



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 29, 2021 – 10:04 AM EDT

PDB ID : 7N56
Title : Crystal Structure of deoxyuridine 5'-triphosphate nucleotidohydrolase from Rickettsia prowazekii str. Madrid E
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2021-06-04
Resolution : 2.15 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 : 2.22
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) : 2.22

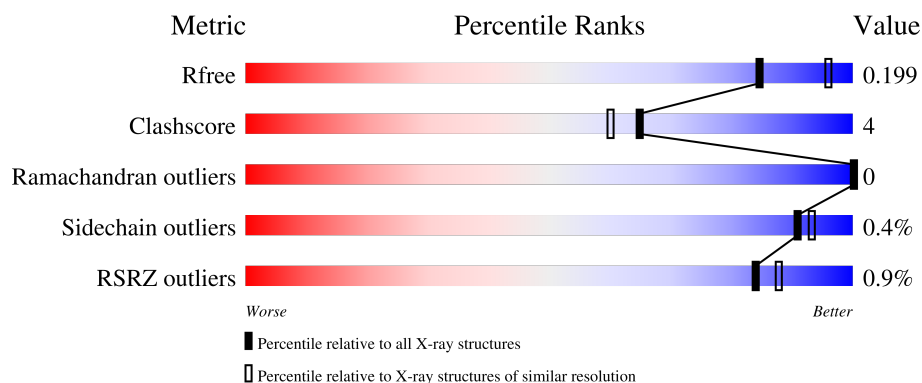
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	152	<div> <div>%</div> <div> <div></div> <div>84%</div> <div>•</div> <div>13%</div> </div> </div>
1	B	152	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>7%</div> <div>13%</div> </div> </div>
1	C	152	<div> <div></div> <div> <div>76%</div> <div>12%</div> <div>11%</div> </div> </div>
1	D	152	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>10%</div> <div>13%</div> </div> </div>
1	E	152	<div> <div>%</div> <div> <div></div> <div>81%</div> <div>6%</div> <div>13%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	152	 % 74% 13% 13%
1	G	152	 % 79% 8% 13%
1	H	152	 3% % 80% 7% 12%
1	I	152	 % 80% 9% 11%
1	J	152	 % 76% 12% 12%
1	K	152	 % 83% 7% 10%
1	L	152	 % 80% 12% 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13371 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 Deoxyuridine 5'-triphosphate nucleotidohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	0	4	0
			1023	660	173	186	4			
1	B	132	Total	C	N	O	S	0	4	0
			1002	646	167	186	3			
1	C	135	Total	C	N	O	S	0	3	0
			1041	668	176	193	4			
1	D	132	Total	C	N	O	S	0	4	0
			1036	667	173	193	3			
1	E	132	Total	C	N	O	S	0	3	0
			1000	644	170	182	4			
1	F	132	Total	C	N	O	S	0	0	0
			989	635	169	182	3			
1	G	132	Total	C	N	O	S	0	3	0
			1008	649	171	185	3			
1	H	133	Total	C	N	O	S	0	2	0
			996	640	168	185	3			
1	I	136	Total	C	N	O	S	0	2	0
			1024	657	173	189	5			
1	J	134	Total	C	N	O	S	0	3	0
			1021	654	175	188	4			
1	K	137	Total	C	N	O	S	0	3	0
			1034	665	176	188	5			
1	L	139	Total	C	N	O	S	0	4	0
			1066	685	178	197	6			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q9ZDD2
A	-2	PRO	-	expression tag	UNP Q9ZDD2
A	-1	GLY	-	expression tag	UNP Q9ZDD2
A	0	SER	-	expression tag	UNP Q9ZDD2
B	-3	GLY	-	expression tag	UNP Q9ZDD2

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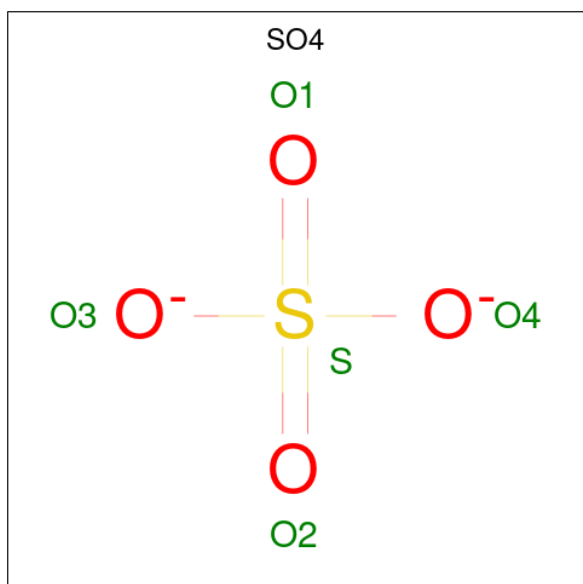
Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	PRO	-	expression tag	UNP Q9ZDD2
B	-1	GLY	-	expression tag	UNP Q9ZDD2
B	0	SER	-	expression tag	UNP Q9ZDD2
C	-3	GLY	-	expression tag	UNP Q9ZDD2
C	-2	PRO	-	expression tag	UNP Q9ZDD2
C	-1	GLY	-	expression tag	UNP Q9ZDD2
C	0	SER	-	expression tag	UNP Q9ZDD2
D	-3	GLY	-	expression tag	UNP Q9ZDD2
D	-2	PRO	-	expression tag	UNP Q9ZDD2
D	-1	GLY	-	expression tag	UNP Q9ZDD2
D	0	SER	-	expression tag	UNP Q9ZDD2
E	-3	GLY	-	expression tag	UNP Q9ZDD2
E	-2	PRO	-	expression tag	UNP Q9ZDD2
E	-1	GLY	-	expression tag	UNP Q9ZDD2
E	0	SER	-	expression tag	UNP Q9ZDD2
F	-3	GLY	-	expression tag	UNP Q9ZDD2
F	-2	PRO	-	expression tag	UNP Q9ZDD2
F	-1	GLY	-	expression tag	UNP Q9ZDD2
F	0	SER	-	expression tag	UNP Q9ZDD2
G	-3	GLY	-	expression tag	UNP Q9ZDD2
G	-2	PRO	-	expression tag	UNP Q9ZDD2
G	-1	GLY	-	expression tag	UNP Q9ZDD2
G	0	SER	-	expression tag	UNP Q9ZDD2
H	-3	GLY	-	expression tag	UNP Q9ZDD2
H	-2	PRO	-	expression tag	UNP Q9ZDD2
H	-1	GLY	-	expression tag	UNP Q9ZDD2
H	0	SER	-	expression tag	UNP Q9ZDD2
I	-3	GLY	-	expression tag	UNP Q9ZDD2
I	-2	PRO	-	expression tag	UNP Q9ZDD2
I	-1	GLY	-	expression tag	UNP Q9ZDD2
I	0	SER	-	expression tag	UNP Q9ZDD2
J	-3	GLY	-	expression tag	UNP Q9ZDD2
J	-2	PRO	-	expression tag	UNP Q9ZDD2
J	-1	GLY	-	expression tag	UNP Q9ZDD2
J	0	SER	-	expression tag	UNP Q9ZDD2
K	-3	GLY	-	expression tag	UNP Q9ZDD2
K	-2	PRO	-	expression tag	UNP Q9ZDD2
K	-1	GLY	-	expression tag	UNP Q9ZDD2
K	0	SER	-	expression tag	UNP Q9ZDD2
L	-3	GLY	-	expression tag	UNP Q9ZDD2
L	-2	PRO	-	expression tag	UNP Q9ZDD2
L	-1	GLY	-	expression tag	UNP Q9ZDD2

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Chain	Residue	Modelled	Actual	Comment	Reference
L	0	SER	-	expression tag	UNP Q9ZDD2

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



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	122	Total	O	0	0
			122	122		
3	B	110	Total	O	0	2
			112	112		
3	C	130	Total	O	0	2
			132	132		
3	D	106	Total	O	0	1
			107	107		
3	E	82	Total	O	0	0
			82	82		
3	F	96	Total	O	0	0
			96	96		

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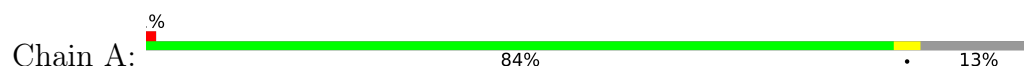
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	85	Total 85	O 85	0	0
3	H	64	Total 64	O 64	0	0
3	I	92	Total 92	O 92	0	0
3	J	75	Total 75	O 75	0	0
3	K	67	Total 67	O 67	0	0
3	L	82	Total 82	O 82	0	0

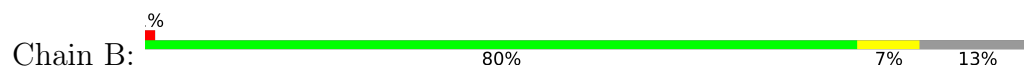
3 Residue-property plots [i](#)

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.

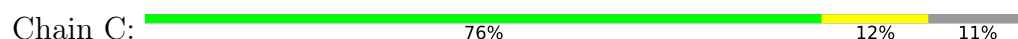
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



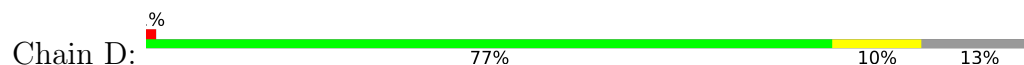
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



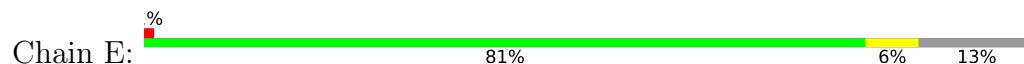
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



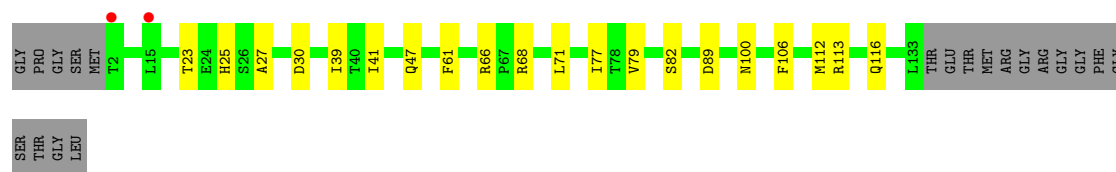
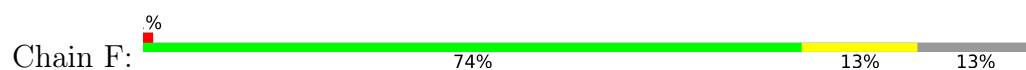
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



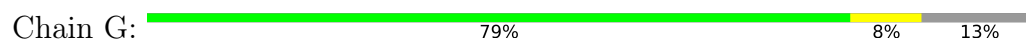
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



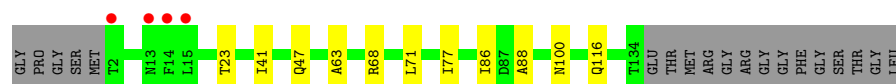
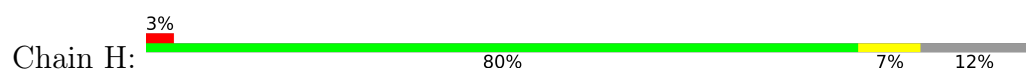
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



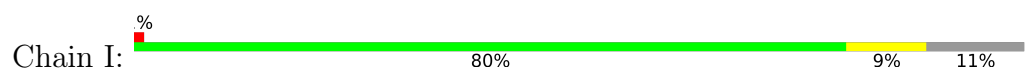
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



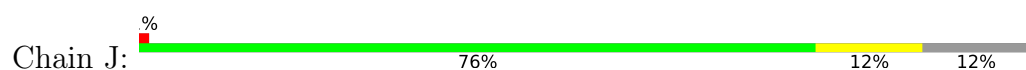
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



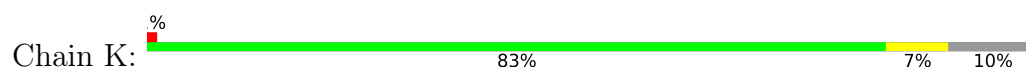
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



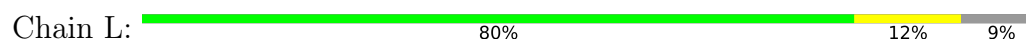
- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



- Molecule 1: Deoxyuridine 5'-triphosphate nucleotidohydrolase



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GLY	PRO	G-1	K9	H25	I41	Q47	E59	A63	R68	L71	H75	G76	I77	P83	I86	D89	F106	M112	Q116	Y122	E123	M137	ARG	GLY	ARG	GLY	GLY	PHE	GLY	SER	THR	GLY	LEU
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	100.37Å 97.39Å 100.54Å 90.00° 118.60° 90.00°	Depositor
Resolution (Å)	48.69 – 2.15 48.70 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.8 (48.69-2.15) 98.9 (48.70-2.15)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.02 (at 2.16Å)	Xtriage
Refinement program	PHENIX 1.18 4224	Depositor
R, R_{free}	0.159 , 0.201 0.159 , 0.199	Depositor DCC
R_{free} test set	1866 reflections (2.04%)	wwPDB-VP
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.389	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h-l,k,h 0.000 for l,k,-h-l 0.000 for h,-k,-h-l 0.000 for -h-l,-k,l 0.012 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13371	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.43 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.7400e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.42	0/1049	0.62	0/1419
1	B	0.37	0/1028	0.59	0/1395
1	C	0.37	0/1064	0.62	0/1438
1	D	0.37	0/1062	0.60	0/1436
1	E	0.35	0/1023	0.58	0/1387
1	F	0.35	0/1003	0.60	0/1360
1	G	0.35	0/1031	0.60	0/1397
1	H	0.35	0/1016	0.58	0/1381
1	I	0.34	0/1045	0.59	0/1415
1	J	0.34	0/1045	0.59	0/1415
1	K	0.34	0/1058	0.62	0/1433
1	L	0.35	0/1092	0.57	0/1476
All	All	0.36	0/12516	0.60	0/16952

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	1023	0	1084	4	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1002	0	1040	8	0
1	C	1041	0	1097	16	1
1	D	1036	0	1096	9	0
1	E	1000	0	1037	6	0
1	F	989	0	1017	16	0
1	G	1008	0	1055	6	0
1	H	996	0	1019	7	0
1	I	1024	0	1063	10	0
1	J	1021	0	1064	13	0
1	K	1034	0	1082	8	0
1	L	1066	0	1121	17	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	F	5	0	0	0	0
3	A	122	0	0	0	0
3	B	112	0	0	1	0
3	C	132	0	0	3	0
3	D	107	0	0	0	0
3	E	82	0	0	0	0
3	F	96	0	0	0	0
3	G	85	0	0	0	0
3	H	64	0	0	0	0
3	I	92	0	0	1	0
3	J	75	0	0	0	0
3	K	67	0	0	2	0
3	L	82	0	0	3	0
All	All	13371	0	12775	102	1

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

All (102) 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:39:ILE:HG21	1:D:47[B]:GLN:HG2	1.65	0.78
1:I:89:ASP:OD2	3:I:201:HOH:O	2.06	0.73
1:C:39:ILE:HG21	1:C:47[B]:GLN:HG2	1.79	0.65
1:A:30:ASP:OD1	1:A:68[A]:ARG:NH1	2.31	0.63
1:F:106:PHE:HE2	1:F:112:MET:HE1	1.62	0.63
1:J:88:ALA:O	1:J:91[B]:ARG:NE	2.32	0.63
1:F:23:THR:OG1	1:L:89:ASP:OD2	2.19	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:68:ARG:HH11	1:I:116:GLN:HE22	1.47	0.60
1:L:106:PHE:HE2	1:L:112[A]:MET:HE1	1.66	0.59
1:H:63:ALA:HB3	1:H:86:ILE:HB	1.85	0.59
1:A:25:HIS:HB3	1:I:88:ALA:HB3	1.84	0.58
1:J:41:ILE:HG12	1:J:47:GLN:HG3	1.86	0.57
1:I:71:LEU:HG	1:I:77:ILE:HB	1.87	0.56
1:K:116:GLN:NE2	3:K:203:HOH:O	2.37	0.56
1:F:106:PHE:CE2	1:F:112:MET:HE1	2.40	0.56
1:C:30:ASP:OD2	1:C:113:ARG:NH1	2.39	0.55
1:G:106:PHE:HE2	1:G:112:MET:HE1	1.71	0.55
1:H:71:LEU:HG	1:H:77:ILE:HB	1.89	0.54
1:C:15:LEU:HD12	1:C:110:LYS:HD3	1.89	0.54
1:K:12:GLU:O	3:K:201:HOH:O	2.18	0.54
1:F:30:ASP:OD2	1:F:113:ARG:NH1	2.40	0.54
1:C:71:LEU:HG	1:C:77:ILE:HB	1.88	0.54
1:C:47[A]:GLN:NE2	3:C:304:HOH:O	2.41	0.54
1:B:106:PHE:HE2	1:B:112:MET:HE1	1.73	0.54
1:G:63:ALA:HB3	1:G:86:ILE:HB	1.90	0.53
1:L:68:ARG:HH11	1:L:116[A]:GLN:HE22	1.57	0.53
1:E:71[A]:LEU:HG	1:E:77:ILE:HB	1.90	0.53
1:L:123:GLU:OE2	3:L:201:HOH:O	2.18	0.53
1:G:71:LEU:HG	1:G:77:ILE:HB	1.91	0.52
1:C:41:ILE:HG12	1:C:47[A]:GLN:HG3	1.92	0.52
1:L:75:HIS:ND1	3:L:202:HOH:O	2.33	0.51
1:L:41:ILE:HG12	1:L:47:GLN:HG3	1.91	0.51
1:I:68:ARG:HH11	1:I:116:GLN:NE2	2.09	0.51
1:K:71:LEU:HG	1:K:77:ILE:HB	1.92	0.51
1:B:88:ALA:HB3	1:D:25:HIS:HB3	1.93	0.50
1:A:88:ALA:HB3	1:C:25:HIS:HB3	1.92	0.50
1:J:66:ARG:HD2	1:J:118:ILE:HD13	1.94	0.50
1:D:61:PHE:CZ	1:D:121:LYS:HE2	2.47	0.49
1:F:41:ILE:HG12	1:F:47:GLN:HG3	1.94	0.49
1:J:106:PHE:HE2	1:J:112:MET:HE1	1.78	0.49
1:C:77:ILE:HD13	1:C:100:ASN:HA	1.94	0.49
1:L:106:PHE:CE2	1:L:112[A]:MET:HE1	2.46	0.49
1:F:71:LEU:HG	1:F:77:ILE:HB	1.95	0.49
1:H:88:ALA:HB3	1:L:25:HIS:HB3	1.95	0.49
1:I:77:ILE:HD13	1:I:100:ASN:HA	1.94	0.49
1:L:71:LEU:HG	1:L:77:ILE:HB	1.93	0.49
1:J:15:LEU:HD23	1:J:15:LEU:HA	1.74	0.48
1:C:3:ILE:HG12	1:J:128:ALA:HB1	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:89:ASP:OD2	1:H:23:THR:OG1	2.20	0.47
1:C:89:ASP:OD2	1:I:23:THR:OG1	2.24	0.47
1:K:106:PHE:CE2	1:K:112:MET:HE1	2.50	0.47
1:K:63:ALA:HB3	1:K:86:ILE:HB	1.97	0.46
1:J:68:ARG:HH11	1:J:116[A]:GLN:HE22	1.62	0.46
1:E:23:THR:OG1	1:K:89:ASP:OD2	2.28	0.45
1:D:39:ILE:CG2	1:D:47[B]:GLN:HG2	2.40	0.45
1:K:106:PHE:HE2	1:K:112:MET:HE1	1.80	0.45
1:D:71[A]:LEU:HG	1:D:77:ILE:HB	1.98	0.45
1:J:68:ARG:HH11	1:J:116[A]:GLN:NE2	2.15	0.45
1:L:9:LYS:NZ	3:L:207:HOH:O	2.50	0.45
1:D:41:ILE:HG12	1:D:47[A]:GLN:HG3	1.98	0.45
1:F:68:ARG:HH11	1:F:116:GLN:HE22	1.65	0.45
1:B:106:PHE:CE2	1:B:112:MET:HE1	2.51	0.44
1:G:38[B]:SER:OG	1:G:109:GLU:HG2	2.16	0.44
1:F:77:ILE:HD13	1:F:100:ASN:HA	1.99	0.44
1:I:66:ARG:HD2	1:I:118:ILE:HD13	1.99	0.44
1:J:106:PHE:CE2	1:J:112:MET:HE1	2.52	0.44
1:F:61:PHE:HE2	1:L:123:GLU:HG3	1.82	0.44
1:E:106:PHE:HE2	1:E:112[A]:MET:HE1	1.81	0.44
1:H:77:ILE:HD13	1:H:100:ASN:HA	2.00	0.44
1:E:14:PHE:O	1:E:15:LEU:HD23	2.17	0.44
1:C:66:ARG:HD2	1:C:118:ILE:HD13	1.98	0.44
1:E:123:GLU:HG3	1:G:61:PHE:CE2	2.53	0.43
1:L:68:ARG:HH11	1:L:116[A]:GLN:NE2	2.14	0.43
1:E:77:ILE:HD13	1:E:100:ASN:HA	2.00	0.43
1:H:41:ILE:HG12	1:H:47:GLN:HG3	2.00	0.43
1:D:106:PHE:HE2	1:D:112:MET:HE1	1.84	0.43
1:B:75:HIS:HB2	1:B:77[B]:ILE:HD12	2.01	0.43
1:C:103[A]:ASN:OD1	3:C:301:HOH:O	2.21	0.43
1:F:61:PHE:CE2	1:L:123:GLU:HG3	2.54	0.43
1:B:25:HIS:HB3	1:J:88:ALA:HB3	2.01	0.43
1:F:27:ALA:HA	1:L:122:TYR:CE1	2.52	0.43
1:C:7:LYS:HB2	1:C:7:LYS:HE3	1.80	0.43
1:B:74:LYS:NZ	3:B:304:HOH:O	2.45	0.42
1:J:14:PHE:HB2	1:J:52:GLY:O	2.19	0.42
1:F:25:HIS:HB2	1:L:89:ASP:OD1	2.19	0.42
1:K:117:MET:HB3	1:K:117:MET:HE2	1.86	0.42
1:L:63:ALA:HB3	1:L:86:ILE:HB	2.01	0.42
1:D:104:LYS:HG2	1:F:39:ILE:HG12	2.01	0.42
1:J:77:ILE:HD13	1:J:100:ASN:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:113:ARG:HB3	3:C:305:HOH:O	2.19	0.42
1:A:122:TYR:CE1	1:C:27:ALA:HA	2.56	0.41
1:F:66:ARG:HD3	1:L:83:PRO:HG2	2.02	0.41
1:I:14:PHE:HZ	1:I:17:ASN:O	2.04	0.41
1:J:63:ALA:HB3	1:J:86:ILE:HB	2.01	0.41
1:D:63:ALA:HB3	1:D:86:ILE:HB	2.02	0.41
1:C:39:ILE:CG2	1:C:47[B]:GLN:HG2	2.49	0.41
1:B:79:VAL:HB	1:B:82:SER:HB2	2.03	0.41
1:G:67:PRO:HD3	1:G:82:SER:OG	2.21	0.41
1:B:41:ILE:HG12	1:B:47:GLN:HG3	2.02	0.41
1:H:68:ARG:HH11	1:H:116[A]:GLN:NE2	2.19	0.40
1:I:63:ALA:HB3	1:I:86:ILE:HB	2.04	0.40
1:F:79:VAL:HB	1:F:82:SER:HB2	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68[A]:ARG:NH2	1:C:36:GLU:OE1[2_656]	2.07	0.13

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	134/152 (88%)	134 (100%)	0	0	100	100
1	B	134/152 (88%)	134 (100%)	0	0	100	100
1	C	136/152 (90%)	136 (100%)	0	0	100	100
1	D	134/152 (88%)	133 (99%)	1 (1%)	0	100	100
1	E	133/152 (88%)	133 (100%)	0	0	100	100
1	F	130/152 (86%)	129 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	133/152 (88%)	133 (100%)	0	0	100	100
1	H	133/152 (88%)	133 (100%)	0	0	100	100
1	I	136/152 (90%)	136 (100%)	0	0	100	100
1	J	135/152 (89%)	135 (100%)	0	0	100	100
1	K	138/152 (91%)	138 (100%)	0	0	100	100
1	L	141/152 (93%)	141 (100%)	0	0	100	100
All	All	1617/1824 (89%)	1615 (100%)	2 (0%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	110/122 (90%)	110 (100%)	0	100	100
1	B	106/122 (87%)	106 (100%)	0	100	100
1	C	112/122 (92%)	112 (100%)	0	100	100
1	D	113/122 (93%)	111 (98%)	2 (2%)	59	63
1	E	104/122 (85%)	104 (100%)	0	100	100
1	F	102/122 (84%)	102 (100%)	0	100	100
1	G	107/122 (88%)	106 (99%)	1 (1%)	78	83
1	H	103/122 (84%)	103 (100%)	0	100	100
1	I	108/122 (88%)	108 (100%)	0	100	100
1	J	108/122 (88%)	108 (100%)	0	100	100
1	K	109/122 (89%)	107 (98%)	2 (2%)	59	63
1	L	114/122 (93%)	113 (99%)	1 (1%)	78	83
All	All	1296/1464 (88%)	1290 (100%)	6 (0%)	91	92

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	91	ARG
1	D	131	SER
1	G	91	ARG
1	K	91[A]	ARG
1	K	91[B]	ARG
1	L	59	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	116	GLN
1	I	116	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](#)

3 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	C	201	-	4,4,4	0.18	0	6,6,6	0.23	0
2	SO4	F	201	-	4,4,4	0.10	0	6,6,6	0.15	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	201	-	4,4,4	0.13	0	6,6,6	0.18	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 ⓘ

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	132/152 (86%)	-0.36	1 (0%) 86 89	16, 24, 46, 74	0
1	B	132/152 (86%)	-0.47	1 (0%) 86 89	17, 26, 46, 63	0
1	C	135/152 (88%)	-0.47	0 100 100	16, 24, 39, 60	0
1	D	132/152 (86%)	-0.47	1 (0%) 86 89	19, 28, 54, 79	0
1	E	132/152 (86%)	-0.45	2 (1%) 73 79	19, 31, 55, 77	0
1	F	132/152 (86%)	-0.46	2 (1%) 73 79	20, 30, 55, 80	0
1	G	132/152 (86%)	-0.52	0 100 100	20, 32, 51, 76	0
1	H	133/152 (87%)	-0.35	4 (3%) 50 59	19, 35, 65, 93	0
1	I	136/152 (89%)	-0.42	1 (0%) 87 91	18, 29, 55, 83	0
1	J	134/152 (88%)	-0.44	1 (0%) 87 91	21, 33, 53, 103	0
1	K	137/152 (90%)	-0.49	1 (0%) 87 91	21, 32, 56, 101	0
1	L	139/152 (91%)	-0.48	0 100 100	19, 32, 49, 64	0
All	All	1606/1824 (88%)	-0.45	14 (0%) 84 88	16, 30, 54, 103	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	132	ILE	3.5
1	H	15	LEU	3.4
1	A	2	THR	3.2
1	H	14	PHE	3.0
1	B	2	THR	2.9
1	E	15	LEU	2.9
1	F	15	LEU	2.6
1	J	-2	PRO	2.6
1	F	2	THR	2.5
1	H	13	ASN	2.4
1	H	2	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	2	THR	2.2
1	K	-2	PRO	2.1
1	I	132	ILE	2.1

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	F	201	5/5	0.62	0.25	27,29,31,51	5
2	SO4	B	201	5/5	0.86	0.14	65,71,73,109	0
2	SO4	C	201	5/5	0.89	0.14	46,47,51,83	0

6.5 Other polymers [i](#)

There are no such residues in this entry.