



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 16, 2023 – 05:12 AM EDT

PDB ID : 1YAR
Title : Structure of Archeobacterial 20S proteasome mutant D9S- PA26 complex
Authors : Forster, A.; Masters, E.I.; Whitby, F.G.; Robinson, H.; Hill, C.P.
Deposited on : 2004-12-17
Resolution : 1.90 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

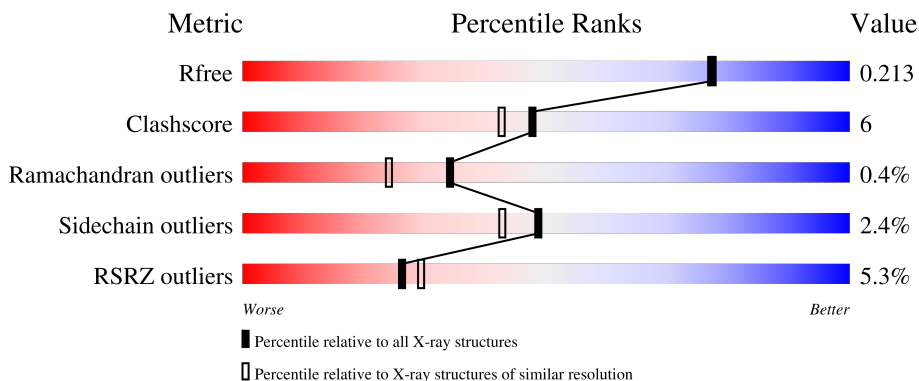
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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.



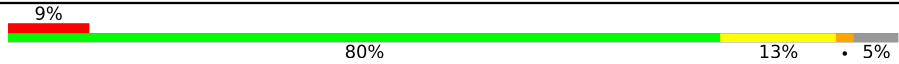
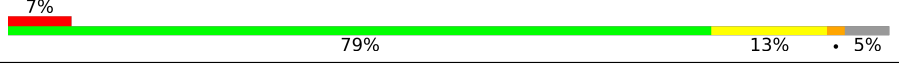
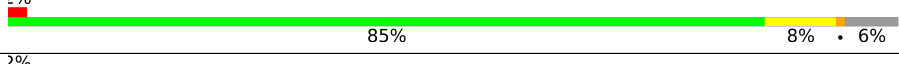


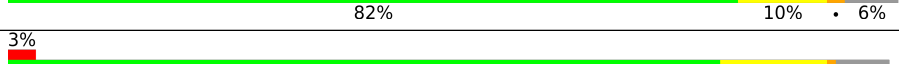
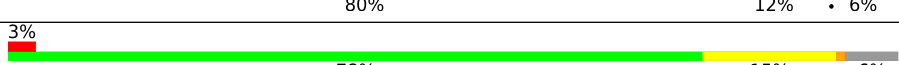
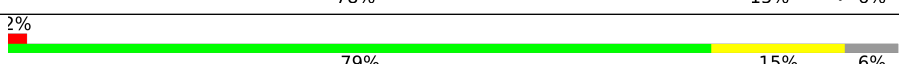
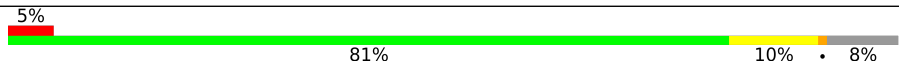


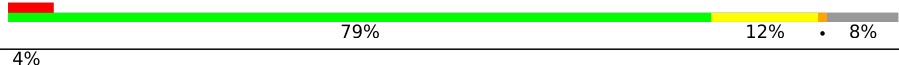
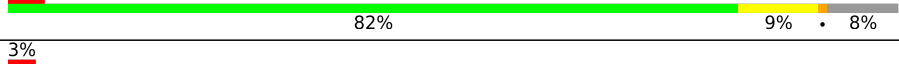

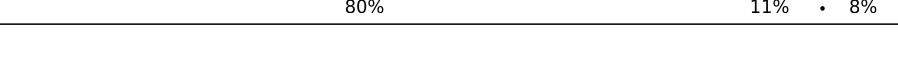

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	233	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 80%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 13%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">80% 13% • 5%</p>
1	B	233	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 85%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">85% 9% • 5%</p>
1	C	233	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 82%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">82% 12% 5%</p>
1	D	233	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 84%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">84% 10% • 5%</p>
1	E	233	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 83%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">83% 10% • 5%</p>

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Mol	Chain	Length	Quality of chain
1	F	233	
1	G	233	
2	H	217	
2	I	217	
2	J	217	
2	K	217	
2	L	217	
2	M	217	
2	N	217	
3	O	237	
3	P	237	
3	Q	237	
3	R	237	
3	S	237	
3	T	237	
3	U	237	

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
4	SO4	G	4007	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 39593 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 Proteasome alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	221	1777	1127	302	345	3	0	7	0
1	B	221	1777	1127	302	345	3	0	7	0
1	C	221	1771	1124	301	343	3	0	6	0
1	D	221	1777	1127	302	345	3	0	7	0
1	E	221	1771	1124	301	343	3	0	6	0
1	F	221	1771	1124	301	343	3	0	6	0
1	G	221	1771	1124	301	343	3	0	6	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	SER	ASP	engineered mutation	UNP P25156
B	9	SER	ASP	engineered mutation	UNP P25156
C	9	SER	ASP	engineered mutation	UNP P25156
D	9	SER	ASP	engineered mutation	UNP P25156
E	9	SER	ASP	engineered mutation	UNP P25156
F	9	SER	ASP	engineered mutation	UNP P25156
G	9	SER	ASP	engineered mutation	UNP P25156

- Molecule 2 is a protein called Proteasome beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	203	1613	1019	271	311	12	0	7	0
2	I	203	1613	1019	271	311	12	0	7	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	J	203	1613	1019	271	311	12	0	7	0
2	K	203	1613	1019	271	311	12	0	7	0
2	L	203	1613	1019	271	311	12	0	7	0
2	M	203	1613	1019	271	311	12	0	7	0
2	N	203	1613	1019	271	311	12	0	7	0

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	204	HIS	-	expression tag	UNP P28061
H	205	HIS	-	expression tag	UNP P28061
H	206	HIS	-	expression tag	UNP P28061
H	207	HIS	-	expression tag	UNP P28061
H	208	HIS	-	expression tag	UNP P28061
H	209	HIS	-	expression tag	UNP P28061
I	204	HIS	-	expression tag	UNP P28061
I	205	HIS	-	expression tag	UNP P28061
I	206	HIS	-	expression tag	UNP P28061
I	207	HIS	-	expression tag	UNP P28061
I	208	HIS	-	expression tag	UNP P28061
I	209	HIS	-	expression tag	UNP P28061
J	204	HIS	-	expression tag	UNP P28061
J	205	HIS	-	expression tag	UNP P28061
J	206	HIS	-	expression tag	UNP P28061
J	207	HIS	-	expression tag	UNP P28061
J	208	HIS	-	expression tag	UNP P28061
J	209	HIS	-	expression tag	UNP P28061
K	204	HIS	-	expression tag	UNP P28061
K	205	HIS	-	expression tag	UNP P28061
K	206	HIS	-	expression tag	UNP P28061
K	207	HIS	-	expression tag	UNP P28061
K	208	HIS	-	expression tag	UNP P28061
K	209	HIS	-	expression tag	UNP P28061
L	204	HIS	-	expression tag	UNP P28061
L	205	HIS	-	expression tag	UNP P28061
L	206	HIS	-	expression tag	UNP P28061
L	207	HIS	-	expression tag	UNP P28061
L	208	HIS	-	expression tag	UNP P28061

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Chain	Residue	Modelled	Actual	Comment	Reference
L	209	HIS	-	expression tag	UNP P28061
M	204	HIS	-	expression tag	UNP P28061
M	205	HIS	-	expression tag	UNP P28061
M	206	HIS	-	expression tag	UNP P28061
M	207	HIS	-	expression tag	UNP P28061
M	208	HIS	-	expression tag	UNP P28061
M	209	HIS	-	expression tag	UNP P28061
N	204	HIS	-	expression tag	UNP P28061
N	205	HIS	-	expression tag	UNP P28061
N	206	HIS	-	expression tag	UNP P28061
N	207	HIS	-	expression tag	UNP P28061
N	208	HIS	-	expression tag	UNP P28061
N	209	HIS	-	expression tag	UNP P28061

- Molecule 3 is a protein called proteasome activator protein PA26.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	O	218	1754	1094	309	343	8	0	10	0
3	P	218	1754	1094	309	343	8	0	10	0
3	Q	218	1754	1094	309	343	8	0	10	0
3	R	218	1754	1094	309	343	8	0	10	0
3	S	218	1754	1094	309	343	8	0	10	0
3	T	218	1746	1089	308	342	7	0	9	0
3	U	218	1754	1094	309	343	8	0	10	0

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
O	-5	MET	-	initiating methionine	GB 5757773
O	-4	HIS	-	expression tag	GB 5757773
O	-3	HIS	-	expression tag	GB 5757773
O	-2	HIS	-	expression tag	GB 5757773
O	-1	HIS	-	expression tag	GB 5757773
O	0	HIS	-	expression tag	GB 5757773
O	1	HIS	-	expression tag	GB 5757773

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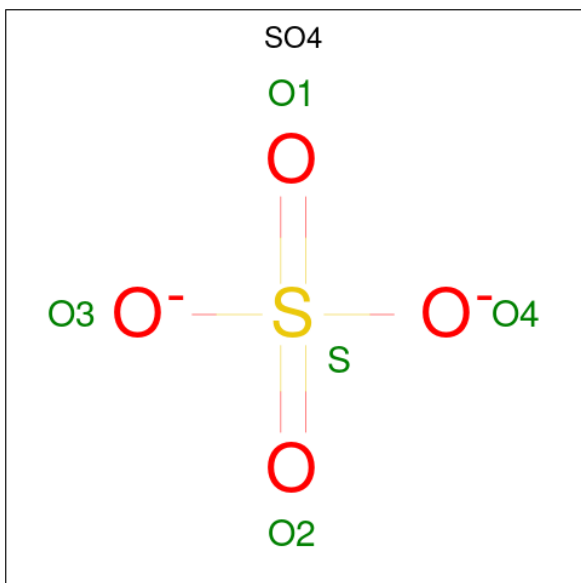
Chain	Residue	Modelled	Actual	Comment	Reference
O	49	VAL	THR	variant	GB 5757773
P	-5	MET	-	initiating methionine	GB 5757773
P	-4	HIS	-	expression tag	GB 5757773
P	-3	HIS	-	expression tag	GB 5757773
P	-2	HIS	-	expression tag	GB 5757773
P	-1	HIS	-	expression tag	GB 5757773
P	0	HIS	-	expression tag	GB 5757773
P	1	HIS	-	expression tag	GB 5757773
P	49	VAL	THR	variant	GB 5757773
Q	-5	MET	-	initiating methionine	GB 5757773
Q	-4	HIS	-	expression tag	GB 5757773
Q	-3	HIS	-	expression tag	GB 5757773
Q	-2	HIS	-	expression tag	GB 5757773
Q	-1	HIS	-	expression tag	GB 5757773
Q	0	HIS	-	expression tag	GB 5757773
Q	1	HIS	-	expression tag	GB 5757773
Q	49	VAL	THR	variant	GB 5757773
R	-5	MET	-	initiating methionine	GB 5757773
R	-4	HIS	-	expression tag	GB 5757773
R	-3	HIS	-	expression tag	GB 5757773
R	-2	HIS	-	expression tag	GB 5757773
R	-1	HIS	-	expression tag	GB 5757773
R	0	HIS	-	expression tag	GB 5757773
R	1	HIS	-	expression tag	GB 5757773
R	49	VAL	THR	variant	GB 5757773
S	-5	MET	-	initiating methionine	GB 5757773
S	-4	HIS	-	expression tag	GB 5757773
S	-3	HIS	-	expression tag	GB 5757773
S	-2	HIS	-	expression tag	GB 5757773
S	-1	HIS	-	expression tag	GB 5757773
S	0	HIS	-	expression tag	GB 5757773
S	1	HIS	-	expression tag	GB 5757773
S	49	VAL	THR	variant	GB 5757773
T	-5	MET	-	initiating methionine	GB 5757773
T	-4	HIS	-	expression tag	GB 5757773
T	-3	HIS	-	expression tag	GB 5757773
T	-2	HIS	-	expression tag	GB 5757773
T	-1	HIS	-	expression tag	GB 5757773
T	0	HIS	-	expression tag	GB 5757773
T	1	HIS	-	expression tag	GB 5757773
T	49	VAL	THR	variant	GB 5757773
U	-5	MET	-	initiating methionine	GB 5757773

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Chain	Residue	Modelled	Actual	Comment	Reference
U	-4	HIS	-	expression tag	GB 5757773
U	-3	HIS	-	expression tag	GB 5757773
U	-2	HIS	-	expression tag	GB 5757773
U	-1	HIS	-	expression tag	GB 5757773
U	0	HIS	-	expression tag	GB 5757773
U	1	HIS	-	expression tag	GB 5757773
U	49	VAL	THR	variant	GB 5757773

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	H	1	Total C O 6 3 3	0	0
5	I	1	Total C O 6 3 3	0	0
5	J	1	Total C O 6 3 3	0	0
5	K	1	Total C O 6 3 3	0	0
5	L	1	Total C O 6 3 3	0	0
5	M	1	Total C O 6 3 3	0	0
5	N	1	Total C O 6 3 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	153	Total O 153 153	0	0
6	B	145	Total O 145 145	0	0
6	C	128	Total O 128 128	0	0
6	D	130	Total O 130 130	0	0
6	E	153	Total O 153 153	0	0

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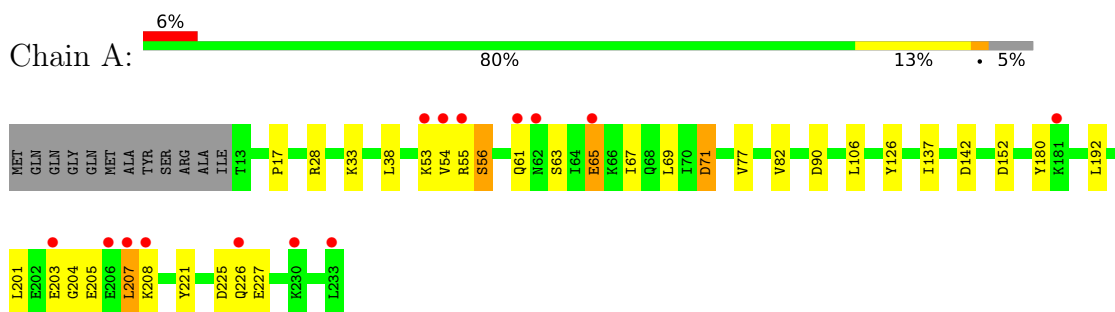
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	F	122	Total 122	O 122	0	0
6	G	133	Total 133	O 133	0	0
6	H	165	Total 165	O 165	0	0
6	I	165	Total 165	O 165	0	0
6	J	159	Total 159	O 159	0	0
6	K	177	Total 177	O 177	0	0
6	L	153	Total 153	O 153	0	0
6	M	180	Total 180	O 180	0	0
6	N	179	Total 179	O 179	0	0
6	O	178	Total 178	O 178	0	0
6	P	162	Total 162	O 162	0	0
6	Q	161	Total 161	O 161	0	0
6	R	189	Total 189	O 189	0	0
6	S	221	Total 221	O 221	0	0
6	T	207	Total 207	O 207	0	0
6	U	210	Total 210	O 210	0	0

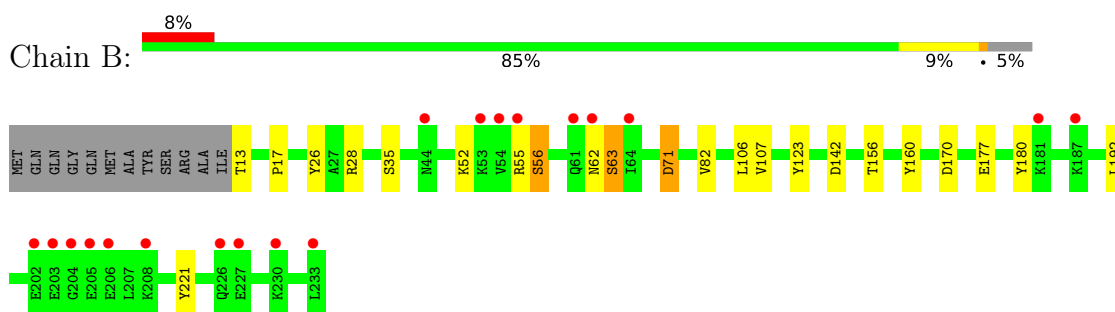
3 Residue-property plots [i](#)

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.

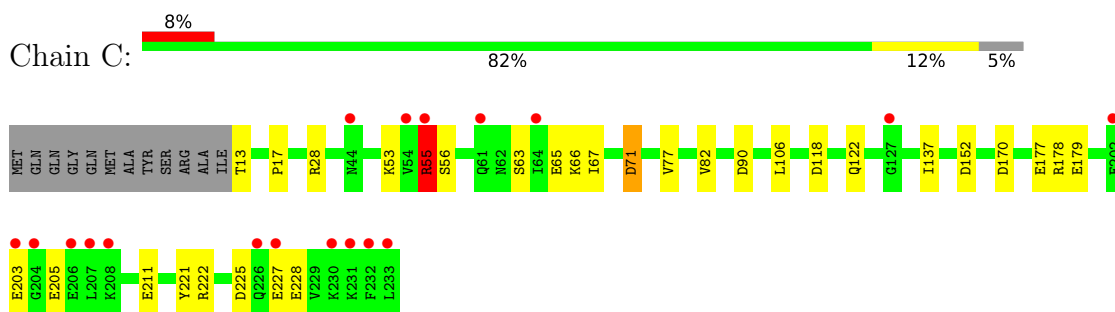
- Molecule 1: Proteasome alpha subunit



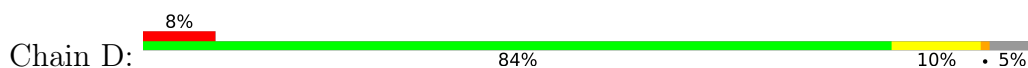
- Molecule 1: Proteasome alpha subunit

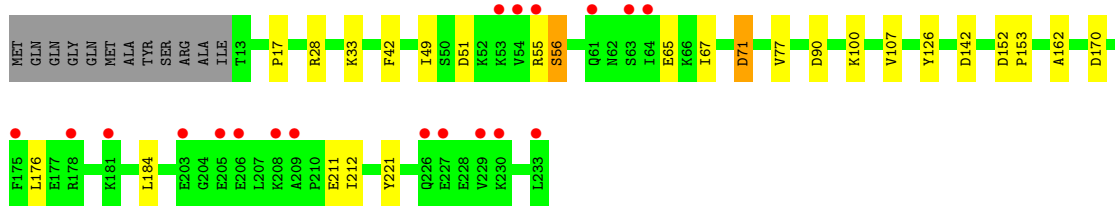


- Molecule 1: Proteasome alpha subunit

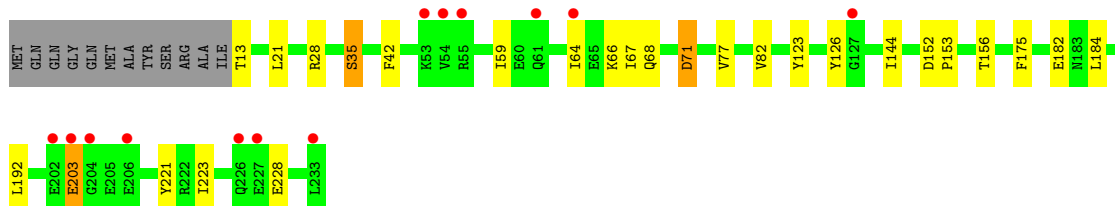
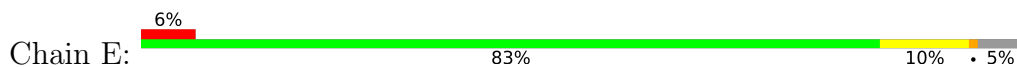


- Molecule 1: Proteasome alpha subunit

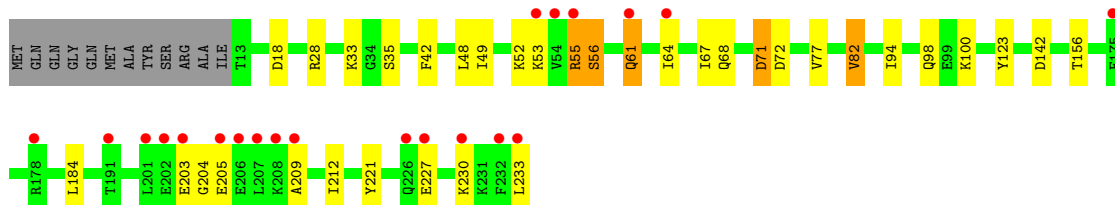
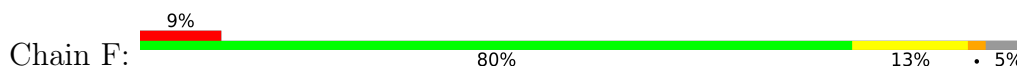




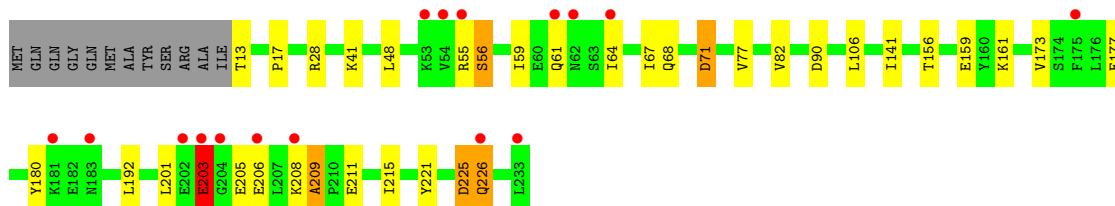
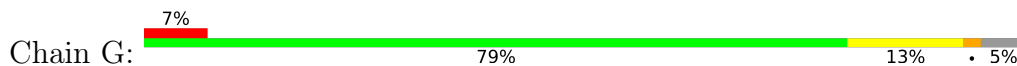
● Molecule 1: Proteasome alpha subunit



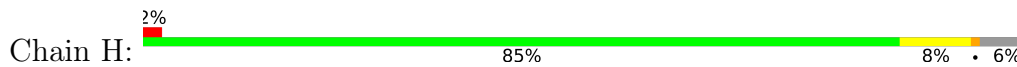
● Molecule 1: Proteasome alpha subunit



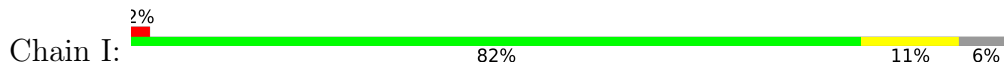
● Molecule 1: Proteasome alpha subunit



● Molecule 2: Proteasome beta subunit

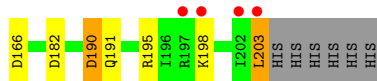
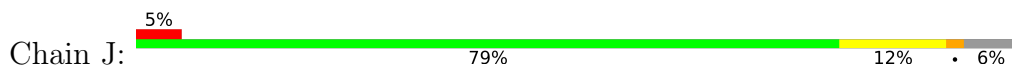


● Molecule 2: Proteasome beta subunit

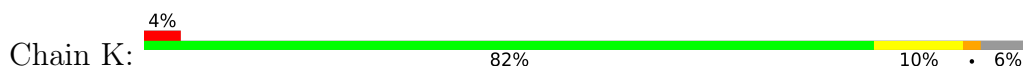




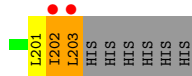
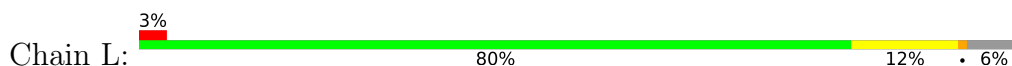
• Molecule 2: Proteasome beta subunit



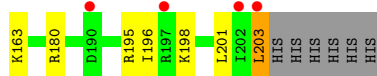
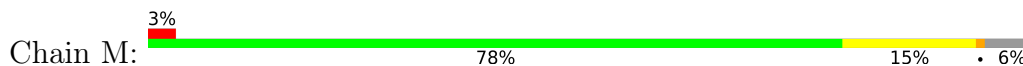
• Molecule 2: Proteasome beta subunit



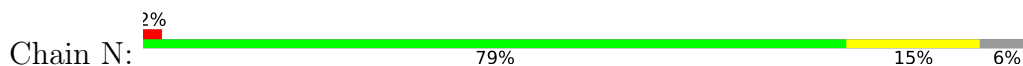
• Molecule 2: Proteasome beta subunit

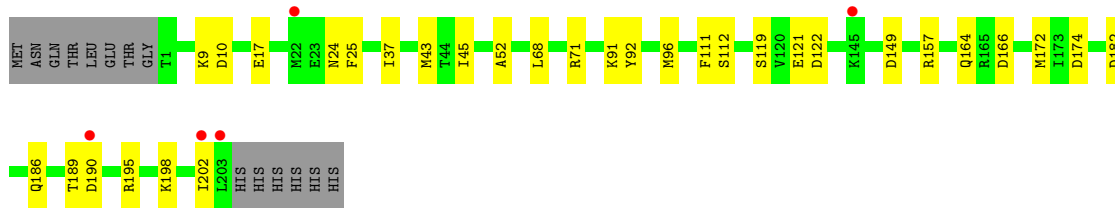


• Molecule 2: Proteasome beta subunit

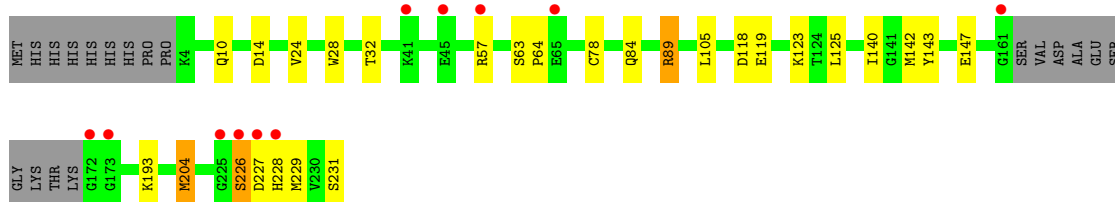
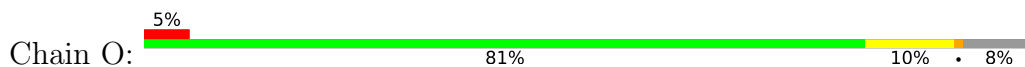


• Molecule 2: Proteasome beta subunit

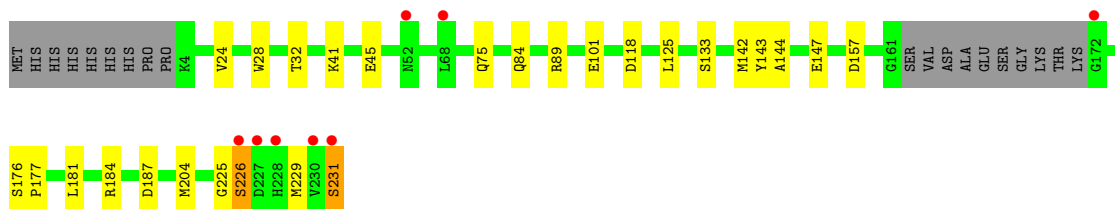
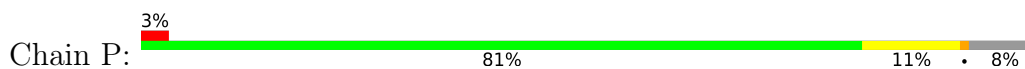




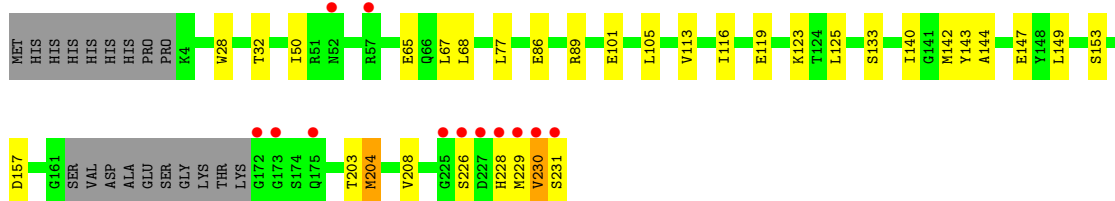
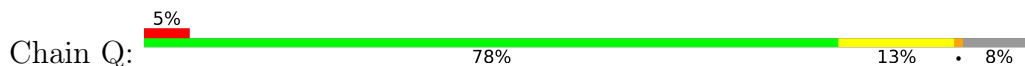
• Molecule 3: proteasome activator protein PA26



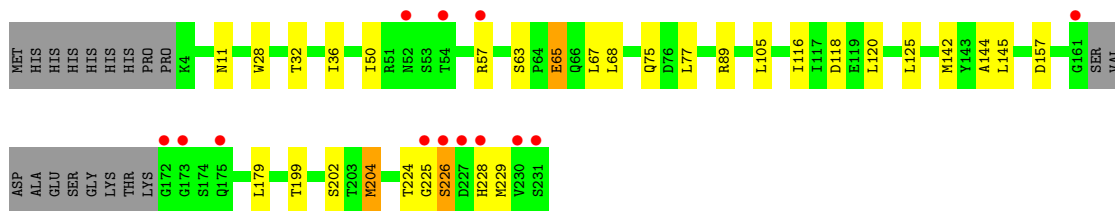
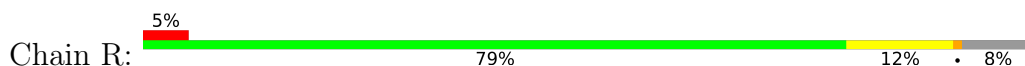
• Molecule 3: proteasome activator protein PA26



• Molecule 3: proteasome activator protein PA26

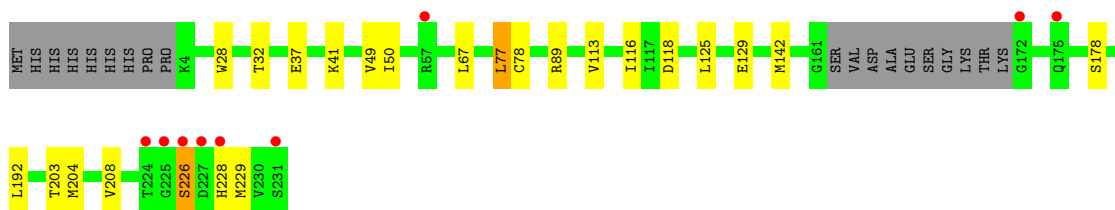


• Molecule 3: proteasome activator protein PA26



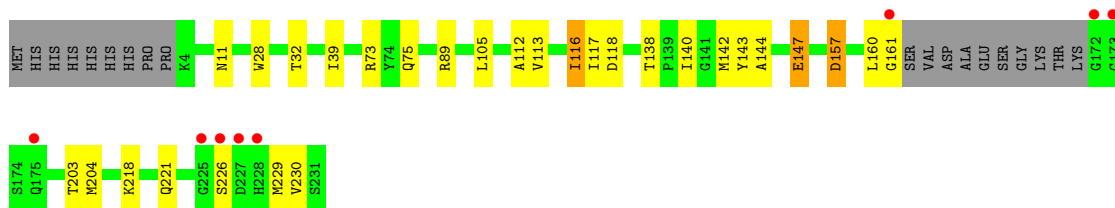
- Molecule 3: proteasome activator protein PA26

Chain S: 4% 82% 9% 8%



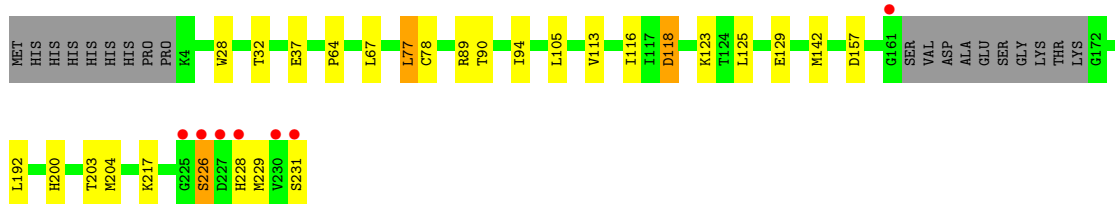
- Molecule 3: proteasome activator protein PA26

Chain T: 3% 80% 11% 8%



- Molecule 3: proteasome activator protein PA26

Chain U: 3% 80% 11% 8%



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	254.87Å 127.51Å 181.18Å 90.00° 92.45° 90.00°	Depositor
Resolution (Å)	47.30 – 1.90 47.30 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.2 (47.30-1.90) 97.2 (47.30-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 1.90Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.182 , 0.216 0.182 , 0.213	Depositor DCC
R_{free} test set	1345 reflections (0.30%)	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtrriage
Anisotropy	0.020	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	0.002 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	39593	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, 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	0.57	0/1800	0.80	4/2425 (0.2%)
1	B	0.57	0/1800	0.75	2/2425 (0.1%)
1	C	0.57	0/1794	0.74	3/2417 (0.1%)
1	D	0.57	0/1800	0.75	3/2425 (0.1%)
1	E	0.59	0/1794	0.76	1/2417 (0.0%)
1	F	0.55	0/1794	0.74	4/2417 (0.2%)
1	G	0.57	0/1794	0.75	2/2417 (0.1%)
2	H	0.64	0/1633	0.83	4/2206 (0.2%)
2	I	0.61	0/1633	0.80	4/2206 (0.2%)
2	J	0.59	0/1633	0.83	6/2206 (0.3%)
2	K	0.59	0/1633	0.83	6/2206 (0.3%)
2	L	0.59	0/1633	0.82	3/2206 (0.1%)
2	M	0.65	0/1633	0.84	2/2206 (0.1%)
2	N	0.61	0/1633	0.84	7/2206 (0.3%)
3	O	0.51	0/1777	0.70	0/2398
3	P	0.50	0/1777	0.68	2/2398 (0.1%)
3	Q	0.50	0/1777	0.69	1/2398 (0.0%)
3	R	0.56	0/1777	0.69	1/2398 (0.0%)
3	S	0.62	0/1777	0.70	0/2398
3	T	0.61	0/1769	0.70	1/2388 (0.0%)
3	U	0.59	0/1777	0.71	2/2398 (0.1%)
All	All	0.58	0/36438	0.76	58/49161 (0.1%)

There are no bond length outliers.

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	71	ASP	CB-CG-OD2	8.91	126.32	118.30
1	D	71	ASP	CB-CG-OD2	8.29	125.76	118.30
1	G	71	ASP	CB-CG-OD2	8.27	125.74	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	71	ASP	CB-CG-OD2	8.18	125.66	118.30
2	N	149	ASP	CB-CG-OD2	7.37	124.93	118.30
1	B	71	ASP	CB-CG-OD2	7.20	124.78	118.30
1	E	71	ASP	CB-CG-OD2	7.12	124.70	118.30
1	F	71	ASP	CB-CG-OD2	6.76	124.38	118.30
2	K	149	ASP	CB-CG-OD2	6.53	124.17	118.30
2	H	114	ASP	CB-CG-OD2	6.44	124.10	118.30
1	A	225	ASP	CB-CG-OD2	6.43	124.09	118.30
2	K	10	ASP	CB-CG-OD2	6.21	123.89	118.30
1	A	142	ASP	CB-CG-OD2	6.13	123.82	118.30
2	N	166	ASP	CB-CG-OD2	5.93	123.63	118.30
2	J	182	ASP	CB-CG-OD2	5.87	123.58	118.30
2	I	10	ASP	CB-CG-OD2	5.85	123.57	118.30
2	H	182	ASP	CB-CG-OD2	5.84	123.56	118.30
2	I	182	ASP	CB-CG-OD2	5.83	123.54	118.30
2	L	114	ASP	CB-CG-OD2	5.79	123.51	118.30
2	M	122	ASP	CB-CG-OD2	5.72	123.45	118.30
3	U	118	ASP	CB-CG-OD2	5.68	123.41	118.30
2	N	182	ASP	CB-CG-OD2	5.62	123.36	118.30
2	N	10	ASP	CB-CG-OD2	5.59	123.33	118.30
1	C	170	ASP	CB-CG-OD2	5.58	123.32	118.30
2	J	114	ASP	CB-CG-OD2	5.58	123.32	118.30
3	Q	157	ASP	CB-CG-OD2	5.56	123.31	118.30
3	P	187	ASP	CB-CG-OD2	5.54	123.28	118.30
3	U	157	ASP	CB-CG-OD2	5.52	123.27	118.30
2	H	190[A]	ASP	CB-CG-OD2	5.48	123.23	118.30
2	H	190[B]	ASP	CB-CG-OD2	5.48	123.23	118.30
2	K	166	ASP	CB-CG-OD2	5.43	123.19	118.30
1	C	225	ASP	CB-CG-OD2	5.43	123.19	118.30
2	J	166	ASP	CB-CG-OD2	5.40	123.16	118.30
1	D	51	ASP	CB-CG-OD2	5.40	123.16	118.30
1	B	170	ASP	CB-CG-OD2	5.37	123.13	118.30
1	D	170	ASP	CB-CG-OD2	5.35	123.11	118.30
2	J	190[A]	ASP	CB-CG-OD2	5.33	123.10	118.30
2	J	190[B]	ASP	CB-CG-OD2	5.33	123.10	118.30
1	F	72	ASP	CB-CG-OD2	5.30	123.07	118.30
2	I	190[A]	ASP	CB-CG-OD2	5.29	123.06	118.30
2	I	190[B]	ASP	CB-CG-OD2	5.29	123.06	118.30
3	T	157	ASP	CB-CG-OD2	5.29	123.06	118.30
1	F	142	ASP	CB-CG-OD2	5.28	123.05	118.30
1	G	225	ASP	CB-CG-OD2	5.25	123.03	118.30
2	J	10	ASP	CB-CG-OD2	5.21	122.99	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	114	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	207	LEU	CA-CB-CG	-5.16	103.42	115.30
2	L	166	ASP	CB-CG-OD2	5.14	122.92	118.30
2	N	190[A]	ASP	CB-CG-OD2	5.10	122.89	118.30
2	N	190[B]	ASP	CB-CG-OD2	5.10	122.89	118.30
3	P	157	ASP	CB-CG-OD2	5.09	122.88	118.30
2	K	182	ASP	CB-CG-OD2	5.08	122.88	118.30
3	R	157	ASP	CB-CG-OD2	5.07	122.87	118.30
2	N	122	ASP	CB-CG-OD2	5.05	122.84	118.30
2	L	10	ASP	CB-CG-OD2	5.03	122.82	118.30
1	F	18	ASP	CB-CG-OD2	5.03	122.82	118.30
2	K	190[A]	ASP	CB-CG-OD2	5.02	122.82	118.30
2	K	190[B]	ASP	CB-CG-OD2	5.02	122.82	118.30

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	1777	0	1809	34	0
1	B	1777	0	1809	21	0
1	C	1771	0	1805	28	0
1	D	1777	0	1809	25	0
1	E	1771	0	1805	27	0
1	F	1771	0	1805	28	0
1	G	1771	0	1805	38	0
2	H	1613	0	1649	14	0
2	I	1613	0	1649	18	0
2	J	1613	0	1649	17	0
2	K	1613	0	1649	21	0
2	L	1613	0	1649	23	0
2	M	1613	0	1649	23	0
2	N	1613	0	1649	16	0
3	O	1754	0	1765	20	0
3	P	1754	0	1765	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	Q	1754	0	1765	30	0
3	R	1754	0	1765	34	0
3	S	1754	0	1765	39	0
3	T	1746	0	1757	30	0
3	U	1754	0	1765	33	0
4	A	5	0	0	0	0
4	B	10	0	0	0	0
4	C	5	0	0	0	0
4	D	5	0	0	0	0
4	F	10	0	0	0	0
4	G	5	0	0	2	0
4	H	10	0	0	1	0
4	I	5	0	0	0	0
4	J	10	0	0	0	0
4	K	10	0	0	0	0
4	L	10	0	0	0	0
4	M	10	0	0	1	0
4	N	10	0	0	0	0
5	H	6	0	8	0	0
5	I	6	0	8	0	0
5	J	6	0	8	0	0
5	K	6	0	8	0	0
5	L	6	0	8	0	0
5	M	6	0	8	0	0
5	N	6	0	8	0	0
6	A	153	0	0	5	0
6	B	145	0	0	4	0
6	C	128	0	0	8	0
6	D	130	0	0	1	0
6	E	153	0	0	6	0
6	F	122	0	0	3	0
6	G	133	0	0	9	0
6	H	165	0	0	5	0
6	I	165	0	0	7	0
6	J	159	0	0	5	0
6	K	177	0	0	4	0
6	L	153	0	0	4	0
6	M	180	0	0	8	0
6	N	179	0	0	6	0
6	O	178	0	0	4	0
6	P	162	0	0	3	0
6	Q	161	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	R	189	0	0	6	0
6	S	221	0	0	1	0
6	T	207	0	0	6	0
6	U	210	0	0	3	0
All	All	39593	0	36593	445	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (445) 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:28[B]:ARG:HH22	3:U:229:MET:CG	1.49	1.23
2:K:96[B]:MET:HE3	2:L:91:LYS:HZ1	1.02	1.10
1:E:28[B]:ARG:HH22	3:O:229:MET:HG3	0.98	1.09
1:F:28[B]:ARG:HH22	3:P:229:MET:HG2	1.10	1.09
2:L:202:ILE:HG22	2:L:203:LEU:H	1.05	1.08
1:C:28[A]:ARG:HH22	3:T:229:MET:CG	1.66	1.08
3:R:89[A]:ARG:HD3	3:S:203:THR:HG21	1.11	1.08
1:D:28[B]:ARG:HH22	3:U:229:MET:HG3	1.14	1.07
2:K:96[B]:MET:HE3	2:L:91:LYS:NZ	1.71	1.05
3:P:89[A]:ARG:HD3	3:Q:203:THR:HG21	1.10	1.05
1:E:28[B]:ARG:NH2	3:O:229:MET:HG3	1.72	1.05
3:S:89[A]:ARG:HD3	3:T:203:THR:HG21	1.07	1.04
1:A:65:GLU:HB3	4:G:4007:SO4:O1	1.57	1.04
2:L:96[B]:MET:HE2	6:M:5041:HOH:O	1.56	1.03
3:P:89[A]:ARG:HD3	3:Q:203:THR:CG2	1.88	1.03
1:B:28[B]:ARG:HH22	3:S:229:MET:CG	1.71	1.03
1:G:28[B]:ARG:HH22	3:Q:229:MET:HG2	1.19	1.02
1:C:28[A]:ARG:HH22	3:T:229:MET:HG3	1.23	1.00
1:F:28[B]:ARG:NH2	3:P:229:MET:HG2	1.79	0.97
1:E:28[B]:ARG:HH22	3:O:229:MET:CG	1.76	0.97
2:L:202:ILE:HG22	2:L:203:LEU:N	1.77	0.97
3:P:89[A]:ARG:CD	3:Q:203:THR:HG21	1.95	0.96
3:Q:101:GLU:HG3	3:R:105:LEU:HD22	1.47	0.95
1:G:90:ASP:HB3	6:M:5136:HOH:O	1.66	0.95
3:S:116[B]:ILE:HD11	3:S:204[B]:MET:HE1	1.47	0.94
3:T:89[A]:ARG:NH1	3:T:118:ASP:OD1	2.01	0.94
1:D:28[B]:ARG:NH2	3:U:229:MET:HG3	1.82	0.94
3:S:89[A]:ARG:HD3	3:T:203:THR:CG2	1.99	0.93
3:S:89[A]:ARG:CD	3:T:203:THR:HG21	1.97	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:U:37[A]:GLU:OE1	6:U:344:HOH:O	1.86	0.93
1:F:68[B]:GLN:NE2	6:F:4108:HOH:O	2.02	0.92
3:S:116[B]:ILE:HD11	3:S:204[B]:MET:CE	1.99	0.92
1:D:28[B]:ARG:HH22	3:U:229:MET:HG2	1.33	0.91
3:T:89[A]:ARG:HD3	3:U:203:THR:HG21	1.52	0.91
3:T:11:ASN:OD1	6:T:310:HOH:O	1.87	0.91
1:D:28[B]:ARG:NH2	3:U:229:MET:CG	2.34	0.91
2:I:67:ARG:NH1	6:I:5168:HOH:O	2.02	0.91
1:F:28[B]:ARG:HH22	3:P:229:MET:CG	1.82	0.90
6:K:5105:HOH:O	2:L:91:LYS:HE2	1.72	0.89
3:R:89[A]:ARG:HD3	3:S:203:THR:CG2	2.00	0.89
3:U:89[A]:ARG:NH1	3:U:118:ASP:OD1	2.05	0.89
6:H:5089:HOH:O	2:I:91:LYS:HE2	1.70	0.89
1:A:28[B]:ARG:HH22	3:R:229:MET:HG2	1.36	0.88
3:R:89[A]:ARG:CD	3:S:203:THR:HG21	2.02	0.88
1:G:28[B]:ARG:NH2	3:Q:229:MET:HG2	1.93	0.83
1:B:28[B]:ARG:HH22	3:S:229:MET:HG2	1.44	0.83
3:T:144:ALA:HA	3:T:147:GLU:HG3	1.61	0.82
6:A:4015:HOH:O	3:Q:230:VAL:HG13	1.79	0.81
2:M:91:LYS:HE3	2:M:92:TYR:CE1	2.15	0.81
1:C:66:LYS:HB2	6:C:4131:HOH:O	1.78	0.81
3:R:75:GLN:OE1	6:R:419:HOH:O	1.99	0.81
1:C:28[A]:ARG:NH2	3:T:229:MET:HG3	1.97	0.78
1:G:68[B]:GLN:OE1	6:G:4058:HOH:O	2.01	0.78
1:A:65:GLU:CB	4:G:4007:SO4:O1	2.32	0.78
1:B:28[B]:ARG:HH22	3:S:229:MET:HG3	1.47	0.76
2:I:22:MET:HB2	2:I:27:MET:SD	2.25	0.76
1:C:28[A]:ARG:NH2	3:T:229:MET:CG	2.48	0.76
1:C:106:LEU:HD21	6:C:4072:HOH:O	1.86	0.75
3:S:116[B]:ILE:HG12	3:S:204[B]:MET:HE2	1.67	0.75
2:H:9:LYS:HG2	6:H:5060:HOH:O	1.85	0.75
1:G:67:ILE:HD12	1:G:211:GLU:HG2	1.67	0.74
3:R:11:ASN:OD1	6:R:276:HOH:O	2.04	0.74
1:A:28[B]:ARG:NH2	6:A:4056:HOH:O	2.18	0.74
3:S:89[A]:ARG:NH1	3:S:118:ASP:OD1	2.21	0.74
1:G:82[A]:VAL:HG11	3:P:231:SER:HB3	1.68	0.73
2:L:202:ILE:CG2	2:L:203:LEU:N	2.46	0.73
3:S:49:VAL:HG21	6:S:412:HOH:O	1.87	0.73
2:I:67:ARG:NE	6:I:5152:HOH:O	2.22	0.73
3:Q:101:GLU:HG3	3:R:105:LEU:CD2	2.20	0.72
3:U:125:LEU:O	3:U:142[A]:MET:HG3	1.89	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:68[B]:GLN:NE2	6:E:371:HOH:O	2.23	0.71
1:F:67:ILE:HG12	1:F:77[B]:VAL:HG12	1.72	0.71
2:J:22:MET:HB2	2:J:27:MET:SD	2.29	0.71
1:A:65:GLU:HG3	1:A:67:ILE:O	1.90	0.71
2:J:67:ARG:HD2	6:J:5151:HOH:O	1.89	0.71
1:D:67:ILE:HG12	1:D:77[A]:VAL:CG1	2.21	0.71
1:A:28[A]:ARG:NH2	3:R:228:HIS:CD2	2.59	0.70
1:G:67:ILE:HG12	1:G:77[A]:VAL:CG1	2.22	0.70
3:S:116[B]:ILE:CG1	3:S:204[B]:MET:HE2	2.22	0.70
1:A:28[B]:ARG:HH22	3:R:229:MET:CG	2.03	0.69
2:I:9:LYS:HG3	6:I:5017:HOH:O	1.92	0.69
6:O:261:HOH:O	3:U:129:GLU:HG2	1.93	0.69
3:O:125:LEU:O	3:O:142[A]:MET:HG3	1.92	0.69
2:K:96[B]:MET:CE	2:L:91:LYS:NZ	2.53	0.68
2:K:157:ARG:HD2	6:K:5018:HOH:O	1.93	0.68
1:A:54:VAL:HG21	1:A:61:GLN:HE22	1.57	0.67
3:T:144:ALA:HA	3:T:147:GLU:CG	2.25	0.67
2:N:91:LYS:HE3	2:N:92:TYR:CE1	2.29	0.67
1:A:90:ASP:HB3	6:N:5079:HOH:O	1.93	0.67
1:G:82[A]:VAL:CG1	3:P:231:SER:HB3	2.23	0.67
2:L:18:ARG:NH1	2:L:30:ASN:OD1	2.26	0.67
2:M:27:MET:HB3	6:N:5104:HOH:O	1.95	0.66
3:O:119:GLU:HG3	3:O:123:LYS:HD2	1.77	0.66
3:P:125:LEU:O	3:P:142[A]:MET:HG3	1.96	0.66
2:J:18:ARG:HH11	2:J:30:ASN:HD21	1.44	0.66
3:S:116[B]:ILE:CD1	3:S:204[B]:MET:CE	2.72	0.66
3:Q:101:GLU:CG	3:R:105:LEU:HD22	2.25	0.65
1:D:90:ASP:HB3	6:J:5096:HOH:O	1.97	0.65
3:S:116[B]:ILE:CD1	3:S:204[B]:MET:HE2	2.27	0.64
1:B:28[B]:ARG:NH2	3:S:229:MET:CG	2.53	0.64
1:D:28[B]:ARG:NH2	3:U:229:MET:HG2	2.09	0.64
1:F:68[B]:GLN:OE1	6:F:4082:HOH:O	2.15	0.64
3:U:89[A]:ARG:CZ	3:U:118:ASP:OD1	2.45	0.64
1:F:67:ILE:HG12	1:F:77[B]:VAL:CG1	2.28	0.63
3:R:144:ALA:N	6:R:419:HOH:O	2.31	0.63
1:B:28[B]:ARG:NH2	3:S:229:MET:HG3	2.12	0.63
2:N:157:ARG:HD2	6:N:5014:HOH:O	1.98	0.63
1:E:228:GLU:OE1	6:E:386:HOH:O	2.16	0.63
1:G:173:VAL:O	1:G:177:GLU:HG3	1.99	0.63
2:K:24:ASN:HD22	2:K:24:ASN:H	1.47	0.63
3:S:204[B]:MET:O	3:S:204[B]:MET:HE3	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:48:LEU:HD22	2:J:96[B]:MET:HE3	1.80	0.62
2:H:163:LYS:HZ1	2:H:203:LEU:HB3	1.63	0.62
3:O:140:ILE:HA	3:U:129:GLU:OE2	1.98	0.62
3:R:125:LEU:O	3:R:142[A]:MET:HG3	2.00	0.62
1:C:67:ILE:HG12	1:C:77[A]:VAL:CG1	2.30	0.61
1:E:28[B]:ARG:NH2	6:E:288:HOH:O	2.21	0.61
1:F:156:THR:HG23	1:G:82[B]:VAL:HG11	1.81	0.61
1:G:141:ILE:HD12	1:G:215:ILE:HG12	1.81	0.61
1:G:67:ILE:HG12	1:G:77[A]:VAL:HG12	1.83	0.60
3:S:113:VAL:O	3:S:116[B]:ILE:HG22	2.01	0.60
3:Q:144:ALA:HB3	6:Q:340:HOH:O	2.00	0.60
1:C:28[A]:ARG:NH2	6:C:4118:HOH:O	2.29	0.60
2:K:27:MET:HB3	6:L:5075:HOH:O	2.02	0.60
1:A:55:ARG:O	1:A:56:SER:CB	2.49	0.60
3:R:226:SER:HB2	3:R:228:HIS:NE2	2.16	0.60
1:G:106:LEU:HD21	6:G:4020:HOH:O	2.01	0.60
3:S:78:CYS:HB3	3:S:142[A]:MET:HG2	1.84	0.60
3:T:89[A]:ARG:CD	3:U:203:THR:HG21	2.30	0.59
3:Q:125:LEU:O	3:Q:142[A]:MET:HG3	2.03	0.59
1:C:28[A]:ARG:HH22	3:T:229:MET:HG2	1.61	0.59
3:R:224:THR:HG23	6:R:315:HOH:O	2.02	0.58
2:L:24:ASN:H	2:L:24:ASN:HD22	1.52	0.58
3:Q:140:ILE:CD1	3:Q:144:ALA:HB1	2.32	0.58
1:F:71:ASP:HA	2:L:68:LEU:HD11	1.86	0.58
2:N:37:ILE:HG21	2:N:43:MET:HE2	1.86	0.58
1:F:100:LYS:NZ	6:L:5117:HOH:O	2.35	0.58
2:H:22:MET:HB2	2:H:27:MET:SD	2.44	0.57
3:Q:140:ILE:HD11	3:Q:144:ALA:HB1	1.86	0.57
1:E:35:SER:HB2	3:U:231:SER:OXT	2.04	0.57
3:P:89[A]:ARG:NH1	3:P:118:ASP:OD1	2.37	0.57
1:E:64:ILE:HG22	1:E:64:ILE:O	2.05	0.57
2:J:18:ARG:HH11	2:J:30:ASN:ND2	2.02	0.57
3:S:226:SER:HB2	3:S:228:HIS:CD2	2.40	0.57
3:Q:116[A]:ILE:HD12	3:Q:208:VAL:HG22	1.85	0.57
1:F:48:LEU:HD13	1:F:77[B]:VAL:HG13	1.86	0.56
4:H:4015:SO4:O4	6:H:5044:HOH:O	2.18	0.56
3:R:89[A]:ARG:NH1	3:R:118:ASP:OD1	2.38	0.56
1:C:13:THR:N	6:C:4077:HOH:O	2.37	0.56
2:N:172:MET:HE2	2:N:189:THR:HG23	1.88	0.56
1:A:55:ARG:O	1:A:55:ARG:HG3	2.05	0.56
3:O:193:LYS:NZ	6:O:384:HOH:O	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:67:ILE:HG12	1:D:77[A]:VAL:HG12	1.86	0.56
1:E:67:ILE:HG12	1:E:77[A]:VAL:CG1	2.35	0.56
2:M:157:ARG:HD2	6:M:5024:HOH:O	2.05	0.56
1:B:63:SER:HB3	6:B:4160:HOH:O	2.06	0.56
3:S:125:LEU:O	3:S:142[A]:MET:HG3	2.05	0.56
2:M:163:LYS:HZ1	2:M:203:LEU:HB3	1.70	0.56
1:F:49:ILE:HG12	1:F:212:ILE:HG12	1.88	0.55
2:I:172:MET:HE2	2:I:189:THR:HG23	1.86	0.55
2:M:91:LYS:HE2	6:M:5038:HOH:O	2.05	0.55
1:D:28[B]:ARG:NH2	6:D:4062:HOH:O	2.37	0.55
1:G:48:LEU:HD13	1:G:77[A]:VAL:HG13	1.89	0.55
2:N:9:LYS:HG3	6:N:5013:HOH:O	2.06	0.55
2:N:195:ARG:HA	2:N:198:LYS:HG2	1.88	0.55
1:E:13:THR:N	6:E:350:HOH:O	2.39	0.55
1:B:177:GLU:OE1	1:C:56:SER:HB2	2.06	0.55
2:H:9:LYS:HE2	6:H:5060:HOH:O	2.06	0.55
2:J:163:LYS:HZ1	2:J:203:LEU:HB3	1.72	0.55
2:I:163:LYS:NZ	2:I:203:LEU:HB3	2.21	0.54
2:N:174[B]:ASP:OD1	2:N:186:GLN:NE2	2.32	0.54
1:C:53:LYS:NZ	6:C:4091:HOH:O	2.41	0.54
1:E:82[B]:VAL:HG12	3:U:229:MET:O	2.08	0.54
2:I:27:MET:HB3	6:J:5084:HOH:O	2.08	0.54
6:P:295:HOH:O	3:Q:204[A]:MET:HE3	2.07	0.54
1:A:28[B]:ARG:NH2	3:R:229:MET:HG2	2.16	0.54
2:L:174[B]:ASP:OD2	6:L:5076:HOH:O	2.18	0.54
1:D:71:ASP:HB2	2:J:64:GLU:OE2	2.08	0.54
1:E:126:TYR:HH	1:F:123:TYR:HH	1.53	0.54
1:D:126:TYR:HH	1:E:123:TYR:HH	1.53	0.54
6:T:330:HOH:O	3:U:200:HIS:CE1	2.61	0.54
2:M:195:ARG:HA	2:M:198:LYS:HG2	1.90	0.53
3:P:143:TYR:O	3:P:147:GLU:HG3	2.08	0.53
3:R:226:SER:O	3:R:228:HIS:CD2	2.60	0.53
1:A:82[B]:VAL:HG11	1:G:156:THR:HG23	1.89	0.53
1:B:17:PRO:HB3	3:T:105:LEU:HD11	1.90	0.53
3:Q:119:GLU:HG3	3:Q:123:LYS:HD2	1.89	0.53
3:S:129:GLU:HG3	3:T:138:THR:O	2.09	0.53
1:D:55:ARG:O	1:D:56:SER:HB3	2.09	0.53
1:G:55:ARG:O	1:G:56:SER:HB2	2.09	0.53
1:G:211:GLU:HG3	6:G:4030:HOH:O	2.07	0.53
1:A:65:GLU:CG	1:A:67:ILE:O	2.56	0.53
1:B:28[B]:ARG:NH2	3:S:229:MET:HG2	2.17	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S:28:TRP:HA	3:S:32:THR:HB	1.91	0.52
1:C:178:ARG:HD2	1:C:179:GLU:OE2	2.09	0.52
1:D:55:ARG:O	1:D:56:SER:CB	2.57	0.52
3:O:28:TRP:HA	3:O:32:THR:HB	1.91	0.52
3:S:116[B]:ILE:CG1	3:S:204[B]:MET:CE	2.86	0.52
1:B:71:ASP:HA	2:H:68:LEU:HD11	1.92	0.52
3:Q:65:GLU:HA	3:Q:68:LEU:HD12	1.92	0.52
1:B:82[B]:VAL:HG12	3:R:229:MET:O	2.09	0.52
1:F:156:THR:HG23	1:G:82[B]:VAL:CG1	2.39	0.52
1:D:49:ILE:HG23	1:D:212:ILE:HD13	1.91	0.51
3:O:89[A]:ARG:NH1	3:O:118:ASP:OD1	2.44	0.51
1:C:66:LYS:O	6:C:4130:HOH:O	2.19	0.51
1:G:68[B]:GLN:NE2	6:G:4063:HOH:O	2.35	0.51
1:F:28[B]:ARG:NH2	6:F:4059:HOH:O	2.43	0.51
3:P:225:GLY:O	3:P:226:SER:HB3	2.09	0.51
3:T:143:TYR:O	3:T:147:GLU:HG2	2.10	0.51
6:R:317:HOH:O	3:S:204[A]:MET:HG2	2.10	0.51
3:R:120:LEU:HD13	3:R:204[B]:MET:HG3	1.93	0.51
1:A:90:ASP:OD1	6:A:4070:HOH:O	2.20	0.50
2:K:24:ASN:H	2:K:24:ASN:ND2	2.09	0.50
2:H:3:THR:HB	2:H:16:THR:HG22	1.93	0.50
2:I:27:MET:HG2	6:J:5154:HOH:O	2.11	0.50
1:G:82[A]:VAL:HG22	3:P:229:MET:O	2.11	0.50
2:I:24:ASN:H	2:I:24:ASN:HD22	1.58	0.50
3:U:64:PRO:HB2	3:U:67:LEU:HG	1.94	0.50
1:G:71:ASP:HA	2:M:68:LEU:HD11	1.93	0.50
2:H:91:LYS:HD3	6:H:5075:HOH:O	2.11	0.50
3:Q:226:SER:HB2	3:Q:228:HIS:NE2	2.26	0.49
3:Q:230:VAL:O	3:Q:231:SER:HB3	2.12	0.49
2:J:111:PHE:CE2	2:J:121:GLU:HB2	2.48	0.49
1:G:28[B]:ARG:NH2	6:G:4066:HOH:O	2.37	0.49
2:H:24:ASN:HD22	2:H:24:ASN:H	1.59	0.49
2:K:9:LYS:HE2	6:K:5172:HOH:O	2.12	0.49
3:U:78:CYS:HB3	3:U:142[A]:MET:HG2	1.94	0.49
2:M:58[B]:TYR:CE2	2:M:86:MET:HE1	2.48	0.49
1:G:17:PRO:HB3	3:R:105:LEU:HD11	1.95	0.49
2:N:24:ASN:HB3	6:N:5109:HOH:O	2.12	0.49
3:U:113:VAL:O	3:U:116[B]:ILE:HG22	2.12	0.49
1:A:54:VAL:O	1:A:208:LYS:NZ	2.46	0.49
1:E:66:LYS:HB2	6:E:318:HOH:O	2.12	0.49
1:G:55:ARG:O	1:G:56:SER:CB	2.61	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:M:4020:SO4:O1	6:M:5097:HOH:O	2.20	0.49
2:N:24:ASN:H	2:N:24:ASN:HD22	1.61	0.49
2:I:157:ARG:CZ	6:I:5068:HOH:O	2.61	0.48
2:M:157:ARG:HD3	6:M:5116:HOH:O	2.13	0.48
1:D:71:ASP:HA	2:J:68:LEU:HD11	1.95	0.48
2:M:22:MET:HE1	6:N:5180:HOH:O	2.13	0.48
1:G:206:GLU:HG2	6:G:4120:HOH:O	2.12	0.48
3:T:28:TRP:HA	3:T:32:THR:HB	1.94	0.48
2:K:163:LYS:NZ	2:K:203:LEU:HB3	2.28	0.48
2:M:111:PHE:CE2	2:M:121:GLU:HB2	2.49	0.48
2:N:111:PHE:CE2	2:N:121:GLU:HB2	2.49	0.48
3:R:28:TRP:HA	3:R:32:THR:HB	1.96	0.48
3:T:160:LEU:O	3:T:161:GLY:C	2.50	0.48
3:P:24:VAL:HG13	3:P:84:GLN:HB3	1.96	0.48
2:I:74:MET:HG2	2:I:78:ALA:HB3	1.96	0.47
2:L:195:ARG:HA	2:L:198:LYS:HG2	1.95	0.47
3:Q:28:TRP:HA	3:Q:32:THR:HB	1.96	0.47
3:U:28:TRP:HA	3:U:32:THR:HB	1.94	0.47
1:A:38:LEU:HD12	1:A:38:LEU:C	2.34	0.47
1:F:82[A]:VAL:HG22	3:O:229:MET:O	2.14	0.47
1:G:13:THR:N	6:G:4083:HOH:O	2.47	0.47
2:J:96[B]:MET:HE3	2:K:91:LYS:HZ2	1.80	0.47
2:L:111:PHE:CE2	2:L:121:GLU:HB2	2.50	0.47
1:A:33:LYS:O	1:A:53:LYS:NZ	2.47	0.47
1:A:180:TYR:HA	1:A:192:LEU:HD21	1.96	0.47
1:E:42:PHE:HB2	1:E:184:LEU:O	2.14	0.47
2:J:45:ILE:HB	2:J:52:ALA:HB1	1.96	0.47
2:M:37:ILE:HG21	2:M:43:MET:HE2	1.97	0.47
2:M:196:ILE:HG23	2:M:201:LEU:HB2	1.95	0.47
3:R:116[B]:ILE:O	3:R:204[B]:MET:HE3	2.15	0.47
3:T:112:ALA:O	3:T:116[A]:ILE:HG23	2.14	0.47
2:K:202:ILE:HG22	2:K:203:LEU:N	2.30	0.47
3:P:41:LYS:O	3:P:45:GLU:HG2	2.15	0.47
1:A:106:LEU:HD21	6:A:4023:HOH:O	2.15	0.47
1:C:65:GLU:HG3	6:I:5134:HOH:O	2.15	0.47
1:G:201:LEU:HB3	1:G:205:GLU:HB3	1.97	0.47
1:C:28[A]:ARG:NH2	3:T:229:MET:HG2	2.25	0.46
1:E:67:ILE:HG12	1:E:77[A]:VAL:HG12	1.97	0.46
1:F:64:ILE:HG22	1:F:64:ILE:O	2.15	0.46
2:N:45:ILE:HB	2:N:52:ALA:HB1	1.97	0.46
2:I:163:LYS:HZ1	2:I:203:LEU:HB3	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:208:LYS:O	1:G:209:ALA:HB2	2.16	0.46
1:A:28[A]:ARG:NH1	1:A:152:ASP:OD2	2.49	0.46
1:A:54:VAL:HG21	1:A:61:GLN:NE2	2.28	0.46
2:L:104:ILE:HD12	2:L:179:THR:HA	1.98	0.46
3:T:39:ILE:HG12	3:T:73:ARG:HG2	1.97	0.46
3:T:113:VAL:O	3:T:116[B]:ILE:HG22	2.16	0.46
1:C:228:GLU:OE1	6:C:4129:HOH:O	2.20	0.46
1:E:144:ILE:N	1:E:144:ILE:HD12	2.31	0.46
2:I:91:LYS:HE3	2:I:92:TYR:CE1	2.50	0.46
2:M:74:MET:HG2	2:M:78:ALA:HB3	1.98	0.46
3:U:123:LYS:HD2	6:U:385:HOH:O	2.14	0.46
2:L:196:ILE:HG23	2:L:201:LEU:HB2	1.97	0.46
1:A:77[A]:VAL:HG12	1:A:137:ILE:HB	1.96	0.46
2:K:3:THR:HB	2:K:16:THR:HG22	1.98	0.46
2:M:24:ASN:HD22	2:M:24:ASN:H	1.63	0.46
3:P:101:GLU:CG	3:Q:105:LEU:HD22	2.46	0.46
3:S:204[B]:MET:CE	3:S:208:VAL:HG23	2.45	0.46
3:U:226:SER:HB2	3:U:228:HIS:CD2	2.51	0.46
6:O:307:HOH:O	3:P:204[A]:MET:HG2	2.15	0.46
1:D:67:ILE:HD12	1:D:211:GLU:HB3	1.96	0.46
2:K:96[B]:MET:CE	2:L:91:LYS:HZ1	1.95	0.46
2:M:96[B]:MET:HA	2:M:96[B]:MET:CE	2.46	0.46
1:F:33:LYS:O	1:F:53:LYS:NZ	2.39	0.45
2:K:91:LYS:HE3	2:K:92:TYR:CE1	2.51	0.45
1:A:201:LEU:HD11	1:A:207:LEU:HB2	1.97	0.45
2:M:58[A]:TYR:CE2	6:M:5099:HOH:O	2.68	0.45
3:P:133:SER:N	3:Q:133:SER:O	2.50	0.45
3:Q:113:VAL:O	3:Q:116[B]:ILE:HG22	2.17	0.45
3:S:37[B]:GLU:OE2	3:S:41:LYS:HE2	2.16	0.45
1:A:126:TYR:OH	1:B:123:TYR:OH	2.35	0.45
2:K:172:MET:HE3	2:K:192[B]:ILE:CG2	2.47	0.45
1:D:28[B]:ARG:CZ	3:U:229:MET:HG3	2.43	0.45
1:E:71:ASP:HA	2:K:68:LEU:HD11	1.98	0.45
1:F:42:PHE:HB2	1:F:184:LEU:O	2.16	0.45
3:U:77:LEU:O	3:U:77:LEU:HG	2.16	0.45
3:U:89[A]:ARG:NH2	3:U:118:ASP:OD1	2.50	0.45
2:I:91:LYS:HE3	2:I:92:TYR:CZ	2.52	0.45
3:P:75:GLN:NE2	6:P:297:HOH:O	2.47	0.45
1:D:33:LYS:HB2	3:T:230:VAL:HG11	1.98	0.45
3:T:89[A]:ARG:NH2	6:T:402:HOH:O	2.49	0.45
3:Q:143:TYR:O	3:Q:147:GLU:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:R:50:ILE:HD11	3:R:67:LEU:HD21	1.99	0.45
1:C:90:ASP:OD1	6:C:4116:HOH:O	2.21	0.44
2:L:45:ILE:HB	2:L:52:ALA:HB1	1.99	0.44
3:Q:153[B]:SER:OG	3:R:179:LEU:HB2	2.18	0.44
3:R:75:GLN:HE21	3:S:192:LEU:CD2	2.31	0.44
1:C:67:ILE:HD12	1:C:211:GLU:HB3	1.99	0.44
2:J:74:MET:HG2	2:J:78:ALA:HB3	1.99	0.44
1:C:222:ARG:HH12	1:C:228:GLU:CD	2.20	0.44
3:S:116[B]:ILE:CD1	3:S:204[B]:MET:HE1	2.34	0.44
3:U:90:THR:O	3:U:94:ILE:HG12	2.17	0.44
3:O:143:TYR:O	3:O:147:GLU:HG3	2.18	0.44
3:O:24:VAL:HG13	3:O:84:GLN:HB3	1.99	0.44
1:F:61:GLN:HE21	1:F:61:GLN:HB3	1.66	0.44
2:K:15:ALA:HA	2:K:174[B]:ASP:O	2.18	0.44
1:D:107:VAL:O	1:D:142:ASP:HB2	2.18	0.44
2:H:163:LYS:NZ	2:H:203:LEU:HB3	2.31	0.44
1:B:107:VAL:O	1:B:142:ASP:HB2	2.18	0.44
1:C:82[A]:VAL:HG22	3:S:229:MET:O	2.18	0.44
1:B:13:THR:HA	6:B:4090:HOH:O	2.17	0.43
1:G:41:LYS:HD3	1:G:161:LYS:HA	2.01	0.43
1:G:208:LYS:O	1:G:209:ALA:CB	2.66	0.43
2:M:24:ASN:H	2:M:24:ASN:ND2	2.15	0.43
3:P:144:ALA:N	6:P:314:HOH:O	2.51	0.43
3:R:65:GLU:HA	3:R:68:LEU:HD12	1.99	0.43
3:S:204[B]:MET:HE2	3:S:208:VAL:HG23	2.00	0.43
1:C:71:ASP:HA	2:I:68:LEU:HD11	2.00	0.43
1:F:55:ARG:O	1:F:56:SER:HB2	2.18	0.43
1:G:64:ILE:HG22	1:G:64:ILE:O	2.19	0.43
1:G:226:GLN:H	1:G:226:GLN:HG3	1.66	0.43
2:L:89:GLN:HG3	6:L:5064:HOH:O	2.18	0.43
3:R:77:LEU:O	3:R:77:LEU:HG	2.15	0.43
1:B:106:LEU:HD21	6:B:4063:HOH:O	2.19	0.43
3:R:75:GLN:HE21	3:S:192:LEU:HD22	1.82	0.43
1:A:82[A]:VAL:HG22	3:Q:229:MET:O	2.19	0.43
1:C:17:PRO:HB3	3:U:105:LEU:HD11	2.00	0.43
1:E:21:LEU:HD12	3:O:229:MET:HE3	1.99	0.43
2:H:91:LYS:HZ1	2:N:96[A]:MET:HG3	1.83	0.43
3:Q:68:LEU:HD21	3:Q:149:LEU:HD21	2.00	0.43
3:Q:140:ILE:HG12	3:Q:144:ALA:HB2	2.01	0.43
1:B:160:TYR:OH	1:C:55:ARG:NH1	2.51	0.43
1:E:223:ILE:N	1:E:223:ILE:HD12	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:PRO:HA	1:B:26:TYR:CD2	2.54	0.43
3:P:28:TRP:HA	3:P:32:THR:HB	2.01	0.43
1:B:28[B]:ARG:NH2	6:B:4091:HOH:O	2.43	0.43
3:Q:86:GLU:HG3	3:R:199[A]:THR:HG22	2.01	0.43
1:C:77[B]:VAL:HG12	1:C:137:ILE:HB	2.00	0.43
2:I:45:ILE:HB	2:I:52:ALA:HB1	1.99	0.43
3:O:226:SER:O	3:O:228:HIS:N	2.50	0.43
1:E:28[A]:ARG:NH1	1:E:152:ASP:OD2	2.52	0.42
3:P:176:SER:HA	3:P:177:PRO:HD2	1.93	0.42
1:B:156:THR:HG23	1:C:82[B]:VAL:HG11	1.99	0.42
1:E:175:PHE:HE1	1:E:192:LEU:HD12	1.84	0.42
1:G:180:TYR:HA	1:G:192:LEU:HD21	2.00	0.42
2:L:24:ASN:H	2:L:24:ASN:ND2	2.17	0.42
1:A:65:GLU:OE2	2:N:71:ARG:NH1	2.53	0.42
2:J:190[B]:ASP:OD1	2:J:191:GLN:N	2.51	0.42
3:Q:50:ILE:HD11	3:Q:67:LEU:HD21	2.00	0.42
6:T:330:HOH:O	3:U:204[A]:MET:HG2	2.19	0.42
2:M:23:GLU:HB3	2:M:24:ASN:H	1.58	0.42
2:N:112[B]:SER:O	2:N:119:SER:HA	2.19	0.42
2:N:164:GLN:HE22	2:N:202:ILE:HD11	1.83	0.42
3:R:11:ASN:HB2	6:R:316:HOH:O	2.18	0.42
3:R:36:ILE:HG21	3:R:202[B]:SER:OG	2.19	0.42
3:T:144:ALA:CA	3:T:147:GLU:HG3	2.39	0.42
3:R:225:GLY:O	3:R:226:SER:HB3	2.19	0.42
1:A:55:ARG:O	1:A:56:SER:HB3	2.20	0.42
3:R:63:SER:HG	3:S:178:SER:HG	1.67	0.42
2:J:24:ASN:HD22	2:J:24:ASN:H	1.66	0.42
3:O:63:SER:HA	3:O:64:PRO:HD3	1.85	0.42
1:A:77[A]:VAL:HG22	6:A:4142:HOH:O	2.19	0.42
1:B:55:ARG:O	1:B:56:SER:CB	2.68	0.42
1:E:156:THR:HG23	1:F:82[B]:VAL:HG11	2.02	0.42
1:F:82[A]:VAL:HG11	3:O:231:SER:HB3	2.02	0.42
2:H:9:LYS:HG2	2:H:9:LYS:H	1.68	0.42
1:F:203:GLU:O	1:F:205:GLU:N	2.53	0.42
1:F:209:ALA:HA	1:F:233:LEU:HD11	2.00	0.42
1:G:56:SER:HB3	1:G:59:ILE:HG12	2.01	0.42
3:P:181:LEU:HD22	3:P:184:ARG:HH12	1.85	0.42
2:K:202:ILE:CG2	2:K:203:LEU:N	2.82	0.41
1:G:13:THR:CA	6:G:4083:HOH:O	2.67	0.41
2:H:24:ASN:H	2:H:24:ASN:ND2	2.18	0.41
1:D:152:ASP:HB2	1:D:153:PRO:CD	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:156:THR:HG23	1:F:82[B]:VAL:CG1	2.51	0.41
2:K:163:LYS:HZ3	2:K:203:LEU:HB3	1.84	0.41
2:L:22:MET:HB2	2:L:27:MET:SD	2.60	0.41
1:D:42:PHE:HB2	1:D:184:LEU:O	2.21	0.41
1:G:13:THR:HA	6:G:4083:HOH:O	2.20	0.41
2:L:163:LYS:HB3	2:L:203:LEU:HD13	2.01	0.41
2:M:20:VAL:HB	2:M:28:HIS:HB2	2.02	0.41
3:U:217:LYS:HD2	6:U:375:HOH:O	2.19	0.41
2:H:54:VAL:HG23	6:I:5043:HOH:O	2.18	0.41
3:S:50:ILE:HD11	3:S:67:LEU:HD21	2.01	0.41
2:K:180:ARG:HD3	6:K:5101:HOH:O	2.20	0.41
3:S:77:LEU:O	3:S:77:LEU:HG	2.19	0.41
2:K:23:GLU:HB3	2:K:24:ASN:HD22	1.85	0.41
2:M:91:LYS:HD2	6:M:5041:HOH:O	2.20	0.41
3:T:75:GLN:HE21	3:U:192:LEU:CD2	2.34	0.41
1:C:118:ASP:O	1:C:122:GLN:HG3	2.21	0.41
1:A:207:LEU:HD12	1:A:207:LEU:HA	1.95	0.41
1:D:100:LYS:NZ	6:J:5095:HOH:O	2.54	0.41
1:E:203:GLU:H	1:E:203:GLU:HG3	1.57	0.41
2:J:3:THR:HB	2:J:16:THR:HG22	2.02	0.41
3:U:78:CYS:CB	3:U:142[B]:MET:HB2	2.51	0.41
2:M:64:GLU:HG3	2:M:67:ARG:HH21	1.86	0.41
1:B:180:TYR:HA	1:B:192:LEU:HD21	2.03	0.40
1:C:28[B]:ARG:NH1	1:C:152:ASP:OD2	2.54	0.40
1:E:152:ASP:HB2	1:E:153:PRO:CD	2.51	0.40
2:J:195:ARG:HA	2:J:198:LYS:HG2	2.02	0.40
3:O:10:GLN:NE2	3:O:14:ASP:OD1	2.49	0.40
6:O:261:HOH:O	3:U:129:GLU:CG	2.60	0.40
3:T:221:GLN:NE2	6:T:304:HOH:O	2.54	0.40
1:A:69:LEU:HD23	1:A:69:LEU:HA	1.91	0.40
1:A:71:ASP:HA	2:N:68:LEU:HD11	2.03	0.40
1:F:55:ARG:O	1:F:56:SER:CB	2.70	0.40
2:M:45:ILE:HB	2:M:52:ALA:HB1	2.02	0.40
3:T:218:LYS:NZ	6:T:428:HOH:O	2.53	0.40
1:A:63:SER:HB2	1:G:159:GLU:OE1	2.21	0.40
1:D:17:PRO:HB3	3:O:105:LEU:HD11	2.03	0.40
1:D:162:ALA:HB1	1:D:176:LEU:HD13	2.02	0.40
1:E:182:GLU:HG2	6:E:341:HOH:O	2.19	0.40
1:F:94:ILE:O	1:F:98:GLN:HG3	2.21	0.40
2:H:74:MET:HG2	2:H:78:ALA:HB3	2.01	0.40
3:O:204[A]:MET:HA	3:O:204[A]:MET:CE	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:203:GLU:H	1:G:203:GLU:HG3	1.58	0.40
2:J:18:ARG:NH1	2:J:30:ASN:HD21	2.15	0.40
3:O:78:CYS:HB3	3:O:142[A]:MET:HG2	2.04	0.40
2:I:71:ARG:NE	6:I:5169:HOH:O	2.54	0.40
2:L:23:GLU:HB3	2:L:24:ASN:H	1.72	0.40
3:T:89[B]:ARG:HG3	3:T:117:ILE:HG21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	226/233 (97%)	219 (97%)	5 (2%)	2 (1%)	17	7
1	B	226/233 (97%)	218 (96%)	7 (3%)	1 (0%)	34	24
1	C	225/233 (97%)	220 (98%)	4 (2%)	1 (0%)	34	24
1	D	226/233 (97%)	219 (97%)	6 (3%)	1 (0%)	34	24
1	E	225/233 (97%)	220 (98%)	5 (2%)	0	100	100
1	F	225/233 (97%)	216 (96%)	7 (3%)	2 (1%)	17	7
1	G	225/233 (97%)	217 (96%)	5 (2%)	3 (1%)	12	4
2	H	208/217 (96%)	204 (98%)	4 (2%)	0	100	100
2	I	208/217 (96%)	202 (97%)	6 (3%)	0	100	100
2	J	208/217 (96%)	204 (98%)	4 (2%)	0	100	100
2	K	208/217 (96%)	204 (98%)	4 (2%)	0	100	100
2	L	208/217 (96%)	201 (97%)	6 (3%)	1 (0%)	29	18
2	M	208/217 (96%)	204 (98%)	4 (2%)	0	100	100
2	N	208/217 (96%)	204 (98%)	4 (2%)	0	100	100
3	O	224/237 (94%)	220 (98%)	2 (1%)	2 (1%)	17	7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	P	224/237 (94%)	218 (97%)	5 (2%)	1 (0%)	34	24
3	Q	224/237 (94%)	221 (99%)	3 (1%)	0	100	100
3	R	224/237 (94%)	220 (98%)	3 (1%)	1 (0%)	34	24
3	S	224/237 (94%)	217 (97%)	6 (3%)	1 (0%)	34	24
3	T	223/237 (94%)	219 (98%)	3 (1%)	1 (0%)	34	24
3	U	224/237 (94%)	219 (98%)	4 (2%)	1 (0%)	34	24
All	All	4601/4809 (96%)	4486 (98%)	97 (2%)	18 (0%)	34	24

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	SER
1	B	56	SER
1	D	56	SER
1	F	56	SER
1	G	56	SER
1	G	209	ALA
3	P	226	SER
3	R	226	SER
3	T	226	SER
1	F	204	GLY
3	O	226	SER
1	C	55	ARG
1	G	203	GLU
3	S	226	SER
3	U	226	SER
2	L	202	ILE
3	O	227	ASP
1	A	204	GLY

5.3.2 Protein sidechains [i](#)

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	191/193 (99%)	185 (97%)	6 (3%)	40	32
1	B	191/193 (99%)	186 (97%)	5 (3%)	46	39
1	C	190/193 (98%)	183 (96%)	7 (4%)	34	25
1	D	191/193 (99%)	189 (99%)	2 (1%)	76	76
1	E	190/193 (98%)	186 (98%)	4 (2%)	53	48
1	F	190/193 (98%)	181 (95%)	9 (5%)	26	16
1	G	190/193 (98%)	185 (97%)	5 (3%)	46	39
2	H	177/183 (97%)	172 (97%)	5 (3%)	43	36
2	I	177/183 (97%)	174 (98%)	3 (2%)	60	57
2	J	177/183 (97%)	171 (97%)	6 (3%)	37	28
2	K	177/183 (97%)	170 (96%)	7 (4%)	31	22
2	L	177/183 (97%)	172 (97%)	5 (3%)	43	36
2	M	177/183 (97%)	172 (97%)	5 (3%)	43	36
2	N	177/183 (97%)	175 (99%)	2 (1%)	73	73
3	O	189/196 (96%)	184 (97%)	5 (3%)	46	39
3	P	189/196 (96%)	188 (100%)	1 (0%)	88	89
3	Q	189/196 (96%)	183 (97%)	6 (3%)	39	30
3	R	189/196 (96%)	184 (97%)	5 (3%)	46	39
3	S	189/196 (96%)	188 (100%)	1 (0%)	88	89
3	T	188/196 (96%)	180 (96%)	8 (4%)	29	19
3	U	189/196 (96%)	188 (100%)	1 (0%)	88	89
All	All	3894/4004 (97%)	3796 (98%)	98 (2%)	49	41

All (98) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	GLU
1	A	203	GLU
1	A	205	GLU
1	A	221	TYR
1	A	226	GLN
1	A	227	GLU
1	B	35	SER
1	B	52	LYS
1	B	62	ASN
1	B	63	SER

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Mol	Chain	Res	Type
1	B	221	TYR
1	C	55	ARG
1	C	63	SER
1	C	177	GLU
1	C	203	GLU
1	C	205	GLU
1	C	221	TYR
1	C	227	GLU
1	D	65	GLU
1	D	221	TYR
1	E	35	SER
1	E	59	ILE
1	E	203	GLU
1	E	221	TYR
1	F	35	SER
1	F	52	LYS
1	F	55	ARG
1	F	61	GLN
1	F	82[A]	VAL
1	F	82[B]	VAL
1	F	221	TYR
1	F	227	GLU
1	F	230	LYS
1	G	61	GLN
1	G	203	GLU
1	G	221	TYR
1	G	225	ASP
1	G	226	GLN
2	H	3	THR
2	H	9	LYS
2	H	17	GLU
2	H	23	GLU
2	H	25	PHE
2	I	17	GLU
2	I	23	GLU
2	I	25	PHE
2	J	3	THR
2	J	17	GLU
2	J	25	PHE
2	J	30	ASN
2	J	105	ASP
2	J	203	LEU

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Mol	Chain	Res	Type
2	K	3	THR
2	K	9	LYS
2	K	17	GLU
2	K	23	GLU
2	K	24	ASN
2	K	25	PHE
2	K	203	LEU
2	L	17	GLU
2	L	24	ASN
2	L	25	PHE
2	L	71	ARG
2	L	203	LEU
2	M	17	GLU
2	M	23	GLU
2	M	25	PHE
2	M	180	ARG
2	M	203	LEU
2	N	17	GLU
2	N	25	PHE
3	O	57	ARG
3	O	89[A]	ARG
3	O	89[B]	ARG
3	O	204[A]	MET
3	O	204[B]	MET
3	P	231	SER
3	Q	77	LEU
3	Q	89[A]	ARG
3	Q	89[B]	ARG
3	Q	204[A]	MET
3	Q	204[B]	MET
3	Q	230	VAL
3	R	57	ARG
3	R	65	GLU
3	R	145	LEU
3	R	204[A]	MET
3	R	204[B]	MET
3	S	77	LEU
3	T	116[A]	ILE
3	T	116[B]	ILE
3	T	140	ILE
3	T	142[A]	MET
3	T	142[B]	MET

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Mol	Chain	Res	Type
3	T	147	GLU
3	T	157	ASP
3	T	204	MET
3	U	77	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	B	143	GLN
1	D	111	ASN
1	F	61	GLN
1	F	143	GLN
1	G	61	GLN
2	H	24	ASN
2	I	24	ASN
2	J	24	ASN
2	J	30	ASN
2	J	191	GLN
2	K	24	ASN
2	L	24	ASN
2	M	24	ASN
2	M	73	ASN
2	N	24	ASN
2	N	164	GLN
3	O	79	HIS
3	P	79	HIS
3	P	228	HIS
3	Q	79	HIS
3	R	228	HIS
3	S	79	HIS
3	S	228	HIS
3	T	79	HIS
3	T	221	GLN
3	T	228	HIS
3	U	79	HIS
3	U	228	HIS

5.3.3 RNA

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

28 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	F	4006	-	4,4,4	0.18	0	6,6,6	0.12	0
4	SO4	I	4011	-	4,4,4	0.11	0	6,6,6	0.38	0
4	SO4	B	4016	-	4,4,4	0.14	0	6,6,6	0.42	0
4	SO4	F	4005	-	4,4,4	0.14	0	6,6,6	0.09	0
4	SO4	H	4015	-	4,4,4	0.20	0	6,6,6	0.39	0
5	GOL	K	5007	-	5,5,5	0.38	0	5,5,5	0.80	0
5	GOL	J	5006	-	5,5,5	0.53	0	5,5,5	0.58	0
4	SO4	G	4007	-	4,4,4	0.12	0	6,6,6	0.30	0
4	SO4	L	4019	-	4,4,4	0.09	0	6,6,6	0.19	0
5	GOL	L	5001	-	5,5,5	0.34	0	5,5,5	0.46	0
4	SO4	J	4017	-	4,4,4	0.16	0	6,6,6	0.13	0
4	SO4	C	4003	-	4,4,4	0.14	0	6,6,6	0.11	0
4	SO4	N	4021	-	4,4,4	0.15	0	6,6,6	0.19	0
4	SO4	D	4004	-	4,4,4	0.15	0	6,6,6	0.11	0
5	GOL	N	5003	-	5,5,5	0.34	0	5,5,5	0.67	0
4	SO4	M	4010	-	4,4,4	0.24	0	6,6,6	0.40	0
4	SO4	J	4012	-	4,4,4	0.10	0	6,6,6	0.41	0
5	GOL	H	5004	-	5,5,5	0.45	0	5,5,5	1.39	1 (20%)
5	GOL	I	5005	-	5,5,5	0.37	0	5,5,5	0.52	0
4	SO4	A	4001	-	4,4,4	0.14	0	6,6,6	0.07	0
4	SO4	N	4009	-	4,4,4	0.18	0	6,6,6	0.28	0
4	SO4	H	4008	-	4,4,4	0.27	0	6,6,6	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	K	4018	-	4,4,4	0.15	0	6,6,6	0.46	0
4	SO4	M	4020	-	4,4,4	0.07	0	6,6,6	0.13	0
4	SO4	B	4002	-	4,4,4	0.16	0	6,6,6	0.18	0
4	SO4	L	4014	-	4,4,4	0.18	0	6,6,6	0.55	0
4	SO4	K	4013	-	4,4,4	0.12	0	6,6,6	0.68	0
5	GOL	M	5002	-	5,5,5	0.26	0	5,5,5	0.44	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	L	5001	-	-	2/4/4/4	-
5	GOL	M	5002	-	-	3/4/4/4	-
5	GOL	N	5003	-	-	1/4/4/4	-
5	GOL	H	5004	-	-	4/4/4/4	-
5	GOL	I	5005	-	-	1/4/4/4	-
5	GOL	K	5007	-	-	0/4/4/4	-
5	GOL	J	5006	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	5004	GOL	O2-C2-C1	2.26	119.06	109.12

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	H	5004	GOL	C1-C2-C3-O3
5	M	5002	GOL	O1-C1-C2-C3
5	H	5004	GOL	O1-C1-C2-C3
5	L	5001	GOL	O1-C1-C2-C3
5	H	5004	GOL	O1-C1-C2-O2
5	H	5004	GOL	O2-C2-C3-O3
5	L	5001	GOL	O1-C1-C2-O2
5	M	5002	GOL	O1-C1-C2-O2
5	M	5002	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
5	J	5006	GOL	O2-C2-C3-O3
5	J	5006	GOL	C1-C2-C3-O3
5	I	5005	GOL	C1-C2-C3-O3
5	N	5003	GOL	C1-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	4015	SO4	1	0
4	G	4007	SO4	2	0
4	M	4020	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	221/233 (94%)	0.27	14 (6%) 20 22	18, 28, 61, 71	0
1	B	221/233 (94%)	0.46	19 (8%) 10 12	16, 29, 60, 71	0
1	C	221/233 (94%)	0.26	18 (8%) 12 13	17, 28, 63, 70	0
1	D	221/233 (94%)	0.53	19 (8%) 10 12	18, 31, 62, 73	0
1	E	221/233 (94%)	0.19	13 (5%) 22 25	18, 27, 55, 65	0
1	F	221/233 (94%)	0.42	21 (9%) 8 9	19, 31, 63, 75	0
1	G	221/233 (94%)	0.41	16 (7%) 15 17	18, 30, 58, 70	0
2	H	203/217 (93%)	-0.02	5 (2%) 57 60	18, 23, 38, 65	0
2	I	203/217 (93%)	0.12	5 (2%) 57 60	19, 24, 40, 61	0
2	J	203/217 (93%)	0.30	11 (5%) 25 29	20, 25, 41, 65	0
2	K	203/217 (93%)	0.07	9 (4%) 34 37	19, 25, 40, 62	0
2	L	203/217 (93%)	0.21	7 (3%) 45 48	19, 25, 40, 65	0
2	M	203/217 (93%)	-0.04	7 (3%) 45 48	17, 22, 37, 60	0
2	N	203/217 (93%)	0.34	5 (2%) 57 60	17, 24, 39, 65	0
3	O	218/237 (91%)	0.05	11 (5%) 28 32	22, 30, 47, 61	0
3	P	218/237 (91%)	0.07	8 (3%) 41 44	24, 32, 50, 67	0
3	Q	218/237 (91%)	0.12	12 (5%) 25 28	22, 32, 51, 79	0
3	R	218/237 (91%)	0.05	13 (5%) 21 24	17, 28, 49, 63	0
3	S	218/237 (91%)	-0.13	9 (4%) 37 40	17, 25, 42, 66	0
3	T	218/237 (91%)	-0.09	8 (3%) 41 44	16, 25, 44, 61	0
3	U	218/237 (91%)	-0.09	7 (3%) 47 50	18, 26, 41, 64	0
All	All	4494/4809 (93%)	0.17	237 (5%) 26 29	16, 27, 54, 79	0

All (237) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	N	203	LEU	14.0
2	L	203	LEU	13.6
2	I	203	LEU	11.8
2	N	202	ILE	11.2
2	K	202	ILE	10.3
2	H	203	LEU	9.9
2	M	203	LEU	9.7
2	K	203	LEU	9.5
3	Q	226	SER	9.3
2	J	203	LEU	9.1
1	G	55	ARG	9.0
2	J	202	ILE	9.0
1	D	54	VAL	8.4
3	U	226	SER	8.3
2	H	202	ILE	7.9
3	S	226	SER	7.9
3	R	226	SER	7.8
2	M	202	ILE	7.6
2	I	202	ILE	7.1
3	Q	227	ASP	7.0
1	F	202	GLU	6.9
1	B	64	ILE	6.6
1	F	55	ARG	6.2
1	B	55	ARG	6.2
3	O	226	SER	6.0
3	U	228	HIS	6.0
1	E	54	VAL	6.0
3	O	225	GLY	5.9
3	P	226	SER	5.8
1	D	55	ARG	5.7
1	B	54	VAL	5.7
1	E	61	GLN	5.6
2	L	202	ILE	5.4
1	A	55	ARG	5.3
3	Q	231	SER	5.2
1	G	203	GLU	5.1
3	O	172	GLY	5.1
1	F	54	VAL	5.0
3	T	226	SER	4.9
3	T	161	GLY	4.8
3	Q	172	GLY	4.8
1	E	203	GLU	4.8
3	T	172	GLY	4.7

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Mol	Chain	Res	Type	RSRZ
1	A	203	GLU	4.7
1	F	233	LEU	4.6
1	F	203	GLU	4.6
3	Q	228	HIS	4.5
3	R	228	HIS	4.5
3	S	227	ASP	4.5
1	C	203	GLU	4.5
1	E	55	ARG	4.4
3	R	225	GLY	4.4
1	F	61	GLN	4.3
3	R	227	ASP	4.3
3	S	228	HIS	4.2
3	T	225	GLY	4.2
1	E	53	LYS	4.2
1	D	233	LEU	4.2
3	P	227	ASP	4.1
1	G	54	VAL	4.1
3	Q	230	VAL	4.1
1	E	64	ILE	4.1
1	G	53	LYS	4.0
3	T	227	ASP	4.0
1	D	230	LYS	4.0
3	Q	52	ASN	4.0
1	D	61	GLN	4.0
1	E	204	GLY	3.9
3	O	227	ASP	3.9
3	U	227	ASP	3.9
1	A	206	GLU	3.9
1	C	226	GLN	3.9
3	T	228	HIS	3.8
1	D	208	LYS	3.8
3	S	172	GLY	3.8
3	O	228	HIS	3.8
1	C	55	ARG	3.8
3	R	172	GLY	3.8
1	D	203	GLU	3.8
1	B	61	GLN	3.8
1	C	233	LEU	3.7
1	G	208	LYS	3.7
3	P	230	VAL	3.7
1	G	206	GLU	3.7
3	Q	225	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	64	ILE	3.7
1	F	64	ILE	3.7
1	B	208	LYS	3.6
1	C	231	LYS	3.6
3	R	52	ASN	3.5
1	B	203	GLU	3.5
1	D	206	GLU	3.4
3	U	231	SER	3.4
1	F	206	GLU	3.4
1	G	61	GLN	3.4
2	K	9	LYS	3.4
3	U	225	GLY	3.4
1	B	206	GLU	3.3
1	C	206	GLU	3.3
1	C	204	GLY	3.3
1	F	207	LEU	3.3
3	O	173	GLY	3.2
1	A	53	LYS	3.2
1	A	62	ASN	3.2
3	O	57	ARG	3.2
3	R	231	SER	3.2
1	G	204	GLY	3.2
1	F	205	GLU	3.2
2	J	22	MET	3.2
3	Q	173	GLY	3.2
3	P	228	HIS	3.1
1	A	65	GLU	3.1
1	G	64	ILE	3.1
2	L	92	TYR	3.1
2	K	93	MET	3.1
1	D	226	GLN	3.0
1	A	226	GLN	3.0
1	G	202	GLU	3.0
1	C	64	ILE	3.0
2	I	92	TYR	3.0
1	A	233	LEU	3.0
1	D	227	GLU	3.0
1	F	232	PHE	3.0
1	G	233	LEU	3.0
2	M	93	MET	3.0
3	Q	229	MET	3.0
1	C	230	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
2	L	27	MET	2.9
1	B	202	GLU	2.9
1	F	226	GLN	2.9
1	C	208	LYS	2.9
1	F	208	LYS	2.9
1	G	175	PHE	2.9
1	B	181	LYS	2.9
1	B	204	GLY	2.9
3	P	172	GLY	2.9
1	B	62	ASN	2.9
1	B	226	GLN	2.8
3	O	161	GLY	2.8
1	B	53	LYS	2.8
2	H	27	MET	2.8
2	N	190[A]	ASP	2.8
2	J	92	TYR	2.8
1	E	226	GLN	2.8
3	P	52	ASN	2.8
3	Q	57	ARG	2.8
2	H	93	MET	2.7
1	E	206	GLU	2.7
1	D	181	LYS	2.7
2	K	197	ARG	2.7
1	D	63	SER	2.7
3	P	231	SER	2.7
2	J	9	LYS	2.6
1	A	207	LEU	2.6
1	C	61	GLN	2.6
2	J	197	ARG	2.6
3	Q	175	GLN	2.6
1	C	202	GLU	2.6
2	I	93	MET	2.6
1	B	205	GLU	2.6
3	R	230	VAL	2.6
1	B	233	LEU	2.5
1	G	226	GLN	2.5
3	R	175	GLN	2.5
1	F	175	PHE	2.5
1	B	230	LYS	2.5
2	L	93	MET	2.5
2	M	197	ARG	2.5
3	U	161	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
3	R	54	THR	2.5
1	C	54	VAL	2.5
2	J	93	MET	2.5
3	T	173	GLY	2.5
1	E	202	GLU	2.5
2	J	26	ILE	2.5
1	E	127	GLY	2.5
1	G	181	LYS	2.4
2	K	27	MET	2.4
1	B	227	GLU	2.4
2	J	27	MET	2.4
2	L	197	ARG	2.4
2	M	22	MET	2.4
1	A	61	GLN	2.4
1	C	232	PHE	2.4
1	D	53	LYS	2.4
1	D	175	PHE	2.4
1	F	209	ALA	2.4
2	I	9	LYS	2.4
1	D	205	GLU	2.4
1	F	201	LEU	2.3
1	A	230	LYS	2.3
1	A	208	LYS	2.3
3	R	161	GLY	2.3
1	F	227	GLU	2.3
2	M	190[A]	ASP	2.3
2	N	145	LYS	2.3
3	S	57	ARG	2.3
1	D	229	VAL	2.3
1	A	181	LYS	2.3
2	J	198	LYS	2.3
2	K	190[A]	ASP	2.3
3	S	225	GLY	2.3
1	D	209	ALA	2.3
3	S	175	GLN	2.3
3	S	231	SER	2.3
3	R	173	GLY	2.2
1	D	178	ARG	2.2
3	R	57	ARG	2.2
3	T	175	GLN	2.2
1	C	44	ASN	2.2
3	S	224	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	207	LEU	2.2
2	L	22	MET	2.2
3	O	45	GLU	2.2
3	O	65	GLU	2.2
2	M	9	LYS	2.2
2	H	197	ARG	2.2
2	N	22	MET	2.2
1	F	191	THR	2.2
1	A	54	VAL	2.2
2	J	58[A]	TYR	2.1
2	K	22	MET	2.1
1	B	187	LYS	2.1
3	P	68	LEU	2.1
1	C	227	GLU	2.1
1	F	178	ARG	2.1
1	B	44	ASN	2.1
1	G	183	ASN	2.1
1	F	53	LYS	2.1
1	C	127	GLY	2.1
2	K	92	TYR	2.0
3	U	230	VAL	2.0
1	E	233	LEU	2.0
1	E	227	GLU	2.0
1	G	62	ASN	2.0
1	F	230	LYS	2.0
3	O	41	LYS	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 monosaccharides 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(\AA^2)	Q<0.9
4	SO4	F	4005	5/5	0.83	0.19	106,106,106,106	0
4	SO4	D	4004	5/5	0.88	0.16	86,86,87,87	0
4	SO4	A	4001	5/5	0.90	0.21	105,106,106,106	0
4	SO4	B	4002	5/5	0.91	0.12	68,69,69,69	0
5	GOL	H	5004	6/6	0.92	0.12	33,37,38,38	0
4	SO4	F	4006	5/5	0.93	0.10	68,69,69,70	0
4	SO4	G	4007	5/5	0.93	0.13	70,71,72,72	0
4	SO4	C	4003	5/5	0.93	0.13	97,98,98,98	0
5	GOL	J	5006	6/6	0.94	0.16	35,37,39,43	0
5	GOL	N	5003	6/6	0.94	0.10	30,33,36,38	0
5	GOL	I	5005	6/6	0.95	0.11	31,34,35,40	0
5	GOL	K	5007	6/6	0.96	0.09	32,36,37,38	0
5	GOL	L	5001	6/6	0.96	0.12	35,36,37,39	0
5	GOL	M	5002	6/6	0.96	0.09	35,36,37,38	0
4	SO4	M	4010	5/5	0.96	0.12	43,46,47,49	0
4	SO4	I	4011	5/5	0.97	0.12	35,40,41,43	0
4	SO4	J	4017	5/5	0.97	0.10	50,51,52,52	0
4	SO4	K	4018	5/5	0.97	0.08	46,49,51,51	0
4	SO4	L	4019	5/5	0.97	0.10	51,54,54,55	0
4	SO4	B	4016	5/5	0.97	0.10	54,55,55,56	0
4	SO4	H	4015	5/5	0.97	0.10	45,47,48,49	0
4	SO4	J	4012	5/5	0.98	0.08	34,38,39,41	0
4	SO4	L	4014	5/5	0.98	0.08	36,40,40,40	0
4	SO4	H	4008	5/5	0.98	0.08	38,40,42,42	0
4	SO4	K	4013	5/5	0.98	0.12	36,39,41,42	0
4	SO4	M	4020	5/5	0.98	0.09	51,52,53,53	0
4	SO4	N	4009	5/5	0.98	0.07	38,39,42,42	0
4	SO4	N	4021	5/5	0.98	0.10	52,53,54,54	0

6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.