



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 09:15 pm BST

PDB ID : 3ST9
Title : Crystal structure of ClpP in heptameric form from *Staphylococcus aureus*
Authors : Zhang, J.; Ye, F.; Lan, L.; Jiang, H.; Luo, C.; Yang, C.-G.
Deposited on : 2011-07-09
Resolution : 2.43 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 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 : 2.11
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) : 2.11

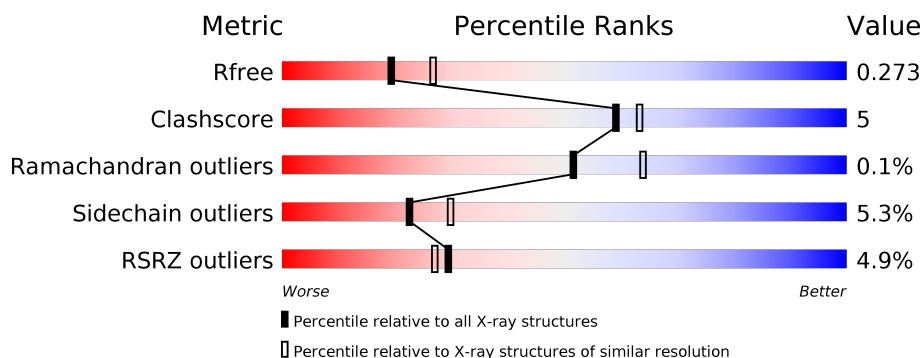
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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	1564 (2.46-2.42)
Clashscore	141614	1631 (2.46-2.42)
Ramachandran outliers	138981	1617 (2.46-2.42)
Sidechain outliers	138945	1617 (2.46-2.42)
RSRZ outliers	127900	1547 (2.46-2.42)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	197	<div> <div>3%</div> <div> <div></div> <div>76%</div> <div>9%</div> <div>•</div> <div>14%</div> </div> </div>
1	B	197	<div> <div>6%</div> <div> <div></div> <div>77%</div> <div>9%</div> <div>•</div> <div>14%</div> </div> </div>
1	C	197	<div> <div>3%</div> <div> <div></div> <div>71%</div> <div>15%</div> <div>•</div> <div>13%</div> </div> </div>
1	D	197	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>11%</div> <div>•</div> <div>13%</div> </div> </div>
1	E	197	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>9%</div> <div>•</div> <div>13%</div> </div> </div>
1	F	197	<div> <div>8%</div> <div> <div></div> <div>73%</div> <div>15%</div> <div>•</div> <div>12%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	197	 <p>3% 78% 10% 12%</p>

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	GOL	C	196	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9499 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 Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	1	0
			1319	831	224	257	7			
1	B	170	Total	C	N	O	S	0	0	0
			1314	827	224	257	6			
1	C	171	Total	C	N	O	S	0	1	0
			1335	842	226	260	7			
1	D	171	Total	C	N	O	S	0	1	0
			1335	842	226	260	7			
1	E	172	Total	C	N	O	S	0	1	0
			1333	839	227	260	7			
1	F	173	Total	C	N	O	S	0	1	0
			1344	845	228	264	7			
1	G	173	Total	C	N	O	S	0	2	0
			1352	851	231	263	7			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP P63786
A	0	SER	-	EXPRESSION TAG	UNP P63786
B	-1	GLY	-	EXPRESSION TAG	UNP P63786
B	0	SER	-	EXPRESSION TAG	UNP P63786
C	-1	GLY	-	EXPRESSION TAG	UNP P63786
C	0	SER	-	EXPRESSION TAG	UNP P63786
D	-1	GLY	-	EXPRESSION TAG	UNP P63786
D	0	SER	-	EXPRESSION TAG	UNP P63786
E	-1	GLY	-	EXPRESSION TAG	UNP P63786
E	0	SER	-	EXPRESSION TAG	UNP P63786
F	-1	GLY	-	EXPRESSION TAG	UNP P63786
F	0	SER	-	EXPRESSION TAG	UNP P63786
G	-1	GLY	-	EXPRESSION TAG	UNP P63786
G	0	SER	-	EXPRESSION TAG	UNP P63786

- 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	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

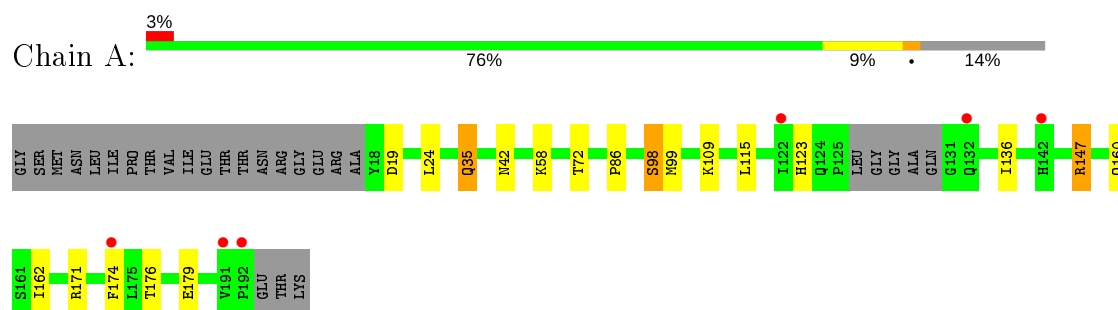
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	25	Total	O	0	0
			25	25		
5	B	18	Total	O	0	0
			18	18		
5	C	35	Total	O	0	0
			35	35		
5	D	17	Total	O	0	0
			17	17		
5	E	14	Total	O	0	0
			14	14		
5	F	24	Total	O	0	0
			24	24		
5	G	17	Total	O	0	0
			17	17		

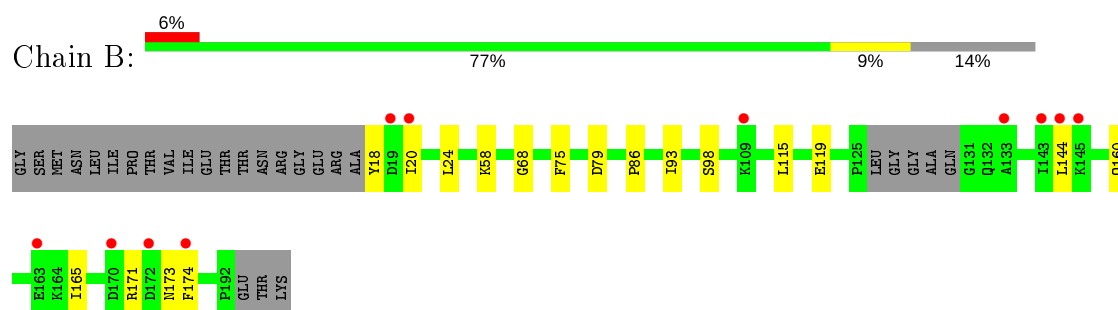
3 Residue-property plots [i](#)

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

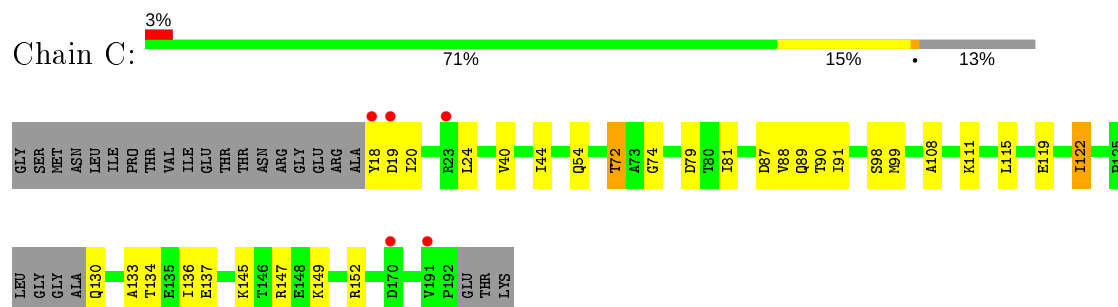
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit

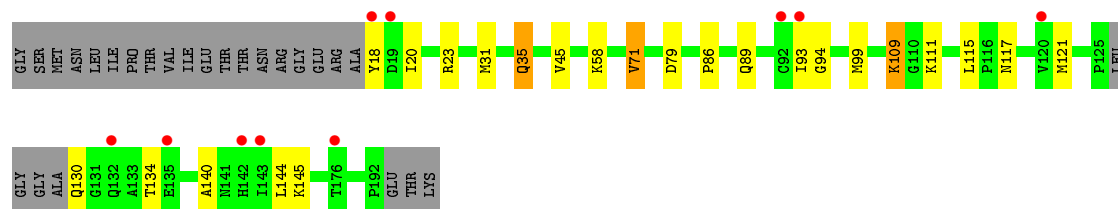


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

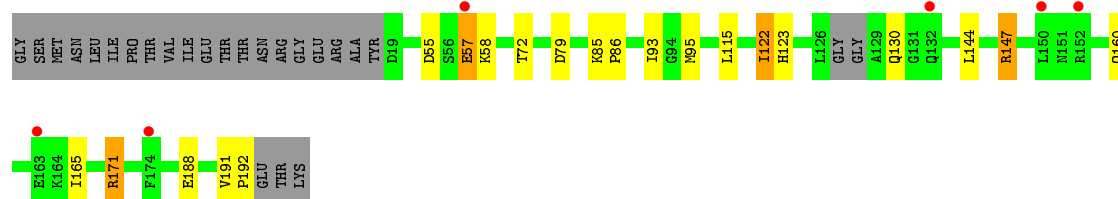
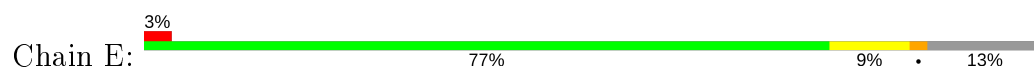


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

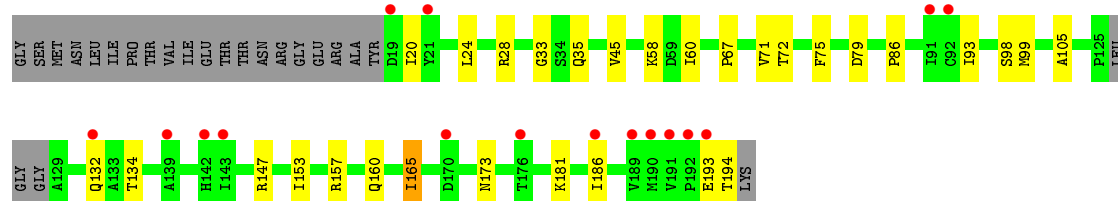




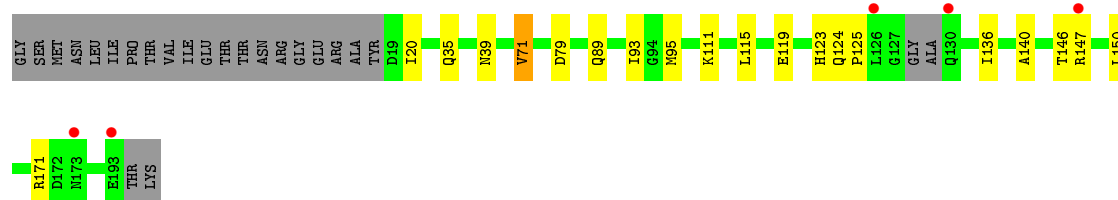
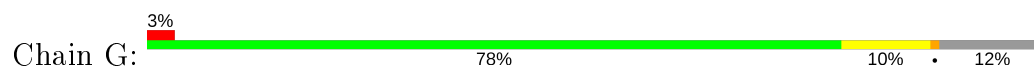
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.29 Å 121.29 Å 404.38 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.43 48.94 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.1 (20.00-2.43) 99.0 (48.94-2.43)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.42 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.239 , 0.271 0.237 , 0.273	Depositor DCC
R_{free} test set	3369 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9499	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CA, 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	0.36	0/1338	0.52	0/1803
1	B	0.35	0/1330	0.48	0/1793
1	C	0.35	0/1355	0.51	0/1826
1	D	0.36	0/1355	0.48	0/1826
1	E	0.34	0/1352	0.50	0/1822
1	F	0.37	0/1363	0.51	0/1837
1	G	0.35	0/1374	0.52	0/1850
All	All	0.35	0/9467	0.50	0/12757

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1319	0	1334	16	0
1	B	1314	0	1325	11	0
1	C	1335	0	1349	22	0
1	D	1335	0	1349	20	0
1	E	1333	0	1347	13	0
1	F	1344	0	1358	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1352	0	1373	11	0
2	A	5	0	0	0	0
2	D	5	0	0	0	0
3	B	1	0	0	0	0
4	C	6	0	8	2	0
5	A	25	0	0	1	0
5	B	18	0	0	0	0
5	C	35	0	0	2	0
5	D	17	0	0	0	0
5	E	14	0	0	0	0
5	F	24	0	0	0	0
5	G	17	0	0	0	0
All	All	9499	0	9443	88	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 5.

All (88) 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:A:147:ARG:HG3	1:A:147:ARG:HH11	1.13	1.07
1:D:109:LYS:HD2	1:D:109:LYS:H	1.31	0.93
1:D:35:GLN:H	1:D:35:GLN:HE21	1.22	0.88
1:D:109:LYS:N	1:D:109:LYS:HD2	1.91	0.83
1:A:147:ARG:CG	1:A:147:ARG:HH11	1.95	0.80
1:C:89:GLN:HG3	1:C:111:LYS:HB3	1.65	0.77
1:D:35:GLN:H	1:D:35:GLN:NE2	1.83	0.77
1:A:147:ARG:HG3	1:A:147:ARG:NH1	1.93	0.72
1:B:115:LEU:HD13	1:D:79:ASP:HB3	1.73	0.69
1:G:89:GLN:HG3	1:G:111:LYS:HB3	1.75	0.68
1:F:160:GLN:HB3	1:F:165:ILE:CD1	2.23	0.68
1:B:93:ILE:HG13	1:B:93:ILE:O	1.94	0.66
1:B:93:ILE:HD13	1:D:45:VAL:HG11	1.78	0.66
1:D:109:LYS:CD	1:D:109:LYS:H	2.09	0.63
1:C:149:LYS:HD3	1:D:117:ASN:ND2	2.15	0.62
1:C:134:THR:HG23	1:C:136:ILE:HG22	1.80	0.62
1:E:93:ILE:HD13	1:F:45:VAL:HG11	1.82	0.61
1:C:136:ILE:HD11	1:D:121:MET:HE2	1.85	0.59
1:A:98:SER:OG	1:A:99:MET:N	2.36	0.58
1:E:122:ILE:H	1:E:122:ILE:HD13	1.68	0.58
1:B:79:ASP:HB3	1:G:115:LEU:HD13	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:TYR:HB3	1:D:20:ILE:HG22	1.87	0.56
1:A:115:LEU:HD13	1:G:79:ASP:HB3	1.88	0.56
1:A:35:GLN:CD	1:A:35:GLN:H	2.10	0.55
1:D:89:GLN:HG3	1:D:111:LYS:HB3	1.89	0.54
1:G:71:VAL:HG21	1:G:146:THR:HG21	1.90	0.54
1:A:162:ILE:H	1:A:162:ILE:HD12	1.73	0.53
1:B:119:GLU:HG2	1:D:144:LEU:HD12	1.91	0.52
1:E:93:ILE:HG13	1:E:93:ILE:O	2.09	0.52
1:F:35:GLN:HG3	1:F:67:PRO:HG2	1.91	0.52
1:F:28:ARG:NH2	1:F:60:ILE:HD11	2.24	0.52
1:C:74:GLY:HA3	1:C:99:MET:HE2	1.91	0.51
1:C:98:SER:OG	1:C:99:MET:N	2.41	0.51
1:F:98:SER:OG	1:F:99:MET:N	2.43	0.51
1:A:176:THR:HG21	5:A:199:HOH:O	2.09	0.51
1:C:149:LYS:NZ	1:D:117:ASN:HD22	2.10	0.50
1:C:87:ASP:HB2	1:C:108:ALA:HB2	1.93	0.50
1:A:176:THR:CG2	1:A:179:GLU:H	2.26	0.49
1:C:136:ILE:HD11	1:D:121:MET:CE	2.42	0.49
1:E:147:ARG:H	1:E:147:ARG:HD3	1.78	0.49
1:E:55:ASP:OD1	1:E:57:GLU:HG2	2.13	0.49
1:F:58:LYS:O	1:F:86:PRO:HB3	2.13	0.49
1:G:123:HIS:HB2	1:G:171:ARG:O	2.14	0.48
4:C:196:GOL:O3	5:C:230:HOH:O	2.20	0.48
1:A:136:ILE:HD12	1:A:136:ILE:H	1.79	0.47
1:B:171:ARG:NE	1:B:173:ASN:HD21	2.12	0.47
1:F:105:ALA:O	1:F:157:ARG:HD3	2.15	0.47
1:G:147[B]:ARG:HA	1:G:150:LEU:HD12	1.96	0.47
1:D:31[B]:MET:HE2	1:D:93:ILE:HD11	1.97	0.47
1:F:20:ILE:HD12	1:F:20:ILE:H	1.79	0.47
1:E:160:GLN:HB2	1:E:165:ILE:CD1	2.46	0.46
1:F:181:LYS:HA	1:F:186:ILE:HG13	1.96	0.46
1:A:147:ARG:NH1	1:A:147:ARG:CG	2.63	0.46
1:A:176:THR:HG22	1:A:179:GLU:H	1.81	0.46
1:E:115:LEU:HD13	1:F:79:ASP:HB3	1.98	0.46
1:A:123:HIS:HB2	1:A:171:ARG:O	2.16	0.46
1:C:152:ARG:HH22	4:C:196:GOL:HO1	1.61	0.45
1:B:144:LEU:HD12	1:G:119:GLU:HG2	1.98	0.45
1:A:174:PHE:CE1	1:G:140:ALA:HB2	2.52	0.45
1:B:18:TYR:C	1:B:20:ILE:H	2.20	0.45
1:C:90:THR:C	1:C:91:ILE:HG13	2.37	0.45
1:B:174:PHE:CE1	1:D:140:ALA:HB2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:GLU:HG2	1:E:144:LEU:HD12	1.99	0.44
1:E:191:VAL:HA	1:E:192:PRO:HD3	1.87	0.44
1:G:147[A]:ARG:HA	1:G:150:LEU:HD12	1.98	0.44
1:E:58:LYS:O	1:E:86:PRO:HB3	2.18	0.44
1:C:72:THR:HG22	5:C:204:HOH:O	2.17	0.43
1:C:115:LEU:HD13	1:E:79:ASP:HB3	1.99	0.43
1:D:71:VAL:HB	1:D:99:MET:HE3	2.00	0.43
1:F:153:ILE:O	1:F:157:ARG:HG2	2.19	0.42
1:G:124:GLN:HA	1:G:125:PRO:HD3	1.76	0.42
1:C:149:LYS:HD3	1:D:117:ASN:HD21	1.81	0.42
1:C:72:THR:HG21	1:C:137:GLU:HB3	2.02	0.42
1:C:79:ASP:HB3	1:D:115:LEU:HD13	2.02	0.42
1:E:123:HIS:HB2	1:E:171:ARG:O	2.19	0.42
1:C:130:GLN:HG3	1:C:133:ALA:HB3	2.02	0.42
1:G:136:ILE:HD12	1:G:136:ILE:H	1.85	0.41
1:C:20:ILE:O	1:C:24:LEU:HB2	2.20	0.41
1:B:58:LYS:O	1:B:86:PRO:HB3	2.20	0.41
1:D:58:LYS:O	1:D:86:PRO:HB3	2.20	0.41
1:B:68:GLY:HA3	1:B:98:SER:HB3	2.03	0.41
1:F:160:GLN:HB3	1:F:165:ILE:HD11	2.01	0.41
1:A:58:LYS:O	1:A:86:PRO:HB3	2.22	0.40
1:E:147:ARG:H	1:E:147:ARG:CD	2.34	0.40
1:A:42:ASN:ND2	1:F:33:GLY:O	2.50	0.40
1:C:122:ILE:H	1:C:122:ILE:HG13	1.68	0.40
1:C:40:VAL:O	1:C:44:ILE:HG12	2.22	0.40
1:C:81:ILE:HG23	1:C:88:VAL:HB	2.04	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	167/197 (85%)	162 (97%)	5 (3%)	0	100	100
1	B	166/197 (84%)	160 (96%)	6 (4%)	0	100	100
1	C	168/197 (85%)	165 (98%)	3 (2%)	0	100	100
1	D	168/197 (85%)	163 (97%)	4 (2%)	1 (1%)	25	29
1	E	169/197 (86%)	166 (98%)	3 (2%)	0	100	100
1	F	170/197 (86%)	167 (98%)	3 (2%)	0	100	100
1	G	171/197 (87%)	164 (96%)	7 (4%)	0	100	100
All	All	1179/1379 (86%)	1147 (97%)	31 (3%)	1 (0%)	51	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	94	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	143/164 (87%)	135 (94%)	8 (6%)	21	27
1	B	142/164 (87%)	138 (97%)	4 (3%)	43	56
1	C	145/164 (88%)	138 (95%)	7 (5%)	25	34
1	D	145/164 (88%)	138 (95%)	7 (5%)	25	34
1	E	144/164 (88%)	135 (94%)	9 (6%)	18	23
1	F	146/164 (89%)	134 (92%)	12 (8%)	11	13
1	G	147/164 (90%)	141 (96%)	6 (4%)	30	40
All	All	1012/1148 (88%)	959 (95%)	53 (5%)	22	31

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ASP

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Mol	Chain	Res	Type
1	A	24	LEU
1	A	35	GLN
1	A	72	THR
1	A	98	SER
1	A	109	LYS
1	A	147	ARG
1	A	160	GLN
1	B	24	LEU
1	B	75	PHE
1	B	160	GLN
1	B	165	ILE
1	C	18	TYR
1	C	19	ASP
1	C	54	GLN
1	C	72	THR
1	C	122	ILE
1	C	145	LYS
1	C	147	ARG
1	D	23	ARG
1	D	35	GLN
1	D	71	VAL
1	D	109	LYS
1	D	130	GLN
1	D	134	THR
1	D	145	LYS
1	E	57	GLU
1	E	72	THR
1	E	85	LYS
1	E	95	MET
1	E	122	ILE
1	E	130	GLN
1	E	147	ARG
1	E	171	ARG
1	E	188	GLU
1	F	24	LEU
1	F	71	VAL
1	F	72	THR
1	F	75	PHE
1	F	93	ILE
1	F	132	GLN
1	F	134	THR
1	F	147	ARG

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Mol	Chain	Res	Type
1	F	165	ILE
1	F	173	ASN
1	F	193	GLU
1	F	194	THR
1	G	20	ILE
1	G	35	GLN
1	G	39	ASN
1	G	71	VAL
1	G	93	ILE
1	G	95	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	39	ASN
1	A	82	GLN
1	A	89	GLN
1	A	132	GLN
1	A	151	ASN
1	B	82	GLN
1	B	89	GLN
1	B	117	ASN
1	B	173	ASN
1	C	130	GLN
1	D	35	GLN
1	D	117	ASN
1	E	89	GLN
1	E	117	ASN
1	E	130	GLN
1	F	82	GLN
1	F	89	GLN
1	F	117	ASN
1	F	173	ASN
1	G	82	GLN
1	G	117	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	196	-	4,4,4	0.22	0	6,6,6	0.14	0
2	SO4	D	196	-	4,4,4	0.24	0	6,6,6	0.22	0
4	GOL	C	196	-	5,5,5	0.28	0	5,5,5	0.63	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	C	196	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	196	GOL	O1-C1-C2-O2
4	C	196	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
4	C	196	GOL	C1-C2-C3-O3
4	C	196	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	196	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	170/197 (86%)	0.26	6 (3%) 44 40	39, 50, 70, 80	0
1	B	170/197 (86%)	0.33	11 (6%) 18 15	42, 53, 75, 82	0
1	C	171/197 (86%)	0.32	5 (2%) 51 47	43, 53, 70, 76	0
1	D	171/197 (86%)	0.54	10 (5%) 23 19	41, 52, 78, 85	0
1	E	172/197 (87%)	0.32	6 (3%) 44 40	48, 62, 77, 88	0
1	F	173/197 (87%)	0.63	16 (9%) 9 6	41, 51, 79, 91	0
1	G	173/197 (87%)	0.23	5 (2%) 51 47	40, 52, 73, 86	0
All	All	1200/1379 (87%)	0.38	59 (4%) 29 27	39, 53, 76, 91	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	19	ASP	4.8
1	A	191	VAL	4.7
1	C	18	TYR	4.4
1	F	191	VAL	4.2
1	F	193	GLU	3.7
1	D	132	GLN	3.6
1	G	126	LEU	3.6
1	D	143	ILE	3.4
1	F	143	ILE	3.4
1	G	173	ASN	3.4
1	B	20	ILE	3.3
1	B	170	ASP	3.1
1	F	142	HIS	3.1
1	B	163	GLU	3.1
1	B	19	ASP	2.9
1	E	174	PHE	2.9
1	D	19	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	E	132	GLN	2.8
1	F	189	VAL	2.8
1	A	122	ILE	2.7
1	B	144	LEU	2.6
1	F	132	GLN	2.6
1	G	193	GLU	2.6
1	D	18	TYR	2.6
1	C	191	VAL	2.5
1	F	139	ALA	2.5
1	E	152	ARG	2.5
1	A	142	HIS	2.5
1	B	109	LYS	2.5
1	G	147[A]	ARG	2.4
1	A	174	PHE	2.4
1	B	145	LYS	2.4
1	F	176	THR	2.3
1	E	163	GLU	2.3
1	D	120	VAL	2.3
1	E	57	GLU	2.3
1	B	174	PHE	2.3
1	D	135	GLU	2.3
1	F	190	MET	2.2
1	F	92	CYS	2.2
1	F	19	ASP	2.2
1	D	93	ILE	2.2
1	B	172	ASP	2.2
1	F	91	ILE	2.2
1	B	133	ALA	2.2
1	F	21	TYR	2.2
1	F	170	ASP	2.1
1	C	170	ASP	2.1
1	G	130	GLN	2.1
1	D	142	HIS	2.1
1	F	186	ILE	2.1
1	C	23	ARG	2.1
1	D	92	CYS	2.1
1	D	176	THR	2.1
1	A	132	GLN	2.1
1	A	192	PRO	2.0
1	B	143	ILE	2.0
1	E	150	LEU	2.0
1	F	192	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	GOL	C	196	6/6	0.71	0.47	49,51,51,53	0
3	CA	B	196	1/1	0.91	0.23	85,85,85,85	0
2	SO4	A	196	5/5	0.93	0.23	83,84,84,84	0
2	SO4	D	196	5/5	0.96	0.22	56,59,59,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.