



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:29 PM UTC

PDB ID : 4P18 / pdb_00004p18
Title : Crystal Structure of frog M ferritin mutant D80K
Authors : Pozzi, C.; Di Pisa, F.; Mangani, S.; Bernacchioni, C.; Ghini, V.; Turano, P.
Deposited on : 2014-02-25
Resolution : 1.91 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

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

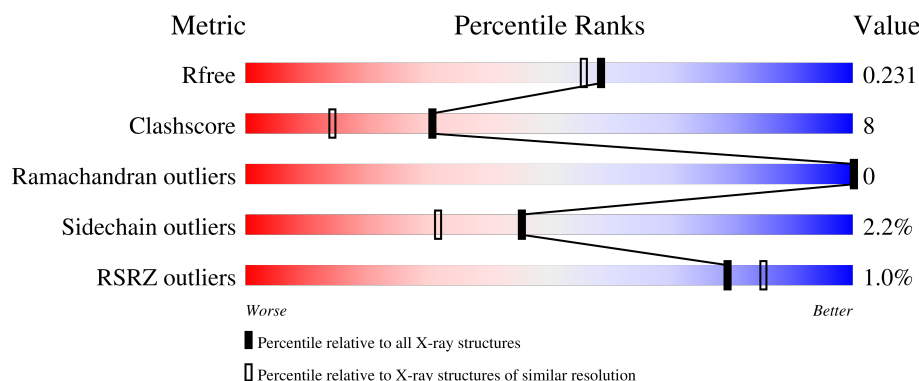
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.91 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



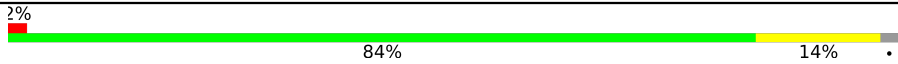
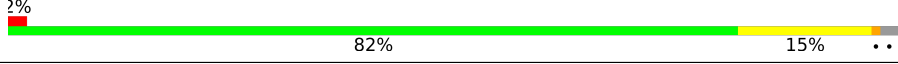
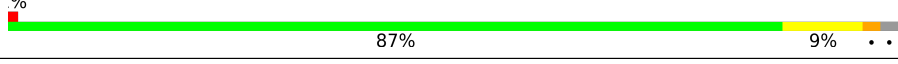
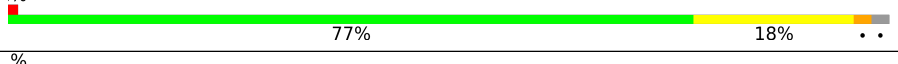
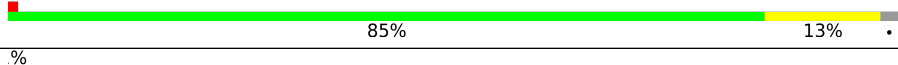
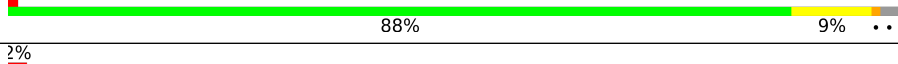
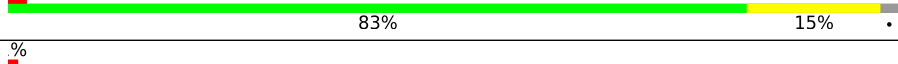
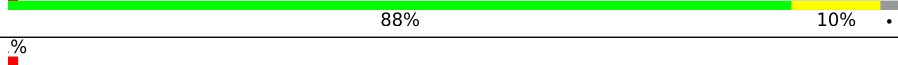
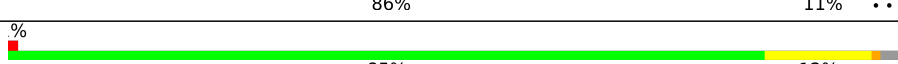
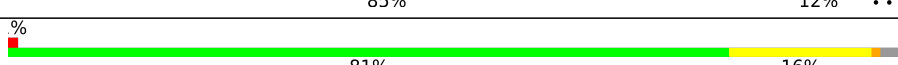
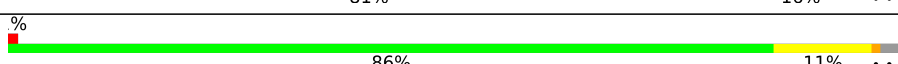
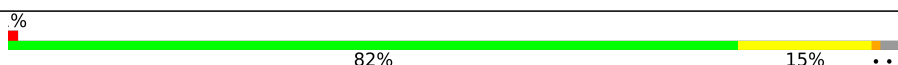
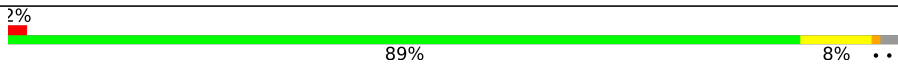
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	176	% 80% 18% ..
1	B	176	82% 15% ..
1	C	176	% 84% 13% ..
1	D	176	2% 88% 10% .
1	E	176	% 88% 10% .

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Mol	Chain	Length	Quality of chain
1	F	176	
1	G	176	
1	H	176	
1	I	176	
1	J	176	
1	K	176	
1	L	176	
1	M	176	
1	N	176	
1	O	176	
1	P	176	
1	Q	176	
1	R	176	
1	S	176	
1	T	176	
1	U	176	
1	V	176	
1	W	176	
1	X	176	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	D	201	-	-	X	-
2	SO4	J	201	-	-	X	-
2	SO4	L	201	-	-	X	-
2	SO4	R	201	-	-	X	-
2	SO4	W	201	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	E	204	-	-	X	-
3	EDO	G	207	-	-	X	-
3	EDO	G	211	-	-	X	-
3	EDO	J	205	-	-	X	-
3	EDO	N	203	-	-	X	-
3	EDO	Q	202	-	-	X	-
3	EDO	R	206	-	-	X	-
3	EDO	S	202	-	-	X	-
3	EDO	W	204	-	-	X	-
4	ACT	A	203	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 40182 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 Ferritin, middle subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	172	Total	C	N	O	S	0	9	0
			1460	922	251	280	7			
1	B	172	Total	C	N	O	S	0	6	0
			1452	915	250	280	7			
1	C	172	Total	C	N	O	S	0	8	0
			1460	922	251	280	7			
1	D	172	Total	C	N	O	S	0	6	0
			1446	912	250	277	7			
1	M	172	Total	C	N	O	S	0	2	0
			1426	899	249	271	7			
1	N	172	Total	C	N	O	S	0	8	0
			1457	919	251	280	7			
1	O	172	Total	C	N	O	S	0	5	0
			1432	903	249	273	7			
1	P	172	Total	C	N	O	S	0	5	0
			1441	908	250	276	7			
1	S	172	Total	C	N	O	S	0	8	0
			1459	922	252	278	7			
1	T	172	Total	C	N	O	S	0	5	0
			1444	912	251	274	7			
1	E	172	Total	C	N	O	S	0	4	0
			1436	904	251	274	7			
1	F	172	Total	C	N	O	S	0	8	0
			1461	923	252	279	7			
1	G	172	Total	C	N	O	S	0	5	0
			1451	913	253	278	7			
1	H	172	Total	C	N	O	S	0	6	0
			1447	913	250	277	7			
1	I	172	Total	C	N	O	S	0	8	0
			1461	923	252	279	7			
1	J	172	Total	C	N	O	S	0	6	0
			1445	912	250	276	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	K	172	Total	C	N	O	S	0	8	0
			1464	923	254	280	7			
1	L	172	Total	C	N	O	S	0	6	0
			1441	909	252	273	7			
1	Q	172	Total	C	N	O	S	0	7	0
			1452	917	252	276	7			
1	R	172	Total	C	N	O	S	0	5	0
			1447	914	252	274	7			
1	U	172	Total	C	N	O	S	0	4	0
			1437	907	250	273	7			
1	V	172	Total	C	N	O	S	0	8	0
			1457	919	252	279	7			
1	W	172	Total	C	N	O	S	0	7	0
			1453	917	253	276	7			
1	X	172	Total	C	N	O	S	0	3	0
			1431	902	249	273	7			

There are 24 discrepancies between the modelled and reference sequences:

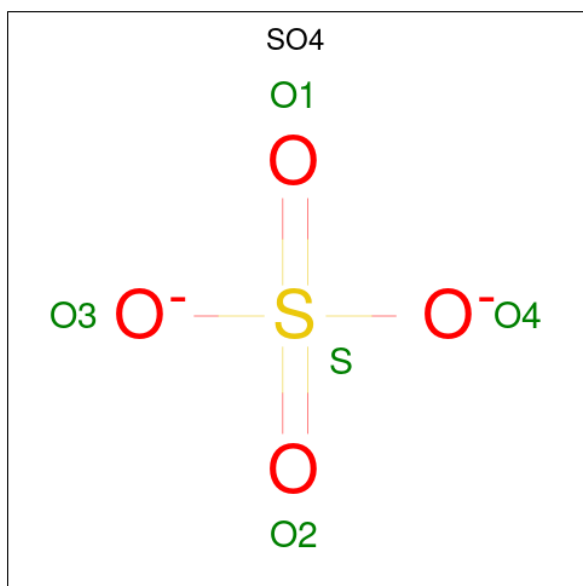
Chain	Residue	Modelled	Actual	Comment	Reference
A	80	LYS	ASP	engineered mutation	UNP P07798
B	80	LYS	ASP	engineered mutation	UNP P07798
C	80	LYS	ASP	engineered mutation	UNP P07798
D	80	LYS	ASP	engineered mutation	UNP P07798
M	80	LYS	ASP	engineered mutation	UNP P07798
N	80	LYS	ASP	engineered mutation	UNP P07798
O	80	LYS	ASP	engineered mutation	UNP P07798
P	80	LYS	ASP	engineered mutation	UNP P07798
S	80	LYS	ASP	engineered mutation	UNP P07798
T	80	LYS	ASP	engineered mutation	UNP P07798
E	80	LYS	ASP	engineered mutation	UNP P07798
F	80	LYS	ASP	engineered mutation	UNP P07798
G	80	LYS	ASP	engineered mutation	UNP P07798
H	80	LYS	ASP	engineered mutation	UNP P07798
I	80	LYS	ASP	engineered mutation	UNP P07798
J	80	LYS	ASP	engineered mutation	UNP P07798
K	80	LYS	ASP	engineered mutation	UNP P07798
L	80	LYS	ASP	engineered mutation	UNP P07798
Q	80	LYS	ASP	engineered mutation	UNP P07798
R	80	LYS	ASP	engineered mutation	UNP P07798
U	80	LYS	ASP	engineered mutation	UNP P07798
V	80	LYS	ASP	engineered mutation	UNP P07798
W	80	LYS	ASP	engineered mutation	UNP P07798

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Chain	Residue	Modelled	Actual	Comment	Reference
X	80	LYS	ASP	engineered mutation	UNP P07798

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	N	1	Total	O	S	0	0
			5	4	1		
2	N	1	Total	O	S	0	0
			5	4	1		
2	O	1	Total	O	S	0	0
			5	4	1		
2	P	1	Total	O	S	0	0
			5	4	1		
2	T	1	Total	O	S	0	0
			5	4	1		
2	T	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	L	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		
2	R	1	Total	O	S	0	0
			5	4	1		
2	U	1	Total	O	S	0	0
			5	4	1		
2	W	1	Total	O	S	0	0
			5	4	1		
2	W	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	M	1	Total	C	O	0	0
			4	2	2		
3	N	1	Total	C	O	0	0
			4	2	2		
3	N	1	Total	C	O	0	0
			4	2	2		
3	N	1	Total	C	O	0	0
			4	2	2		
3	O	1	Total	C	O	0	0
			4	2	2		
3	P	1	Total	C	O	0	0
			4	2	2		

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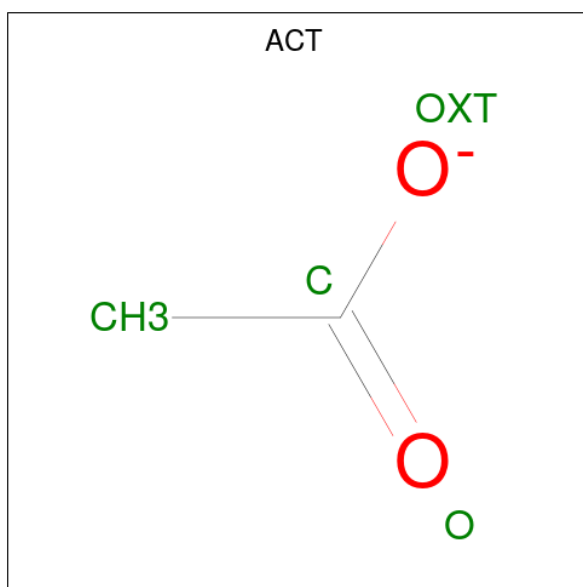
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	P	1	Total 4	C 2	O 2	0	0
3	P	1	Total 4	C 2	O 2	0	0
3	S	1	Total 4	C 2	O 2	0	0
3	S	1	Total 4	C 2	O 2	0	0
3	S	1	Total 4	C 2	O 2	0	0
3	T	1	Total 4	C 2	O 2	0	0
3	E	1	Total 4	C 2	O 2	0	0
3	E	1	Total 4	C 2	O 2	0	0
3	F	1	Total 4	C 2	O 2	0	0
3	F	1	Total 4	C 2	O 2	0	0
3	G	1	Total 4	C 2	O 2	0	0
3	G	1	Total 4	C 2	O 2	0	0
3	G	1	Total 4	C 2	O 2	0	0
3	G	1	Total 4	C 2	O 2	0	0
3	G	1	Total 4	C 2	O 2	0	0
3	G	1	Total 4	C 2	O 2	0	0
3	G	1	Total 4	C 2	O 2	0	0
3	I	1	Total 4	C 2	O 2	0	0
3	I	1	Total 4	C 2	O 2	0	0
3	J	1	Total 4	C 2	O 2	0	0
3	J	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	K	1	Total 4	C 2	O 2	0	0
3	K	1	Total 4	C 2	O 2	0	0
3	L	1	Total 4	C 2	O 2	0	0
3	L	1	Total 4	C 2	O 2	0	0
3	Q	1	Total 4	C 2	O 2	0	0
3	Q	1	Total 4	C 2	O 2	0	0
3	R	1	Total 4	C 2	O 2	0	0
3	R	1	Total 4	C 2	O 2	0	0
3	R	1	Total 4	C 2	O 2	0	0
3	R	1	Total 4	C 2	O 2	0	0
3	U	1	Total 4	C 2	O 2	0	0
3	V	1	Total 4	C 2	O 2	0	0
3	V	1	Total 4	C 2	O 2	0	0
3	W	1	Total 4	C 2	O 2	0	0
3	X	1	Total 4	C 2	O 2	0	0
3	X	1	Total 4	C 2	O 2	0	0

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2$).



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

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Cl	0	0
			2	2		
5	C	1	Total	Cl	0	0
			1	1		
5	D	2	Total	Cl	0	0
			2	2		
5	M	1	Total	Cl	0	0
			1	1		
5	O	2	Total	Cl	0	0
			2	2		
5	S	1	Total	Cl	0	0
			1	1		
5	T	2	Total	Cl	0	0
			2	2		
5	E	3	Total	Cl	0	0
			3	3		
5	F	2	Total	Cl	0	0
			2	2		
5	G	3	Total	Cl	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	1	Total 1	Cl 1	0	0
5	I	1	Total 1	Cl 1	0	0
5	J	1	Total 1	Cl 1	0	0
5	L	2	Total 2	Cl 2	0	0
5	Q	1	Total 1	Cl 1	0	0
5	U	2	Total 2	Cl 2	0	0
5	W	1	Total 1	Cl 1	0	0
5	X	3	Total 3	Cl 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	214	Total 214	O 214	0	0
6	B	211	Total 211	O 211	0	0
6	C	203	Total 203	O 203	0	0
6	D	193	Total 193	O 193	0	0
6	M	208	Total 208	O 208	0	0
6	N	194	Total 194	O 194	0	0
6	O	210	Total 210	O 210	0	0
6	P	191	Total 191	O 191	0	0
6	S	226	Total 226	O 226	0	0
6	T	207	Total 207	O 207	0	0
6	E	205	Total 205	O 205	0	0

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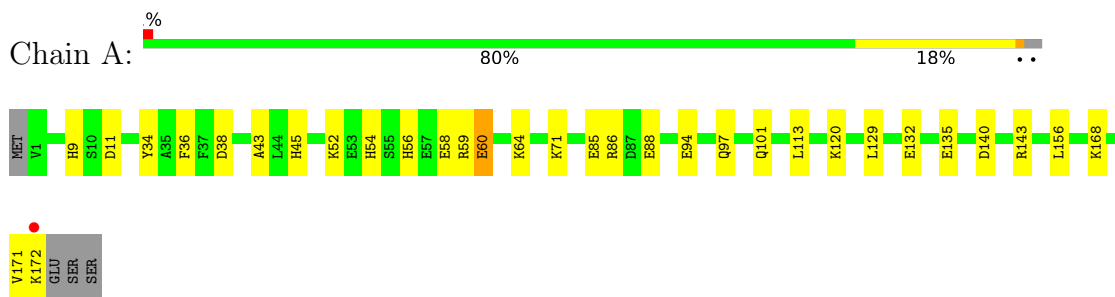
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	F	229	Total 229	O 229	0	0
6	G	210	Total 210	O 210	0	0
6	H	194	Total 194	O 194	0	0
6	I	224	Total 224	O 224	0	0
6	J	231	Total 231	O 231	0	0
6	K	199	Total 199	O 199	0	0
6	L	209	Total 209	O 209	0	0
6	Q	203	Total 203	O 203	0	0
6	R	207	Total 207	O 207	0	0
6	U	204	Total 204	O 204	0	0
6	V	245	Total 245	O 245	0	0
6	W	227	Total 227	O 227	0	0
6	X	210	Total 210	O 210	0	0

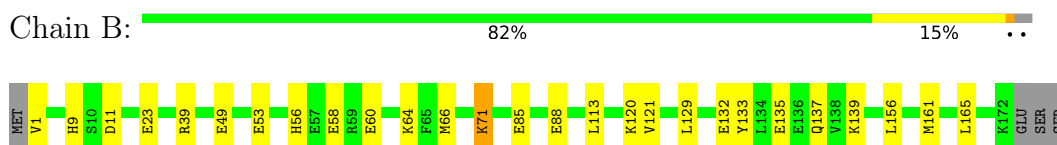
3 Residue-property plots [i](#)

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

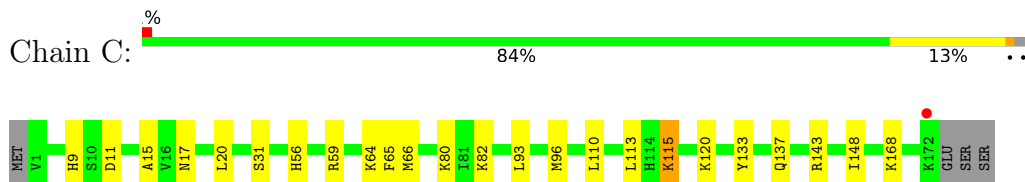
- Molecule 1: Ferritin, middle subunit



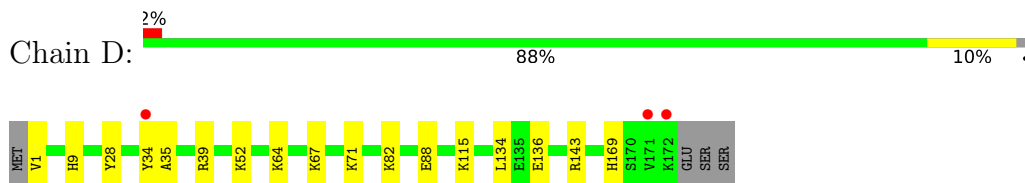
- Molecule 1: Ferritin, middle subunit



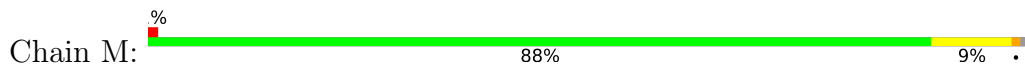
- Molecule 1: Ferritin, middle subunit



- Molecule 1: Ferritin, middle subunit

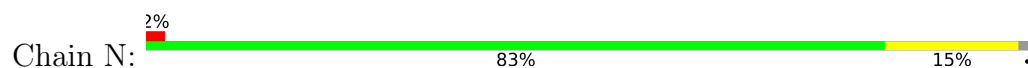


- Molecule 1: Ferritin, middle subunit

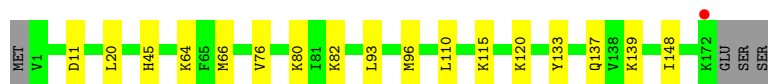
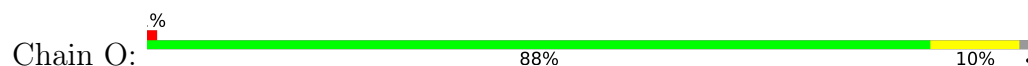




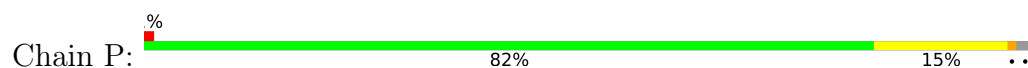
- Molecule 1: Ferritin, middle subunit



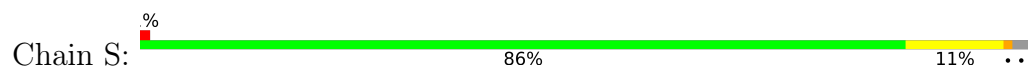
- Molecule 1: Ferritin, middle subunit



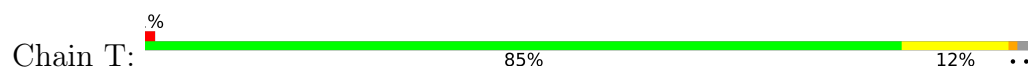
- Molecule 1: Ferritin, middle subunit



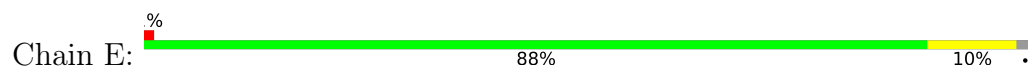
- Molecule 1: Ferritin, middle subunit



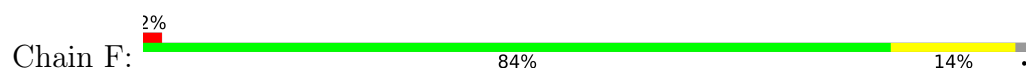
- Molecule 1: Ferritin, middle subunit



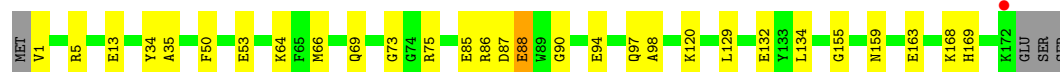
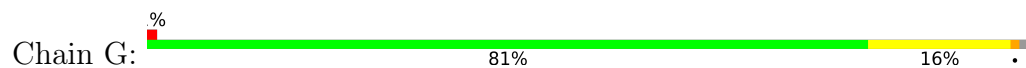
- Molecule 1: Ferritin, middle subunit



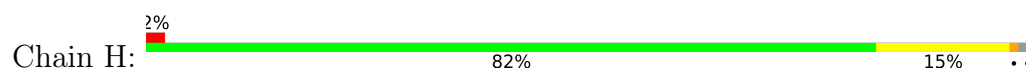
- Molecule 1: Ferritin, middle subunit



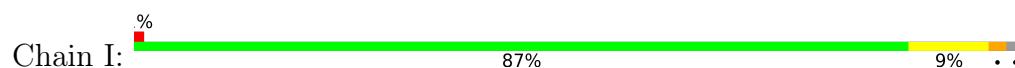
- Molecule 1: Ferritin, middle subunit



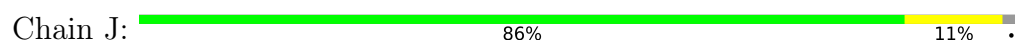
- Molecule 1: Ferritin, middle subunit



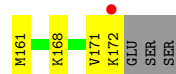
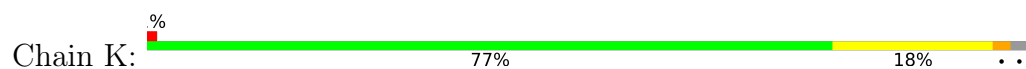
- Molecule 1: Ferritin, middle subunit



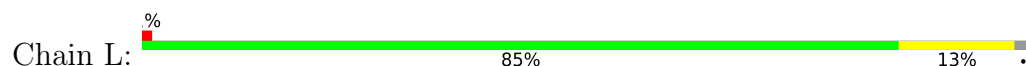
- Molecule 1: Ferritin, middle subunit

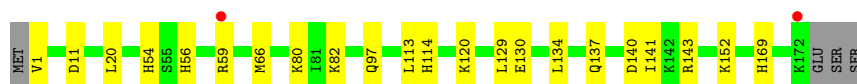


- Molecule 1: Ferritin, middle subunit

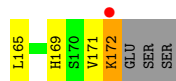
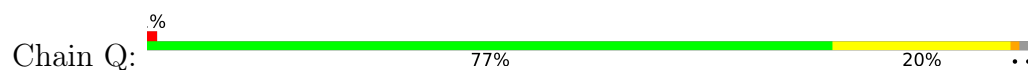


- Molecule 1: Ferritin, middle subunit

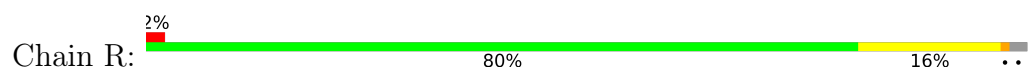




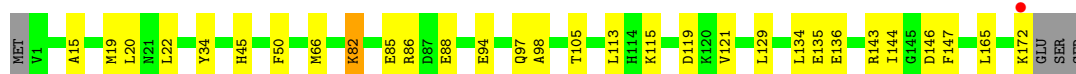
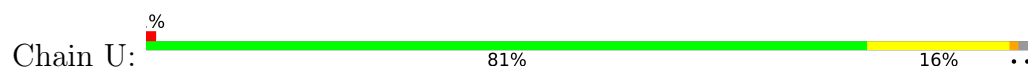
- Molecule 1: Ferritin, middle subunit



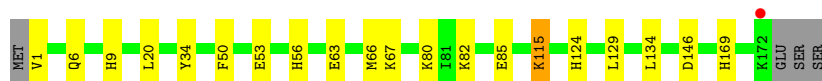
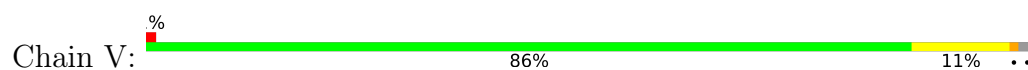
- Molecule 1: Ferritin, middle subunit



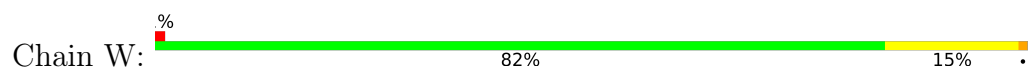
- Molecule 1: Ferritin, middle subunit



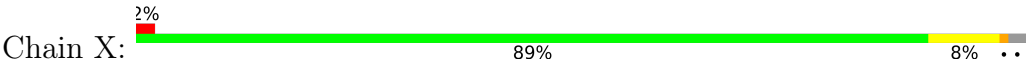
- Molecule 1: Ferritin, middle subunit



- Molecule 1: Ferritin, middle subunit



- Molecule 1: Ferritin, middle subunit



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	238.93Å 238.43Å 119.69Å 90.00° 94.33° 90.00°	Depositor
Resolution (Å)	75.35 – 1.91 75.35 – 1.91	Depositor EDS
% Data completeness (in resolution range)	96.7 (75.35-1.91) 96.8 (75.35-1.91)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.196 , 0.231 0.196 , 0.231	Depositor DCC
R_{free} test set	25123 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	9.7	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 59.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	40182	wwPDB-VP
Average B, all atoms (Å ²)	8.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.26% 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: CL, EDO, ACT, SO4

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	0.65	0/1513	0.80	0/2033
1	B	0.65	0/1496	0.83	0/2011
1	C	0.69	0/1510	0.81	0/2030
1	D	0.66	0/1490	0.85	0/2003
1	E	0.65	0/1471	0.82	0/1978
1	F	0.66	0/1511	0.82	0/2029
1	G	0.65	0/1488	0.84	0/1998
1	H	0.64	0/1494	0.79	0/2007
1	I	0.65	0/1514	0.83	0/2033
1	J	0.68	0/1492	0.83	0/2005
1	K	0.65	0/1511	0.79	0/2029
1	L	0.68	0/1488	0.84	0/2000
1	M	0.67	0/1461	0.82	0/1964
1	N	0.64	0/1507	0.77	0/2025
1	O	0.67	0/1476	0.83	0/1985
1	P	0.65	0/1484	0.83	0/1994
1	Q	0.65	0/1502	0.80	0/2017
1	R	0.71	0/1485	0.84	0/1996
1	S	0.64	0/1510	0.78	0/2029
1	T	0.65	0/1485	0.81	0/1996
1	U	0.67	0/1477	0.80	0/1983
1	V	0.67	0/1504	0.83	0/2021
1	W	0.65	0/1501	0.81	0/2017
1	X	0.64	0/1469	0.81	0/1975
All	All	0.66	0/35839	0.82	0/48158

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	1460	0	1425	30	0
1	B	1452	0	1407	19	0
1	C	1460	0	1424	37	0
1	D	1446	0	1403	18	0
1	E	1436	0	1389	19	0
1	F	1461	0	1432	30	0
1	G	1451	0	1415	28	0
1	H	1447	0	1416	34	0
1	I	1461	0	1436	19	0
1	J	1445	0	1411	21	0
1	K	1464	0	1433	44	0
1	L	1441	0	1408	23	0
1	M	1426	0	1385	15	0
1	N	1457	0	1421	16	0
1	O	1432	0	1393	11	0
1	P	1441	0	1406	32	0
1	Q	1452	0	1429	50	0
1	R	1447	0	1408	41	0
1	S	1459	0	1421	19	0
1	T	1444	0	1410	21	0
1	U	1437	0	1413	29	0
1	V	1457	0	1418	22	0
1	W	1453	0	1418	27	0
1	X	1431	0	1389	16	0
2	A	5	0	0	0	0
2	C	5	0	0	0	0
2	D	10	0	0	2	0
2	F	5	0	0	0	0
2	G	5	0	0	0	0
2	H	10	0	0	0	0
2	I	5	0	0	0	0
2	J	10	0	0	2	0
2	K	10	0	0	1	0
2	L	5	0	0	2	0
2	N	10	0	0	2	0
2	O	5	0	0	0	0
2	P	5	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	R	10	0	0	3	0
2	T	10	0	0	1	0
2	U	5	0	0	0	0
2	W	10	0	0	3	0
3	A	4	0	6	2	0
3	B	12	0	18	4	0
3	C	12	0	18	0	0
3	D	4	0	6	1	0
3	E	8	0	12	4	0
3	F	8	0	12	3	0
3	G	28	0	42	12	0
3	I	8	0	12	4	0
3	J	8	0	12	4	0
3	K	8	0	12	2	0
3	L	8	0	12	1	0
3	M	4	0	6	0	0
3	N	12	0	18	7	0
3	O	4	0	6	0	0
3	P	12	0	18	5	0
3	Q	8	0	12	5	0
3	R	16	0	24	9	0
3	S	12	0	18	5	0
3	T	4	0	6	2	0
3	U	4	0	6	0	0
3	V	8	0	12	3	0
3	W	4	0	6	7	0
3	X	8	0	12	1	0
4	A	4	0	3	5	0
4	Q	4	0	3	1	0
5	B	2	0	0	0	0
5	C	1	0	0	0	0
5	D	2	0	0	0	0
5	E	3	0	0	0	0
5	F	2	0	0	0	0
5	G	3	0	0	1	0
5	H	1	0	0	1	0
5	I	1	0	0	0	0
5	J	1	0	0	0	0
5	L	2	0	0	0	0
5	M	1	0	0	0	0
5	O	2	0	0	0	0
5	Q	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	S	1	0	0	1	0
5	T	2	0	0	0	0
5	U	2	0	0	1	0
5	W	1	0	0	1	0
5	X	3	0	0	0	0
6	A	214	0	0	12	0
6	B	211	0	0	5	0
6	C	203	0	0	8	0
6	D	193	0	0	10	0
6	E	205	0	0	7	0
6	F	229	0	0	14	0
6	G	210	0	0	7	0
6	H	194	0	0	11	0
6	I	224	0	0	3	0
6	J	231	0	0	5	0
6	K	199	0	0	15	0
6	L	209	0	0	11	0
6	M	208	0	0	5	0
6	N	194	0	0	7	0
6	O	210	0	0	4	0
6	P	191	0	0	19	0
6	Q	203	0	0	23	0
6	R	207	0	0	3	0
6	S	226	0	0	6	0
6	T	207	0	0	3	0
6	U	204	0	0	11	0
6	V	245	0	0	5	0
6	W	227	0	0	3	0
6	X	210	0	0	6	0
All	All	40182	0	34222	578	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 8.

All (578) 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:C:20[B]:LEU:CD2	1:C:65:PHE:HB3	1.78	1.11
1:F:171:VAL:HG11	6:F:524:HOH:O	1.50	1.09
1:C:148:ILE:HG12	6:C:451:HOH:O	1.52	1.07
1:J:115:LYS:HD2	6:J:525:HOH:O	1.52	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:11[B]:ASP:OD2	1:L:120:LYS:HE3	1.57	1.04
1:F:108:GLN:HG2	6:F:526:HOH:O	1.59	1.02
1:L:137:GLN:HB3	6:L:503:HOH:O	1.57	1.02
1:H:66:MET:HE3	1:J:35:ALA:HB2	1.44	0.99
1:R:115:LYS:HD3	6:R:489:HOH:O	1.60	0.99
1:P:144:ILE:HB	6:P:489:HOH:O	1.63	0.99
1:C:20[B]:LEU:HD23	1:C:65:PHE:HB3	1.44	0.97
1:G:73:GLY:O	3:G:211:EDO:H12	1.65	0.97
1:S:73:GLY:O	3:S:202:EDO:H22	1.65	0.97
1:V:80:LYS:HE3	1:V:82[A]:LYS:HE2	1.47	0.94
1:V:80:LYS:HE3	1:V:82[A]:LYS:CE	1.99	0.92
1:C:66:MET:HE3	1:D:35:ALA:HB2	1.50	0.91
1:E:11[B]:ASP:OD2	1:E:120:LYS:HE3	1.70	0.91
1:N:171:VAL:HG11	6:N:488:HOH:O	1.71	0.91
1:P:50:PHE:HE2	6:P:488:HOH:O	1.51	0.91
1:W:11[B]:ASP:OD2	1:W:120:LYS:HE3	1.70	0.90
1:C:20[B]:LEU:HD22	1:C:65:PHE:HB3	1.54	0.90
1:P:161:MET:HE3	1:F:161:MET:SD	2.12	0.90
1:L:130:GLU:HG3	6:L:447:HOH:O	1.69	0.89
1:T:9[B]:HIS:HD2	6:T:505:HOH:O	1.54	0.89
1:F:35:ALA:HB2	1:G:66:MET:HE3	1.53	0.88
1:Q:147:PHE:HB3	6:Q:449:HOH:O	1.72	0.88
1:P:12:CYS:HB3	6:P:442:HOH:O	1.73	0.88
1:K:96:MET:HE1	6:K:497:HOH:O	1.75	0.87
1:Q:35:ALA:HB3	6:Q:446:HOH:O	1.75	0.86
1:H:159:ASN:HD21	1:K:161:MET:CE	1.89	0.85
6:F:417:HOH:O	1:G:87:ASP:HB3	1.75	0.85
1:H:164:TYR:HE1	6:Q:484:HOH:O	1.57	0.85
1:Q:148:ILE:HG12	6:Q:464:HOH:O	1.76	0.83
1:V:9[B]:HIS:CD2	6:V:474:HOH:O	2.32	0.83
1:X:82:LYS:HD3	6:X:378:HOH:O	1.79	0.82
1:Q:66:MET:HE3	1:R:35:ALA:HB2	1.59	0.82
1:U:82:LYS:CE	1:U:82:LYS:HA	2.10	0.82
1:W:171:VAL:O	1:W:172:LYS:HB2	1.77	0.82
1:R:5:ARG:HH11	3:R:206:EDO:H21	1.44	0.82
1:U:82:LYS:HA	1:U:82:LYS:HE2	1.63	0.81
1:K:20:LEU:HD11	1:K:66:MET:HG3	1.63	0.81
1:S:11[B]:ASP:OD2	1:S:120:LYS:HE3	1.80	0.80
3:N:203:EDO:H11	1:W:135:GLU:HB3	1.62	0.80
1:O:148:ILE:HG12	6:O:483:HOH:O	1.82	0.80
3:V:201:EDO:H12	1:W:45:HIS:HB2	1.64	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:9[B]:HIS:CD2	6:M:464:HOH:O	2.35	0.79
1:G:13:GLU:HG2	6:G:500:HOH:O	1.83	0.78
6:S:333:HOH:O	3:J:205:EDO:H11	1.83	0.78
1:Q:96:MET:HG3	6:Q:449:HOH:O	1.82	0.78
1:Q:93:LEU:HD12	6:Q:464:HOH:O	1.83	0.77
1:W:2:SER:H	3:W:204:EDO:H11	1.48	0.77
1:U:20:LEU:HD11	1:U:66:MET:HG2	1.65	0.76
1:P:100:LEU:HD13	6:P:489:HOH:O	1.86	0.76
1:P:62:ALA:HA	6:P:370:HOH:O	1.83	0.76
3:E:204:EDO:H11	6:E:428:HOH:O	1.85	0.76
1:R:70:ASN:HA	6:R:462:HOH:O	1.86	0.76
1:K:11[B]:ASP:OD2	1:K:120:LYS:HE3	1.87	0.75
1:E:135:GLU:HG3	1:E:139:LYS:CE	2.17	0.74
1:N:153:ARG:HD2	6:N:428:HOH:O	1.88	0.74
1:P:97:GLN:HG2	6:P:465:HOH:O	1.87	0.74
1:F:11[B]:ASP:OD2	1:F:120:LYS:HE2	1.88	0.74
1:H:159:ASN:HD21	1:K:161:MET:HE3	1.50	0.74
1:Q:171:VAL:O	1:Q:172:LYS:HB2	1.86	0.74
1:P:65:PHE:HB2	6:P:370:HOH:O	1.87	0.73
1:A:9[B]:HIS:HD2	6:A:472:HOH:O	1.71	0.73
1:F:9[B]:HIS:CD2	6:F:523:HOH:O	2.41	0.73
1:J:164:TYR:O	1:J:168[B]:LYS:HE2	1.88	0.73
1:K:171:VAL:O	1:K:172:LYS:HB2	1.85	0.73
1:A:11[B]:ASP:OD2	1:A:120:LYS:HE3	1.87	0.73
6:S:333:HOH:O	3:J:205:EDO:C1	2.35	0.73
1:R:16:VAL:HG12	6:R:417:HOH:O	1.89	0.73
1:K:9[B]:HIS:CD2	6:K:386:HOH:O	2.41	0.72
3:N:203:EDO:O2	1:W:139:LYS:HD2	1.88	0.72
1:P:20:LEU:HA	6:P:370:HOH:O	1.89	0.71
1:H:171:VAL:O	1:H:172:LYS:HB2	1.90	0.71
1:C:11[B]:ASP:OD2	1:C:120:LYS:HE3	1.90	0.71
1:B:9[B]:HIS:CD2	6:B:434:HOH:O	2.44	0.71
6:D:465:HOH:O	1:N:165:LEU:HD23	1.90	0.71
1:R:115:LYS:HE2	1:R:115:LYS:HA	1.71	0.71
1:F:9[B]:HIS:HD2	6:F:523:HOH:O	1.72	0.71
1:F:82:LYS:HG3	6:F:474:HOH:O	1.91	0.70
1:T:9[B]:HIS:CD2	6:T:505:HOH:O	2.34	0.70
1:E:135:GLU:HG3	1:E:139:LYS:HE3	1.73	0.70
1:R:5:ARG:NH1	3:R:206:EDO:H21	2.06	0.70
1:V:9[B]:HIS:HD2	6:V:474:HOH:O	1.73	0.70
1:P:96:MET:SD	6:P:488:HOH:O	2.48	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:80:LYS:CE	1:V:82[A]:LYS:HE2	2.20	0.70
1:U:136:GLU:HG2	6:U:343:HOH:O	1.93	0.68
1:A:9[B]:HIS:CD2	6:A:472:HOH:O	2.44	0.68
1:I:11[B]:ASP:OD2	1:I:120:LYS:HE3	1.92	0.68
1:A:85[A]:GLU:HG3	1:A:86:ARG:HG3	1.76	0.68
1:R:56:HIS:HD2	1:R:59:ARG:HH22	1.41	0.68
1:F:75:ARG:NE	3:F:205:EDO:H21	2.08	0.67
1:M:67:LYS:HG3	6:N:484:HOH:O	1.93	0.67
1:Q:66:MET:HE1	1:R:34:TYR:HD2	1.59	0.67
1:S:5:ARG:NH1	3:S:202:EDO:H21	2.10	0.67
1:W:143:ARG:HD3	6:W:397:HOH:O	1.93	0.67
1:T:121:VAL:HG11	1:J:115:LYS:HE3	1.76	0.67
1:T:5:ARG:HD2	3:T:205:EDO:O2	1.94	0.67
1:L:114:HIS:CD2	6:L:447:HOH:O	2.47	0.67
1:J:115:LYS:HE2	1:J:115:LYS:HA	1.77	0.67
1:V:6:GLN:HA	3:V:202:EDO:H11	1.76	0.67
1:L:114:HIS:HD2	6:L:447:HOH:O	1.78	0.67
1:A:38:ASP:OD1	3:A:202:EDO:H22	1.95	0.66
1:S:171:VAL:O	1:S:172:LYS:HB2	1.94	0.66
1:G:73:GLY:O	3:G:211:EDO:C1	2.42	0.66
1:U:144:ILE:HG12	6:U:500:HOH:O	1.95	0.66
1:N:73:GLY:O	3:N:205:EDO:H12	1.96	0.66
1:I:101[B]:GLN:HG2	6:I:408:HOH:O	1.95	0.66
1:N:64[A]:LYS:HE3	6:N:436:HOH:O	1.96	0.65
1:P:161:MET:CE	1:F:161:MET:SD	2.84	0.65
1:U:82:LYS:HE2	6:U:334:HOH:O	1.96	0.65
1:E:169:HIS:NE2	2:J:201:SO4:O1	2.29	0.65
1:C:93:LEU:HD12	6:C:451:HOH:O	1.97	0.65
3:D:205:EDO:H21	1:N:45:HIS:HB2	1.78	0.65
1:K:45:HIS:CD2	6:K:498:HOH:O	2.50	0.65
1:W:2:SER:HB3	3:W:204:EDO:H11	1.79	0.65
1:Q:165:LEU:HB3	6:Q:484:HOH:O	1.97	0.65
1:C:20[B]:LEU:HD22	1:C:65:PHE:CB	2.26	0.64
1:R:5:ARG:HH11	3:R:206:EDO:C2	2.09	0.64
1:S:5:ARG:HH11	3:S:202:EDO:H21	1.61	0.64
1:K:158:GLU:CD	1:K:158:GLU:H	2.06	0.64
1:I:5:ARG:HH11	3:I:203:EDO:H12	1.63	0.64
1:E:9[B]:HIS:CD2	6:E:502:HOH:O	2.51	0.64
1:G:155:GLY:O	3:G:210:EDO:H12	1.97	0.64
1:K:143:ARG:HD3	6:K:363:HOH:O	1.97	0.64
1:C:66:MET:HE1	1:D:34:TYR:HD2	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:444:HOH:O	1:R:139:LYS:HE3	1.98	0.63
1:C:20[B]:LEU:CD2	1:C:65:PHE:CB	2.68	0.63
1:A:43:ALA:HA	6:A:455:HOH:O	1.98	0.63
1:V:80:LYS:HE3	1:V:82[A]:LYS:HE3	1.80	0.63
1:F:67:LYS:HE2	1:F:71:LYS:HE3	1.81	0.63
1:Q:43:ALA:HB1	3:Q:202:EDO:C2	2.28	0.62
1:W:2:SER:H	3:W:204:EDO:C1	2.12	0.62
1:M:56:HIS:CD2	1:M:59:ARG:HH12	2.17	0.62
1:Q:31[A]:SER:HA	6:Q:503:HOH:O	1.99	0.62
1:C:17:ASN:HA	1:C:20[B]:LEU:HG	1.81	0.62
1:P:50:PHE:CE2	6:P:488:HOH:O	2.36	0.62
1:X:143:ARG:NH1	6:X:469:HOH:O	2.32	0.62
1:E:9[B]:HIS:HD2	6:E:502:HOH:O	1.80	0.62
1:I:80:LYS:NZ	1:I:80:LYS:HB3	2.14	0.62
1:B:11[B]:ASP:OD2	1:B:120:LYS:HE3	1.99	0.61
1:Q:31[B]:SER:HA	6:Q:503:HOH:O	2.00	0.61
1:O:11[B]:ASP:OD2	1:O:120:LYS:HE3	2.00	0.61
1:G:50:PHE:O	1:G:53[B]:GLU:HG2	2.01	0.61
1:H:159:ASN:ND2	1:K:161:MET:CE	2.62	0.61
1:Q:9[B]:HIS:CD2	6:Q:445:HOH:O	2.53	0.61
1:Q:172:LYS:HA	6:Q:483:HOH:O	2.01	0.61
1:S:5:ARG:HH11	3:S:202:EDO:C2	2.14	0.61
1:K:136:GLU:HA	1:K:139:LYS:HE2	1.83	0.61
1:Q:66:MET:CE	1:R:34:TYR:CD2	2.85	0.60
1:A:45:HIS:ND1	3:A:202:EDO:H21	2.16	0.60
1:R:7:ASN:H	3:R:204:EDO:H11	1.67	0.60
1:O:93:LEU:HD12	6:O:483:HOH:O	2.01	0.60
1:K:76:VAL:O	3:L:205:EDO:H22	2.01	0.60
1:O:96:MET:HB2	6:O:483:HOH:O	2.00	0.60
1:K:85:GLU:HG3	1:K:86:ARG:HG3	1.84	0.60
1:A:56:HIS:CD2	1:A:59:ARG:HH22	2.20	0.60
1:H:143:ARG:HD3	6:H:386:HOH:O	2.02	0.60
1:I:36:PHE:CE1	1:I:88:GLU:HG2	2.37	0.59
3:N:203:EDO:H11	1:W:135:GLU:CB	2.32	0.59
1:G:86:ARG:HD2	3:G:207:EDO:H12	1.83	0.59
1:L:54:HIS:CE1	1:L:143[A]:ARG:NH1	2.70	0.59
1:Q:169:HIS:HD2	6:Q:484:HOH:O	1.85	0.59
1:J:82:LYS:HE2	6:J:361:HOH:O	2.02	0.59
1:Q:11[B]:ASP:OD2	1:Q:120:LYS:HE3	2.01	0.59
1:Q:43:ALA:HB1	3:Q:202:EDO:H21	1.85	0.59
1:E:64:LYS:HG3	3:E:204:EDO:H21	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:HIS:NE2	1:B:60:GLU:OE1	2.35	0.59
1:L:152:LYS:HE3	6:L:499:HOH:O	2.02	0.59
1:R:56:HIS:CD2	1:R:59:ARG:HH22	2.21	0.59
1:R:73:GLY:O	3:R:206:EDO:C2	2.51	0.59
1:B:49:GLU:O	1:B:53[A]:GLU:HG3	2.03	0.58
1:D:115:LYS:HG3	6:D:402:HOH:O	2.01	0.58
1:I:80:LYS:NZ	1:I:80:LYS:CB	2.66	0.58
1:B:9[B]:HIS:HD2	6:B:434:HOH:O	1.82	0.58
1:Q:9[B]:HIS:HD2	6:Q:445:HOH:O	1.87	0.58
1:Q:34:TYR:CD2	6:Q:503:HOH:O	2.52	0.58
1:T:121:VAL:CG1	1:J:115:LYS:HE3	2.34	0.57
1:H:20:LEU:HD11	1:H:66:MET:CG	2.34	0.57
1:Q:85:GLU:HG3	1:Q:86:ARG:HG3	1.85	0.57
1:H:159:ASN:HD21	1:K:161:MET:HE1	1.68	0.57
1:H:20:LEU:HD11	1:H:66:MET:HG2	1.87	0.57
1:U:105:THR:HG21	6:U:466:HOH:O	2.04	0.57
1:O:80:LYS:HE3	1:O:82:LYS:HD3	1.85	0.57
1:G:86:ARG:CD	3:G:207:EDO:H12	2.35	0.57
1:I:80:LYS:HB3	1:I:80:LYS:HZ3	1.69	0.57
1:U:50:PHE:CZ	6:U:500:HOH:O	2.53	0.57
1:H:169:HIS:NE2	2:K:201:SO4:O3	2.38	0.56
1:I:136[A]:GLU:OE1	1:I:139:LYS:NZ	2.37	0.56
2:L:201:SO4:O1	1:X:169:HIS:NE2	2.35	0.56
1:V:85:GLU:HG2	6:V:461:HOH:O	2.05	0.56
5:S:201:CL:CL	1:G:163:GLU:OE1	2.60	0.56
1:B:161:MET:HB2	3:B:204:EDO:H22	1.85	0.56
1:S:73:GLY:O	3:S:202:EDO:C2	2.47	0.56
1:F:80:LYS:HE3	1:F:82:LYS:HD3	1.87	0.56
4:A:203:ACT:O	1:M:139:LYS:NZ	2.37	0.56
1:B:135[B]:GLU:OE2	1:B:139:LYS:NZ	2.35	0.56
1:Q:66:MET:CE	1:R:34:TYR:HD2	2.19	0.56
3:Q:202:EDO:H22	1:U:146:ASP:HB3	1.87	0.56
1:W:20:LEU:HD11	1:W:66:MET:HG2	1.88	0.56
1:K:143:ARG:NH1	6:K:424:HOH:O	2.38	0.56
1:B:85:GLU:HG2	6:B:465:HOH:O	2.06	0.55
1:Q:11[B]:ASP:OD2	1:Q:120:LYS:CE	2.55	0.55
1:M:67:LYS:NZ	2:N:202:SO4:S	2.73	0.55
1:F:64[A]:LYS:HE3	6:F:449:HOH:O	2.06	0.55
1:G:13:GLU:CG	6:G:500:HOH:O	2.47	0.55
1:P:46:ASN:HD21	3:P:203:EDO:H21	1.70	0.55
1:J:71:LYS:HE2	6:J:473:HOH:O	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:168:LYS:HG2	6:M:503:HOH:O	2.07	0.55
1:H:113:LEU:HG	1:H:129:LEU:HD11	1.89	0.55
1:A:71:LYS:HZ1	4:A:203:ACT:C	2.20	0.54
1:T:88:GLU:HG3	2:T:201:SO4:O1	2.07	0.54
1:J:18:ARG:HD3	6:J:445:HOH:O	2.07	0.54
1:L:80:LYS:HE3	1:L:82:LYS:NZ	2.22	0.54
1:D:1:VAL:HB	6:D:362:HOH:O	2.08	0.54
1:E:53:GLU:HG3	6:E:487:HOH:O	2.07	0.54
1:D:143:ARG:NH2	6:D:492:HOH:O	2.35	0.54
1:D:9[B]:HIS:ND1	6:D:424:HOH:O	2.32	0.54
3:F:204:EDO:C1	6:F:497:HOH:O	2.54	0.54
1:K:80:LYS:NZ	1:K:82:LYS:HE2	2.23	0.54
1:N:66:MET:HE2	1:N:78:LEU:HD21	1.88	0.54
1:P:52:LYS:HE3	2:P:201:SO4:O4	2.08	0.54
1:U:136:GLU:HG3	6:U:407:HOH:O	2.07	0.54
1:Q:63:GLU:OE1	1:R:52:LYS:NZ	2.40	0.54
1:Q:66:MET:HE2	1:R:34:TYR:CE2	2.43	0.54
1:V:115:LYS:HE3	1:V:115:LYS:O	2.08	0.54
1:W:143:ARG:HD2	6:W:468:HOH:O	2.07	0.54
1:D:169:HIS:NE2	2:D:201:SO4:O3	2.40	0.53
1:H:66:MET:HE1	1:J:34:TYR:HD2	1.72	0.53
1:V:50:PHE:O	1:V:53[B]:GLU:HG2	2.08	0.53
1:A:168:LYS:HE2	6:A:511:HOH:O	2.09	0.53
1:Q:169:HIS:CD2	6:Q:484:HOH:O	2.59	0.53
1:A:36:PHE:CE1	1:A:88:GLU:HG2	2.42	0.53
1:H:143:ARG:NH1	6:H:413:HOH:O	2.40	0.53
1:W:9[B]:HIS:CD2	1:W:120:LYS:HB3	2.44	0.53
1:T:64:LYS:NZ	1:T:132:GLU:OE1	2.41	0.53
1:G:69:GLN:NE2	6:G:500:HOH:O	2.41	0.53
1:K:9[B]:HIS:HD2	6:K:386:HOH:O	1.83	0.53
1:K:20:LEU:CD1	1:K:66:MET:HG3	2.37	0.53
1:T:9[A]:HIS:CD2	1:T:120:LYS:HE2	2.44	0.52
1:W:64:LYS:HD3	6:W:467:HOH:O	2.09	0.52
1:Q:94:GLU:HA	1:Q:97:GLN:HE21	1.74	0.52
1:F:108:GLN:CG	6:F:526:HOH:O	2.36	0.52
1:C:66:MET:CE	1:D:34:TYR:HD2	2.23	0.52
1:K:97:GLN:HG2	6:K:405:HOH:O	2.08	0.52
1:K:101[B]:GLN:CD	1:K:104[B]:LYS:HD2	2.35	0.52
1:G:88:GLU:OE1	1:G:90:GLY:N	2.42	0.52
1:R:121:VAL:HG11	1:U:115[B]:LYS:HD3	1.91	0.52
1:X:94:GLU:HA	1:X:97:GLN:HE21	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:169:HIS:NE2	2:R:201:SO4:O2	2.43	0.52
1:J:85:GLU:HG2	1:J:86:ARG:HG3	1.92	0.52
1:U:82:LYS:CE	1:U:82:LYS:CA	2.85	0.52
1:B:23:GLU:OE1	1:B:58:GLU:OE1	2.28	0.52
1:F:23:GLU:OE1	1:F:58:GLU:OE1	2.28	0.52
1:K:23:GLU:OE1	1:K:58:GLU:OE1	2.28	0.52
1:R:63:GLU:HA	1:R:66:MET:HE2	1.91	0.52
1:W:66:MET:HE1	1:X:34:TYR:HD2	1.75	0.52
1:A:71:LYS:NZ	4:A:203:ACT:O	2.40	0.52
1:Q:67[A]:LYS:HD2	2:R:202:SO4:O1	2.10	0.52
4:A:203:ACT:H1	6:A:486:HOH:O	2.10	0.51
1:C:80:LYS:HE2	1:C:82:LYS:HE3	1.92	0.51
1:M:163:GLU:OE1	5:W:203:CL:CL	2.65	0.51
1:Q:20:LEU:HD11	1:Q:66:MET:HG3	1.91	0.51
1:Q:34:TYR:HB3	6:Q:503:HOH:O	2.09	0.51
1:Q:143:ARG:HD3	6:Q:429:HOH:O	2.09	0.51
1:S:80:LYS:HE3	6:S:526:HOH:O	2.10	0.51
1:J:67:LYS:HD3	1:J:71:LYS:HE3	1.92	0.51
1:W:2:SER:CB	3:W:204:EDO:H11	2.39	0.51
1:S:143:ARG:NH1	6:S:445:HOH:O	2.40	0.51
1:F:18:ARG:HD3	6:F:384:HOH:O	2.09	0.51
1:C:66:MET:CE	1:D:34:TYR:CD2	2.94	0.51
1:D:136[A]:GLU:HG2	6:D:328:HOH:O	2.09	0.51
1:R:56:HIS:CD2	1:R:59:ARG:NH2	2.79	0.51
1:A:64[A]:LYS:HD2	6:A:446:HOH:O	2.10	0.51
1:E:113:LEU:HG	1:E:129:LEU:HD11	1.92	0.51
1:F:66:MET:HE3	1:G:35:ALA:HB2	1.93	0.51
1:Q:43:ALA:HB1	3:Q:202:EDO:H22	1.91	0.51
1:Q:43:ALA:C	3:Q:202:EDO:H21	2.36	0.51
1:X:101:GLN:HG3	6:X:481:HOH:O	2.10	0.51
1:L:56:HIS:CD2	1:L:59:ARG:HH12	2.29	0.50
1:P:125:LEU:HD22	6:P:442:HOH:O	2.11	0.50
1:I:143:ARG:NH1	6:I:503:HOH:O	2.43	0.50
1:Q:66:MET:HE2	1:R:34:TYR:HE2	1.75	0.50
1:H:53:GLU:HG3	6:H:385:HOH:O	2.11	0.50
1:H:154:LEU:HG	6:H:445:HOH:O	2.10	0.50
1:C:9[B]:HIS:CD2	1:C:120:LYS:HB3	2.45	0.50
1:C:110:LEU:HD12	1:C:137:GLN:HG3	1.93	0.50
1:P:42:VAL:HG13	3:P:204:EDO:H22	1.93	0.50
1:V:115:LYS:HD3	6:V:414:HOH:O	2.11	0.50
1:A:56:HIS:CE1	1:A:60:GLU:OE1	2.65	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:101[B]:GLN:NE2	1:K:104[B]:LYS:HD2	2.26	0.50
1:S:80:LYS:HD2	1:T:80:LYS:HD2	1.93	0.50
1:S:85:GLU:HG3	1:S:86:ARG:HG3	1.92	0.50
1:G:94:GLU:HA	1:G:97:GLN:HE21	1.77	0.50
1:H:159:ASN:ND2	1:K:161:MET:HE3	2.22	0.50
1:Q:67[A]:LYS:HG2	1:R:34:TYR:OH	2.11	0.50
1:X:143:ARG:HD3	6:X:409:HOH:O	2.11	0.50
1:D:28:TYR:HB3	6:D:391:HOH:O	2.12	0.50
1:E:64:LYS:CG	3:E:204:EDO:H21	2.42	0.50
1:N:11[B]:ASP:OD2	1:N:120:LYS:HE3	2.12	0.50
1:F:101:GLN:CD	6:F:381:HOH:O	2.55	0.50
1:H:31[B]:SER:OG	1:J:66:MET:HE1	2.11	0.50
1:K:121:VAL:HG11	1:X:115:LYS:HE3	1.93	0.49
1:X:91:ASN:HA	3:X:204:EDO:H11	1.94	0.49
1:C:56:HIS:CD2	1:C:59:ARG:HH12	2.31	0.49
1:T:59:ARG:HB3	1:T:59:ARG:HH11	1.77	0.49
1:E:20:LEU:HD11	1:E:66:MET:HG2	1.94	0.49
1:V:115:LYS:HD2	1:X:121:VAL:HG11	1.93	0.49
1:H:92:THR:HA	6:H:467:HOH:O	2.12	0.49
1:I:11[B]:ASP:OD2	1:I:120:LYS:CE	2.59	0.49
1:B:66:MET:HB3	6:B:471:HOH:O	2.13	0.49
1:P:45:HIS:HE1	6:P:364:HOH:O	1.96	0.49
1:P:59:ARG:NH1	6:P:456:HOH:O	2.42	0.49
1:I:9[B]:HIS:HD2	1:I:120:LYS:HD2	1.78	0.49
1:J:142:LYS:NZ	3:J:205:EDO:H22	2.28	0.49
1:V:146:ASP:HB3	3:V:201:EDO:H22	1.95	0.49
1:S:9[B]:HIS:CD2	1:S:120:LYS:HB3	2.47	0.49
1:E:135:GLU:CG	1:E:139:LYS:HE3	2.41	0.49
1:R:2:SER:HB2	3:R:206:EDO:H11	1.95	0.49
1:T:80:LYS:HE2	1:T:82:LYS:HD3	1.95	0.49
1:E:135:GLU:OE2	1:H:71:LYS:NZ	2.42	0.49
1:Q:20:LEU:HD11	1:Q:66:MET:CG	2.43	0.49
1:Q:113:LEU:HG	1:Q:129:LEU:HD11	1.95	0.48
1:K:171:VAL:CG1	6:K:497:HOH:O	2.61	0.48
1:L:169:HIS:NE2	2:L:201:SO4:O2	2.46	0.48
1:W:7:ASN:N	2:W:202:SO4:O2	2.39	0.48
1:O:66:MET:HE2	1:O:76:VAL:HG11	1.96	0.48
1:M:66:MET:HE2	1:M:78:LEU:HD21	1.95	0.48
1:C:20[A]:LEU:HD11	1:C:66:MET:CG	2.44	0.48
1:Q:34:TYR:HD2	6:Q:503:HOH:O	1.90	0.48
1:S:9[B]:HIS:HD2	6:S:334:HOH:O	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:115:LYS:NZ	1:R:119:ASP:OD1	2.46	0.48
1:R:120:LYS:HD2	1:R:120:LYS:N	2.28	0.48
1:V:56:HIS:HB3	6:V:538:HOH:O	2.12	0.48
1:E:66:MET:HE2	1:E:78:LEU:HD21	1.96	0.48
1:C:66:MET:HE2	1:D:34:TYR:CE2	2.49	0.47
1:M:9[B]:HIS:HD2	1:M:120:LYS:HB3	1.80	0.47
1:T:66:MET:HE2	1:T:78:LEU:HD21	1.96	0.47
1:U:19:MET:HE1	1:U:22:LEU:HD23	1.96	0.47
1:A:59:ARG:NH2	6:A:468:HOH:O	2.46	0.47
1:W:73:GLY:O	3:W:204:EDO:H12	2.14	0.47
1:A:156:LEU:CD2	6:A:460:HOH:O	2.61	0.47
1:K:101[B]:GLN:HG2	6:K:410:HOH:O	2.14	0.47
1:Q:96:MET:CG	6:Q:449:HOH:O	2.53	0.47
1:S:168:LYS:HE2	6:S:519:HOH:O	2.15	0.47
1:F:171:VAL:O	1:F:172:LYS:HB3	2.14	0.47
1:K:96:MET:CE	6:K:497:HOH:O	2.45	0.47
1:U:82:LYS:HE2	1:U:82:LYS:CA	2.34	0.47
1:U:115[B]:LYS:HE2	1:U:119:ASP:OD1	2.15	0.47
1:E:168:LYS:HE2	6:E:503:HOH:O	2.14	0.47
1:G:168:LYS:HG2	6:G:509:HOH:O	2.14	0.47
1:J:115:LYS:NZ	6:J:477:HOH:O	2.41	0.47
1:Q:97:GLN:HG2	6:Q:486:HOH:O	2.15	0.47
1:F:171:VAL:O	1:F:172:LYS:CB	2.62	0.47
1:L:130:GLU:N	6:L:447:HOH:O	2.48	0.47
1:A:64[A]:LYS:NZ	1:A:132:GLU:OE1	2.37	0.46
1:C:64:LYS:HB3	1:C:133:TYR:OH	2.15	0.46
1:P:136[A]:GLU:HG2	6:P:321:HOH:O	2.13	0.46
1:B:88:GLU:HG2	3:B:203:EDO:H21	1.96	0.46
1:C:143:ARG:NH1	6:C:365:HOH:O	2.48	0.46
1:M:9[B]:HIS:CD2	1:M:120:LYS:HB3	2.50	0.46
1:I:36:PHE:HE1	1:I:88:GLU:HG2	1.79	0.46
1:U:34:TYR:OH	1:V:67:LYS:HG2	2.15	0.46
1:N:113:LEU:HG	1:N:129:LEU:HD11	1.96	0.46
3:N:205:EDO:H11	6:N:333:HOH:O	2.15	0.46
1:T:113:LEU:HG	1:T:129:LEU:HD11	1.97	0.46
1:Q:80:LYS:HE3	1:Q:82:LYS:HD3	1.98	0.46
1:R:73:GLY:O	3:R:206:EDO:H22	2.12	0.46
1:U:82:LYS:HA	1:U:82:LYS:HE3	1.93	0.46
1:T:1:VAL:HA	3:T:205:EDO:H21	1.97	0.46
1:X:97:GLN:NE2	6:X:435:HOH:O	2.48	0.46
1:C:143:ARG:HD3	6:C:400:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:172:LYS:CB	6:T:368:HOH:O	2.63	0.46
1:W:66:MET:HE1	1:X:34:TYR:CD2	2.51	0.46
1:P:64:LYS:HG3	3:P:202:EDO:C1	2.45	0.46
1:P:144:ILE:HG12	6:P:488:HOH:O	2.15	0.46
1:K:71:LYS:HD3	3:K:203:EDO:O1	2.16	0.46
1:C:66:MET:HE1	1:D:34:TYR:CD2	2.49	0.46
1:F:66:MET:HE1	1:G:34:TYR:HD2	1.79	0.46
1:C:20[B]:LEU:HD23	1:C:65:PHE:CB	2.30	0.45
1:C:143:ARG:NE	6:C:372:HOH:O	2.48	0.45
1:G:155:GLY:HA3	3:G:210:EDO:H21	1.98	0.45
1:R:23:GLU:OE1	1:R:58:GLU:OE1	2.34	0.45
1:A:171:VAL:O	1:A:172:LYS:HB3	2.16	0.45
1:G:5:ARG:HH11	3:G:211:EDO:C2	2.29	0.45
1:K:104[B]:LYS:NZ	6:K:468:HOH:O	2.47	0.45
1:W:5:ARG:HH11	3:W:204:EDO:H12	1.82	0.45
1:H:171:VAL:O	1:H:172:LYS:CB	2.63	0.45
1:I:67[A]:LYS:HB3	1:I:67[A]:LYS:HE3	1.66	0.45
1:L:152:LYS:CE	6:L:499:HOH:O	2.63	0.45
1:Q:67[A]:LYS:HB3	1:Q:67[A]:LYS:HE2	1.75	0.45
1:V:169:HIS:NE2	2:W:201:SO4:O1	2.49	0.45
1:N:121:VAL:HG13	3:N:204:EDO:H22	1.99	0.45
1:H:136[A]:GLU:HG2	6:H:397:HOH:O	2.16	0.45
1:W:67:LYS:HB3	1:W:67:LYS:HE3	1.53	0.45
1:A:101:GLN:HG2	6:A:490:HOH:O	2.17	0.45
1:C:9[A]:HIS:CD2	1:C:120:LYS:HD2	2.52	0.45
1:P:45:HIS:CE1	6:P:364:HOH:O	2.68	0.45
1:G:97:GLN:NE2	6:G:415:HOH:O	2.48	0.45
1:H:92:THR:HG23	6:H:467:HOH:O	2.17	0.45
1:C:80:LYS:HE2	1:C:82:LYS:HD3	1.98	0.45
1:H:80:LYS:HE3	6:H:406:HOH:O	2.17	0.45
1:Q:124:HIS:HE1	6:Q:376:HOH:O	1.99	0.45
1:R:127:ASP:OD2	1:U:135:GLU:OE2	2.34	0.45
1:V:20:LEU:HD11	1:V:66:MET:HG2	1.99	0.45
1:C:80:LYS:HE2	1:C:82:LYS:CD	2.47	0.45
1:C:96:MET:HB2	6:C:451:HOH:O	2.17	0.45
1:F:104:LYS:HE2	6:F:403:HOH:O	2.17	0.45
1:G:75:ARG:HG2	3:G:211:EDO:O1	2.17	0.45
1:H:164:TYR:CE1	6:Q:484:HOH:O	2.46	0.45
1:Q:34:TYR:OH	4:Q:204:ACT:H3	2.18	0.44
1:M:113:LEU:HG	1:M:129:LEU:HD11	1.97	0.44
1:N:168:LYS:HE2	6:N:482:HOH:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:45:HIS:HD2	6:O:475:HOH:O	2.00	0.44
1:S:67:LYS:HE3	1:S:67:LYS:HB3	1.61	0.44
1:H:9[B]:HIS:HD2	1:H:120:LYS:HD2	1.82	0.44
1:I:121:VAL:HA	3:I:204:EDO:H11	1.99	0.44
1:J:142:LYS:HE3	3:J:205:EDO:H12	1.99	0.44
1:K:94:GLU:HA	1:K:97:GLN:HE21	1.81	0.44
1:B:71:LYS:HB2	1:B:71:LYS:HE3	1.79	0.44
1:O:66:MET:CE	1:O:76:VAL:HG11	2.48	0.44
1:G:85:GLU:OE2	3:G:207:EDO:O2	2.26	0.44
1:L:80:LYS:HE3	1:L:82:LYS:HZ3	1.82	0.44
1:L:113:LEU:HG	1:L:129:LEU:HD11	1.98	0.44
1:C:15:ALA:HB1	1:C:113:LEU:HD13	1.98	0.44
1:F:80:LYS:CE	1:F:82:LYS:HD3	2.48	0.44
1:J:113:LEU:HG	1:J:129:LEU:HD11	1.98	0.44
1:U:147:PHE:HD2	6:U:500:HOH:O	2.00	0.44
1:B:113:LEU:HG	1:B:129:LEU:HD11	1.98	0.44
1:I:86:ARG:NH1	1:I:94:GLU:OE2	2.42	0.44
1:B:39:ARG:HH12	3:B:203:EDO:H22	1.83	0.44
1:P:82:LYS:HE2	1:P:82:LYS:HB3	1.81	0.44
1:P:113:LEU:HG	1:P:129:LEU:HD11	2.00	0.44
1:H:9[B]:HIS:HE1	6:H:458:HOH:O	1.99	0.44
1:K:101[B]:GLN:HA	1:K:104[B]:LYS:HG2	1.99	0.44
1:M:67:LYS:HB2	1:M:67:LYS:HE2	1.86	0.44
1:P:64:LYS:O	3:P:202:EDO:H11	2.18	0.44
1:V:63:GLU:HG2	1:V:66:MET:CE	2.47	0.44
1:K:101[B]:GLN:OE1	1:K:104[B]:LYS:HD2	2.17	0.43
1:I:113:LEU:HG	1:I:129:LEU:HD11	1.99	0.43
1:L:56:HIS:CD2	1:L:59:ARG:HH22	2.36	0.43
1:L:143[A]:ARG:NH1	6:L:422:HOH:O	2.50	0.43
1:P:115:LYS:HG3	6:G:430:HOH:O	2.18	0.43
1:S:50:PHE:HE1	1:S:143:ARG:NH1	2.16	0.43
1:I:5:ARG:HH11	3:I:203:EDO:C1	2.30	0.43
1:K:11[B]:ASP:OD2	1:K:120:LYS:CE	2.62	0.43
1:Q:66:MET:CE	1:R:34:TYR:CE2	3.00	0.43
1:A:71:LYS:NZ	4:A:203:ACT:C	2.81	0.43
1:H:20:LEU:HD11	1:H:66:MET:HG3	2.00	0.43
1:K:67:LYS:HG2	3:K:203:EDO:H21	2.00	0.43
1:L:54:HIS:CE1	1:L:143[A]:ARG:HH12	2.36	0.43
1:A:36:PHE:HE1	1:A:88:GLU:HG2	1.84	0.43
1:A:140:ASP:OD1	1:A:143:ARG:NH1	2.51	0.43
1:M:56:HIS:HD2	6:M:496:HOH:O	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:66:MET:HB3	6:E:400:HOH:O	2.18	0.43
1:F:28:TYR:OH	5:G:203:CL:CL	2.74	0.43
1:A:59:ARG:NH1	6:A:494:HOH:O	2.41	0.43
3:I:203:EDO:H21	6:I:380:HOH:O	2.17	0.43
1:K:18:ARG:HD2	6:K:406:HOH:O	2.17	0.43
1:Q:27:SER:HB2	1:Q:58:GLU:HB2	2.01	0.43
1:X:168:LYS:HE2	6:X:365:HOH:O	2.17	0.43
1:A:94:GLU:HA	1:A:97:GLN:HE21	1.83	0.43
1:L:20:LEU:HD11	1:L:66:MET:HG3	2.00	0.43
1:R:82[B]:LYS:HE3	1:R:82[B]:LYS:HB3	1.25	0.43
1:T:56:HIS:O	1:T:60:GLU:HG3	2.19	0.43
1:H:168:LYS:HG2	6:H:492:HOH:O	2.18	0.43
1:L:140:ASP:OD1	1:L:143[A]:ARG:NH1	2.51	0.43
1:B:121:VAL:HA	3:R:205:EDO:H22	2.01	0.43
1:P:143[A]:ARG:CZ	6:P:389:HOH:O	2.66	0.43
1:K:136:GLU:HA	1:K:139:LYS:CE	2.46	0.43
1:K:138:VAL:HG21	1:V:124:HIS:CD2	2.53	0.43
1:X:9[B]:HIS:HD2	1:X:120:LYS:CG	2.32	0.43
1:A:54:HIS:O	1:A:58:GLU:HG2	2.19	0.43
1:C:115:LYS:HA	1:C:115:LYS:HD3	1.84	0.43
1:F:75:ARG:CZ	3:F:205:EDO:H21	2.48	0.43
1:J:115:LYS:O	1:J:115:LYS:HD3	2.18	0.43
1:R:2:SER:HB3	3:R:206:EDO:H22	2.01	0.43
1:M:168:LYS:HE2	6:M:503:HOH:O	2.19	0.42
1:P:71:LYS:HD3	3:P:202:EDO:C2	2.49	0.42
1:R:71:LYS:HG2	6:U:449:HOH:O	2.19	0.42
1:O:110:LEU:HD12	1:O:137:GLN:HG3	2.01	0.42
1:P:71:LYS:HB3	1:P:71:LYS:HE3	1.78	0.42
1:T:94:GLU:HA	1:T:97:GLN:HE21	1.84	0.42
1:G:64:LYS:NZ	1:G:132:GLU:OE1	2.48	0.42
1:F:168:LYS:HE2	6:F:388:HOH:O	2.19	0.42
1:Q:136:GLU:HA	1:Q:139:LYS:CE	2.49	0.42
1:V:63:GLU:HA	1:V:66:MET:CE	2.49	0.42
1:D:9[B]:HIS:CE1	6:D:424:HOH:O	2.71	0.42
1:K:20:LEU:C	1:K:20:LEU:HD23	2.44	0.42
1:L:134:LEU:CD1	6:L:447:HOH:O	2.67	0.42
1:U:113:LEU:HG	1:U:129:LEU:HD11	2.00	0.42
1:B:64:LYS:NZ	1:B:132:GLU:OE1	2.46	0.42
1:X:113:LEU:HG	1:X:129:LEU:HD11	2.00	0.42
6:C:402:HOH:O	1:F:104:LYS:CE	2.68	0.42
1:O:64:LYS:HB3	1:O:133:TYR:OH	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:LEU:HD22	2:W:201:SO4:O2	2.19	0.42
1:C:20[A]:LEU:HD11	1:C:66:MET:HG3	2.02	0.42
1:Q:66:MET:HE1	1:R:34:TYR:CD2	2.43	0.42
1:S:56[A]:HIS:CE1	1:T:59:ARG:HH21	2.37	0.42
1:G:168:LYS:HE2	6:G:509:HOH:O	2.18	0.42
1:K:97:GLN:NE2	6:K:417:HOH:O	2.53	0.42
1:A:113:LEU:HG	1:A:129:LEU:HD11	2.02	0.42
1:T:116:LEU:O	1:T:120:LYS:HG2	2.20	0.42
1:F:82:LYS:HE3	6:F:340:HOH:O	2.19	0.42
1:H:43:ALA:HA	5:H:203:CL:CL	2.57	0.42
1:H:56:HIS:CD2	1:H:59:ARG:HH22	2.38	0.42
1:K:171:VAL:HG11	6:K:497:HOH:O	2.20	0.42
1:H:85:GLU:HG3	1:H:86:ARG:HG3	2.02	0.42
1:L:134:LEU:HD12	6:L:447:HOH:O	2.18	0.42
1:V:63:GLU:HA	1:V:66:MET:HE3	2.02	0.42
1:E:23:GLU:OE1	1:E:58:GLU:OE1	2.38	0.41
1:U:98:ALA:O	5:U:202:CL:CL	2.75	0.41
1:W:23:GLU:OE1	1:W:58:GLU:OE1	2.37	0.41
1:B:39:ARG:HH22	3:B:203:EDO:H12	1.85	0.41
1:N:49:GLU:OE1	1:N:52:LYS:HE3	2.21	0.41
1:G:159:ASN:HB2	3:G:210:EDO:H12	2.03	0.41
1:K:104[B]:LYS:HB3	1:K:104[B]:LYS:HE2	1.70	0.41
1:K:151:LEU:HD22	1:K:156:LEU:HD22	2.01	0.41
1:R:113:LEU:HG	1:R:129:LEU:HD11	2.01	0.41
1:W:31[B]:SER:OG	1:X:66:MET:HE1	2.20	0.41
1:B:64:LYS:HB3	1:B:133:TYR:OH	2.20	0.41
1:C:15:ALA:CB	1:C:113:LEU:HD13	2.50	0.41
1:C:168:LYS:HE2	6:C:503:HOH:O	2.20	0.41
1:N:6:GLN:HA	2:N:201:SO4:O1	2.21	0.41
1:R:115:LYS:HE2	1:R:115:LYS:CA	2.45	0.41
1:M:71:LYS:HE2	6:M:428:HOH:O	2.21	0.41
1:E:34:TYR:OH	1:I:67[A]:LYS:HG2	2.20	0.41
1:A:34:TYR:CD1	1:A:52:LYS:HB2	2.55	0.41
1:G:5:ARG:HH11	3:G:211:EDO:H22	1.85	0.41
1:U:86:ARG:NH2	1:U:88:GLU:OE2	2.49	0.41
1:W:5:ARG:HH11	3:W:204:EDO:C1	2.34	0.41
1:A:38:ASP:C	6:A:455:HOH:O	2.63	0.41
1:C:115:LYS:HE3	1:S:121:VAL:HG11	2.03	0.41
1:U:45:HIS:HD2	6:U:496:HOH:O	2.03	0.41
1:G:98:ALA:HB2	3:G:207:EDO:H11	2.03	0.41
1:L:54:HIS:HE1	1:L:143[A]:ARG:NH1	2.13	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:64:LYS:NZ	1:W:132:GLU:HB3	2.35	0.41
1:C:66:MET:HE2	1:D:34:TYR:HE2	1.85	0.41
1:O:115:LYS:O	1:O:115:LYS:HD3	2.20	0.41
1:J:151:LEU:HD22	1:J:156:LEU:HD22	2.02	0.41
1:R:11[B]:ASP:OD2	1:R:120:LYS:NZ	2.48	0.41
1:R:20[A]:LEU:HD11	1:R:66:MET:HG2	2.02	0.41
1:U:94:GLU:HA	1:U:97:GLN:HE21	1.85	0.41
1:A:135[A]:GLU:OE2	1:D:71:LYS:NZ	2.47	0.41
1:A:156:LEU:HD21	6:A:460:HOH:O	2.19	0.41
1:D:64:LYS:HD3	6:D:454:HOH:O	2.20	0.41
1:P:28:TYR:HB3	6:P:399:HOH:O	2.20	0.41
1:L:141:ILE:HD12	6:L:503:HOH:O	2.20	0.41
1:S:135:GLU:OE2	1:F:127:ASP:OD2	2.39	0.41
1:F:64[A]:LYS:HE2	1:F:64[A]:LYS:HB2	1.66	0.41
1:U:66:MET:HE3	1:V:34:TYR:CE2	2.56	0.41
1:D:39:ARG:NH1	1:D:88[B]:GLU:HG2	2.36	0.40
1:N:57:GLU:OE1	1:N:140:ASP:OD2	2.39	0.40
1:E:135:GLU:CD	1:E:139:LYS:HE3	2.46	0.40
1:H:167:ASP:HB3	6:H:492:HOH:O	2.21	0.40
1:J:23:GLU:OE1	1:J:58:GLU:OE1	2.38	0.40
1:Q:161:MET:HE2	1:U:165:LEU:HD12	2.03	0.40
1:U:15:ALA:HB1	1:U:113:LEU:HD13	2.03	0.40
3:N:203:EDO:H12	6:N:453:HOH:O	2.21	0.40
1:P:20:LEU:HG	6:P:370:HOH:O	2.20	0.40
1:W:113:LEU:HG	1:W:129:LEU:HD11	2.02	0.40
1:P:123:PRO:HB3	1:I:114:HIS:CE1	2.55	0.40
1:T:20:LEU:HD11	1:T:66:MET:CG	2.51	0.40
3:E:204:EDO:C1	6:E:504:HOH:O	2.69	0.40
1:U:115[A]:LYS:HD2	6:U:473:HOH:O	2.22	0.40
1:U:143:ARG:HD3	6:U:483:HOH:O	2.20	0.40
1:C:20[A]:LEU:HD11	1:C:66:MET:HG2	2.02	0.40
2:D:201:SO4:O2	1:N:169:HIS:NE2	2.54	0.40
1:H:66:MET:HE2	1:J:34:TYR:CE2	2.57	0.40
1:K:168:LYS:HE2	6:K:496:HOH:O	2.20	0.40
1:Q:147:PHE:CB	6:Q:449:HOH:O	2.47	0.40
1:R:161:MET:HE3	1:R:161:MET:HB2	1.95	0.40
1:R:169:HIS:NE2	2:R:201:SO4:O1	2.54	0.40
6:D:465:HOH:O	1:N:165:LEU:CD2	2.60	0.40
1:G:169:HIS:NE2	2:J:201:SO4:O3	2.49	0.40
1:W:61:HIS:HB3	1:W:133:TYR:HE1	1.85	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	179/176 (102%)	176 (98%)	3 (2%)	0	100	100
1	B	176/176 (100%)	172 (98%)	4 (2%)	0	100	100
1	C	178/176 (101%)	175 (98%)	3 (2%)	0	100	100
1	D	176/176 (100%)	172 (98%)	4 (2%)	0	100	100
1	E	174/176 (99%)	170 (98%)	4 (2%)	0	100	100
1	F	178/176 (101%)	175 (98%)	3 (2%)	0	100	100
1	G	175/176 (99%)	173 (99%)	2 (1%)	0	100	100
1	H	176/176 (100%)	172 (98%)	4 (2%)	0	100	100
1	I	178/176 (101%)	174 (98%)	4 (2%)	0	100	100
1	J	176/176 (100%)	172 (98%)	4 (2%)	0	100	100
1	K	178/176 (101%)	174 (98%)	4 (2%)	0	100	100
1	L	176/176 (100%)	173 (98%)	3 (2%)	0	100	100
1	M	172/176 (98%)	169 (98%)	3 (2%)	0	100	100
1	N	178/176 (101%)	175 (98%)	3 (2%)	0	100	100
1	O	175/176 (99%)	171 (98%)	4 (2%)	0	100	100
1	P	175/176 (99%)	171 (98%)	4 (2%)	0	100	100
1	Q	177/176 (101%)	175 (99%)	2 (1%)	0	100	100
1	R	175/176 (99%)	173 (99%)	2 (1%)	0	100	100
1	S	178/176 (101%)	175 (98%)	3 (2%)	0	100	100
1	T	175/176 (99%)	170 (97%)	5 (3%)	0	100	100
1	U	174/176 (99%)	171 (98%)	3 (2%)	0	100	100
1	V	178/176 (101%)	174 (98%)	4 (2%)	0	100	100
1	W	177/176 (101%)	172 (97%)	5 (3%)	0	100	100
1	X	173/176 (98%)	169 (98%)	4 (2%)	0	100	100
All	All	4227/4224 (100%)	4143 (98%)	84 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	160/157 (102%)	159 (99%)	1 (1%)	78	75
1	B	158/157 (101%)	154 (98%)	4 (2%)	42	26
1	C	160/157 (102%)	158 (99%)	2 (1%)	61	53
1	D	157/157 (100%)	153 (98%)	4 (2%)	42	26
1	E	155/157 (99%)	154 (99%)	1 (1%)	78	75
1	F	160/157 (102%)	158 (99%)	2 (1%)	61	53
1	G	158/157 (101%)	153 (97%)	5 (3%)	34	18
1	H	159/157 (101%)	156 (98%)	3 (2%)	50	37
1	I	161/157 (102%)	156 (97%)	5 (3%)	35	19
1	J	158/157 (101%)	157 (99%)	1 (1%)	78	75
1	K	161/157 (102%)	153 (95%)	8 (5%)	22	8
1	L	157/157 (100%)	155 (99%)	2 (1%)	61	53
1	M	154/157 (98%)	149 (97%)	5 (3%)	34	18
1	N	160/157 (102%)	154 (96%)	6 (4%)	29	14
1	O	156/157 (99%)	154 (99%)	2 (1%)	61	53
1	P	157/157 (100%)	154 (98%)	3 (2%)	50	37
1	Q	160/157 (102%)	156 (98%)	4 (2%)	42	26
1	R	156/157 (99%)	152 (97%)	4 (3%)	40	24
1	S	159/157 (101%)	154 (97%)	5 (3%)	35	19
1	T	157/157 (100%)	155 (99%)	2 (1%)	61	53
1	U	157/157 (100%)	152 (97%)	5 (3%)	34	18
1	V	159/157 (101%)	155 (98%)	4 (2%)	42	26
1	W	159/157 (101%)	154 (97%)	5 (3%)	35	19
1	X	155/157 (99%)	153 (99%)	2 (1%)	61	53

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	3793/3768 (101%)	3708 (98%)	85 (2%)	45 32

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	GLU
1	B	1	VAL
1	B	71	LYS
1	B	137	GLN
1	B	156	LEU
1	C	31	SER
1	C	115	LYS
1	D	52	LYS
1	D	67	LYS
1	D	82	LYS
1	D	134	LEU
1	M	20	LEU
1	M	101	GLN
1	M	121	VAL
1	M	134	LEU
1	M	139	LYS
1	N	20	LEU
1	N	53	GLU
1	N	82	LYS
1	N	134	LEU
1	N	137	GLN
1	N	156	LEU
1	O	20	LEU
1	O	139	LYS
1	P	53	GLU
1	P	115	LYS
1	P	121	VAL
1	S	1	VAL
1	S	67	LYS
1	S	71	LYS
1	S	139	LYS
1	S	156	LEU
1	T	1	VAL
1	T	59	ARG
1	E	1	VAL
1	F	121	VAL
1	F	134	LEU

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Mol	Chain	Res	Type
1	G	1	VAL
1	G	88	GLU
1	G	120	LYS
1	G	129	LEU
1	G	134	LEU
1	H	101	GLN
1	H	134	LEU
1	H	171	VAL
1	I	63	GLU
1	I	67[A]	LYS
1	I	67[B]	LYS
1	I	80	LYS
1	I	139	LYS
1	J	97	GLN
1	K	60	GLU
1	K	85	GLU
1	K	101[A]	GLN
1	K	101[B]	GLN
1	K	113	LEU
1	K	121	VAL
1	K	137	GLN
1	K	158	GLU
1	L	1	VAL
1	L	97	GLN
1	Q	101[A]	GLN
1	Q	101[B]	GLN
1	Q	137	GLN
1	Q	172	LYS
1	R	115	LYS
1	R	120	LYS
1	R	137	GLN
1	R	156	LEU
1	U	82	LYS
1	U	85	GLU
1	U	121	VAL
1	U	134	LEU
1	U	172	LYS
1	V	1	VAL
1	V	115	LYS
1	V	129	LEU
1	V	134	LEU
1	W	67	LYS

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Mol	Chain	Res	Type
1	W	71	LYS
1	W	101	GLN
1	W	134	LEU
1	W	172	LYS
1	X	115	LYS
1	X	120	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (78) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	56	HIS
1	A	79	GLN
1	A	97	GLN
1	A	107	ASN
1	A	108	GLN
1	B	21	ASN
1	B	46	ASN
1	C	56	HIS
1	C	69	GLN
1	C	97	GLN
1	C	107	ASN
1	D	3	GLN
1	D	21	ASN
1	D	97	GLN
1	D	107	ASN
1	M	56	HIS
1	M	97	GLN
1	N	97	GLN
1	N	137	GLN
1	O	45	HIS
1	O	97	GLN
1	O	107	ASN
1	O	108	GLN
1	P	45	HIS
1	P	46	ASN
1	P	56	HIS
1	P	97	GLN
1	P	107	ASN
1	P	137	GLN
1	S	3	GLN
1	S	46	ASN

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Mol	Chain	Res	Type
1	S	97	GLN
1	S	108	GLN
1	T	97	GLN
1	E	3	GLN
1	E	56	HIS
1	E	97	GLN
1	E	101	GLN
1	E	108	GLN
1	F	56	HIS
1	F	97	GLN
1	F	107	ASN
1	G	97	GLN
1	G	107	ASN
1	H	46	ASN
1	H	56	HIS
1	H	79	GLN
1	H	97	GLN
1	H	108	GLN
1	H	159	ASN
1	I	3	GLN
1	I	45	HIS
1	I	97	GLN
1	J	97	GLN
1	K	45	HIS
1	K	97	GLN
1	K	107	ASN
1	K	137	GLN
1	L	45	HIS
1	L	97	GLN
1	L	101	GLN
1	L	137	GLN
1	Q	79	GLN
1	Q	97	GLN
1	Q	107	ASN
1	R	45	HIS
1	R	56	HIS
1	R	137	GLN
1	U	45	HIS
1	U	97	GLN
1	V	45	HIS
1	W	45	HIS
1	W	97	GLN

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Mol	Chain	Res	Type
1	W	101	GLN
1	X	3	GLN
1	X	97	GLN
1	X	101	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 109 ligands modelled in this entry, 31 are monoatomic - leaving 78 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	$\# Z > 2$
2	SO4	K	201	-	4,4,4	0.24	0	6,6,6	0.23	0
2	SO4	L	201	-	4,4,4	0.27	0	6,6,6	0.37	0
3	EDO	P	203	-	3,3,3	0.49	0	2,2,2	0.28	0
3	EDO	B	203	-	3,3,3	0.30	0	2,2,2	0.65	0
3	EDO	X	205	-	3,3,3	0.40	0	2,2,2	0.47	0
4	ACT	Q	204	-	3,3,3	0.90	0	3,3,3	0.76	0
3	EDO	M	202	-	3,3,3	0.52	0	2,2,2	0.24	0
3	EDO	J	205	-	3,3,3	0.37	0	2,2,2	0.49	0
3	EDO	F	205	-	3,3,3	0.47	0	2,2,2	0.44	0
3	EDO	W	204	-	3,3,3	0.26	0	2,2,2	0.65	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	L	205	-	3,3,3	0.51	0	2,2,2	0.05	0
3	EDO	C	205	-	3,3,3	0.37	0	2,2,2	0.54	0
3	EDO	G	207	-	3,3,3	0.31	0	2,2,2	0.63	0
3	EDO	A	202	-	3,3,3	0.44	0	2,2,2	0.36	0
2	SO4	J	202	-	4,4,4	0.48	0	6,6,6	0.11	0
3	EDO	G	208	-	3,3,3	0.39	0	2,2,2	0.47	0
2	SO4	N	202	-	4,4,4	0.39	0	6,6,6	0.19	0
3	EDO	J	204	-	3,3,3	0.31	0	2,2,2	0.67	0
2	SO4	U	201	-	4,4,4	0.40	0	6,6,6	0.14	0
2	SO4	R	202	-	4,4,4	0.43	0	6,6,6	0.13	0
3	EDO	G	211	-	3,3,3	0.27	0	2,2,2	0.57	0
3	EDO	I	204	-	3,3,3	0.45	0	2,2,2	0.59	0
3	EDO	G	205	-	3,3,3	0.47	0	2,2,2	0.76	0
2	SO4	W	201	-	4,4,4	0.23	0	6,6,6	0.14	0
3	EDO	I	203	-	3,3,3	0.24	0	2,2,2	0.89	0
3	EDO	B	204	-	3,3,3	0.36	0	2,2,2	0.21	0
3	EDO	Q	203	-	3,3,3	0.52	0	2,2,2	0.31	0
3	EDO	U	204	-	3,3,3	0.40	0	2,2,2	0.58	0
2	SO4	H	201	-	4,4,4	0.44	0	6,6,6	0.13	0
3	EDO	E	205	-	3,3,3	0.32	0	2,2,2	0.59	0
3	EDO	G	210	-	3,3,3	0.40	0	2,2,2	0.61	0
2	SO4	N	201	-	4,4,4	0.41	0	6,6,6	0.20	0
4	ACT	A	203	-	3,3,3	0.66	0	3,3,3	1.21	0
2	SO4	J	201	-	4,4,4	0.37	0	6,6,6	0.20	0
2	SO4	K	202	-	4,4,4	0.49	0	6,6,6	0.12	0
2	SO4	W	202	-	4,4,4	0.43	0	6,6,6	0.12	0
3	EDO	R	204	-	3,3,3	0.32	0	2,2,2	0.68	0
2	SO4	O	201	-	4,4,4	0.44	0	6,6,6	0.12	0
3	EDO	E	204	-	3,3,3	0.50	0	2,2,2	0.12	0
2	SO4	D	202	-	4,4,4	0.43	0	6,6,6	0.14	0
3	EDO	Q	202	-	3,3,3	0.35	0	2,2,2	0.56	0
3	EDO	K	204	-	3,3,3	0.52	0	2,2,2	0.32	0
3	EDO	F	204	-	3,3,3	0.41	0	2,2,2	0.35	0
3	EDO	N	203	-	3,3,3	0.27	0	2,2,2	0.36	0
3	EDO	R	203	-	3,3,3	0.51	0	2,2,2	0.27	0
2	SO4	T	202	-	4,4,4	0.39	0	6,6,6	0.25	0
3	EDO	N	204	-	3,3,3	0.55	0	2,2,2	0.22	0
2	SO4	H	202	-	4,4,4	0.39	0	6,6,6	0.13	0
3	EDO	C	203	-	3,3,3	0.48	0	2,2,2	0.47	0
2	SO4	A	201	-	4,4,4	0.44	0	6,6,6	0.16	0
3	EDO	P	202	-	3,3,3	0.45	0	2,2,2	0.25	0
3	EDO	C	204	-	3,3,3	0.53	0	2,2,2	0.15	0
3	EDO	V	201	-	3,3,3	0.38	0	2,2,2	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	R	201	-	4,4,4	0.28	0	6,6,6	0.25	0
3	EDO	S	204	-	3,3,3	0.37	0	2,2,2	0.56	0
3	EDO	S	202	-	3,3,3	0.37	0	2,2,2	0.39	0
3	EDO	O	204	-	3,3,3	0.46	0	2,2,2	0.33	0
3	EDO	X	204	-	3,3,3	0.51	0	2,2,2	0.36	0
2	SO4	F	201	-	4,4,4	0.48	0	6,6,6	0.21	0
2	SO4	D	201	-	4,4,4	0.38	0	6,6,6	0.31	0
3	EDO	V	202	-	3,3,3	0.47	0	2,2,2	0.32	0
2	SO4	I	201	-	4,4,4	0.45	0	6,6,6	0.14	0
2	SO4	T	201	-	4,4,4	0.45	0	6,6,6	0.17	0
3	EDO	G	206	-	3,3,3	0.37	0	2,2,2	0.53	0
3	EDO	S	203	-	3,3,3	0.37	0	2,2,2	0.63	0
3	EDO	N	205	-	3,3,3	0.38	0	2,2,2	0.40	0
3	EDO	R	205	-	3,3,3	0.46	0	2,2,2	0.41	0
3	EDO	T	205	-	3,3,3	0.33	0	2,2,2	0.63	0
2	SO4	C	201	-	4,4,4	0.47	0	6,6,6	0.23	0
3	EDO	D	205	-	3,3,3	0.40	0	2,2,2	0.31	0
3	EDO	L	204	-	3,3,3	0.60	0	2,2,2	0.25	0
2	SO4	G	201	-	4,4,4	0.36	0	6,6,6	0.14	0
2	SO4	P	201	-	4,4,4	0.42	0	6,6,6	0.15	0
3	EDO	K	203	-	3,3,3	0.34	0	2,2,2	0.85	0
3	EDO	G	209	-	3,3,3	0.44	0	2,2,2	0.23	0
3	EDO	P	204	-	3,3,3	0.45	0	2,2,2	0.32	0
3	EDO	B	205	-	3,3,3	0.34	0	2,2,2	0.54	0
3	EDO	R	206	-	3,3,3	0.32	0	2,2,2	0.54	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	O	204	-	-	0/1/1/1	-
3	EDO	R	204	-	-	1/1/1/1	-
3	EDO	X	204	-	-	0/1/1/1	-
3	EDO	J	204	-	-	1/1/1/1	-
3	EDO	E	204	-	-	1/1/1/1	-
3	EDO	G	211	-	-	1/1/1/1	-
3	EDO	Q	202	-	-	0/1/1/1	-
3	EDO	V	202	-	-	1/1/1/1	-
3	EDO	I	204	-	-	0/1/1/1	-
3	EDO	G	206	-	-	0/1/1/1	-
3	EDO	F	204	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	G	205	-	-	0/1/1/1	-
3	EDO	K	204	-	-	0/1/1/1	-
3	EDO	S	203	-	-	1/1/1/1	-
3	EDO	N	203	-	-	1/1/1/1	-
3	EDO	P	203	-	-	0/1/1/1	-
3	EDO	I	203	-	-	0/1/1/1	-
3	EDO	B	204	-	-	1/1/1/1	-
3	EDO	R	203	-	-	0/1/1/1	-
3	EDO	B	203	-	-	0/1/1/1	-
3	EDO	N	205	-	-	1/1/1/1	-
3	EDO	R	205	-	-	1/1/1/1	-
3	EDO	X	205	-	-	1/1/1/1	-
3	EDO	Q	203	-	-	1/1/1/1	-
3	EDO	T	205	-	-	0/1/1/1	-
3	EDO	N	204	-	-	1/1/1/1	-
3	EDO	C	203	-	-	0/1/1/1	-
3	EDO	U	204	-	-	0/1/1/1	-
3	EDO	M	202	-	-	0/1/1/1	-
3	EDO	P	202	-	-	0/1/1/1	-
3	EDO	J	205	-	-	1/1/1/1	-
3	EDO	C	204	-	-	0/1/1/1	-
3	EDO	D	205	-	-	1/1/1/1	-
3	EDO	L	204	-	-	0/1/1/1	-
3	EDO	E	205	-	-	1/1/1/1	-
3	EDO	F	205	-	-	1/1/1/1	-
3	EDO	G	210	-	-	1/1/1/1	-
3	EDO	W	204	-	-	1/1/1/1	-
3	EDO	A	202	-	-	1/1/1/1	-
3	EDO	K	203	-	-	0/1/1/1	-
3	EDO	G	209	-	-	1/1/1/1	-
3	EDO	L	205	-	-	1/1/1/1	-
3	EDO	V	201	-	-	0/1/1/1	-
3	EDO	C	205	-	-	0/1/1/1	-
3	EDO	G	207	-	-	1/1/1/1	-
3	EDO	P	204	-	-	1/1/1/1	-
3	EDO	B	205	-	-	0/1/1/1	-
3	EDO	S	204	-	-	0/1/1/1	-
3	EDO	R	206	-	-	1/1/1/1	-
3	EDO	G	208	-	-	1/1/1/1	-
3	EDO	S	202	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	R	206	EDO	O1-C1-C2-O2
3	F	205	EDO	O1-C1-C2-O2
3	G	211	EDO	O1-C1-C2-O2
3	Q	203	EDO	O1-C1-C2-O2
3	V	202	EDO	O1-C1-C2-O2
3	W	204	EDO	O1-C1-C2-O2
3	A	202	EDO	O1-C1-C2-O2
3	D	205	EDO	O1-C1-C2-O2
3	R	204	EDO	O1-C1-C2-O2
3	N	204	EDO	O1-C1-C2-O2
3	N	205	EDO	O1-C1-C2-O2
3	F	204	EDO	O1-C1-C2-O2
3	L	205	EDO	O1-C1-C2-O2
3	G	210	EDO	O1-C1-C2-O2
3	J	204	EDO	O1-C1-C2-O2
3	B	204	EDO	O1-C1-C2-O2
3	P	204	EDO	O1-C1-C2-O2
3	G	209	EDO	O1-C1-C2-O2
3	S	203	EDO	O1-C1-C2-O2
3	E	205	EDO	O1-C1-C2-O2
3	G	207	EDO	O1-C1-C2-O2
3	G	208	EDO	O1-C1-C2-O2
3	R	205	EDO	O1-C1-C2-O2
3	N	203	EDO	O1-C1-C2-O2
3	E	204	EDO	O1-C1-C2-O2
3	S	202	EDO	O1-C1-C2-O2
3	J	205	EDO	O1-C1-C2-O2
3	X	205	EDO	O1-C1-C2-O2

There are no ring outliers.

45 monomers are involved in 104 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	201	SO4	1	0
2	L	201	SO4	2	0
3	P	203	EDO	1	0
3	B	203	EDO	3	0
4	Q	204	ACT	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	205	EDO	4	0
3	F	205	EDO	2	0
3	W	204	EDO	7	0
3	L	205	EDO	1	0
3	G	207	EDO	4	0
3	A	202	EDO	2	0
2	N	202	SO4	1	0
2	R	202	SO4	1	0
3	G	211	EDO	5	0
3	I	204	EDO	1	0
2	W	201	SO4	2	0
3	I	203	EDO	3	0
3	B	204	EDO	1	0
3	G	210	EDO	3	0
2	N	201	SO4	1	0
4	A	203	ACT	5	0
2	J	201	SO4	2	0
2	W	202	SO4	1	0
3	R	204	EDO	1	0
3	E	204	EDO	4	0
3	Q	202	EDO	5	0
3	F	204	EDO	1	0
3	N	203	EDO	4	0
3	N	204	EDO	1	0
3	P	202	EDO	3	0
3	V	201	EDO	2	0
2	R	201	SO4	2	0
3	S	202	EDO	5	0
3	X	204	EDO	1	0
2	D	201	SO4	2	0
3	V	202	EDO	1	0
2	T	201	SO4	1	0
3	N	205	EDO	2	0
3	R	205	EDO	1	0
3	T	205	EDO	2	0
3	D	205	EDO	1	0
2	P	201	SO4	1	0
3	K	203	EDO	2	0
3	P	204	EDO	1	0
3	R	206	EDO	7	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	172/176 (97%)	-0.10	1 (0%) 85 89	2, 5, 12, 29	16 (9%)
1	B	172/176 (97%)	-0.14	0 100 100	2, 5, 11, 27	9 (5%)
1	C	172/176 (97%)	0.03	1 (0%) 85 89	3, 6, 12, 23	16 (9%)
1	D	172/176 (97%)	0.14	3 (1%) 69 75	4, 7, 14, 28	12 (6%)
1	E	172/176 (97%)	-0.17	1 (0%) 85 89	3, 5, 11, 22	10 (5%)
1	F	172/176 (97%)	0.10	3 (1%) 69 75	3, 7, 12, 36	18 (10%)
1	G	172/176 (97%)	-0.04	1 (0%) 85 89	3, 6, 12, 31	11 (6%)
1	H	172/176 (97%)	0.01	3 (1%) 69 75	3, 6, 14, 34	12 (6%)
1	I	172/176 (97%)	-0.02	1 (0%) 85 89	2, 6, 12, 31	13 (7%)
1	J	172/176 (97%)	-0.13	0 100 100	2, 5, 11, 22	12 (6%)
1	K	172/176 (97%)	0.14	1 (0%) 85 89	3, 6, 14, 29	12 (6%)
1	L	172/176 (97%)	-0.03	2 (1%) 76 82	3, 6, 11, 23	11 (6%)
1	M	172/176 (97%)	-0.10	2 (1%) 76 82	2, 6, 12, 19	12 (6%)
1	N	172/176 (97%)	-0.01	3 (1%) 69 75	4, 7, 14, 32	13 (7%)
1	O	172/176 (97%)	-0.07	1 (0%) 85 89	3, 6, 12, 26	12 (6%)
1	P	172/176 (97%)	0.18	2 (1%) 76 82	4, 7, 16, 32	16 (9%)
1	Q	172/176 (97%)	0.04	1 (0%) 85 89	4, 6, 13, 30	12 (6%)
1	R	172/176 (97%)	0.04	3 (1%) 69 75	3, 6, 11, 20	12 (6%)
1	S	172/176 (97%)	-0.13	2 (1%) 76 82	3, 6, 12, 29	11 (6%)
1	T	172/176 (97%)	-0.08	2 (1%) 76 82	2, 5, 12, 24	11 (6%)
1	U	172/176 (97%)	0.06	1 (0%) 85 89	3, 6, 13, 32	12 (6%)
1	V	172/176 (97%)	-0.06	1 (0%) 85 89	2, 6, 11, 27	15 (8%)
1	W	172/176 (97%)	-0.18	2 (1%) 76 82	3, 5, 11, 29	14 (8%)
1	X	172/176 (97%)	-0.07	3 (1%) 69 75	3, 6, 12, 25	11 (6%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
All	All	4128/4224 (97%)	-0.03	40 (0%) 79 85	2, 6, 12, 36	303 (7%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	T	172	LYS	4.3
1	H	172	LYS	4.3
1	L	172	LYS	4.2
1	C	172	LYS	3.7
1	O	172	LYS	3.7
1	A	172	LYS	3.7
1	U	172	LYS	3.7
1	I	172	LYS	3.6
1	E	172	LYS	3.3
1	D	172	LYS	3.2
1	F	172	LYS	3.1
1	W	172	LYS	3.1
1	T	1	VAL	3.1
1	G	172	LYS	3.0
1	V	172	LYS	2.9
1	D	171	VAL	2.9
1	S	172	LYS	2.8
1	N	1	VAL	2.8
1	X	1	VAL	2.8
1	N	171	VAL	2.7
1	K	172	LYS	2.7
1	Q	172	LYS	2.6
1	S	1	VAL	2.6
1	M	171	VAL	2.5
1	N	172	LYS	2.5
1	X	172	LYS	2.5
1	P	172	LYS	2.4
1	F	1	VAL	2.4
1	F	9[A]	HIS	2.4
1	R	172	LYS	2.4
1	H	60	GLU	2.3
1	D	34	TYR	2.3
1	R	56	HIS	2.2
1	X	171	VAL	2.2
1	L	59	ARG	2.2
1	M	1	VAL	2.1
1	W	56[A]	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	P	171	VAL	2.1
1	H	171	VAL	2.0
1	R	157	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	T	201	5/5	0.67	0.21	58,59,61,62	0
3	EDO	R	204	4/4	0.70	0.20	28,29,29,30	0
3	EDO	Q	202	4/4	0.71	0.23	22,22,24,26	0
3	EDO	G	208	4/4	0.75	0.21	29,32,32,34	0
5	CL	G	204	1/1	0.75	0.18	46,46,46,46	0
3	EDO	I	203	4/4	0.78	0.22	20,22,25,27	0
3	EDO	P	202	4/4	0.79	0.17	17,17,18,21	0
3	EDO	B	204	4/4	0.79	0.19	20,20,20,23	0
3	EDO	E	204	4/4	0.80	0.17	25,28,28,29	0
3	EDO	Q	203	4/4	0.80	0.18	23,24,24,25	0
3	EDO	C	205	4/4	0.80	0.18	23,24,24,27	0
4	ACT	A	203	4/4	0.80	0.17	17,18,19,20	0
3	EDO	B	203	4/4	0.80	0.17	27,28,28,32	0
3	EDO	N	205	4/4	0.81	0.19	26,28,28,30	0
3	EDO	S	203	4/4	0.81	0.18	18,21,22,25	0
3	EDO	S	204	4/4	0.81	0.20	30,32,34,39	0
3	EDO	K	203	4/4	0.81	0.22	25,25,26,26	0
3	EDO	K	204	4/4	0.81	0.17	23,24,24,24	0
3	EDO	I	204	4/4	0.82	0.16	20,20,20,21	0
3	EDO	A	202	4/4	0.82	0.17	27,27,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	G	207	4/4	0.83	0.19	27,27,27,27	0
3	EDO	G	206	4/4	0.84	0.16	31,33,34,35	0
3	EDO	R	205	4/4	0.84	0.17	29,29,30,30	0
3	EDO	V	202	4/4	0.84	0.14	20,20,21,21	0
3	EDO	J	205	4/4	0.84	0.17	26,26,27,27	0
3	EDO	F	205	4/4	0.84	0.15	24,25,25,25	0
3	EDO	N	204	4/4	0.85	0.12	21,21,21,21	0
3	EDO	P	203	4/4	0.85	0.14	22,22,23,24	0
2	SO4	A	201	5/5	0.85	0.14	45,45,47,47	0
5	CL	X	201	1/1	0.85	0.17	45,45,45,45	0
3	EDO	X	204	4/4	0.86	0.14	19,20,20,21	0
3	EDO	B	205	4/4	0.86	0.15	23,25,26,26	0
3	EDO	L	205	4/4	0.87	0.16	17,17,18,20	0
3	EDO	W	204	4/4	0.87	0.19	21,23,24,25	0
3	EDO	S	202	4/4	0.87	0.17	16,19,19,23	0
3	EDO	X	205	4/4	0.87	0.13	29,30,31,31	0
3	EDO	C	204	4/4	0.87	0.12	14,14,15,15	0
5	CL	D	203	1/1	0.87	0.12	36,36,36,36	0
5	CL	E	202	1/1	0.87	0.13	39,39,39,39	0
3	EDO	G	210	4/4	0.87	0.13	14,14,14,15	0
2	SO4	N	202	5/5	0.87	0.17	45,45,47,48	0
3	EDO	L	204	4/4	0.88	0.12	11,12,12,12	0
3	EDO	N	203	4/4	0.88	0.14	17,19,20,21	0
3	EDO	T	205	4/4	0.88	0.17	23,24,24,25	0
3	EDO	G	211	4/4	0.88	0.22	24,25,25,26	0
3	EDO	P	204	4/4	0.88	0.15	25,25,26,27	0
4	ACT	Q	204	4/4	0.89	0.12	24,24,25,25	0
3	EDO	E	205	4/4	0.89	0.13	25,26,27,27	0
3	EDO	F	204	4/4	0.89	0.11	24,25,25,25	0
5	CL	F	203	1/1	0.89	0.12	40,40,40,40	0
2	SO4	G	201	5/5	0.89	0.18	40,40,41,41	0
2	SO4	T	202	5/5	0.89	0.17	31,34,34,35	0
3	EDO	R	206	4/4	0.90	0.17	16,17,18,19	0
3	EDO	J	204	4/4	0.90	0.18	14,15,15,15	0
2	SO4	C	201	5/5	0.90	0.16	36,37,39,39	0
2	SO4	I	201	5/5	0.90	0.17	42,42,43,44	0
3	EDO	D	205	4/4	0.90	0.18	22,22,23,25	0
5	CL	Q	201	1/1	0.90	0.14	34,34,34,34	0
5	CL	U	203	1/1	0.90	0.14	38,38,38,38	0
2	SO4	F	201	5/5	0.90	0.15	30,31,32,32	0
5	CL	F	202	1/1	0.91	0.13	35,35,35,35	0
3	EDO	G	209	4/4	0.91	0.11	28,30,30,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	K	202	5/5	0.91	0.14	38,39,40,40	0
2	SO4	L	201	5/5	0.91	0.19	14,15,16,16	5
5	CL	M	201	1/1	0.91	0.12	32,32,32,32	0
2	SO4	W	202	5/5	0.91	0.15	34,34,36,37	0
2	SO4	H	201	5/5	0.92	0.13	39,39,41,41	0
5	CL	O	203	1/1	0.92	0.17	38,38,38,38	0
3	EDO	G	205	4/4	0.92	0.09	7,7,7,8	0
2	SO4	R	201	5/5	0.92	0.21	15,16,17,17	5
2	SO4	R	202	5/5	0.92	0.12	39,40,41,42	0
2	SO4	O	201	5/5	0.92	0.14	38,38,40,40	0
2	SO4	K	201	5/5	0.92	0.18	13,14,14,14	5
2	SO4	P	201	5/5	0.92	0.13	38,41,41,42	0
5	CL	D	204	1/1	0.92	0.15	38,38,38,38	0
5	CL	X	203	1/1	0.92	0.16	37,37,37,37	0
3	EDO	C	203	4/4	0.93	0.09	10,11,11,12	0
2	SO4	W	201	5/5	0.93	0.15	22,25,26,26	0
3	EDO	U	204	4/4	0.93	0.12	25,26,26,26	0
5	CL	C	202	1/1	0.93	0.12	31,31,31,31	0
3	EDO	V	201	4/4	0.93	0.12	17,18,19,19	0
2	SO4	J	202	5/5	0.93	0.12	35,36,38,38	0
3	EDO	R	203	4/4	0.93	0.10	12,13,13,13	0
2	SO4	D	202	5/5	0.93	0.11	32,32,33,33	0
3	EDO	M	202	4/4	0.94	0.09	9,10,10,11	0
2	SO4	H	202	5/5	0.94	0.15	32,33,34,35	0
5	CL	E	203	1/1	0.94	0.12	31,31,31,31	0
2	SO4	U	201	5/5	0.94	0.14	30,32,32,32	0
5	CL	O	202	1/1	0.94	0.09	36,36,36,36	0
5	CL	L	203	1/1	0.95	0.21	38,38,38,38	0
2	SO4	J	201	5/5	0.95	0.14	22,24,25,26	0
2	SO4	D	201	5/5	0.95	0.18	16,18,19,19	5
5	CL	T	204	1/1	0.95	0.18	40,40,40,40	0
5	CL	X	202	1/1	0.95	0.13	32,32,32,32	0
2	SO4	N	201	5/5	0.95	0.12	31,31,31,31	0
5	CL	I	202	1/1	0.96	0.13	31,31,31,31	0
5	CL	L	202	1/1	0.96	0.15	32,32,32,32	0
5	CL	E	201	1/1	0.96	0.17	29,29,29,29	1
5	CL	H	203	1/1	0.96	0.09	38,38,38,38	0
3	EDO	O	204	4/4	0.97	0.05	11,11,11,11	0
5	CL	U	202	1/1	0.97	0.11	27,27,27,27	0
5	CL	B	201	1/1	0.97	0.11	29,29,29,29	0
5	CL	S	201	1/1	0.97	0.15	25,25,25,25	0
5	CL	G	202	1/1	0.97	0.14	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	CL	G	203	1/1	0.97	0.12	24,24,24,24	0
5	CL	W	203	1/1	0.98	0.16	24,24,24,24	0
5	CL	J	203	1/1	0.98	0.16	30,30,30,30	0
5	CL	T	203	1/1	0.99	0.14	23,23,23,23	0
5	CL	B	202	1/1	0.99	0.18	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.