



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 06:47 am GMT

PDB ID : 2V67
Title : CRYSTAL STRUCTURE OF CHLAMYDOMONAS REINHARDTII RUBISCO WITH A LARGE-SUBUNIT SUPPRESSOR MUTATION T342I
Authors : Karkehabadi, S.; Satagopan, S.; Taylor, T.C.; Spreitzer, R.J.; Andersson, I.
Deposited on : 2007-07-13
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

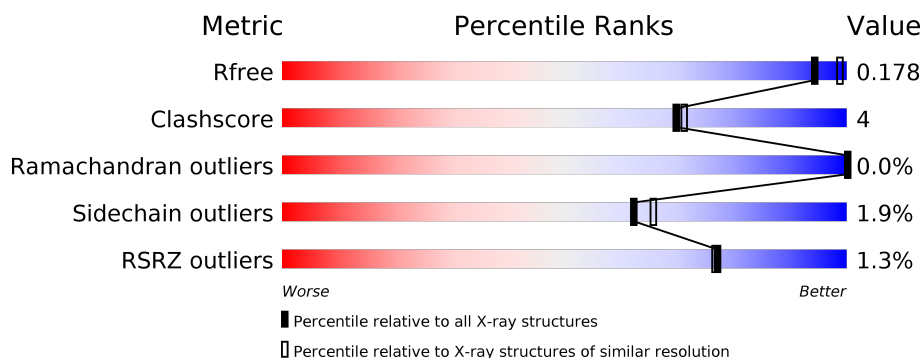
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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.






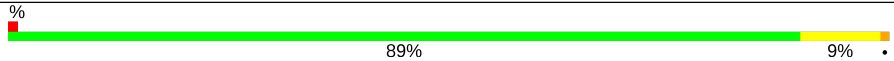
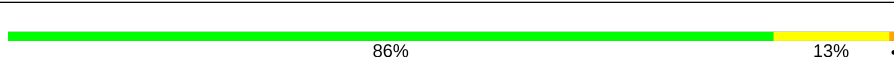
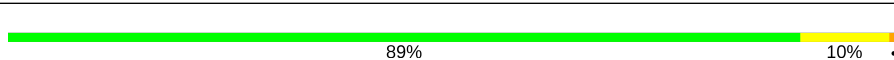
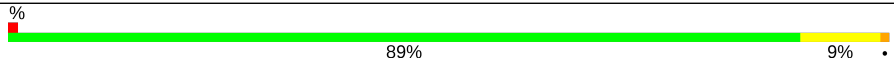
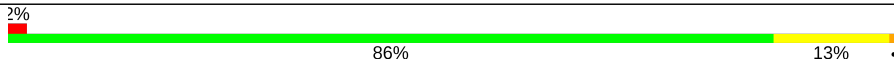
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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 ≥ 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	475	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 89%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 89% 9% </div> </div>
1	B	475	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 1%, yellow 1%, green 89%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 89% 9% </div> </div>
1	C	475	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 89%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 89% 9% </div> </div>
1	D	475	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 88%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 88% 10% </div> </div>
1	E	475	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 90%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 90% 8% </div> </div>
1	F	475	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 1%, yellow 1%, green 89%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 89% 9% </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	475	
1	H	475	
2	I	140	
2	J	140	
2	K	140	
2	L	140	
2	M	140	
2	N	140	
2	O	140	
2	P	140	

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
5	EDO	A	1478	-	-	-	X
5	EDO	A	1479	-	-	-	X
5	EDO	B	1476	-	-	-	X
5	EDO	B	1478	-	-	-	X
5	EDO	B	1480	-	-	-	X
5	EDO	C	1476	-	-	-	X
5	EDO	C	1479	-	-	-	X
5	EDO	C	1480	-	-	-	X
5	EDO	D	1477	-	-	-	X
5	EDO	D	1480	-	-	-	X
5	EDO	E	1476	-	-	-	X
5	EDO	E	1479	-	-	-	X
5	EDO	E	1481	-	-	-	X
5	EDO	F	1477	-	-	-	X
5	EDO	F	1480	-	-	-	X
5	EDO	G	1476	-	-	-	X
5	EDO	G	1480	-	-	-	X
5	EDO	G	1481	-	-	-	X
5	EDO	H	1478	-	-	-	X
5	EDO	H	1480	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	EDO	I	1141	-	-	-	X
5	EDO	N	1142	-	-	-	X
5	EDO	P	1141	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 41543 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 RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	0	4	0
			3646	2305	642	673	26			
1	B	467	Total	C	N	O	S	0	6	0
			3654	2308	643	676	27			
1	C	467	Total	C	N	O	S	0	3	0
			3649	2307	643	674	25			
1	D	465	Total	C	N	O	S	0	4	0
			3642	2303	641	672	26			
1	E	465	Total	C	N	O	S	0	4	0
			3641	2303	641	671	26			
1	F	467	Total	C	N	O	S	0	3	0
			3646	2306	640	674	26			
1	G	469	Total	C	N	O	S	0	3	0
			3664	2316	646	677	25			
1	H	466	Total	C	N	O	S	0	3	0
			3644	2304	642	673	25			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
A	342	ILE	THR	ENGINEERED MUTATION	UNP P00877
B	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
B	342	ILE	THR	ENGINEERED MUTATION	UNP P00877
C	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
C	342	ILE	THR	ENGINEERED MUTATION	UNP P00877
D	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
D	342	ILE	THR	ENGINEERED MUTATION	UNP P00877
E	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
E	342	ILE	THR	ENGINEERED MUTATION	UNP P00877
F	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
F	342	ILE	THR	ENGINEERED MUTATION	UNP P00877

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Chain	Residue	Modelled	Actual	Comment	Reference
G	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
G	342	ILE	THR	ENGINEERED MUTATION	UNP P00877
H	46	PRO	LEU	CONFLICT SEE REMARK 9	UNP P00877
H	342	ILE	THR	ENGINEERED MUTATION	UNP P00877

- Molecule 2 is a protein called RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	140	Total	C	N	O	S	0	0	0
			1143	739	190	203	11			
2	J	140	Total	C	N	O	S	0	1	0
			1145	740	190	203	12			
2	K	140	Total	C	N	O	S	0	1	0
			1145	740	190	203	12			
2	L	140	Total	C	N	O	S	0	1	0
			1147	740	193	203	11			
2	M	140	Total	C	N	O	S	0	3	0
			1152	743	193	204	12			
2	N	140	Total	C	N	O	S	0	0	0
			1143	739	190	203	11			
2	O	140	Total	C	N	O	S	0	1	0
			1145	740	190	203	12			
2	P	140	Total	C	N	O	S	0	0	0
			1143	739	190	203	11			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

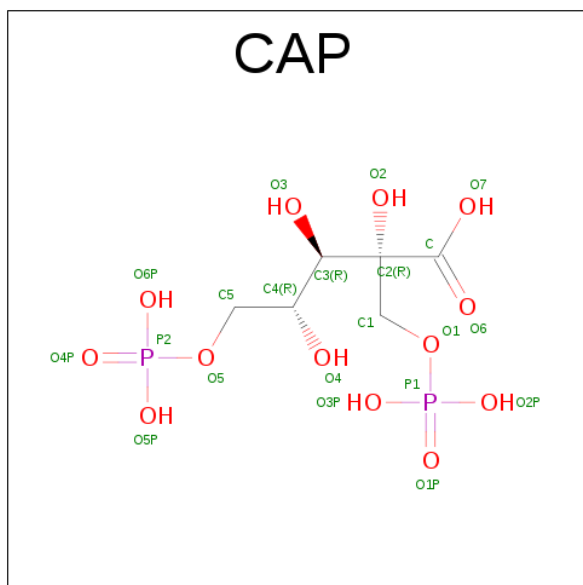
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	H	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	F	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 2-CARBOXYARABINITOL-1,5-DIPHOSPHATE (three-letter code: CAP) (formula: $C_6H_{14}O_{13}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			21	6	13	2		
4	B	1	Total	C	O	P	0	0
			21	6	13	2		
4	C	1	Total	C	O	P	0	0
			21	6	13	2		
4	D	1	Total	C	O	P	0	0
			21	6	13	2		
4	E	1	Total	C	O	P	0	0
			21	6	13	2		
4	F	1	Total	C	O	P	0	0
			21	6	13	2		
4	G	1	Total	C	O	P	0	0
			21	6	13	2		
4	H	1	Total	C	O	P	0	0
			21	6	13	2		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	G	1	Total 4	C 2	O 2	0	0
5	G	1	Total 4	C 2	O 2	0	0
5	G	1	Total 4	C 2	O 2	0	0
5	G	1	Total 4	C 2	O 2	0	0
5	H	1	Total 4	C 2	O 2	0	0
5	H	1	Total 4	C 2	O 2	0	0
5	H	1	Total 4	C 2	O 2	0	0
5	H	1	Total 4	C 2	O 2	0	0
5	H	1	Total 4	C 2	O 2	0	0
5	I	1	Total 4	C 2	O 2	0	0
5	J	1	Total 4	C 2	O 2	0	0
5	J	1	Total 4	C 2	O 2	0	0
5	K	1	Total 4	C 2	O 2	0	0
5	K	1	Total 4	C 2	O 2	0	0
5	L	1	Total 4	C 2	O 2	0	0
5	L	1	Total 4	C 2	O 2	0	0
5	M	1	Total 4	C 2	O 2	0	0
5	M	1	Total 4	C 2	O 2	0	0
5	N	1	Total 4	C 2	O 2	0	0
5	N	1	Total 4	C 2	O 2	0	0
5	O	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	O	1	Total	C	O	0	0
			4	2	2		
5	P	1	Total	C	O	0	0
			4	2	2		

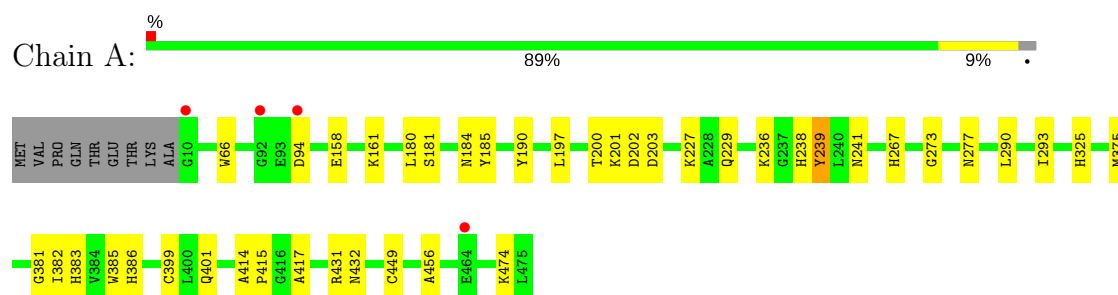
- Molecule 6 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	287	Total	O		0	0
			287	287			
6	B	272	Total	O		0	0
			272	272			
6	C	274	Total	O		0	0
			274	274			
6	D	269	Total	O		0	0
			269	269			
6	E	266	Total	O		0	0
			266	266			
6	F	241	Total	O		0	0
			241	241			
6	G	271	Total	O		0	0
			271	271			
6	H	273	Total	O		0	0
			273	273			
6	I	77	Total	O		0	0
			77	77			
6	J	78	Total	O		0	0
			78	78			
6	K	90	Total	O		0	0
			90	90			
6	L	84	Total	O		0	0
			84	84			
6	M	98	Total	O		0	0
			98	98			
6	N	64	Total	O		0	0
			64	64			
6	O	71	Total	O		0	0
			71	71			
6	P	71	Total	O		0	0
			71	71			

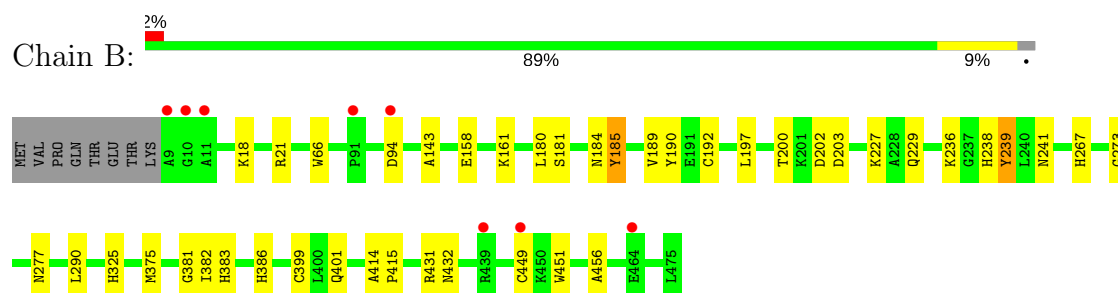
3 Residue-property plots

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

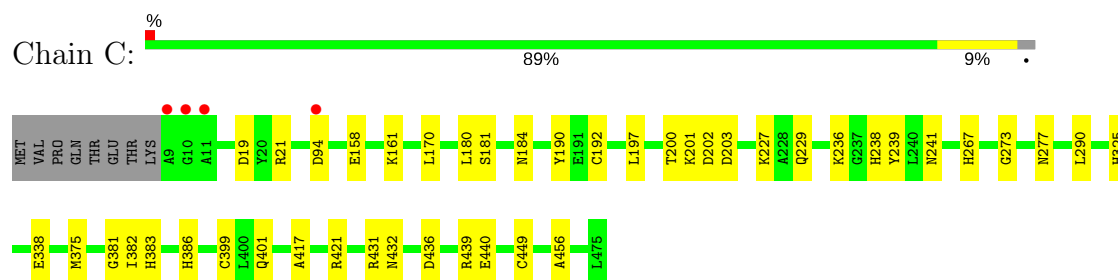
• Molecule 1: RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN



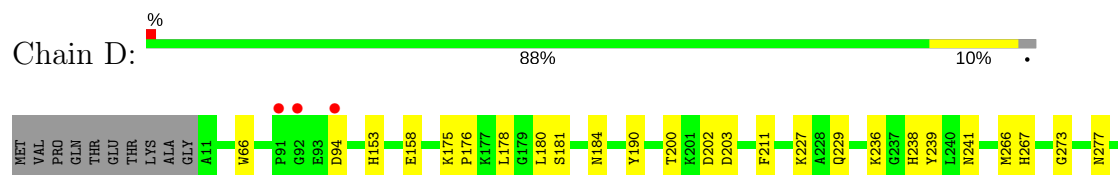
• Molecule 1: RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN



• Molecule 1: RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN

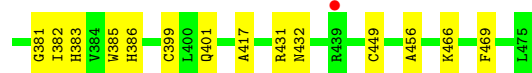
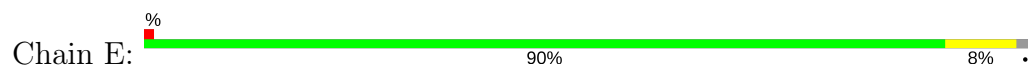


• Molecule 1: RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN

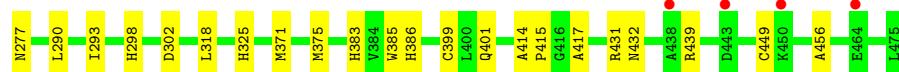
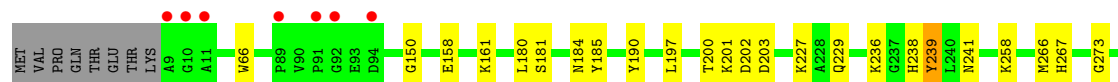
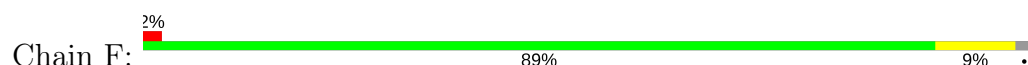




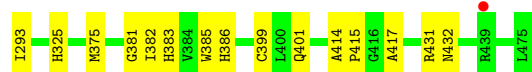
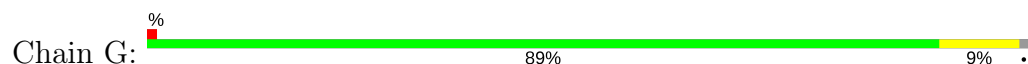
• Molecule 1: RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN



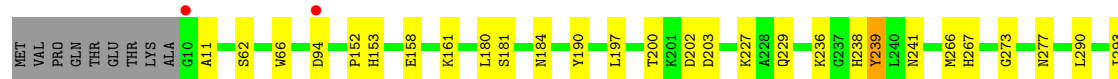
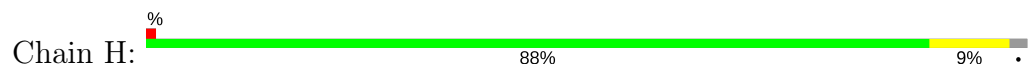
• Molecule 1: RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN



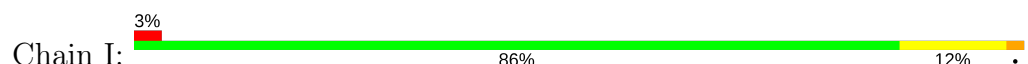
• Molecule 1: RIBULOSE BISPHOSPHATE CARBOXYLASE LARGE CHAIN



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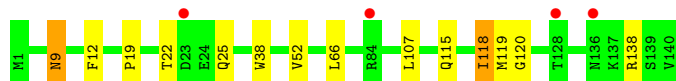
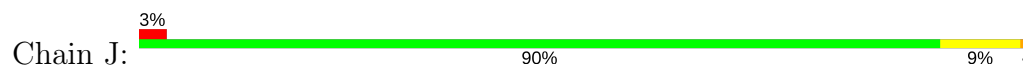


• Molecule 2: RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1

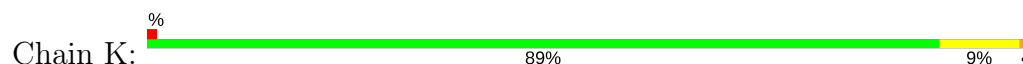




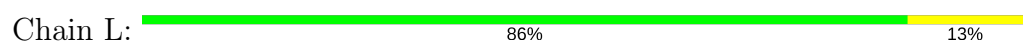
- Molecule 2: RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1



- Molecule 2: RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1



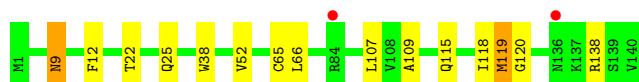
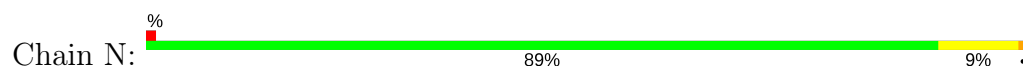
- Molecule 2: RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1



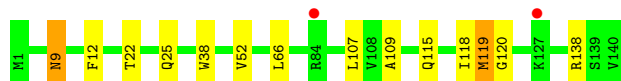
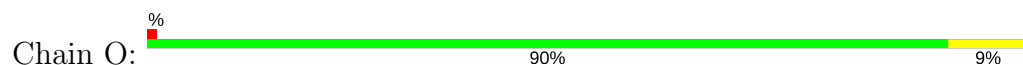
- Molecule 2: RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1



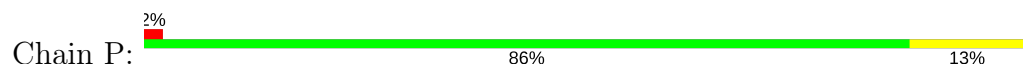
- Molecule 2: RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1

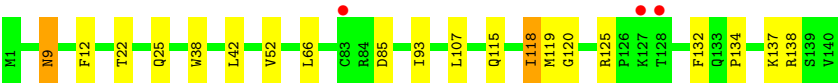


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- Molecule 2: RIBULOSE BISPHOSPHATE CARBOXYLASE SMALL CHAIN 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	120.09Å 178.56Å 122.50Å 90.00° 117.83° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 19.88 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.0 (30.00-2.00) 89.1 (19.88-2.00)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.174 , 0.208 0.177 , 0.178	Depositor DCC
R_{free} test set	13773 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	17.1	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.000 for -h-l,k,h 0.000 for l,k,-h-l 0.005 for h,-k,-h-l 0.009 for -h-l,-k,l 0.108 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	41543	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, CAP, EDO, HYP, MME, SMC, KCX

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	0.41	0/3707	0.54	0/5009
1	B	0.46	0/3727	0.56	0/5036
1	C	0.46	0/3704	0.57	0/5005
1	D	0.45	0/3703	0.56	0/5004
1	E	0.45	0/3702	0.55	0/5004
1	F	0.45	0/3701	0.55	0/5002
1	G	0.45	0/3720	0.56	0/5026
1	H	0.46	0/3699	0.56	0/4998
2	I	0.43	0/1166	0.52	0/1584
2	J	0.44	0/1174	0.55	0/1594
2	K	0.45	0/1174	0.55	0/1594
2	L	0.45	0/1177	0.55	0/1598
2	M	0.44	0/1191	0.54	0/1616
2	N	0.43	0/1166	0.53	0/1584
2	O	0.46	0/1174	0.54	0/1594
2	P	0.46	0/1166	0.56	0/1584
All	All	0.45	0/39051	0.55	0/52832

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3646	0	3553	32	0
1	B	3654	0	3553	37	0
1	C	3649	0	3557	37	0
1	D	3642	0	3550	35	0
1	E	3641	0	3550	28	0
1	F	3646	0	3553	35	0
1	G	3664	0	3573	33	0
1	H	3644	0	3552	32	0
2	I	1143	0	1122	13	0
2	J	1145	0	1123	10	0
2	K	1145	0	1123	11	0
2	L	1147	0	1123	16	0
2	M	1152	0	1133	12	0
2	N	1143	0	1122	11	0
2	O	1145	0	1123	10	0
2	P	1143	0	1122	13	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	21	0	7	0	0
4	B	21	0	8	0	0
4	C	21	0	8	0	0
4	D	21	0	7	0	0
4	E	21	0	7	0	0
4	F	21	0	7	0	0
4	G	21	0	7	0	0
4	H	21	0	8	0	0
5	A	24	0	36	0	0
5	B	20	0	30	2	0
5	C	24	0	36	0	0
5	D	20	0	30	1	0
5	E	24	0	36	2	0
5	F	20	0	30	0	0
5	G	24	0	36	0	0
5	H	20	0	30	0	0
5	I	4	0	6	0	0
5	J	8	0	12	0	0
5	K	8	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	L	8	0	12	2	0
5	M	8	0	12	0	0
5	N	8	0	12	0	0
5	O	8	0	12	0	0
5	P	4	0	6	0	0
6	A	287	0	0	3	0
6	B	272	0	0	2	0
6	C	274	0	0	3	0
6	D	269	0	0	6	0
6	E	266	0	0	3	0
6	F	241	0	0	1	0
6	G	271	0	0	2	0
6	H	273	0	0	1	0
6	I	77	0	0	1	0
6	J	78	0	0	0	0
6	K	90	0	0	1	0
6	L	84	0	0	3	0
6	M	98	0	0	0	0
6	N	64	0	0	0	0
6	O	71	0	0	0	0
6	P	71	0	0	1	0
All	All	41543	0	37839	319	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (319) 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:D:375[B]:MET:HE3	1:D:399:CYS:HB2	1.23	1.16
1:H:375[B]:MET:HE3	1:H:399:CYS:HB2	1.28	1.15
1:B:375[B]:MET:HE3	1:B:399:CYS:HB2	1.28	1.15
1:A:375[B]:MET:HE3	1:A:399:CYS:HB2	1.26	1.12
1:C:375[B]:MET:HE3	1:C:399:CYS:HB2	1.24	1.12
1:F:375[B]:MET:HE3	1:F:399:CYS:HB2	1.32	1.11
1:E:375[B]:MET:HE3	1:E:399:CYS:HB2	1.28	1.10
1:G:375[B]:MET:HE3	1:G:399:CYS:HB2	1.24	1.10
1:E:267:HIS:HD2	1:E:277:ASN:HD22	1.08	1.01
1:A:267:HIS:HD2	1:A:277:ASN:HD22	1.04	1.01
1:D:267:HIS:HD2	1:D:277:ASN:HD22	1.05	0.99
1:G:267:HIS:HD2	1:G:277:ASN:HD22	1.02	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:HIS:HD2	1:B:277:ASN:HD22	1.07	0.98
1:H:267:HIS:HD2	1:H:277:ASN:HD22	1.07	0.97
1:G:184:ASN:HD22	2:M:115:GLN:HE21	1.11	0.97
1:C:267:HIS:HD2	1:C:277:ASN:HD22	1.00	0.95
1:F:267:HIS:HD2	1:F:277:ASN:HD22	1.06	0.93
1:B:184:ASN:HD22	2:L:115:GLN:HE21	1.19	0.89
1:E:184:ASN:HD22	2:K:115:GLN:HE21	1.18	0.89
1:A:184:ASN:HD22	2:O:115:GLN:HE21	1.23	0.87
1:C:267:HIS:CD2	1:C:277:ASN:HD22	1.92	0.87
1:H:184:ASN:HD22	2:J:115:GLN:HE21	1.23	0.86
2:I:134:PRO:HG2	2:I:137:LYS:HB2	1.57	0.85
1:F:184:ASN:HD22	2:P:115:GLN:HE21	1.25	0.85
1:G:267:HIS:CD2	1:G:277:ASN:HD22	1.93	0.84
1:B:267:HIS:CD2	1:B:277:ASN:HD22	1.94	0.84
1:H:267:HIS:CD2	1:H:277:ASN:HD22	1.96	0.83
2:P:134:PRO:HG2	2:P:137:LYS:HB2	1.59	0.83
1:A:161:LYS:HE2	6:C:2131:HOH:O	1.77	0.82
1:C:184:ASN:HD22	2:I:115:GLN:HE21	1.26	0.82
1:E:431:ARG:HH21	1:E:432:ASN:HD21	1.29	0.81
1:G:431:ARG:HH21	1:G:432:ASN:HD21	1.27	0.80
1:A:267:HIS:CD2	1:A:277:ASN:HD22	1.95	0.79
1:A:431:ARG:HH21	1:A:432:ASN:HD21	1.30	0.79
1:F:431:ARG:HH21	1:F:432:ASN:HD21	1.29	0.79
1:D:184:ASN:HD22	2:N:115:GLN:HE21	1.30	0.78
1:D:431:ARG:HH21	1:D:432:ASN:HD21	1.32	0.78
1:C:431:ARG:HH21	1:C:432:ASN:HD21	1.32	0.77
2:M:22:THR:H	2:M:25:GLN:HE21	1.34	0.75
1:H:431:ARG:HH21	1:H:432:ASN:HD21	1.31	0.75
1:E:267:HIS:CD2	1:E:277:ASN:HD22	1.99	0.75
2:P:22:THR:H	2:P:25:GLN:HE21	1.35	0.73
1:D:383:HIS:H	1:D:386:HIS:HD2	1.36	0.73
1:G:93:GLU:O	6:G:2056:HOH:O	2.07	0.71
1:F:267:HIS:CD2	1:F:277:ASN:HD22	1.99	0.71
2:I:22:THR:H	2:I:25:GLN:HE21	1.38	0.71
2:L:130[B]:ARG:NH1	6:L:2080:HOH:O	2.24	0.71
1:C:383:HIS:H	1:C:386:HIS:HD2	1.36	0.71
1:D:267:HIS:CD2	1:D:277:ASN:HD22	1.97	0.71
1:F:383:HIS:H	1:F:386:HIS:HD2	1.39	0.71
2:L:22:THR:H	2:L:25:GLN:HE21	1.38	0.70
1:B:431:ARG:HH21	1:B:432:ASN:HD21	1.36	0.70
1:B:383:HIS:H	1:B:386:HIS:HD2	1.40	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:383:HIS:H	1:G:386:HIS:HD2	1.41	0.68
2:N:22:THR:H	2:N:25:GLN:HE21	1.41	0.68
1:C:436:ASP:OD2	1:C:439:ARG:HD2	1.95	0.67
2:O:22:THR:H	2:O:25:GLN:HE21	1.39	0.67
1:H:383:HIS:H	1:H:386:HIS:HD2	1.41	0.67
1:A:383:HIS:H	1:A:386:HIS:HD2	1.41	0.67
1:G:290:LEU:HG	2:O:66:LEU:HD11	1.76	0.67
1:H:290:LEU:HG	2:P:66:LEU:HD11	1.77	0.66
1:H:200:THR:OG1	1:H:238:HIS:HD2	1.79	0.66
1:B:202:ASP:OD1	1:B:238:HIS:HE1	1.79	0.66
2:J:22:THR:H	2:J:25:GLN:HE21	1.44	0.65
1:E:161:LYS:HE2	6:G:2136:HOH:O	1.97	0.65
1:D:290:LEU:HG	2:L:66:LEU:HD11	1.78	0.65
1:F:290:LEU:HG	2:N:66:LEU:HD11	1.79	0.65
1:D:375[B]:MET:CE	1:D:399:CYS:HB2	2.15	0.65
1:F:200:THR:OG1	1:F:238:HIS:HD2	1.79	0.64
1:A:290:LEU:HG	2:I:66:LEU:HD11	1.80	0.64
1:F:202:ASP:OD1	1:F:238:HIS:HE1	1.80	0.63
1:H:229:GLN:HE21	1:H:236:LYS:H	1.47	0.63
1:E:202:ASP:OD1	1:E:238:HIS:HE1	1.81	0.63
1:G:375[B]:MET:HE3	1:G:399:CYS:CB	2.16	0.63
1:C:202:ASP:OD1	1:C:238:HIS:HE1	1.81	0.62
1:C:290:LEU:HG	2:K:66:LEU:HD11	1.82	0.62
1:B:375[B]:MET:CE	1:B:399:CYS:HB2	2.19	0.62
1:A:200:THR:OG1	1:A:238:HIS:HD2	1.83	0.61
1:B:229:GLN:HE21	1:B:236:LYS:H	1.48	0.61
1:F:375[B]:MET:CE	1:F:399:CYS:HB2	2.20	0.61
1:G:202:ASP:OD1	1:G:238:HIS:HE1	1.82	0.61
1:G:181:SER:H	2:M:115:GLN:NE2	1.99	0.60
1:E:383:HIS:H	1:E:386:HIS:HD2	1.47	0.60
1:A:202:ASP:OD1	1:A:238:HIS:HE1	1.84	0.60
1:A:375[B]:MET:CE	1:A:399:CYS:HB2	2.18	0.60
2:K:22:THR:H	2:K:25:GLN:HE21	1.48	0.60
1:E:200:THR:OG1	1:E:238:HIS:HD2	1.85	0.59
1:A:229:GLN:HE21	1:A:236:LYS:H	1.51	0.59
1:G:229:GLN:HE21	1:G:236:LYS:H	1.51	0.59
6:D:2130:HOH:O	1:F:161:LYS:HE2	2.01	0.59
1:H:202:ASP:OD1	1:H:238:HIS:HE1	1.86	0.58
1:B:181:SER:H	2:L:115:GLN:NE2	2.01	0.58
1:A:181:SER:H	2:O:115:GLN:NE2	2.02	0.58
1:D:383:HIS:H	1:D:386:HIS:CD2	2.20	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:290:LEU:HG	2:M:66:LEU:HD11	1.86	0.58
1:C:383:HIS:H	1:C:386:HIS:CD2	2.19	0.57
1:H:180:LEU:HA	2:J:115:GLN:HE22	1.70	0.57
2:P:85:ASP:HB3	6:P:2047:HOH:O	2.04	0.57
1:C:180:LEU:HA	2:I:115:GLN:HE22	1.70	0.57
2:K:9:ASN:HD21	2:K:138:ARG:HG2	1.70	0.57
1:H:181:SER:H	2:J:115:GLN:NE2	2.03	0.56
1:B:200:THR:OG1	1:B:238:HIS:HD2	1.88	0.56
1:D:202:ASP:OD1	1:D:238:HIS:HE1	1.88	0.56
1:D:229:GLN:HE21	1:D:236:LYS:H	1.51	0.56
1:B:267:HIS:HD2	1:B:277:ASN:ND2	1.90	0.56
1:C:200:THR:OG1	1:C:238:HIS:HD2	1.89	0.56
1:E:469:PHE:CE2	5:E:1480:EDO:H21	2.41	0.56
1:H:383:HIS:H	1:H:386:HIS:CD2	2.24	0.55
1:G:200:THR:OG1	1:G:238:HIS:HD2	1.90	0.55
1:C:181:SER:H	2:I:115:GLN:NE2	2.03	0.55
1:F:158:GLU:CD	1:F:325:HIS:HE2	2.09	0.55
6:F:2116:HOH:O	1:H:161:LYS:HE2	2.06	0.55
1:C:229:GLN:HE21	1:C:236:LYS:H	1.53	0.55
2:J:9:ASN:HD21	2:J:138:ARG:HG2	1.71	0.55
1:E:229:GLN:HE21	1:E:236:LYS:H	1.54	0.55
1:E:466:LYS:O	5:E:1480:EDO:H22	2.06	0.55
1:F:180:LEU:HA	2:P:115:GLN:HE22	1.72	0.54
1:D:180:LEU:HA	2:N:115:GLN:HE22	1.73	0.54
1:B:180:LEU:HA	2:L:115:GLN:HE22	1.72	0.54
1:F:267:HIS:HD2	1:F:277:ASN:ND2	1.91	0.54
2:I:85:ASP:HB3	6:I:2048:HOH:O	2.07	0.54
2:N:22:THR:H	2:N:25:GLN:NE2	2.05	0.54
1:B:383:HIS:H	1:B:386:HIS:CD2	2.24	0.53
1:B:290:LEU:HG	2:J:66:LEU:HD11	1.89	0.53
1:A:180:LEU:HA	2:O:115:GLN:HE22	1.72	0.53
1:F:239:TYR:HE2	1:F:401:GLN:HE22	1.57	0.53
2:O:9:ASN:HD21	2:O:138:ARG:HG2	1.74	0.53
1:C:375[B]:MET:CE	1:C:399:CYS:HB2	2.17	0.53
1:E:239:TYR:HE2	1:E:401:GLN:HE22	1.57	0.53
1:G:375[B]:MET:CE	1:G:399:CYS:HB2	2.16	0.52
2:M:9:ASN:HD21	2:M:138:ARG:HG2	1.74	0.52
1:A:383:HIS:H	1:A:386:HIS:CD2	2.24	0.52
1:D:375[B]:MET:HE3	1:D:399:CYS:CB	2.16	0.52
2:L:9:ASN:HD21	2:L:138:ARG:HG2	1.75	0.52
2:L:87:MET:HG2	6:L:2055:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:22:THR:H	2:M:25:GLN:NE2	2.05	0.52
1:C:338:GLU:HG3	6:C:2259:HOH:O	2.10	0.52
1:B:18:LYS:O	5:B:1477:EDO:H21	2.10	0.52
1:G:158:GLU:CD	1:G:325:HIS:HE2	2.12	0.51
1:E:180:LEU:HA	2:K:115:GLN:HE22	1.75	0.51
2:I:22:THR:H	2:I:25:GLN:NE2	2.06	0.51
1:C:267:HIS:HE1	6:D:2159:HOH:O	1.93	0.51
1:E:383:HIS:H	1:E:386:HIS:CD2	2.27	0.51
2:I:9:ASN:HD21	2:I:138:ARG:HG2	1.75	0.51
2:P:22:THR:H	2:P:25:GLN:NE2	2.06	0.51
1:G:383:HIS:H	1:G:386:HIS:CD2	2.25	0.51
2:P:125:ARG:HD2	2:P:132:PHE:CE2	2.46	0.51
1:C:161:LYS:HE2	6:E:2116:HOH:O	2.09	0.51
1:F:181:SER:H	2:P:115:GLN:NE2	2.09	0.50
2:I:125:ARG:HD2	2:I:132:PHE:CE2	2.46	0.50
1:B:158:GLU:CD	1:B:325:HIS:HE2	2.15	0.50
1:D:158:GLU:CD	1:D:325:HIS:HE2	2.15	0.50
1:G:180:LEU:HA	2:M:115:GLN:HE22	1.76	0.50
1:B:185:TYR:O	1:B:189:VAL:HG23	2.12	0.50
1:C:382:ILE:HA	1:C:386:HIS:CD2	2.46	0.50
2:P:9:ASN:HD21	2:P:138:ARG:HG2	1.76	0.50
1:D:435:ARG:NH2	6:D:2239:HOH:O	2.43	0.49
1:F:298:HIS:ND1	1:F:302:ASP:OD2	2.35	0.49
1:D:239:TYR:HE2	1:D:401:GLN:HE22	1.59	0.49
1:D:341:VAL:HG11	5:D:1480:EDO:H22	1.95	0.49
1:D:200:THR:OG1	1:D:238:HIS:HD2	1.94	0.49
1:C:436:ASP:O	1:C:440:GLU:HB2	2.12	0.49
1:F:229:GLN:HE21	1:F:236:LYS:H	1.59	0.49
1:G:414:ALA:HB3	1:G:415:PRO:HD3	1.94	0.49
1:E:181:SER:H	2:K:115:GLN:NE2	2.11	0.49
1:H:375[B]:MET:CE	1:H:399:CYS:HB2	2.20	0.49
2:L:22:THR:H	2:L:25:GLN:NE2	2.08	0.49
2:K:24:GLU:HB2	6:K:2024:HOH:O	2.12	0.48
1:B:449:CYS:HB3	1:B:456:ALA:HA	1.94	0.48
2:K:22:THR:H	2:K:25:GLN:NE2	2.09	0.48
1:F:383:HIS:H	1:F:386:HIS:CD2	2.26	0.48
1:A:239:TYR:HE2	1:A:401:GLN:HE22	1.61	0.48
1:B:239:TYR:HE2	1:B:401:GLN:HE22	1.61	0.48
1:E:158:GLU:CD	1:E:325:HIS:HE2	2.14	0.48
2:O:22:THR:H	2:O:25:GLN:NE2	2.09	0.48
2:P:107:LEU:O	2:P:120:GLY:HA2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:273:GLY:HA3	1:H:273:GLY:HA3	1.95	0.48
1:F:449:CYS:HB3	1:F:456:ALA:HA	1.96	0.47
1:C:201:KCX:HB2	1:C:239:TYR:CD2	2.49	0.47
1:B:382:ILE:HA	1:B:386:HIS:CD2	2.49	0.47
1:E:382:ILE:HA	1:E:386:HIS:CD2	2.49	0.47
1:E:377:VAL:HG22	1:E:399:CYS:HB3	1.96	0.47
1:A:474:LYS:HD2	6:A:2277:HOH:O	2.14	0.47
1:C:449:CYS:HB3	1:C:456:ALA:HA	1.96	0.47
1:D:153:HIS:HE1	6:D:2165:HOH:O	1.95	0.47
1:D:446:ARG:NH1	6:D:2243:HOH:O	2.44	0.47
6:E:2148:HOH:O	1:F:267:HIS:HE1	1.97	0.47
1:H:293:ILE:HG13	1:H:318:LEU:HD21	1.95	0.47
1:C:192:CYS:HB3	1:C:197:LEU:HD12	1.96	0.47
1:A:158:GLU:CD	1:A:325:HIS:HE2	2.15	0.47
1:F:190:TYR:CZ	1:F:227:LYS:HE3	2.50	0.47
5:B:1480:EDO:O1	6:B:2270:HOH:O	2.18	0.47
6:D:2138:HOH:O	2:L:10:LYS:HE3	2.14	0.47
1:D:181:SER:H	2:N:115:GLN:NE2	2.14	0.46
1:D:449:CYS:O	1:D:456:ALA:HB2	2.16	0.46
1:G:239:TYR:HE2	1:G:401:GLN:HE22	1.62	0.46
1:H:190:TYR:CZ	1:H:227:LYS:HE3	2.50	0.46
1:A:382:ILE:HA	1:A:386:HIS:CD2	2.50	0.46
2:N:9:ASN:HD21	2:N:138:ARG:HG2	1.79	0.46
1:A:273:GLY:HA3	1:B:273:GLY:HA3	1.98	0.46
1:C:375[B]:MET:HE3	1:C:399:CYS:CB	2.18	0.46
1:A:381:GLY:HA2	1:B:66:TRP:CD1	2.51	0.46
1:D:190:TYR:CZ	1:D:227:LYS:HE3	2.51	0.46
1:C:381:GLY:HA2	1:D:66:TRP:CD1	2.51	0.46
1:F:201:KCX:HB2	1:F:239:TYR:CD2	2.51	0.46
1:G:382:ILE:HA	1:G:386:HIS:CD2	2.51	0.46
2:J:22:THR:H	2:J:25:GLN:NE2	2.11	0.46
2:O:109:ALA:HB3	2:O:119:MET:HG2	1.98	0.46
1:H:239:TYR:HE2	1:H:401:GLN:HE22	1.63	0.46
1:A:414:ALA:HB3	1:A:415:PRO:HD3	1.98	0.46
1:E:382:ILE:HA	1:E:386:HIS:HD2	1.81	0.46
1:A:267:HIS:HE1	6:B:2164:HOH:O	1.98	0.45
1:D:239:TYR:HB3	1:D:266:MET:HB3	1.98	0.45
1:D:382:ILE:HA	1:D:386:HIS:CD2	2.51	0.45
1:F:375[B]:MET:HE3	1:F:399:CYS:CB	2.23	0.45
1:B:190:TYR:CZ	1:B:227:LYS:HE3	2.51	0.45
1:F:239:TYR:HE2	1:F:401:GLN:NE2	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:449:CYS:O	1:H:456:ALA:HB2	2.16	0.45
2:O:107:LEU:O	2:O:120:GLY:HA2	2.17	0.45
1:C:158:GLU:CD	1:C:325:HIS:HE2	2.17	0.45
1:H:382:ILE:HA	1:H:386:HIS:CD2	2.51	0.45
1:H:302:ASP:OD2	1:H:311:PHE:HB2	2.17	0.45
1:G:197:LEU:HG	1:G:417:ALA:HB1	1.99	0.45
2:P:38:TRP:CD2	2:P:118:ILE:HG21	2.52	0.45
1:C:239:TYR:HE2	1:C:401:GLN:HE22	1.64	0.44
1:C:273:GLY:HA3	1:D:273:GLY:HA3	1.99	0.44
2:P:42:LEU:HD21	2:P:93:ILE:HG12	1.99	0.44
1:A:375[B]:MET:HE3	1:A:399:CYS:CB	2.19	0.44
1:A:190:TYR:CZ	1:A:227:LYS:HE3	2.52	0.44
1:C:383:HIS:N	1:C:386:HIS:HD2	2.09	0.44
1:F:383:HIS:CE1	1:F:385:TRP:HB2	2.52	0.44
1:H:197:LEU:HG	1:H:417:ALA:HB1	1.99	0.44
1:G:177:LYS:HB2	1:H:62:SER:O	2.18	0.44
1:B:414:ALA:HB3	1:B:415:PRO:HD3	2.00	0.44
6:A:2172:HOH:O	1:B:267:HIS:HE1	1.99	0.44
1:A:449:CYS:HB3	1:A:456:ALA:HA	2.00	0.44
1:C:197:LEU:HG	1:C:417:ALA:HB1	2.00	0.44
1:E:197:LEU:HG	1:E:417:ALA:HB1	2.00	0.44
1:H:239:TYR:HB3	1:H:266:MET:HB3	2.00	0.44
1:E:381:GLY:HA2	1:F:66:TRP:CD1	2.53	0.44
1:H:152:PRO:HB2	1:H:153:HIS:CD2	2.52	0.44
2:I:107:LEU:O	2:I:120:GLY:HA2	2.18	0.43
1:F:150:GLY:HA3	1:F:371[B]:MET:SD	2.58	0.43
2:M:107:LEU:O	2:M:120:GLY:HA2	2.18	0.43
1:H:158:GLU:CD	1:H:325:HIS:HE2	2.17	0.43
2:K:42:LEU:HD21	2:K:93:ILE:HG12	1.99	0.43
2:J:107:LEU:O	2:J:120:GLY:HA2	2.19	0.43
1:C:190:TYR:CZ	1:C:227:LYS:HE3	2.53	0.43
1:H:449:CYS:HB3	1:H:456:ALA:HA	2.01	0.43
1:F:258:LYS:HA	2:N:65:CYS:SG	2.59	0.43
1:A:200:THR:OG1	1:A:238:HIS:CD2	2.68	0.43
1:D:277:ASN:HD21	1:D:293:ILE:HD12	1.83	0.43
1:F:200:THR:OG1	1:F:238:HIS:CD2	2.66	0.43
2:I:38:TRP:CD2	2:I:118:ILE:HG21	2.54	0.43
1:C:19:ASP:HB3	1:C:21:ARG:HG2	1.99	0.43
1:D:295:ARG:HG2	1:D:327:HIS:HB2	2.00	0.43
6:A:2145:HOH:O	1:G:161:LYS:HE2	2.19	0.43
1:G:190:TYR:CZ	1:G:227:LYS:HE3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:414:ALA:HB3	1:D:415:PRO:HD3	2.01	0.42
2:L:107:LEU:O	2:L:120:GLY:HA2	2.19	0.42
1:G:383:HIS:CE1	1:G:385:TRP:HB2	2.55	0.42
2:L:125:ARG:HD2	2:L:132:PHE:CE2	2.54	0.42
1:F:293:ILE:HG13	1:F:318:LEU:HD21	2.01	0.42
1:G:381:GLY:HA2	1:H:66:TRP:CD1	2.55	0.42
1:H:414:ALA:HB3	1:H:415:PRO:HD3	2.02	0.42
1:B:184:ASN:HD22	2:L:115:GLN:NE2	2.01	0.42
1:C:267:HIS:HD2	1:C:277:ASN:ND2	1.85	0.42
1:G:181:SER:H	2:M:115:GLN:HE22	1.65	0.42
2:J:38:TRP:CD2	2:J:118:ILE:HG21	2.54	0.42
2:K:38:TRP:CD2	2:K:118:ILE:HG21	2.54	0.42
1:B:192:CYS:HB3	1:B:197:LEU:HD12	2.01	0.42
1:B:143:ALA:HA	1:G:143:ALA:HA	2.02	0.42
1:B:161:LYS:HE2	6:H:2129:HOH:O	2.19	0.42
1:E:277:ASN:HD21	1:E:293:ILE:HD12	1.84	0.42
1:E:383:HIS:CE1	1:E:385:TRP:HB2	2.55	0.42
1:G:201:KCX:HB2	1:G:239:TYR:CD2	2.55	0.42
1:G:277:ASN:HD21	1:G:293:ILE:HD12	1.84	0.42
1:C:200:THR:OG1	1:C:238:HIS:CD2	2.72	0.42
1:D:383:HIS:CE1	1:D:385:TRP:HB2	2.55	0.42
1:E:273:GLY:HA3	1:F:273:GLY:HA3	2.01	0.41
2:L:42:LEU:HD21	2:L:93:ILE:HG12	2.01	0.41
2:M:125:ARG:HD2	2:M:132:PHE:CE2	2.54	0.41
2:N:38:TRP:CD2	2:N:118:ILE:HG21	2.55	0.41
6:C:2152:HOH:O	1:D:267:HIS:HE1	2.03	0.41
1:F:197:LEU:HG	1:F:417:ALA:HB1	2.02	0.41
2:O:38:TRP:CD2	2:O:118:ILE:HG21	2.54	0.41
1:A:197:LEU:HG	1:A:417:ALA:HB1	2.01	0.41
1:B:181:SER:H	2:L:115:GLN:HE22	1.69	0.41
1:F:239:TYR:HB3	1:F:266:MET:HB3	2.02	0.41
1:G:185:TYR:O	1:G:189:VAL:HG23	2.19	0.41
1:H:200:THR:OG1	1:H:238:HIS:CD2	2.67	0.41
1:H:382:ILE:HA	1:H:386:HIS:HD2	1.86	0.41
1:A:201:KCX:HB2	1:A:239:TYR:CD2	2.55	0.41
1:A:277:ASN:HD21	1:A:293:ILE:HD12	1.85	0.41
1:E:338:GLU:HG3	6:E:2256:HOH:O	2.20	0.41
1:E:449:CYS:O	1:E:456:ALA:HB2	2.20	0.41
1:A:383:HIS:CE1	1:A:385:TRP:HB2	2.55	0.41
2:K:125:ARG:HD2	2:K:132:PHE:CE2	2.55	0.41
1:B:227:LYS:HA	5:L:1141:EDO:H12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:451:TRP:CE2	2:J:19:PRO:HG3	2.55	0.41
1:C:382:ILE:HA	1:C:386:HIS:HD2	1.85	0.41
1:C:449:CYS:O	1:C:456:ALA:HB2	2.21	0.41
1:F:414:ALA:HB3	1:F:415:PRO:HD3	2.03	0.41
1:B:200:THR:OG1	1:B:238:HIS:CD2	2.72	0.41
1:D:457:ALA:O	1:D:461:VAL:HG23	2.21	0.41
2:I:109:ALA:HB3	2:I:119:MET:HG2	2.03	0.41
1:A:66:TRP:CD1	1:B:381:GLY:HA2	2.56	0.41
1:D:178:LEU:HD22	1:D:211:PHE:HZ	1.86	0.41
2:M:42:LEU:HD21	2:M:93:ILE:HG12	2.03	0.41
2:L:38:TRP:CD2	2:L:118:ILE:HG21	2.55	0.40
2:N:107:LEU:O	2:N:120:GLY:HA2	2.21	0.40
1:C:170:LEU:HD11	1:C:421:ARG:HA	2.03	0.40
5:L:1141:EDO:H11	6:L:2044:HOH:O	2.21	0.40
2:N:109:ALA:HB3	2:N:119:MET:HG2	2.03	0.40
1:B:383:HIS:N	1:B:386:HIS:HD2	2.13	0.40
1:D:175:LYS:HA	1:D:176:PRO:C	2.41	0.40
1:G:239:TYR:HB3	1:G:266:MET:HB3	2.03	0.40
1:B:449:CYS:O	1:B:456:ALA:HB2	2.21	0.40
2:M:109:ALA:HB3	2:M:119:MET:HG2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	463/475 (98%)	451 (97%)	12 (3%)	0	100	100
1	B	466/475 (98%)	454 (97%)	12 (3%)	0	100	100
1	C	463/475 (98%)	450 (97%)	13 (3%)	0	100	100
1	D	462/475 (97%)	450 (97%)	12 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	462/475 (97%)	448 (97%)	14 (3%)	0	100	100
1	F	463/475 (98%)	449 (97%)	14 (3%)	0	100	100
1	G	465/475 (98%)	451 (97%)	14 (3%)	0	100	100
1	H	462/475 (97%)	450 (97%)	11 (2%)	1 (0%)	51	48
2	I	138/140 (99%)	132 (96%)	6 (4%)	0	100	100
2	J	139/140 (99%)	134 (96%)	5 (4%)	0	100	100
2	K	139/140 (99%)	133 (96%)	6 (4%)	0	100	100
2	L	139/140 (99%)	133 (96%)	6 (4%)	0	100	100
2	M	141/140 (101%)	135 (96%)	6 (4%)	0	100	100
2	N	138/140 (99%)	132 (96%)	6 (4%)	0	100	100
2	O	139/140 (99%)	134 (96%)	5 (4%)	0	100	100
2	P	138/140 (99%)	132 (96%)	6 (4%)	0	100	100
All	All	4817/4920 (98%)	4668 (97%)	148 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	11	ALA

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	372/376 (99%)	367 (99%)	5 (1%)	73	78
1	B	374/376 (100%)	368 (98%)	6 (2%)	68	72
1	C	371/376 (99%)	368 (99%)	3 (1%)	85	88
1	D	372/376 (99%)	369 (99%)	3 (1%)	85	88
1	E	372/376 (99%)	368 (99%)	4 (1%)	78	82
1	F	371/376 (99%)	366 (99%)	5 (1%)	73	78

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	373/376 (99%)	366 (98%)	7 (2%)	62	66
1	H	371/376 (99%)	366 (99%)	5 (1%)	73	78
2	I	122/122 (100%)	116 (95%)	6 (5%)	29	24
2	J	123/122 (101%)	118 (96%)	5 (4%)	35	31
2	K	123/122 (101%)	119 (97%)	4 (3%)	43	41
2	L	123/122 (101%)	118 (96%)	5 (4%)	35	31
2	M	125/122 (102%)	121 (97%)	4 (3%)	44	42
2	N	122/122 (100%)	118 (97%)	4 (3%)	43	41
2	O	123/122 (101%)	119 (97%)	4 (3%)	43	41
2	P	122/122 (100%)	117 (96%)	5 (4%)	35	31
All	All	3959/3984 (99%)	3884 (98%)	75 (2%)	62	66

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	ASP
1	A	185	TYR
1	A	203	ASP
1	A	239	TYR
1	A	241	ASN
1	B	21	ARG
1	B	94	ASP
1	B	185	TYR
1	B	203	ASP
1	B	239	TYR
1	B	241	ASN
1	C	94	ASP
1	C	203	ASP
1	C	241	ASN
1	D	94	ASP
1	D	203	ASP
1	D	241	ASN
1	E	185	TYR
1	E	203	ASP
1	E	239	TYR
1	E	241	ASN
1	F	185	TYR
1	F	203	ASP

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Mol	Chain	Res	Type
1	F	239	TYR
1	F	241	ASN
1	F	439	ARG
1	G	8	LYS
1	G	94	ASP
1	G	172	CYS
1	G	185	TYR
1	G	203	ASP
1	G	239	TYR
1	G	241	ASN
1	H	94	ASP
1	H	203	ASP
1	H	239	TYR
1	H	241	ASN
1	H	439	ARG
2	I	9	ASN
2	I	12	PHE
2	I	52	VAL
2	I	84	ARG
2	I	118	ILE
2	I	119	MET
2	J	9	ASN
2	J	12	PHE
2	J	52	VAL
2	J	118	ILE
2	J	119	MET
2	K	9	ASN
2	K	12	PHE
2	K	118	ILE
2	K	119	MET
2	L	9	ASN
2	L	12	PHE
2	L	52	VAL
2	L	118	ILE
2	L	119	MET
2	M	9	ASN
2	M	12	PHE
2	M	52	VAL
2	M	119	MET
2	N	9	ASN
2	N	12	PHE
2	N	52	VAL

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Mol	Chain	Res	Type
2	N	119	MET
2	O	9	ASN
2	O	12	PHE
2	O	52	VAL
2	O	119	MET
2	P	9	ASN
2	P	12	PHE
2	P	52	VAL
2	P	118	ILE
2	P	119	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (125) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	HIS
1	A	229	GLN
1	A	238	HIS
1	A	241	ASN
1	A	267	HIS
1	A	277	ASN
1	A	304	GLN
1	A	386	HIS
1	A	401	GLN
1	A	432	ASN
1	B	153	HIS
1	B	229	GLN
1	B	238	HIS
1	B	241	ASN
1	B	267	HIS
1	B	277	ASN
1	B	304	GLN
1	B	386	HIS
1	B	401	GLN
1	B	432	ASN
1	C	153	HIS
1	C	163	ASN
1	C	229	GLN
1	C	238	HIS
1	C	241	ASN
1	C	267	HIS
1	C	277	ASN
1	C	304	GLN

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Mol	Chain	Res	Type
1	C	386	HIS
1	C	401	GLN
1	C	432	ASN
1	D	153	HIS
1	D	229	GLN
1	D	238	HIS
1	D	241	ASN
1	D	267	HIS
1	D	277	ASN
1	D	304	GLN
1	D	386	HIS
1	D	401	GLN
1	D	432	ASN
1	E	153	HIS
1	E	229	GLN
1	E	238	HIS
1	E	241	ASN
1	E	267	HIS
1	E	277	ASN
1	E	304	GLN
1	E	386	HIS
1	E	401	GLN
1	E	432	ASN
1	F	153	HIS
1	F	229	GLN
1	F	238	HIS
1	F	241	ASN
1	F	267	HIS
1	F	277	ASN
1	F	304	GLN
1	F	386	HIS
1	F	401	GLN
1	F	432	ASN
1	G	153	HIS
1	G	229	GLN
1	G	238	HIS
1	G	241	ASN
1	G	267	HIS
1	G	277	ASN
1	G	304	GLN
1	G	386	HIS
1	G	401	GLN

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Mol	Chain	Res	Type
1	G	420	ASN
1	G	432	ASN
1	H	153	HIS
1	H	207	ASN
1	H	229	GLN
1	H	238	HIS
1	H	241	ASN
1	H	267	HIS
1	H	277	ASN
1	H	304	GLN
1	H	386	HIS
1	H	401	GLN
1	H	432	ASN
2	I	8	ASN
2	I	9	ASN
2	I	25	GLN
2	I	29	GLN
2	I	115	GLN
2	I	133	GLN
2	J	9	ASN
2	J	25	GLN
2	J	29	GLN
2	J	115	GLN
2	J	133	GLN
2	K	9	ASN
2	K	25	GLN
2	K	29	GLN
2	K	115	GLN
2	K	133	GLN
2	L	8	ASN
2	L	9	ASN
2	L	25	GLN
2	L	29	GLN
2	L	115	GLN
2	L	133	GLN
2	M	8	ASN
2	M	9	ASN
2	M	25	GLN
2	M	29	GLN
2	M	115	GLN
2	N	8	ASN
2	N	9	ASN

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Mol	Chain	Res	Type
2	N	25	GLN
2	N	29	GLN
2	N	115	GLN
2	O	8	ASN
2	O	9	ASN
2	O	25	GLN
2	O	29	GLN
2	O	115	GLN
2	P	9	ASN
2	P	25	GLN
2	P	29	GLN
2	P	115	GLN
2	P	133	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

48 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	HYP	A	104	1	7,8,9	0.63	0	5,10,12	1.13	1 (20%)
1	HYP	A	151	1	7,8,9	0.69	0	5,10,12	1.26	1 (20%)
1	KCX	A	201	1,3	8,11,12	1.05	0	6,12,14	0.82	0
1	SMC	A	256	1	6,6,7	0.81	0	3,6,8	0.94	0
1	SMC	A	369	1	6,6,7	0.88	0	3,6,8	1.13	0
1	HYP	B	104	1	7,8,9	0.67	0	5,10,12	1.09	1 (20%)
1	HYP	B	151	1	7,8,9	0.51	0	5,10,12	1.54	1 (20%)
1	KCX	B	201	1,3	8,11,12	1.12	1 (12%)	6,12,14	0.83	0
1	SMC	B	256	1	6,6,7	0.69	0	3,6,8	1.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	SMC	B	369	1	6,6,7	0.84	0	3,6,8	1.75	1 (33%)
1	HYP	C	104	1	7,8,9	0.50	0	5,10,12	1.22	1 (20%)
1	HYP	C	151	1	7,8,9	0.82	0	5,10,12	1.38	1 (20%)
1	KCX	C	201	1,3	8,11,12	0.91	0	6,12,14	0.82	0
1	SMC	C	256	1	6,6,7	0.69	0	3,6,8	1.17	0
1	SMC	C	369	1	6,6,7	0.92	0	3,6,8	1.25	0
1	HYP	D	104	1	7,8,9	0.52	0	5,10,12	1.20	1 (20%)
1	HYP	D	151	1	7,8,9	0.67	0	5,10,12	1.23	1 (20%)
1	KCX	D	201	1,3	8,11,12	1.07	1 (12%)	6,12,14	1.17	0
1	SMC	D	256	1	6,6,7	1.07	1 (16%)	3,6,8	1.23	0
1	SMC	D	369	1	6,6,7	0.92	0	3,6,8	1.10	0
1	HYP	E	104	1	7,8,9	0.60	0	5,10,12	1.07	1 (20%)
1	HYP	E	151	1	7,8,9	1.00	1 (14%)	5,10,12	1.31	1 (20%)
1	KCX	E	201	1,3	8,11,12	0.92	1 (12%)	6,12,14	1.26	1 (16%)
1	SMC	E	256	1	6,6,7	0.70	0	3,6,8	1.04	0
1	SMC	E	369	1	6,6,7	0.63	0	3,6,8	1.42	0
1	HYP	F	104	1	7,8,9	0.75	0	5,10,12	1.14	1 (20%)
1	HYP	F	151	1	7,8,9	0.57	0	5,10,12	1.25	1 (20%)
1	KCX	F	201	1,3	8,11,12	0.87	0	6,12,14	1.12	0
1	SMC	F	256	1	6,6,7	1.07	0	3,6,8	1.02	0
1	SMC	F	369	1	6,6,7	1.28	1 (16%)	3,6,8	1.26	0
1	HYP	G	104	1	7,8,9	0.68	0	5,10,12	1.20	1 (20%)
1	HYP	G	151	1	7,8,9	0.56	0	5,10,12	1.42	1 (20%)
1	KCX	G	201	1,3	8,11,12	0.89	0	6,12,14	1.00	0
1	SMC	G	256	1	6,6,7	0.62	0	3,6,8	1.14	0
1	SMC	G	369	1	6,6,7	0.81	0	3,6,8	0.88	0
1	HYP	H	104	1	7,8,9	0.67	0	5,10,12	1.23	1 (20%)
1	HYP	H	151	1	7,8,9	0.61	0	5,10,12	1.56	1 (20%)
1	KCX	H	201	1,3	8,11,12	0.97	0	6,12,14	0.84	0
1	SMC	H	256	1	6,6,7	0.81	0	3,6,8	1.28	0
1	SMC	H	369	1	6,6,7	1.00	1 (16%)	3,6,8	1.05	0
2	MME	I	1	2	8,8,9	2.59	1 (12%)	7,8,10	1.44	1 (14%)
2	MME	J	1	2	8,8,9	2.65	1 (12%)	7,8,10	1.30	1 (14%)
2	MME	K	1	2	8,8,9	2.54	1 (12%)	7,8,10	1.40	2 (28%)
2	MME	L	1	2	8,8,9	2.54	1 (12%)	7,8,10	1.31	1 (14%)
2	MME	M	1	2	8,8,9	2.58	1 (12%)	7,8,10	1.41	2 (28%)
2	MME	N	1	2	8,8,9	2.58	1 (12%)	7,8,10	1.48	2 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MME	O	1	2	8,8,9	2.61	1 (12%)	7,8,10	1.36	2 (28%)
2	MME	P	1	2	8,8,9	2.55	1 (12%)	7,8,10	1.78	2 (28%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	A	104	1	-	0/0/11/13	0/1/1/1
1	HYP	A	151	1	-	0/0/11/13	0/1/1/1
1	KCX	A	201	1,3	-	0/6/10/12	0/0/0/0
1	SMC	A	256	1	-	0/3/5/7	0/0/0/0
1	SMC	A	369	1	-	0/3/5/7	0/0/0/0
1	HYP	B	104	1	-	0/0/11/13	0/1/1/1
1	HYP	B	151	1	-	0/0/11/13	0/1/1/1
1	KCX	B	201	1,3	-	0/6/10/12	0/0/0/0
1	SMC	B	256	1	-	0/3/5/7	0/0/0/0
1	SMC	B	369	1	-	0/3/5/7	0/0/0/0
1	HYP	C	104	1	-	0/0/11/13	0/1/1/1
1	HYP	C	151	1	-	0/0/11/13	0/1/1/1
1	KCX	C	201	1,3	-	0/6/10/12	0/0/0/0
1	SMC	C	256	1	-	0/3/5/7	0/0/0/0
1	SMC	C	369	1	-	0/3/5/7	0/0/0/0
1	HYP	D	104	1	-	0/0/11/13	0/1/1/1
1	HYP	D	151	1	-	0/0/11/13	0/1/1/1
1	KCX	D	201	1,3	-	0/6/10/12	0/0/0/0
1	SMC	D	256	1	-	0/3/5/7	0/0/0/0
1	SMC	D	369	1	-	0/3/5/7	0/0/0/0
1	HYP	E	104	1	-	0/0/11/13	0/1/1/1
1	HYP	E	151	1	-	0/0/11/13	0/1/1/1
1	KCX	E	201	1,3	-	0/6/10/12	0/0/0/0
1	SMC	E	256	1	-	0/3/5/7	0/0/0/0
1	SMC	E	369	1	-	0/3/5/7	0/0/0/0
1	HYP	F	104	1	-	0/0/11/13	0/1/1/1
1	HYP	F	151	1	-	0/0/11/13	0/1/1/1
1	KCX	F	201	1,3	-	0/6/10/12	0/0/0/0
1	SMC	F	256	1	-	0/3/5/7	0/0/0/0
1	SMC	F	369	1	-	0/3/5/7	0/0/0/0
1	HYP	G	104	1	-	0/0/11/13	0/1/1/1
1	HYP	G	151	1	-	0/0/11/13	0/1/1/1
1	KCX	G	201	1,3	-	0/6/10/12	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SMC	G	256	1	-	0/3/5/7	0/0/0/0
1	SMC	G	369	1	-	0/3/5/7	0/0/0/0
1	HYP	H	104	1	-	0/0/11/13	0/1/1/1
1	HYP	H	151	1	-	0/0/11/13	0/1/1/1
1	KCX	H	201	1,3	-	0/6/10/12	0/0/0/0
1	SMC	H	256	1	-	0/3/5/7	0/0/0/0
1	SMC	H	369	1	-	0/3/5/7	0/0/0/0
2	MME	I	1	2	-	0/4/8/10	0/0/0/0
2	MME	J	1	2	-	0/4/8/10	0/0/0/0
2	MME	K	1	2	-	0/4/8/10	0/0/0/0
2	MME	L	1	2	-	0/4/8/10	0/0/0/0
2	MME	M	1	2	-	0/4/8/10	0/0/0/0
2	MME	N	1	2	-	0/4/8/10	0/0/0/0
2	MME	O	1	2	-	0/4/8/10	0/0/0/0
2	MME	P	1	2	-	0/4/8/10	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	1	MME	CM-N	-7.17	1.26	1.46
2	I	1	MME	CM-N	-7.07	1.27	1.46
2	K	1	MME	CM-N	-7.03	1.27	1.46
2	N	1	MME	CM-N	-7.03	1.27	1.46
2	O	1	MME	CM-N	-7.00	1.27	1.46
2	P	1	MME	CM-N	-6.99	1.27	1.46
2	M	1	MME	CM-N	-6.90	1.27	1.46
2	L	1	MME	CM-N	-6.88	1.27	1.46
1	E	201	KCX	CE-NZ	2.01	1.50	1.46
1	H	369	SMC	CB-SG	2.12	1.83	1.80
1	B	201	KCX	CA-C	2.13	1.53	1.50
1	D	201	KCX	CE-NZ	2.21	1.51	1.46
1	E	151	HYP	CA-C	2.24	1.53	1.50
1	D	256	SMC	CB-SG	2.36	1.83	1.80
1	F	369	SMC	CB-SG	2.52	1.83	1.80

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	1	MME	CB-CA-C	-3.52	105.85	111.65
1	H	151	HYP	O-C-CA	-2.59	119.10	125.15
2	N	1	MME	CB-CA-C	-2.56	107.44	111.65
1	B	151	HYP	O-C-CA	-2.56	119.18	125.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	201	KCX	CE-NZ-CX	-2.48	120.31	123.35
1	G	151	HYP	O-C-CA	-2.46	119.41	125.15
2	I	1	MME	CB-CA-C	-2.46	107.60	111.65
1	B	369	SMC	CA-CB-SG	-2.42	110.89	114.22
1	A	151	HYP	O-C-CA	-2.37	119.62	125.15
1	F	151	HYP	O-C-CA	-2.35	119.66	125.15
1	C	151	HYP	O-C-CA	-2.33	119.72	125.15
1	D	104	HYP	O-C-CA	-2.31	119.77	125.15
1	C	104	HYP	O-C-CA	-2.27	119.86	125.15
1	E	151	HYP	O-C-CA	-2.26	119.87	125.15
1	E	104	HYP	O-C-CA	-2.17	120.09	125.15
2	K	1	MME	O-C-CA	-2.16	120.12	125.15
1	D	151	HYP	O-C-CA	-2.15	120.13	125.15
1	A	104	HYP	O-C-CA	-2.14	120.16	125.15
1	G	104	HYP	O-C-CA	-2.12	120.20	125.15
1	H	104	HYP	O-C-CA	-2.12	120.20	125.15
2	M	1	MME	O-C-CA	-2.07	120.32	125.15
2	J	1	MME	CB-CA-C	-2.06	108.26	111.65
1	B	104	HYP	O-C-CA	-2.04	120.39	125.15
2	O	1	MME	CB-CA-C	-2.02	108.31	111.65
1	F	104	HYP	O-C-CA	-2.00	120.48	125.15
2	N	1	MME	CM-N-CA	2.01	120.15	113.60
2	O	1	MME	CM-N-CA	2.01	120.16	113.60
2	L	1	MME	CM-N-CA	2.09	120.40	113.60
2	K	1	MME	CM-N-CA	2.13	120.53	113.60
2	M	1	MME	CM-N-CA	2.29	121.08	113.60
2	P	1	MME	CM-N-CA	2.29	121.08	113.60

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	201	KCX	1	0
1	C	201	KCX	1	0
1	F	201	KCX	1	0
1	G	201	KCX	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 74 ligands modelled in this entry, 8 are monoatomic - leaving 66 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	EDO	A	1476	-	3,3,3	0.57	0	2,2,2	0.27	0
5	EDO	A	1477	-	3,3,3	0.40	0	2,2,2	0.45	0
5	EDO	A	1478	-	3,3,3	0.52	0	2,2,2	0.31	0
5	EDO	A	1479	-	3,3,3	0.43	0	2,2,2	0.36	0
5	EDO	A	1480	-	3,3,3	0.57	0	2,2,2	0.26	0
5	EDO	A	1481	-	3,3,3	0.46	0	2,2,2	0.36	0
4	CAP	A	477	3	14,20,20	0.75	0	17,31,31	0.58	0
5	EDO	B	1476	-	3,3,3	0.64	0	2,2,2	0.12	0
5	EDO	B	1477	-	3,3,3	0.44	0	2,2,2	0.31	0
5	EDO	B	1478	-	3,3,3	0.49	0	2,2,2	0.25	0
5	EDO	B	1479	-	3,3,3	0.50	0	2,2,2	0.32	0
5	EDO	B	1480	-	3,3,3	0.46	0	2,2,2	0.45	0
4	CAP	B	477	3	14,20,20	0.70	0	17,31,31	0.54	0
5	EDO	C	1476	-	3,3,3	0.49	0	2,2,2	0.31	0
5	EDO	C	1477	-	3,3,3	0.56	0	2,2,2	0.28	0
5	EDO	C	1478	-	3,3,3	0.54	0	2,2,2	0.24	0
5	EDO	C	1479	-	3,3,3	0.54	0	2,2,2	0.07	0
5	EDO	C	1480	-	3,3,3	0.48	0	2,2,2	0.39	0
5	EDO	C	1481	-	3,3,3	0.52	0	2,2,2	0.12	0
4	CAP	C	477	3	14,20,20	0.78	0	17,31,31	0.60	0
5	EDO	D	1476	-	3,3,3	0.48	0	2,2,2	0.26	0
5	EDO	D	1477	-	3,3,3	0.50	0	2,2,2	0.26	0
5	EDO	D	1478	-	3,3,3	0.54	0	2,2,2	0.09	0
5	EDO	D	1479	-	3,3,3	0.60	0	2,2,2	0.10	0
5	EDO	D	1480	-	3,3,3	0.41	0	2,2,2	0.51	0
4	CAP	D	477	3	14,20,20	0.81	0	17,31,31	0.68	0
5	EDO	E	1476	-	3,3,3	0.50	0	2,2,2	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	E	1477	-	3,3,3	0.61	0	2,2,2	0.08	0
5	EDO	E	1478	-	3,3,3	0.36	0	2,2,2	0.56	0
5	EDO	E	1479	-	3,3,3	0.47	0	2,2,2	0.41	0
5	EDO	E	1480	-	3,3,3	0.43	0	2,2,2	0.49	0
5	EDO	E	1481	-	3,3,3	0.45	0	2,2,2	0.40	0
4	CAP	E	477	3	14,20,20	0.73	0	17,31,31	0.65	0
5	EDO	F	1476	-	3,3,3	0.42	0	2,2,2	0.54	0
5	EDO	F	1477	-	3,3,3	0.49	0	2,2,2	0.32	0
5	EDO	F	1478	-	3,3,3	0.53	0	2,2,2	0.25	0
5	EDO	F	1479	-	3,3,3	0.56	0	2,2,2	0.14	0
5	EDO	F	1480	-	3,3,3	0.48	0	2,2,2	0.34	0
4	CAP	F	477	3	14,20,20	0.79	0	17,31,31	0.66	0
5	EDO	G	1476	-	3,3,3	0.45	0	2,2,2	0.35	0
5	EDO	G	1477	-	3,3,3	0.52	0	2,2,2	0.23	0
5	EDO	G	1478	-	3,3,3	0.37	0	2,2,2	0.56	0
5	EDO	G	1479	-	3,3,3	0.50	0	2,2,2	0.38	0
5	EDO	G	1480	-	3,3,3	0.46	0	2,2,2	0.15	0
5	EDO	G	1481	-	3,3,3	0.55	0	2,2,2	0.29	0
4	CAP	G	477	3	14,20,20	0.75	0	17,31,31	0.63	0
5	EDO	H	1476	-	3,3,3	0.64	0	2,2,2	0.05	0
5	EDO	H	1477	-	3,3,3	0.41	0	2,2,2	0.45	0
5	EDO	H	1478	-	3,3,3	0.54	0	2,2,2	0.35	0
5	EDO	H	1479	-	3,3,3	0.51	0	2,2,2	0.31	0
5	EDO	H	1480	-	3,3,3	0.44	0	2,2,2	0.34	0
4	CAP	H	477	3	14,20,20	0.84	1 (7%)	17,31,31	0.55	0
5	EDO	I	1141	-	3,3,3	0.46	0	2,2,2	0.26	0
5	EDO	J	1141	-	3,3,3	0.44	0	2,2,2	0.26	0
5	EDO	J	1142	-	3,3,3	0.47	0	2,2,2	0.24	0
5	EDO	K	1141	-	3,3,3	0.49	0	2,2,2	0.22	0
5	EDO	K	1142	-	3,3,3	0.46	0	2,2,2	0.34	0
5	EDO	L	1141	-	3,3,3	0.45	0	2,2,2	0.28	0
5	EDO	L	1142	-	3,3,3	0.48	0	2,2,2	0.34	0
5	EDO	M	1141	-	3,3,3	0.46	0	2,2,2	0.30	0
5	EDO	M	1142	-	3,3,3	0.50	0	2,2,2	0.18	0
5	EDO	N	1141	-	3,3,3	0.48	0	2,2,2	0.39	0
5	EDO	N	1142	-	3,3,3	0.52	0	2,2,2	0.23	0
5	EDO	O	1141	-	3,3,3	0.53	0	2,2,2	0.17	0
5	EDO	O	1142	-	3,3,3	0.42	0	2,2,2	0.44	0
5	EDO	P	1141	-	3,3,3	0.44	0	2,2,2	0.38	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. '-' means

no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1478	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1480	-	-	0/1/1/1	0/0/0/0
5	EDO	A	1481	-	-	0/1/1/1	0/0/0/0
4	CAP	A	477	3	-	0/23/29/29	0/0/0/0
5	EDO	B	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	B	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	B	1478	-	-	0/1/1/1	0/0/0/0
5	EDO	B	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	B	1480	-	-	0/1/1/1	0/0/0/0
4	CAP	B	477	3	-	0/23/29/29	0/0/0/0
5	EDO	C	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	C	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	C	1478	-	-	0/1/1/1	0/0/0/0
5	EDO	C	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	C	1480	-	-	0/1/1/1	0/0/0/0
5	EDO	C	1481	-	-	0/1/1/1	0/0/0/0
4	CAP	C	477	3	-	0/23/29/29	0/0/0/0
5	EDO	D	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	D	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	D	1478	-	-	0/1/1/1	0/0/0/0
5	EDO	D	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	D	1480	-	-	0/1/1/1	0/0/0/0
4	CAP	D	477	3	-	0/23/29/29	0/0/0/0
5	EDO	E	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	E	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	E	1478	-	-	0/1/1/1	0/0/0/0
5	EDO	E	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	E	1480	-	-	0/1/1/1	0/0/0/0
5	EDO	E	1481	-	-	0/1/1/1	0/0/0/0
4	CAP	E	477	3	-	0/23/29/29	0/0/0/0
5	EDO	F	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	F	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	F	1478	-	-	0/1/1/1	0/0/0/0
5	EDO	F	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	F	1480	-	-	0/1/1/1	0/0/0/0
4	CAP	F	477	3	-	0/23/29/29	0/0/0/0
5	EDO	G	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	G	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	G	1478	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	G	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	G	1480	-	-	0/1/1/1	0/0/0/0
5	EDO	G	1481	-	-	0/1/1/1	0/0/0/0
4	CAP	G	477	3	-	0/23/29/29	0/0/0/0
5	EDO	H	1476	-	-	0/1/1/1	0/0/0/0
5	EDO	H	1477	-	-	0/1/1/1	0/0/0/0
5	EDO	H	1478	-	-	0/1/1/1	0/0/0/0
5	EDO	H	1479	-	-	0/1/1/1	0/0/0/0
5	EDO	H	1480	-	-	0/1/1/1	0/0/0/0
4	CAP	H	477	3	-	0/23/29/29	0/0/0/0
5	EDO	I	1141	-	-	0/1/1/1	0/0/0/0
5	EDO	J	1141	-	-	0/1/1/1	0/0/0/0
5	EDO	J	1142	-	-	0/1/1/1	0/0/0/0
5	EDO	K	1141	-	-	0/1/1/1	0/0/0/0
5	EDO	K	1142	-	-	0/1/1/1	0/0/0/0
5	EDO	L	1141	-	-	0/1/1/1	0/0/0/0
5	EDO	L	1142	-	-	0/1/1/1	0/0/0/0
5	EDO	M	1141	-	-	0/1/1/1	0/0/0/0
5	EDO	M	1142	-	-	0/1/1/1	0/0/0/0
5	EDO	N	1141	-	-	0/1/1/1	0/0/0/0
5	EDO	N	1142	-	-	0/1/1/1	0/0/0/0
5	EDO	O	1141	-	-	0/1/1/1	0/0/0/0
5	EDO	O	1142	-	-	0/1/1/1	0/0/0/0
5	EDO	P	1141	-	-	0/1/1/1	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	477	CAP	O2-C2	2.03	1.46	1.43

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1477	EDO	1	0
5	B	1480	EDO	1	0
5	D	1480	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	1480	EDO	2	0
5	L	1141	EDO	2	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	461/475 (97%)	-0.55	4 (0%) 84 83	9, 14, 30, 40	0
1	B	462/475 (97%)	-0.58	8 (1%) 70 69	9, 14, 30, 40	0
1	C	462/475 (97%)	-0.57	4 (0%) 84 83	9, 14, 29, 38	0
1	D	460/475 (96%)	-0.59	6 (1%) 77 77	9, 14, 30, 40	0
1	E	460/475 (96%)	-0.58	5 (1%) 80 80	9, 15, 30, 41	0
1	F	462/475 (97%)	-0.53	11 (2%) 59 59	9, 15, 30, 40	1 (0%)
1	G	464/475 (97%)	-0.58	5 (1%) 80 80	9, 14, 30, 51	0
1	H	461/475 (97%)	-0.60	5 (1%) 80 80	9, 14, 30, 42	0
2	I	139/140 (99%)	-0.22	4 (2%) 52 52	12, 20, 33, 36	1 (0%)
2	J	139/140 (99%)	-0.26	4 (2%) 52 52	12, 20, 31, 35	1 (0%)
2	K	139/140 (99%)	-0.31	1 (0%) 87 87	12, 20, 31, 37	1 (0%)
2	L	139/140 (99%)	-0.35	0 100 100	12, 20, 33, 36	0
2	M	139/140 (99%)	-0.36	0 100 100	12, 19, 30, 33	0
2	N	139/140 (99%)	-0.29	2 (1%) 75 75	12, 20, 35, 41	1 (0%)
2	O	139/140 (99%)	-0.31	2 (1%) 75 75	12, 20, 31, 35	1 (0%)
2	P	139/140 (99%)	-0.26	3 (2%) 62 61	12, 20, 32, 36	1 (0%)
All	All	4804/4920 (97%)	-0.51	64 (1%) 77 77	9, 16, 31, 51	7 (0%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	10	GLY	5.4
1	E	92	GLY	5.1
1	B	9	ALA	4.8
1	G	8	LYS	4.8
1	B	10	GLY	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	10	GLY	4.1
1	C	10	GLY	4.1
1	F	10	GLY	4.0
1	C	11	ALA	3.8
1	F	91	PRO	3.8
1	G	7	THR	3.7
1	G	9	ALA	3.7
1	F	9	ALA	3.6
1	F	94	ASP	3.5
1	H	10	GLY	3.5
1	B	94	ASP	3.4
1	C	94	ASP	3.4
1	H	94	ASP	3.4
1	E	11	ALA	3.3
1	E	94	ASP	3.3
2	P	128	THR	3.3
1	F	11	ALA	3.2
2	I	84	ARG	3.1
2	P	83	CYS	3.1
1	D	91	PRO	3.1
1	G	439	ARG	3.1
1	C	9	ALA	3.0
1	F	92	GLY	2.9
1	D	92	GLY	2.9
1	B	11	ALA	2.9
1	D	94	ASP	2.8
2	P	127	LYS	2.8
1	A	94	ASP	2.8
1	B	439	ARG	2.7
1	H	464	GLU	2.7
1	D	439	ARG	2.7
1	E	91	PRO	2.6
1	F	443	ASP	2.6
1	A	92	GLY	2.6
1	E	439	ARG	2.5
1	H	439	ARG	2.5
1	F	464	GLU	2.5
2	I	128	THR	2.5
2	I	83	CYS	2.4
1	A	464	GLU	2.4
1	B	464	GLU	2.3
1	H	449	CYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	438	ALA	2.3
1	F	89	PRO	2.3
2	J	23	ASP	2.3
2	J	84	ARG	2.3
2	J	128	THR	2.3
2	O	84	ARG	2.2
2	N	136	ASN	2.2
1	B	449	CYS	2.2
1	B	91	PRO	2.2
1	F	450	LYS	2.2
2	N	84	ARG	2.2
2	I	127	LYS	2.1
1	D	449	CYS	2.1
2	J	136	ASN	2.0
1	D	450	LYS	2.0
2	O	127	LYS	2.0
2	K	84	ARG	2.0

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

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	A	104	8/9	0.97	0.06	-	11,11,12,13	0
1	HYP	F	104	8/9	0.96	0.07	-	11,12,12,13	0
1	HYP	E	104	8/9	0.97	0.06	-	11,12,12,13	0
1	KCX	F	201	12/13	0.97	0.07	-	10,11,11,12	0
1	SMC	D	256	7/8	0.98	0.08	-	9,10,10,10	0
1	SMC	D	369	7/8	0.98	0.06	-	14,14,16,17	0
1	SMC	A	256	7/8	0.99	0.06	-	10,10,10,11	0
1	SMC	F	256	7/8	0.98	0.06	-	10,10,11,12	0
1	SMC	H	369	7/8	0.97	0.10	-	14,15,16,18	0
1	KCX	A	201	12/13	0.98	0.07	-	9,11,11,11	0
1	HYP	E	151	8/9	0.97	0.07	-	11,12,12,12	0
2	MME	O	1	9/10	0.90	0.16	-	24,25,32,33	0
2	MME	N	1	9/10	0.94	0.20	-	25,26,33,33	0
2	MME	I	1	9/10	0.95	0.13	-	24,25,32,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	C	151	8/9	0.97	0.07	-	11,12,12,12	0
1	KCX	B	201	12/13	0.98	0.07	-	9,10,11,12	0
1	HYP	H	151	8/9	0.95	0.08	-	11,11,12,12	0
1	HYP	D	104	8/9	0.97	0.08	-	11,11,12,12	0
1	SMC	F	369	7/8	0.96	0.09	-	14,15,16,18	0
1	SMC	C	369	7/8	0.98	0.07	-	14,15,16,18	0
1	HYP	F	151	8/9	0.97	0.06	-	11,12,12,12	0
1	SMC	B	369	7/8	0.97	0.06	-	14,14,16,17	0
1	SMC	C	256	7/8	0.99	0.06	-	9,10,10,11	0
1	HYP	B	151	8/9	0.97	0.06	-	11,12,12,12	0
1	SMC	H	256	7/8	0.98	0.07	-	10,10,10,11	0
1	SMC	G	369	7/8	0.98	0.06	-	14,14,16,17	0
1	KCX	E	201	12/13	0.98	0.06	-	9,10,11,11	0
2	MME	K	1	9/10	0.89	0.15	-	24,25,33,33	0
1	SMC	E	256	7/8	0.99	0.06	-	9,10,10,11	0
1	SMC	G	256	7/8	0.99	0.06	-	9,10,10,11	0
1	HYP	H	104	8/9	0.97	0.07	-	11,12,12,13	0
1	KCX	D	201	12/13	0.97	0.08	-	10,11,11,11	0
2	MME	J	1	9/10	0.90	0.17	-	24,26,32,33	0
2	MME	L	1	9/10	0.91	0.17	-	24,25,32,33	0
1	SMC	A	369	7/8	0.98	0.08	-	14,14,16,18	0
2	MME	M	1	9/10	0.94	0.14	-	24,25,32,32	0
1	HYP	A	151	8/9	0.96	0.07	-	11,12,12,12	0
1	KCX	G	201	12/13	0.97	0.07	-	10,10,11,11	0
1	HYP	B	104	8/9	0.98	0.06	-	11,11,12,13	0
1	HYP	G	104	8/9	0.97	0.07	-	11,11,12,12	0
2	MME	P	1	9/10	0.91	0.14	-	24,25,32,33	0
1	HYP	D	151	8/9	0.98	0.06	-	11,12,12,12	0
1	KCX	H	201	12/13	0.98	0.06	-	9,11,11,11	0
1	SMC	E	369	7/8	0.97	0.07	-	14,14,16,18	0
1	SMC	B	256	7/8	0.99	0.05	-	10,10,10,11	0
1	KCX	C	201	12/13	0.97	0.07	-	10,10,11,12	0
1	HYP	G	151	8/9	0.97	0.07	-	11,12,12,12	0
1	HYP	C	104	8/9	0.97	0.07	-	11,11,12,13	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	EDO	G	1480	4/4	0.92	0.32	24.73	8,8,8,8	4
5	EDO	E	1476	4/4	0.97	0.32	22.14	8,8,9,9	4
5	EDO	C	1479	4/4	0.90	0.32	17.54	9,10,10,11	4
5	EDO	D	1477	4/4	0.67	0.21	10.15	42,42,42,44	0
5	EDO	C	1480	4/4	0.73	0.19	7.36	48,48,48,48	0
5	EDO	F	1480	4/4	0.72	0.28	6.70	49,49,49,49	0
5	EDO	A	1478	4/4	0.79	0.17	6.40	40,41,42,43	0
5	EDO	C	1476	4/4	0.79	0.28	5.71	45,46,46,47	0
5	EDO	B	1478	4/4	0.90	0.11	5.15	35,35,35,36	0
5	EDO	F	1477	4/4	0.86	0.13	4.39	40,40,41,42	0
5	EDO	G	1481	4/4	0.77	0.18	4.27	37,38,38,38	0
5	EDO	G	1476	4/4	0.85	0.20	4.13	41,42,42,43	0
5	EDO	D	1480	4/4	0.83	0.23	4.02	47,47,47,48	0
5	EDO	H	1478	4/4	0.78	0.14	3.77	38,38,38,38	0
5	EDO	H	1480	4/4	0.93	0.22	3.48	55,55,55,55	0
5	EDO	E	1481	4/4	0.83	0.21	3.02	45,47,47,48	0
5	EDO	E	1479	4/4	0.67	0.21	2.82	49,50,50,50	0
5	EDO	P	1141	4/4	0.86	0.16	2.73	38,38,38,40	0
5	EDO	I	1141	4/4	0.86	0.16	2.72	35,35,36,36	0
5	EDO	N	1142	4/4	0.22	0.42	2.38	54,55,57,57	0
5	EDO	B	1480	4/4	0.89	0.17	2.20	44,45,45,45	0
5	EDO	A	1479	4/4	0.88	0.20	2.13	49,50,50,50	0
5	EDO	B	1476	4/4	0.92	0.12	2.01	18,21,22,22	0
5	EDO	M	1141	4/4	0.83	0.15	1.76	47,47,48,48	0
5	EDO	L	1141	4/4	0.88	0.18	1.62	46,46,46,47	0
5	EDO	E	1478	4/4	0.95	0.13	1.47	24,25,26,26	0
5	EDO	F	1479	4/4	0.88	0.14	1.40	27,28,29,29	0
5	EDO	J	1141	4/4	0.90	0.13	1.11	32,32,32,33	0
5	EDO	C	1478	4/4	0.94	0.10	0.90	22,22,22,24	0
5	EDO	L	1142	4/4	0.87	0.19	0.88	49,50,51,51	0
5	EDO	M	1142	4/4	0.87	0.18	0.84	53,53,53,53	0
5	EDO	O	1141	4/4	0.94	0.12	0.72	33,33,33,34	0
5	EDO	E	1477	4/4	0.89	0.12	0.71	25,26,27,27	0
5	EDO	H	1476	4/4	0.93	0.10	0.61	19,20,21,22	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	EDO	D	1479	4/4	0.95	0.09	0.60	15,16,16,17	0
5	EDO	K	1142	4/4	0.92	0.10	0.45	33,33,34,34	0
5	EDO	G	1478	4/4	0.95	0.09	0.43	17,18,18,19	0
5	EDO	C	1477	4/4	0.93	0.09	0.37	19,20,21,22	0
5	EDO	D	1476	4/4	0.95	0.09	0.28	21,21,23,23	0
5	EDO	N	1141	4/4	0.94	0.10	0.25	27,27,27,27	0
5	EDO	K	1141	4/4	0.79	0.17	0.18	31,33,34,37	0
5	EDO	A	1477	4/4	0.97	0.08	0.02	18,18,19,20	0
5	EDO	J	1142	4/4	0.90	0.13	0.01	33,33,34,36	0
5	EDO	A	1476	4/4	0.96	0.09	-0.07	17,18,18,19	0
5	EDO	O	1142	4/4	0.93	0.11	-0.09	29,29,30,31	0
5	EDO	F	1476	4/4	0.95	0.09	-0.22	26,26,27,28	0
4	CAP	E	477	21/21	0.98	0.07	-0.43	12,15,16,17	0
5	EDO	G	1477	4/4	0.98	0.08	-0.58	18,20,20,21	0
5	EDO	B	1477	4/4	0.96	0.07	-0.67	24,25,25,26	0
4	CAP	F	477	21/21	0.98	0.07	-0.88	13,15,16,18	0
4	CAP	D	477	21/21	0.98	0.06	-1.09	10,15,15,17	0
4	CAP	H	477	21/21	0.98	0.06	-1.09	11,15,16,17	0
4	CAP	A	477	21/21	0.99	0.06	-1.10	12,15,16,18	0
5	EDO	H	1477	4/4	0.97	0.07	-1.14	15,17,18,20	0
4	CAP	G	477	21/21	0.98	0.06	-1.20	12,15,16,18	0
4	CAP	B	477	21/21	0.98	0.07	-1.22	12,15,16,18	0
3	MG	A	476	1/1	0.98	0.06	-1.29	11,11,11,11	0
5	EDO	A	1480	4/4	0.97	0.07	-1.32	18,19,20,20	0
3	MG	G	476	1/1	0.97	0.06	-1.37	11,11,11,11	0
4	CAP	C	477	21/21	0.98	0.06	-1.47	11,15,16,18	0
3	MG	C	476	1/1	0.98	0.05	-1.74	10,10,10,10	0
3	MG	E	476	1/1	0.98	0.03	-2.50	11,11,11,11	0
3	MG	D	476	1/1	0.99	0.03	-2.93	12,12,12,12	0
3	MG	F	476	1/1	0.97	0.04	-4.36	11,11,11,11	0
3	MG	B	476	1/1	0.99	0.03	-4.90	10,10,10,10	0
3	MG	H	476	1/1	0.99	0.02	-6.74	11,11,11,11	0
5	EDO	A	1481	4/4	0.97	0.09	-	24,25,26,26	0
5	EDO	E	1480	4/4	0.77	0.19	-	37,38,39,41	0
5	EDO	G	1479	4/4	0.95	0.15	-	21,23,23,24	0
5	EDO	B	1479	4/4	0.91	0.15	-	28,30,31,31	0
5	EDO	F	1478	4/4	0.94	0.16	-	29,30,30,31	0
5	EDO	H	1479	4/4	0.95	0.15	-	19,21,21,23	0
5	EDO	C	1481	4/4	0.96	0.15	-	22,23,24,24	0
5	EDO	D	1478	4/4	0.90	0.15	-	31,31,31,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.