



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2023 – 04:20 AM JST

PDB ID : 6KR1
Title : ATP dependent protease HslV from Staphylococcus aureus
Authors : Ha, N.-C.; Jeong, S.
Deposited on : 2019-08-20
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

<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 : 2.36
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) : 2.36

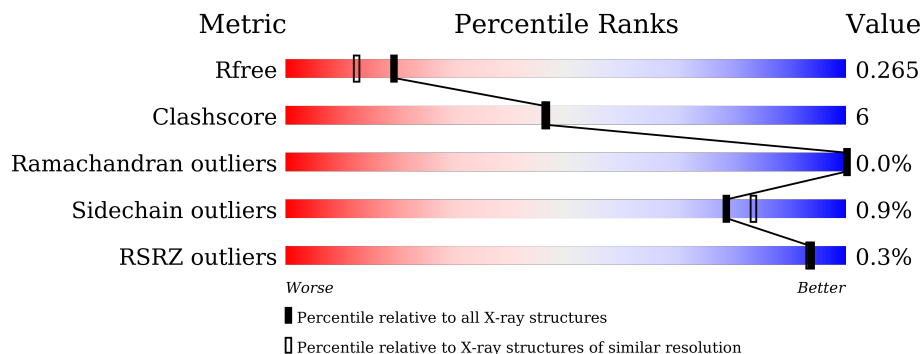
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	191	
1	B	191	
1	C	191	
1	D	191	
1	E	191	
1	F	191	

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Mol	Chain	Length	Quality of chain
1	G	191	<p>% 76% 16% 8%</p>
1	H	191	<p>82% 6% 12%</p>
1	I	191	<p>69% 16% 14%</p>
1	J	191	<p>% 80% 12% 8%</p>
1	K	191	<p>75% 13% 13%</p>
1	L	191	<p>% 75% 15% 9%</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16584 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 ATP-dependent protease subunit HslV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	176	1336	839	236	255	6	0	0	0
1	B	168	1275	802	224	243	6	0	0	0
1	C	168	1272	798	225	243	6	0	0	0
1	D	177	1343	843	237	257	6	0	0	0
1	E	168	1276	803	225	242	6	0	0	0
1	F	169	1280	802	226	246	6	0	0	0
1	G	176	1336	839	236	255	6	0	0	0
1	H	168	1272	798	225	243	6	0	0	0
1	I	164	1243	782	216	239	6	0	0	0
1	J	176	1336	839	236	255	6	0	0	0
1	K	167	1266	795	224	242	5	0	0	0
1	L	173	1313	824	231	252	6	0	0	0

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	MET	-	expression tag	UNP P65796
A	-8	HIS	-	expression tag	UNP P65796
A	-7	HIS	-	expression tag	UNP P65796
A	-6	HIS	-	expression tag	UNP P65796
A	-5	HIS	-	expression tag	UNP P65796

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	HIS	-	expression tag	UNP P65796
A	-3	GLY	-	expression tag	UNP P65796
A	-2	ALA	-	expression tag	UNP P65796
A	-1	ALA	-	expression tag	UNP P65796
A	0	SER	-	expression tag	UNP P65796
B	-9	MET	-	expression tag	UNP P65796
B	-8	HIS	-	expression tag	UNP P65796
B	-7	HIS	-	expression tag	UNP P65796
B	-6	HIS	-	expression tag	UNP P65796
B	-5	HIS	-	expression tag	UNP P65796
B	-4	HIS	-	expression tag	UNP P65796
B	-3	GLY	-	expression tag	UNP P65796
B	-2	ALA	-	expression tag	UNP P65796
B	-1	ALA	-	expression tag	UNP P65796
B	0	SER	-	expression tag	UNP P65796
C	-9	MET	-	expression tag	UNP P65796
C	-8	HIS	-	expression tag	UNP P65796
C	-7	HIS	-	expression tag	UNP P65796
C	-6	HIS	-	expression tag	UNP P65796
C	-5	HIS	-	expression tag	UNP P65796
C	-4	HIS	-	expression tag	UNP P65796
C	-3	GLY	-	expression tag	UNP P65796
C	-2	ALA	-	expression tag	UNP P65796
C	-1	ALA	-	expression tag	UNP P65796
C	0	SER	-	expression tag	UNP P65796
D	-9	MET	-	expression tag	UNP P65796
D	-8	HIS	-	expression tag	UNP P65796
D	-7	HIS	-	expression tag	UNP P65796
D	-6	HIS	-	expression tag	UNP P65796
D	-5	HIS	-	expression tag	UNP P65796
D	-4	HIS	-	expression tag	UNP P65796
D	-3	GLY	-	expression tag	UNP P65796
D	-2	ALA	-	expression tag	UNP P65796
D	-1	ALA	-	expression tag	UNP P65796
D	0	SER	-	expression tag	UNP P65796
E	-9	MET	-	expression tag	UNP P65796
E	-8	HIS	-	expression tag	UNP P65796
E	-7	HIS	-	expression tag	UNP P65796
E	-6	HIS	-	expression tag	UNP P65796
E	-5	HIS	-	expression tag	UNP P65796
E	-4	HIS	-	expression tag	UNP P65796
E	-3	GLY	-	expression tag	UNP P65796

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	ALA	-	expression tag	UNP P65796
E	-1	ALA	-	expression tag	UNP P65796
E	0	SER	-	expression tag	UNP P65796
F	-9	MET	-	expression tag	UNP P65796
F	-8	HIS	-	expression tag	UNP P65796
F	-7	HIS	-	expression tag	UNP P65796
F	-6	HIS	-	expression tag	UNP P65796
F	-5	HIS	-	expression tag	UNP P65796
F	-4	HIS	-	expression tag	UNP P65796
F	-3	GLY	-	expression tag	UNP P65796
F	-2	ALA	-	expression tag	UNP P65796
F	-1	ALA	-	expression tag	UNP P65796
F	0	SER	-	expression tag	UNP P65796
G	-9	MET	-	expression tag	UNP P65796
G	-8	HIS	-	expression tag	UNP P65796
G	-7	HIS	-	expression tag	UNP P65796
G	-6	HIS	-	expression tag	UNP P65796
G	-5	HIS	-	expression tag	UNP P65796
G	-4	HIS	-	expression tag	UNP P65796
G	-3	GLY	-	expression tag	UNP P65796
G	-2	ALA	-	expression tag	UNP P65796
G	-1	ALA	-	expression tag	UNP P65796
G	0	SER	-	expression tag	UNP P65796
H	-9	MET	-	expression tag	UNP P65796
H	-8	HIS	-	expression tag	UNP P65796
H	-7	HIS	-	expression tag	UNP P65796
H	-6	HIS	-	expression tag	UNP P65796
H	-5	HIS	-	expression tag	UNP P65796
H	-4	HIS	-	expression tag	UNP P65796
H	-3	GLY	-	expression tag	UNP P65796
H	-2	ALA	-	expression tag	UNP P65796
H	-1	ALA	-	expression tag	UNP P65796
H	0	SER	-	expression tag	UNP P65796
I	-9	MET	-	expression tag	UNP P65796
I	-8	HIS	-	expression tag	UNP P65796
I	-7	HIS	-	expression tag	UNP P65796
I	-6	HIS	-	expression tag	UNP P65796
I	-5	HIS	-	expression tag	UNP P65796
I	-4	HIS	-	expression tag	UNP P65796
I	-3	GLY	-	expression tag	UNP P65796
I	-2	ALA	-	expression tag	UNP P65796
I	-1	ALA	-	expression tag	UNP P65796

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Chain	Residue	Modelled	Actual	Comment	Reference
I	0	SER	-	expression tag	UNP P65796
J	-9	MET	-	expression tag	UNP P65796
J	-8	HIS	-	expression tag	UNP P65796
J	-7	HIS	-	expression tag	UNP P65796
J	-6	HIS	-	expression tag	UNP P65796
J	-5	HIS	-	expression tag	UNP P65796
J	-4	HIS	-	expression tag	UNP P65796
J	-3	GLY	-	expression tag	UNP P65796
J	-2	ALA	-	expression tag	UNP P65796
J	-1	ALA	-	expression tag	UNP P65796
J	0	SER	-	expression tag	UNP P65796
K	-9	MET	-	expression tag	UNP P65796
K	-8	HIS	-	expression tag	UNP P65796
K	-7	HIS	-	expression tag	UNP P65796
K	-6	HIS	-	expression tag	UNP P65796
K	-5	HIS	-	expression tag	UNP P65796
K	-4	HIS	-	expression tag	UNP P65796
K	-3	GLY	-	expression tag	UNP P65796
K	-2	ALA	-	expression tag	UNP P65796
K	-1	ALA	-	expression tag	UNP P65796
K	0	SER	-	expression tag	UNP P65796
L	-9	MET	-	expression tag	UNP P65796
L	-8	HIS	-	expression tag	UNP P65796
L	-7	HIS	-	expression tag	UNP P65796
L	-6	HIS	-	expression tag	UNP P65796
L	-5	HIS	-	expression tag	UNP P65796
L	-4	HIS	-	expression tag	UNP P65796
L	-3	GLY	-	expression tag	UNP P65796
L	-2	ALA	-	expression tag	UNP P65796
L	-1	ALA	-	expression tag	UNP P65796
L	0	SER	-	expression tag	UNP P65796

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	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	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	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		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	94	Total	O	0	0
			94	94		
3	B	80	Total	O	0	0
			80	80		
3	C	77	Total	O	0	0
			77	77		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	101	Total 101	O 101	0	0
3	E	60	Total 60	O 60	0	0
3	F	85	Total 85	O 85	0	0
3	G	103	Total 103	O 103	0	0
3	H	79	Total 79	O 79	0	0
3	I	72	Total 72	O 72	0	0
3	J	89	Total 89	O 89	0	0
3	K	86	Total 86	O 86	0	0
3	L	65	Total 65	O 65	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.

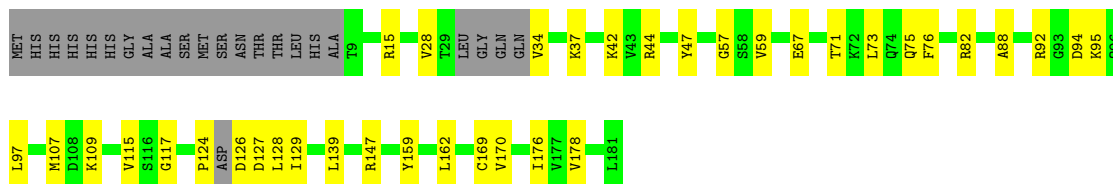
- Molecule 1: ATP-dependent protease subunit HslV

Chain A:  82% 10% 8%



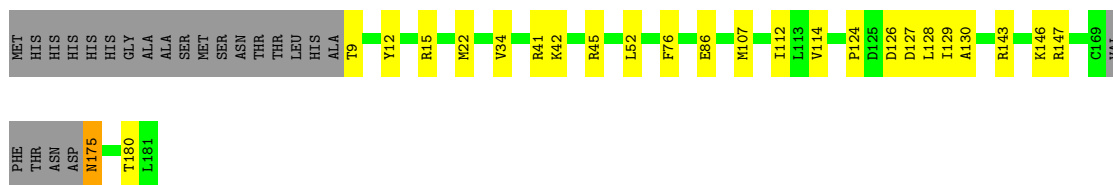
- Molecule 1: ATP-dependent protease subunit HslV

Chain B:  69% 19% 12%




- Molecule 1: ATP-dependent protease subunit HslV

Chain C:  75% 13% 12%




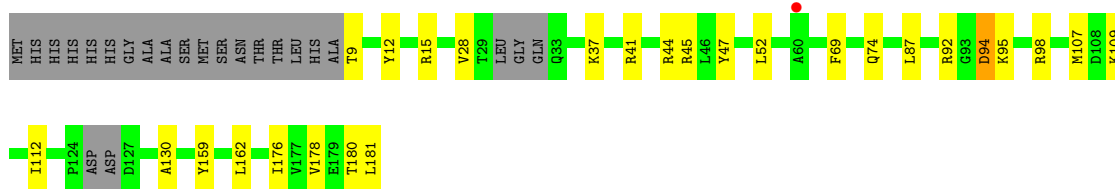
- Molecule 1: ATP-dependent protease subunit HslV

Chain D:  85% 7% 7%



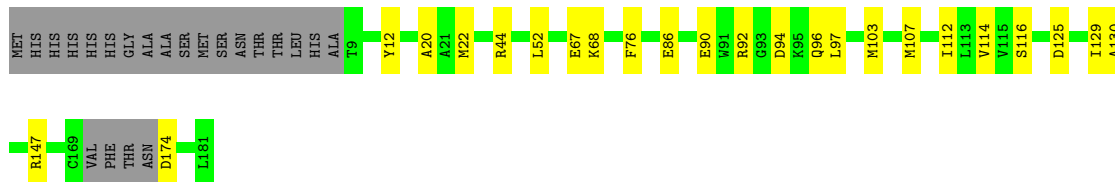
- Molecule 1: ATP-dependent protease subunit HslV

Chain E:  74% 14% 12%



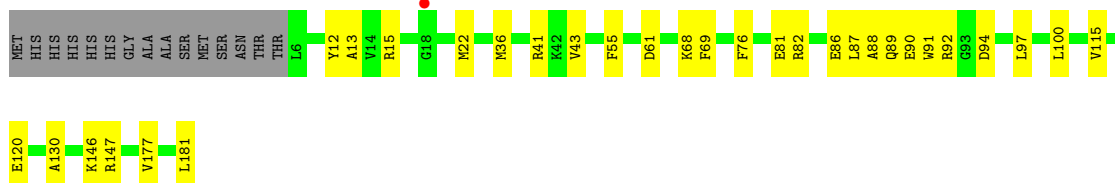
- Molecule 1: ATP-dependent protease subunit HslV

Chain F: 76% 13% 12%



- Molecule 1: ATP-dependent protease subunit HslV

Chain G: 76% 16% 8%



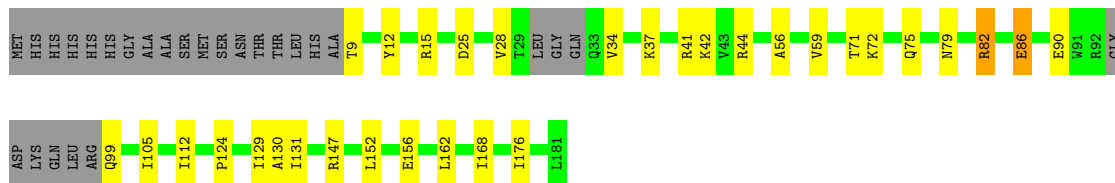
- Molecule 1: ATP-dependent protease subunit HslV

Chain H: 82% 6% 12%



- Molecule 1: ATP-dependent protease subunit HslV

Chain I: 69% 16% 14%



- Molecule 1: ATP-dependent protease subunit HslV

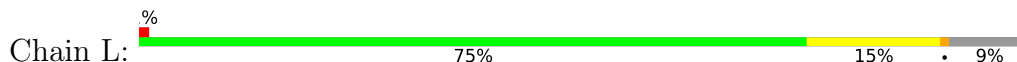
Chain J: 80% 12% 8%



• Molecule 1: ATP-dependent protease subunit HslV



• Molecule 1: ATP-dependent protease subunit HslV



4 Data and refinement statistics i

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	94.30Å 94.30Å 227.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.16 – 2.00 33.16 – 2.00	Depositor EDS
% Data completeness (in resolution range)	82.3 (33.16-2.00) 82.3 (33.16-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.16_3549, PHENIX 1.16_3549	Depositor
R, R_{free}	0.216 , 0.266 0.216 , 0.265	Depositor DCC
R_{free} test set	6166 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.480 for -h,-k,l 0.487 for h,-h-k,-l 0.487 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16584	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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: 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	0.24	0/1352	0.42	0/1822
1	B	0.24	0/1288	0.42	0/1732
1	C	0.23	0/1285	0.41	0/1728
1	D	0.23	0/1359	0.43	0/1832
1	E	0.24	0/1289	0.42	0/1733
1	F	0.23	0/1293	0.41	0/1739
1	G	0.24	0/1352	0.41	0/1822
1	H	0.24	0/1285	0.41	0/1728
1	I	0.24	0/1256	0.41	0/1691
1	J	0.24	0/1352	0.42	0/1822
1	K	0.24	0/1279	0.42	0/1720
1	L	0.23	0/1328	0.41	0/1789
All	All	0.24	0/15718	0.42	0/21158

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1336	0	1350	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1275	0	1291	27	0
1	C	1272	0	1291	17	0
1	D	1343	0	1357	10	0
1	E	1276	0	1295	21	0
1	F	1280	0	1295	15	0
1	G	1336	0	1350	23	0
1	H	1272	0	1291	9	0
1	I	1243	0	1251	21	0
1	J	1336	0	1350	17	0
1	K	1266	0	1286	15	0
1	L	1313	0	1327	21	0
2	A	10	0	0	0	0
2	D	10	0	0	0	0
2	G	10	0	0	0	0
2	H	5	0	0	0	0
2	J	5	0	0	0	0
2	K	5	0	0	0	0
3	A	94	0	0	2	0
3	B	80	0	0	3	0
3	C	77	0	0	5	0
3	D	101	0	0	2	0
3	E	60	0	0	5	0
3	F	85	0	0	3	0
3	G	103	0	0	5	0
3	H	79	0	0	4	0
3	I	72	0	0	3	0
3	J	89	0	0	2	0
3	K	86	0	0	4	0
3	L	65	0	0	5	0
All	All	16584	0	15734	197	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 (197) 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:L:15:ARG:HB2	1:L:127:ASP:HB3	1.64	0.80
1:G:120:GLU:HG2	1:L:30:LEU:HD21	1.65	0.78
1:L:71:THR:HG22	1:L:75:GLN:HE21	1.48	0.77
1:L:107:MET:SD	3:L:202:HOH:O	2.43	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:95:LYS:HD3	1:J:98:ARG:HD3	1.71	0.72
1:K:37:LYS:NZ	3:K:302:HOH:O	2.22	0.71
1:K:19:LYS:NZ	3:K:301:HOH:O	2.18	0.71
1:E:162:LEU:HD23	1:E:176:ILE:HG12	1.75	0.68
1:H:22:MET:SD	3:H:358:HOH:O	2.53	0.67
1:G:147:ARG:NH1	3:G:302:HOH:O	2.27	0.66
1:C:22:MET:SD	3:C:253:HOH:O	2.52	0.66
1:L:108:ASP:O	3:L:201:HOH:O	2.15	0.65
1:C:45:ARG:NH2	1:C:180:THR:O	2.29	0.64
1:F:44:ARG:NH1	3:F:201:HOH:O	2.31	0.64
1:C:22:MET:HE1	1:C:52:LEU:HB3	1.79	0.63
1:L:28:VAL:HB	1:L:37:LYS:HB2	1.79	0.63
1:B:34:VAL:N	3:B:205:HOH:O	2.32	0.63
1:I:99:GLN:N	3:I:204:HOH:O	2.32	0.62
1:B:15:ARG:HB2	1:B:127:ASP:HB3	1.81	0.62
1:D:122:ILE:HD12	1:E:37:LYS:HD2	1.80	0.62
1:F:103:MET:HG2	1:F:116:SER:HB3	1.81	0.62
1:B:82:ARG:HG2	1:C:34:VAL:HG21	1.82	0.61
1:E:95:LYS:HA	1:E:98:ARG:HD3	1.82	0.61
1:A:44:ARG:HA	1:A:44:ARG:HE	1.65	0.61
1:F:22:MET:HE1	1:F:52:LEU:HB3	1.81	0.60
1:H:22:MET:HE1	1:H:52:LEU:HB3	1.84	0.60
1:B:169:CYS:SG	3:G:314:HOH:O	2.56	0.60
1:I:15:ARG:HG3	1:I:112:ILE:HD11	1.83	0.60
1:C:114:VAL:HG21	1:C:129:ILE:HG12	1.85	0.59
1:G:13:ALA:HB2	1:G:22:MET:HG3	1.86	0.58
1:G:41:ARG:HH21	1:G:177:VAL:HG11	1.69	0.58
1:J:15:ARG:O	1:J:146:LYS:NZ	2.36	0.58
1:E:44:ARG:NH2	3:E:206:HOH:O	2.37	0.58
1:H:41:ARG:NH2	3:H:302:HOH:O	2.36	0.57
1:C:175:ASN:ND2	3:C:203:HOH:O	2.33	0.57
1:E:41:ARG:NH2	3:E:201:HOH:O	2.37	0.57
1:A:13:ALA:HB2	1:A:22:MET:HG3	1.85	0.57
1:E:41:ARG:NH1	3:E:205:HOH:O	2.37	0.56
1:G:15:ARG:O	1:G:146:LYS:NZ	2.37	0.56
1:K:45:ARG:HG2	1:K:49:GLY:HA2	1.86	0.56
1:C:15:ARG:HB2	1:C:127:ASP:HB3	1.87	0.55
1:K:114:VAL:HG21	1:K:129:ILE:HG12	1.88	0.55
1:B:124:PRO:O	1:B:126:ASP:N	2.39	0.55
1:I:41:ARG:NH2	3:I:208:HOH:O	2.39	0.55
1:I:71:THR:HG22	1:I:75:GLN:OE1	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:94:ASP:HB3	1:G:97:LEU:HB2	1.89	0.55
1:A:147:ARG:NH1	3:A:307:HOH:O	2.40	0.55
1:F:174:ASP:N	3:F:207:HOH:O	2.40	0.55
1:D:128:LEU:HD21	1:D:143:ARG:HE	1.72	0.54
1:E:52:LEU:HG	1:E:181:LEU:HD13	1.87	0.54
1:L:143:ARG:NH1	3:L:203:HOH:O	2.40	0.54
1:G:88:ALA:HB2	1:G:115:VAL:HG13	1.88	0.54
1:F:114:VAL:HG21	1:F:129:ILE:HG12	1.88	0.54
1:B:94:ASP:HB3	1:B:97:LEU:HB2	1.90	0.53
1:E:28:VAL:HB	1:E:37:LYS:HB2	1.90	0.53
1:B:44:ARG:NH2	3:B:211:HOH:O	2.41	0.52
1:K:33:GLN:HE22	1:K:37:LYS:NZ	2.07	0.52
1:I:162:LEU:HD23	1:I:176:ILE:HG12	1.91	0.52
1:J:7:HIS:HB2	1:J:29:THR:HB	1.91	0.52
1:E:159:TYR:HB2	1:E:178:VAL:HG21	1.91	0.52
1:I:124:PRO:HG2	1:I:129:ILE:HG21	1.91	0.52
1:I:152:LEU:HB3	1:I:156:GLU:HG3	1.92	0.52
1:L:50:LYS:HE3	1:L:109:LYS:HE2	1.92	0.52
1:G:81:GLU:OE2	3:G:301:HOH:O	2.18	0.52
1:B:159:TYR:HB2	1:B:178:VAL:HG21	1.91	0.51
1:D:5:THR:N	3:D:312:HOH:O	2.43	0.51
1:G:12:TYR:HD1	1:G:130:ALA:HB2	1.74	0.51
1:H:179:GLU:OE2	3:H:301:HOH:O	2.19	0.51
1:L:19:LYS:NZ	3:L:204:HOH:O	2.43	0.51
1:L:70:GLU:O	1:L:74:GLN:HG2	2.11	0.51
1:C:147:ARG:NH1	3:C:208:HOH:O	2.44	0.50
1:H:105:ILE:HG12	1:H:114:VAL:HG22	1.93	0.50
1:I:72:LYS:HE2	1:I:86:GLU:HG2	1.94	0.50
1:G:89:GLN:HB3	1:L:96:GLN:OE1	2.11	0.50
1:E:15:ARG:HD3	1:E:112:ILE:HG13	1.93	0.50
1:F:94:ASP:HB3	1:F:97:LEU:HB2	1.94	0.50
1:D:15:ARG:O	1:D:146:LYS:NZ	2.45	0.50
1:G:92:ARG:HD2	1:L:99:GLN:HE21	1.76	0.50
3:C:207:HOH:O	1:I:147:ARG:NH1	2.44	0.49
1:D:81:GLU:OE2	3:D:301:HOH:O	2.19	0.49
1:K:44:ARG:NH2	3:K:309:HOH:O	2.45	0.49
1:G:41:ARG:NH2	3:G:313:HOH:O	2.46	0.49
1:J:61:ASP:OD1	1:K:92:ARG:NH1	2.46	0.49
1:L:162:LEU:HD23	1:L:176:ILE:HG12	1.94	0.49
1:K:9:THR:HG21	1:K:42:LYS:HD2	1.95	0.48
1:I:99:GLN:NE2	3:I:213:HOH:O	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:ASP:N	1:B:126:ASP:OD1	2.46	0.48
1:K:155:GLU:HG3	1:K:178:VAL:HG11	1.95	0.48
1:G:97:LEU:HA	1:G:100:LEU:HD13	1.95	0.48
1:J:67:GLU:OE2	3:J:301:HOH:O	2.19	0.48
1:A:22:MET:HE2	1:A:43:VAL:HG22	1.96	0.48
1:I:72:LYS:NZ	1:I:90:GLU:OE1	2.37	0.48
1:J:28:VAL:HG22	1:J:37:LYS:HB3	1.95	0.48
1:J:76:PHE:CE2	1:J:86:GLU:HG3	2.49	0.47
1:I:28:VAL:HG22	1:I:37:LYS:HB2	1.96	0.47
1:L:42:LYS:NZ	1:L:57:GLY:O	2.37	0.47
1:F:147:ARG:NE	3:F:210:HOH:O	2.42	0.47
1:K:105:ILE:HG12	1:K:114:VAL:HG22	1.96	0.47
1:B:147:ARG:NH1	3:H:305:HOH:O	2.49	0.46
1:E:52:LEU:HB2	1:E:107:MET:HG2	1.97	0.46
1:J:95:LYS:HD3	1:J:98:ARG:CD	2.45	0.46
1:A:37:LYS:HG3	1:F:125:ASP:OD1	2.15	0.46
1:F:76:PHE:CZ	1:F:86:GLU:HG3	2.51	0.46
1:F:68:LYS:NZ	1:F:90:GLU:OE1	2.48	0.46
1:B:42:LYS:NZ	1:B:57:GLY:O	2.29	0.46
1:D:95:LYS:HA	1:D:98:ARG:HG2	1.96	0.45
1:G:61:ASP:HB3	1:G:100:LEU:HD21	1.98	0.45
1:B:128:LEU:HD23	1:B:139:LEU:HD12	1.98	0.45
1:C:107:MET:HE3	1:C:112:ILE:HG12	1.99	0.45
1:G:91:TRP:CD1	1:G:97:LEU:HB3	2.52	0.45
1:L:159:TYR:HB2	1:L:178:VAL:HG21	1.97	0.45
1:C:124:PRO:HG3	1:C:129:ILE:HB	1.99	0.45
1:E:181:LEU:HD12	1:E:181:LEU:HA	1.84	0.45
1:F:12:TYR:HD2	1:F:130:ALA:HB2	1.81	0.45
1:E:92:ARG:NH1	1:F:67:GLU:OE2	2.49	0.45
1:E:94:ASP:O	1:E:98:ARG:HG3	2.16	0.45
1:I:42:LYS:HE3	1:I:59:VAL:HG22	1.99	0.45
1:E:12:TYR:HD1	1:E:130:ALA:HB2	1.81	0.45
1:E:69:PHE:CD1	1:E:87:LEU:HD12	2.52	0.45
1:J:12:TYR:HD1	1:J:130:ALA:HB2	1.82	0.44
1:J:95:LYS:HA	1:J:98:ARG:HD3	1.98	0.44
1:A:45:ARG:NH1	1:A:179:GLU:OE1	2.50	0.44
1:A:76:PHE:CE2	1:A:86:GLU:HG3	2.53	0.44
1:I:9:THR:HG22	1:I:56:ALA:HA	1.99	0.44
1:F:76:PHE:CE2	1:F:86:GLU:HG3	2.52	0.44
1:B:128:LEU:CD2	1:B:139:LEU:HD12	2.48	0.44
1:D:168:ILE:HG23	1:I:34:VAL:HG13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:LYS:HA	1:A:109:LYS:HD2	1.84	0.44
1:B:42:LYS:HE2	1:B:59:VAL:HG12	1.98	0.44
1:B:47:TYR:HB2	1:B:73:LEU:HD13	2.00	0.44
1:E:9:THR:N	3:E:213:HOH:O	2.51	0.44
1:C:76:PHE:CE1	1:C:86:GLU:HG3	2.53	0.44
1:G:55:PHE:HZ	1:G:100:LEU:HD23	1.83	0.44
1:K:76:PHE:CE2	1:K:86:GLU:HG3	2.53	0.44
1:I:9:THR:OG1	1:I:25:ASP:OD1	2.28	0.43
1:B:92:ARG:NH2	3:B:204:HOH:O	2.32	0.43
1:I:105:ILE:HG13	1:I:131:ILE:HD13	1.99	0.43
1:B:92:ARG:HH21	1:B:117:GLY:C	2.22	0.43
1:J:58:SER:OG	1:J:61:ASP:OD2	2.36	0.43
1:B:162:LEU:HB3	1:B:176:ILE:HD13	2.00	0.43
1:L:129:ILE:HD12	1:L:129:ILE:HA	1.86	0.43
1:B:88:ALA:HB2	1:B:115:VAL:HG13	2.00	0.43
1:J:129:ILE:HD12	1:J:129:ILE:HA	1.89	0.43
1:L:41:ARG:NH2	3:L:207:HOH:O	2.51	0.43
3:A:303:HOH:O	1:B:37:LYS:NZ	2.41	0.43
1:G:76:PHE:CE2	1:G:86:GLU:HG3	2.54	0.43
1:G:36:MET:SD	1:H:122:ILE:HD12	2.59	0.43
1:B:124:PRO:HG3	1:B:129:ILE:HG21	2.01	0.42
1:F:20:ALA:HB1	1:F:107:MET:CE	2.49	0.42
1:C:143:ARG:NH1	1:I:168:ILE:HG12	2.34	0.42
1:K:12:TYR:HD2	1:K:130:ALA:HB2	1.85	0.42
1:A:44:ARG:HA	1:A:44:ARG:NE	2.34	0.42
1:D:76:PHE:CE2	1:D:86:GLU:HG3	2.55	0.42
1:B:170:VAL:N	3:G:314:HOH:O	2.50	0.42
1:A:61:ASP:HB3	1:A:100:LEU:HD21	2.01	0.42
1:C:12:TYR:HD2	1:C:130:ALA:HB2	1.85	0.42
1:C:126:ASP:OD1	1:C:126:ASP:N	2.45	0.42
1:E:45:ARG:NH1	1:E:180:THR:O	2.53	0.42
1:D:44:ARG:HD3	1:D:44:ARG:HA	1.86	0.42
1:G:76:PHE:CD1	1:G:82:ARG:HG2	2.54	0.42
1:H:105:ILE:HG13	1:H:131:ILE:HD13	2.01	0.42
1:A:162:LEU:HD23	1:A:162:LEU:HA	1.94	0.42
1:B:67:GLU:O	1:B:71:THR:HG23	2.20	0.41
1:G:181:LEU:HD12	1:G:181:LEU:HA	1.81	0.41
1:H:114:VAL:HG21	1:H:129:ILE:HG12	2.03	0.41
1:J:122:ILE:HD13	1:J:122:ILE:HA	1.89	0.41
1:H:91:TRP:CD1	1:H:97:LEU:HB3	2.55	0.41
1:J:120:GLU:HG3	1:J:122:ILE:HG12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:15:ARG:HB2	1:K:127:ASP:HB3	2.02	0.41
1:F:107:MET:HG3	1:F:112:ILE:HG12	2.02	0.41
1:G:22:MET:HE2	1:G:43:VAL:HG22	2.02	0.41
1:A:168:ILE:HG23	1:L:34:VAL:HG13	2.03	0.41
1:B:28:VAL:HB	1:B:37:LYS:HB2	2.03	0.41
1:G:68:LYS:HD3	1:G:90:GLU:OE2	2.19	0.41
1:I:79:ASN:OD1	1:I:82:ARG:HB2	2.20	0.41
1:I:162:LEU:HB3	1:I:176:ILE:HD13	2.02	0.41
1:K:22:MET:HE3	1:K:52:LEU:HB3	2.02	0.41
1:L:30:LEU:HD23	1:L:30:LEU:HA	1.87	0.41
1:B:15:ARG:NH1	1:B:109:LYS:O	2.54	0.41
1:C:41:ARG:HG2	1:C:42:LYS:HG3	2.02	0.41
1:E:44:ARG:NH1	3:E:218:HOH:O	2.54	0.41
1:E:109:LYS:H	1:E:109:LYS:HG2	1.72	0.41
1:L:15:ARG:HD3	1:L:112:ILE:HG13	2.03	0.41
1:L:76:PHE:CD2	1:L:82:ARG:HD2	2.56	0.41
1:A:105:ILE:HD13	1:A:129:ILE:HG23	2.02	0.41
1:A:122:ILE:HD12	1:B:37:LYS:HD3	2.01	0.41
1:C:128:LEU:HD12	1:C:146:LYS:HD2	2.03	0.41
1:A:23:ALA:HB2	1:A:178:VAL:HG22	2.03	0.40
1:D:136:ASN:HB3	1:J:137:TYR:CE2	2.56	0.40
1:J:68:LYS:O	1:J:72:LYS:HG2	2.21	0.40
1:J:82:ARG:NH2	3:J:308:HOH:O	2.43	0.40
1:K:175:ASN:N	3:K:314:HOH:O	2.55	0.40
1:C:9:THR:N	3:C:214:HOH:O	2.54	0.40
1:G:69:PHE:HB2	1:G:87:LEU:HD22	2.03	0.40
1:B:15:ARG:HG3	1:B:107:MET:HE2	2.02	0.40
1:E:47:TYR:HE2	1:E:74:GLN:HA	1.87	0.40
1:I:12:TYR:HD1	1:I:130:ALA:HB2	1.86	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	174/191 (91%)	167 (96%)	7 (4%)	0	100	100
1	B	162/191 (85%)	154 (95%)	7 (4%)	1 (1%)	25	19
1	C	164/191 (86%)	160 (98%)	4 (2%)	0	100	100
1	D	175/191 (92%)	167 (95%)	8 (5%)	0	100	100
1	E	162/191 (85%)	156 (96%)	6 (4%)	0	100	100
1	F	165/191 (86%)	162 (98%)	3 (2%)	0	100	100
1	G	174/191 (91%)	168 (97%)	6 (3%)	0	100	100
1	H	164/191 (86%)	160 (98%)	4 (2%)	0	100	100
1	I	158/191 (83%)	154 (98%)	4 (2%)	0	100	100
1	J	174/191 (91%)	166 (95%)	8 (5%)	0	100	100
1	K	163/191 (85%)	161 (99%)	2 (1%)	0	100	100
1	L	171/191 (90%)	166 (97%)	5 (3%)	0	100	100
All	All	2006/2292 (88%)	1941 (97%)	64 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	75	GLN

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	136/148 (92%)	136 (100%)	0	100	100
1	B	130/148 (88%)	128 (98%)	2 (2%)	65	69
1	C	129/148 (87%)	128 (99%)	1 (1%)	81	86
1	D	137/148 (93%)	137 (100%)	0	100	100
1	E	130/148 (88%)	129 (99%)	1 (1%)	81	86
1	F	130/148 (88%)	128 (98%)	2 (2%)	65	69
1	G	136/148 (92%)	136 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	129/148 (87%)	129 (100%)	0	100	100
1	I	127/148 (86%)	124 (98%)	3 (2%)	49	51
1	J	136/148 (92%)	134 (98%)	2 (2%)	65	69
1	K	128/148 (86%)	127 (99%)	1 (1%)	81	86
1	L	134/148 (90%)	132 (98%)	2 (2%)	65	69
All	All	1582/1776 (89%)	1568 (99%)	14 (1%)	78	83

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	76	PHE
1	B	95	LYS
1	C	175	ASN
1	E	94	ASP
1	F	92	ARG
1	F	96	GLN
1	I	44	ARG
1	I	82	ARG
1	I	86	GLU
1	J	36	MET
1	J	82	ARG
1	K	25	ASP
1	L	33	GLN
1	L	109	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	99	GLN
1	C	175	ASN
1	D	32	GLN
1	D	99	GLN
1	L	74	GLN
1	L	75	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	202	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	D	202	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	G	202	-	4,4,4	0.14	0	6,6,6	0.04	0
2	SO4	A	201	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	D	201	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	G	201	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	K	201	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	H	201	-	4,4,4	0.15	0	6,6,6	0.05	0
2	SO4	J	201	-	4,4,4	0.14	0	6,6,6	0.05	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	176/191 (92%)	-0.26	0 100 100	10, 25, 48, 65	0
1	B	168/191 (87%)	-0.16	0 100 100	15, 30, 60, 74	0
1	C	168/191 (87%)	-0.27	0 100 100	14, 28, 52, 70	0
1	D	177/191 (92%)	-0.20	1 (0%) 89 88	8, 25, 51, 66	0
1	E	168/191 (87%)	-0.13	1 (0%) 89 88	14, 31, 57, 65	0
1	F	169/191 (88%)	-0.24	0 100 100	14, 28, 56, 69	0
1	G	176/191 (92%)	-0.23	1 (0%) 89 88	9, 26, 50, 67	0
1	H	168/191 (87%)	-0.23	0 100 100	14, 29, 51, 67	0
1	I	164/191 (85%)	-0.14	0 100 100	16, 31, 55, 66	0
1	J	176/191 (92%)	-0.21	2 (1%) 80 79	9, 26, 51, 64	0
1	K	167/191 (87%)	-0.23	0 100 100	12, 28, 53, 67	0
1	L	173/191 (90%)	-0.14	1 (0%) 89 88	16, 32, 60, 70	0
All	All	2050/2292 (89%)	-0.20	6 (0%) 94 93	8, 28, 56, 74	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	31	GLY	3.1
1	D	18	GLY	2.2
1	E	60	ALA	2.2
1	J	93	GLY	2.2
1	L	75	GLN	2.1
1	G	18	GLY	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(Å ²)	Q<0.9
2	SO4	A	201	5/5	0.94	0.11	40,59,70,77	0
2	SO4	D	202	5/5	0.94	0.09	44,64,70,76	0
2	SO4	J	201	5/5	0.94	0.10	41,52,67,71	0
2	SO4	G	201	5/5	0.96	0.10	48,59,67,75	0
2	SO4	K	201	5/5	0.96	0.09	56,63,75,79	0
2	SO4	G	202	5/5	0.97	0.13	39,42,53,58	0
2	SO4	H	201	5/5	0.98	0.15	49,50,52,62	0
2	SO4	D	201	5/5	0.98	0.10	43,50,58,61	0
2	SO4	A	202	5/5	0.98	0.11	38,40,55,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.