33	120.68	111.55
18	G	101	CDL	OB6-CB5-C51	4.43	120.90	111.55
11	P	411	PEE	O2-C10-C11	4.47	120.97	111.55
18	Q	502	CDL	OB6-CB5-C51	4.47	120.98	111.55
11	C	505	PEE	O2-C10-C11	4.56	121.17	111.55
18	T	101	CDL	OB6-CB5-C51	4.57	121.18	111.55



There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

23 monomers are involved in 48 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	A	501	PEE	1	0
12	C	501	HEM	4	0
12	C	502	HEM	3	0
13	C	503	Y52	2	0
14	C	504	U10	1	0
16	C	508	GOL	1	0
11	C	509	PEE	2	0
18	D	502	CDL	1	0
11	E	502	PEE	4	0
11	E	503	PEE	2	0
18	G	101	CDL	1	0
11	P	401	PEE	1	0
12	P	402	HEM	5	0
12	P	403	HEM	3	0
13	P	404	Y52	1	0
14	P	405	U10	3	0
19	P	406	BOG	3	0
11	P	407	PEE	2	0
16	P	408	GOL	1	0
11	P	409	PEE	3	0
11	P	411	PEE	1	0
11	R	502	PEE	5	0
18	T	101	CDL	3	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	443/446 (99%)	0.05	8 (1%) 68 58	28, 53, 78, 101	0
1	N	442/446 (99%)	0.15	9 (2%) 65 54	36, 57, 81, 99	0
2	B	421/441 (95%)	0.24	7 (1%) 70 60	47, 67, 102, 123	0
2	O	422/441 (95%)	0.17	8 (1%) 66 56	35, 63, 86, 104	0
3	C	379/380 (99%)	-0.17	4 (1%) 80 72	22, 33, 70, 102	0
3	P	379/380 (99%)	0.03	2 (0%) 90 87	27, 51, 78, 104	0
4	D	241/241 (100%)	-0.14	3 (1%) 79 69	25, 37, 73, 95	0
4	Q	241/241 (100%)	0.14	5 (2%) 63 52	39, 59, 85, 117	0
5	E	196/196 (100%)	1.61	75 (38%) 0 0	28, 114, 166, 182	9 (4%)
5	R	196/196 (100%)	1.00	41 (20%) 1 1	36, 81, 131, 141	8 (4%)
6	F	101/110 (91%)	-0.17	0 100 100	25, 37, 53, 78	0
6	S	101/110 (91%)	0.22	2 (1%) 65 54	46, 59, 86, 102	0
7	G	80/81 (98%)	0.18	1 (1%) 77 68	28, 45, 73, 89	0
7	T	79/81 (97%)	0.65	9 (11%) 5 4	40, 70, 122, 127	0
8	H	68/77 (88%)	0.14	1 (1%) 73 64	35, 50, 68, 101	0
8	U	68/77 (88%)	1.45	20 (29%) 0 0	66, 91, 113, 117	0
9	I	30/76 (39%)	1.01	7 (23%) 0 1	55, 77, 116, 120	0
9	V	30/76 (39%)	1.22	7 (23%) 0 1	47, 78, 119, 130	0
10	J	61/61 (100%)	0.81	6 (9%) 7 6	34, 46, 79, 129	0
10	W	60/61 (98%)	0.74	3 (5%) 29 19	41, 55, 92, 117	0
All	All	4038/4218 (95%)	0.26	218 (5%) 26 16	22, 56, 107, 182	17 (0%)

All (218) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	E	163	SER	14.2
5	R	163	SER	7.5
5	E	162	GLY	7.4
10	J	62	SER	7.2
5	E	172	ARG	6.6
5	E	115	SER	6.6
5	E	82	PRO	6.3
10	J	64	GLU	6.2
5	E	84	GLY	6.0
5	E	120	PRO	5.9
5	E	121	GLN	5.6
5	R	102	THR	5.6
5	E	85	LYS	5.5
5	E	157	TYR	5.3
5	E	102	THR	5.3
9	I	51	CYS	5.2
5	R	125	ASP	5.1
5	E	83	GLU	5.1
5	R	114	VAL	5.0
5	E	117	LEU	5.0
9	V	57	GLY	5.0
9	V	51	CYS	5.0
5	E	171	ILE	4.9
5	E	104	ALA	4.8
5	E	107	ASN	4.7
5	E	176	ALA	4.7
5	E	106	ILE	4.6
5	R	117	LEU	4.4
5	E	103	GLN	4.4
5	E	174	GLY	4.2
5	R	157	TYR	4.2
5	E	150	SER	4.2
5	E	114	VAL	4.2
5	R	165	TYR	4.1
8	U	50	THR	4.0
8	U	44	VAL	4.0
5	R	120	PRO	4.0
5	E	173	LYS	4.0
5	E	116	LYS	3.9
5	E	126	ARG	3.8
10	J	63	GLU	3.8
8	U	13	LEU	3.8
5	E	164	HIS	3.8

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Mol	Chain	Res	Type	RSRZ
5	E	119	ASP	3.8
3	P	156	TYR	3.8
5	E	123	ASP	3.8
5	E	112	VAL	3.7
5	E	152	ASP	3.7
8	U	39	LEU	3.7
5	R	171	ILE	3.7
7	T	77	TYR	3.7
5	E	99	ARG	3.7
5	E	179	ASN	3.7
1	A	218	GLY	3.6
5	E	113	ASP	3.6
5	E	178	TYR	3.6
9	I	57	GLY	3.6
5	R	162	GLY	3.6
5	E	167	ALA	3.6
5	E	140	THR	3.6
5	E	118	ARG	3.5
5	E	175	PRO	3.5
5	R	113	ASP	3.5
5	R	164	HIS	3.5
5	E	125	ASP	3.5
1	N	217	SER	3.5
5	E	146	PRO	3.5
10	J	61	ALA	3.4
5	R	177	PRO	3.4
5	E	168	SER	3.4
5	E	149	ASN	3.4
3	C	8	SER	3.4
7	T	80	ASP	3.3
5	E	165	TYR	3.3
5	R	115	SER	3.3
5	R	124	LEU	3.3
1	A	217	SER	3.3
3	P	371	GLY	3.3
5	E	151	GLY	3.3
5	E	79	SER	3.2
5	E	158	CYS	3.2
5	R	119	ASP	3.2
8	U	46	SER	3.2
9	I	52	ARG	3.2
9	V	50	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
5	R	122	HIS	3.2
5	E	187	PHE	3.1
5	E	156	TYR	3.1
5	E	81	ILE	3.1
9	I	53	GLU	3.1
6	S	13	MET	3.0
5	R	156	TYR	3.0
5	E	191	ASP	3.0
5	R	170	ARG	3.0
5	R	158	CYS	3.0
8	U	35	GLU	3.0
5	E	108	GLN	3.0
5	R	112	VAL	3.0
5	R	155	GLY	3.0
5	R	84	GLY	3.0
2	B	31	ASN	3.0
8	U	37	LEU	3.0
5	E	177	PRO	3.0
9	V	48	PRO	3.0
5	E	190	ASP	2.9
9	V	54	SER	2.9
9	V	52	ARG	2.9
5	R	127	VAL	2.9
5	R	140	THR	2.9
5	R	172	ARG	2.9
2	B	230	ALA	2.9
2	B	226	ILE	2.9
5	E	185	TYR	2.9
5	E	122	HIS	2.9
5	R	105	GLU	2.8
8	U	41	ASP	2.8
2	B	227	ARG	2.8
1	N	68	LYS	2.8
2	B	229	GLY	2.8
8	U	52	GLU	2.8
8	U	28	GLU	2.7
9	I	54	SER	2.7
2	O	355	GLU	2.7
5	E	184	THR	2.7
5	E	105	GLU	2.7
7	T	78	GLU	2.7
10	W	4	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
5	E	76	ILE	2.7
8	U	45	SER	2.6
5	E	100	HIS	2.6
5	E	109	GLU	2.6
4	D	80	LEU	2.6
5	E	192	LEU	2.6
5	R	118	ARG	2.6
5	R	82	PRO	2.6
5	E	148	ALA	2.6
10	J	33	ARG	2.6
5	R	174	GLY	2.6
5	R	167	ALA	2.5
5	E	132	TRP	2.5
1	N	66	GLY	2.5
5	E	86	ASN	2.5
5	E	101	ARG	2.5
6	S	11	ARG	2.5
5	E	153	PHE	2.5
1	N	69	LYS	2.5
5	R	152	ASP	2.5
5	E	141	HIS	2.5
1	N	71	PRO	2.4
5	E	124	LEU	2.4
9	I	48	PRO	2.4
5	R	85	LYS	2.4
4	Q	1	GLY	2.4
1	N	178	THR	2.4
8	U	11	GLU	2.4
10	W	60	GLU	2.4
2	O	296	TYR	2.4
9	V	49	LEU	2.4
5	E	154	GLY	2.4
8	U	40	CYS	2.4
1	A	219	VAL	2.3
7	T	74	PRO	2.3
8	U	49	HIS	2.3
1	A	226	ASP	2.3
10	W	63	GLU	2.3
1	A	4	TYR	2.3
4	Q	81	PHE	2.3
5	R	100	HIS	2.3
4	D	79	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
7	G	80	ASP	2.3
3	C	285	ILE	2.3
1	A	223	TYR	2.3
5	R	176	ALA	2.3
1	A	69	LYS	2.2
1	N	404	ALA	2.2
5	E	194	VAL	2.2
5	E	143	GLY	2.2
1	A	222	THR	2.2
8	U	23	HIS	2.2
4	Q	82	MET	2.2
5	R	83	GLU	2.2
2	O	19	PRO	2.2
3	C	156	TYR	2.2
4	Q	77	ASN	2.2
5	R	187	PHE	2.2
5	R	109	GLU	2.2
8	H	11	GLU	2.2
8	U	43	ARG	2.2
5	E	110	ALA	2.2
2	O	350	GLY	2.2
7	T	2	ILE	2.2
5	R	143	GLY	2.2
1	N	204	SER	2.2
5	E	111	GLU	2.2
8	U	12	GLU	2.2
8	U	33	ALA	2.1
2	O	222	GLN	2.1
2	O	349	GLN	2.1
5	R	121	GLN	2.1
7	T	68	ARG	2.1
7	T	75	ALA	2.1
7	T	79	ASN	2.1
2	O	218	GLN	2.1
5	E	186	GLN	2.1
1	N	67	THR	2.1
5	E	188	VAL	2.1
10	J	60	GLU	2.1
8	U	27	THR	2.1
9	I	56	SER	2.1
4	D	78	GLY	2.1
8	U	26	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
5	R	103	GLN	2.1
4	Q	79	GLU	2.1
2	B	228	SER	2.1
2	O	200	THR	2.0
3	C	155	PRO	2.0
7	T	66	PHE	2.0
2	B	80	ALA	2.0

## 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	FME	C	1	9/11	0.70	0.41	72,88,96,99	0

## 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
11	PEE	N	501	8/51	0.72	0.28	57,79,96,112	0
11	PEE	A	501	26/51	0.72	0.33	27,87,106,123	0
11	PEE	P	409	41/51	0.72	0.34	54,69,112,128	0
18	CDL	Q	502	42/100	0.73	0.36	40,99,121,136	0
11	PEE	P	411	25/51	0.74	0.41	36,60,75,80	25
11	PEE	C	509	47/51	0.76	0.32	22,48,88,107	0
11	PEE	R	502	49/51	0.76	0.39	33,64,79,88	0
11	PEE	C	507	11/51	0.76	0.38	25,51,62,63	11
11	PEE	P	401	15/51	0.78	0.37	41,58,88,89	0
19	BOG	P	406	20/20	0.79	0.27	39,74,96,98	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
14	U10	P	405	19/63	0.80	0.39	60,75,88,91	0
18	CDL	T	101	40/100	0.81	0.32	44,85,97,98	0
11	PEE	E	503	15/51	0.81	0.33	22,49,63,68	0
11	PEE	E	502	48/51	0.81	0.33	36,55,80,86	0
18	CDL	D	502	42/100	0.83	0.30	46,80,99,127	0
15	MES	C	506	12/12	0.85	0.33	45,69,109,122	0
11	PEE	P	410	12/51	0.86	0.26	30,37,46,47	0
14	U10	C	504	19/63	0.86	0.32	45,59,67,70	0
11	PEE	P	407	49/51	0.86	0.28	45,68,84,89	0
18	CDL	G	101	40/100	0.86	0.27	24,56,69,73	0
11	PEE	C	510	15/51	0.87	0.25	25,54,81,84	0
19	BOG	D	503	20/20	0.88	0.31	29,46,54,57	0
16	GOL	C	508	6/6	0.91	0.24	34,35,38,43	0
11	PEE	C	505	49/51	0.91	0.22	22,43,54,57	0
19	BOG	Q	503	20/20	0.92	0.24	34,54,62,63	0
16	GOL	P	408	6/6	0.93	0.26	42,43,49,51	0
13	Y52	P	404	27/27	0.95	0.22	38,46,58,63	0
12	HEM	C	501	43/43	0.96	0.23	18,24,32,35	0
20	FES	R	501	4/4	0.96	0.13	71,84,92,107	0
17	HEC	D	501	43/43	0.96	0.21	26,34,40,42	0
17	HEC	Q	501	43/43	0.96	0.23	44,51,63,71	0
12	HEM	P	402	43/43	0.96	0.26	26,32,43,53	0
12	HEM	P	403	43/43	0.96	0.25	27,33,50,53	0
13	Y52	C	503	27/27	0.96	0.18	25,33,45,51	0
12	HEM	C	502	43/43	0.97	0.24	19,24,32,37	0
20	FES	E	501	4/4	0.98	0.11	92,103,103,106	4

## 6.5 Other polymers [i](#)

There are no such residues in this entry.