



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 05:30 PM BST

PDB ID : 6KJA  
Title : E. coli ATCase holoenzyme mutant - G128/130A (catalytic chain)  
Authors : Lei, Z.; Zheng, J.; Jia, Z.C.  
Deposited on : 2019-07-22  
Resolution : 3.06 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](mailto: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.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.13  
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) : 2.13

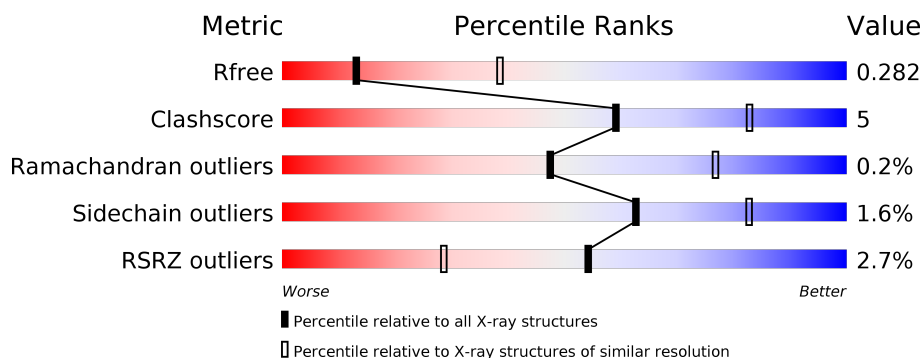
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.06 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	1754 (3.10-3.02)
Clashscore	141614	1864 (3.10-3.02)
Ramachandran outliers	138981	1794 (3.10-3.02)
Sidechain outliers	138945	1793 (3.10-3.02)
RSRZ outliers	127900	1713 (3.10-3.02)

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  $\geq 3$ , 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	310	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>8%</div> <div>..</div> </div> </div>
1	C	310	<div> <div>85%</div> <div>12%</div> <div>..</div> </div>
1	E	310	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>14%</div> <div>..</div> </div> </div>
2	B	153	<div> <div>8%</div> <div> <div></div> <div>75%</div> <div>12%</div> <div>12%</div> </div> </div>
2	D	153	<div> <div>8%</div> <div> <div></div> <div>78%</div> <div>9%</div> <div>13%</div> </div> </div>
2	F	153	<div> <div>3%</div> <div> <div></div> <div>74%</div> <div>14%</div> <div>12%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10019 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 Aspartate carbamoyltransferase catalytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2354	1491	411	443	9			
1	C	303	Total	C	N	O	S	0	0	0
			2355	1493	409	444	9			
1	E	304	Total	C	N	O	S	0	0	0
			2356	1495	409	443	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	128	ALA	GLY	engineered mutation	UNP P0A786
A	130	ALA	GLY	engineered mutation	UNP P0A786
C	128	ALA	GLY	engineered mutation	UNP P0A786
C	130	ALA	GLY	engineered mutation	UNP P0A786
E	128	ALA	GLY	engineered mutation	UNP P0A786
E	130	ALA	GLY	engineered mutation	UNP P0A786

- Molecule 2 is a protein called Aspartate carbamoyltransferase regulatory chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	134	Total	C	N	O	S	0	0	0
			976	616	168	188	4			
2	D	133	Total	C	N	O	S	0	0	0
			955	602	163	186	4			
2	F	134	Total	C	N	O	S	0	0	0
			1009	638	174	193	4			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total 1	Zn 1	0	0
3	F	1	Total 1	Zn 1	0	0

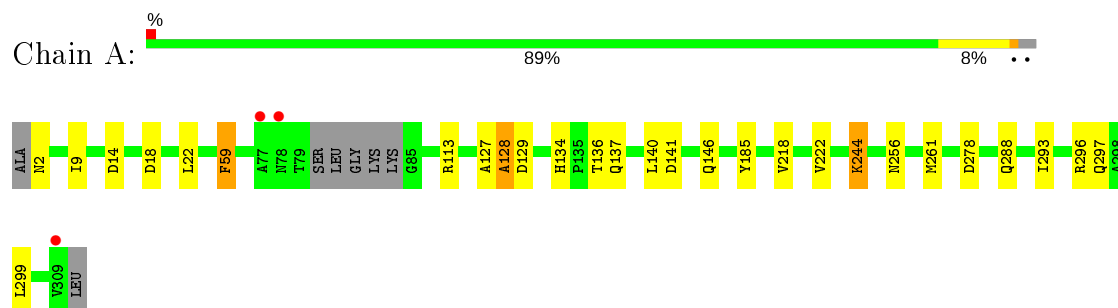
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total 4	O 4	0	0
4	B	2	Total 2	O 2	0	0
4	C	2	Total 2	O 2	0	0
4	D	2	Total 2	O 2	0	0
4	E	1	Total 1	O 1	0	0

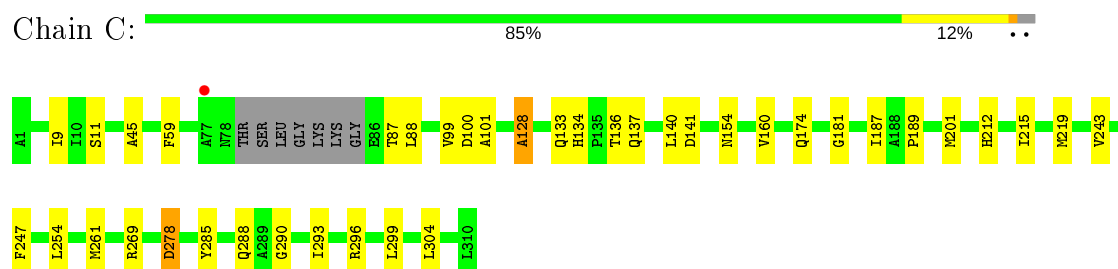
### 3 Residue-property plots

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.

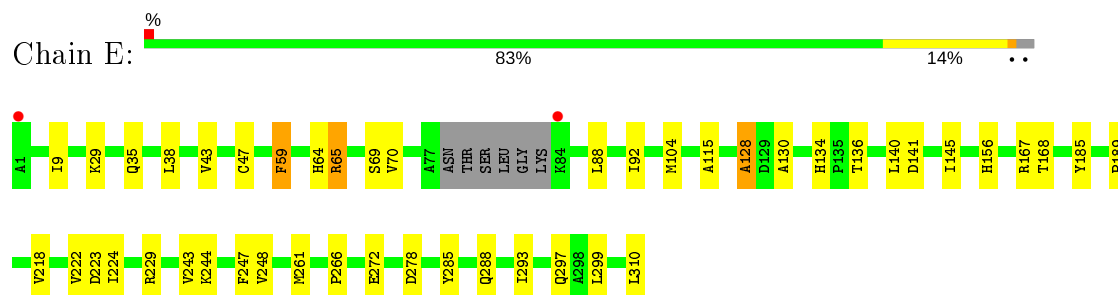
- Molecule 1: Aspartate carbamoyltransferase catalytic subunit



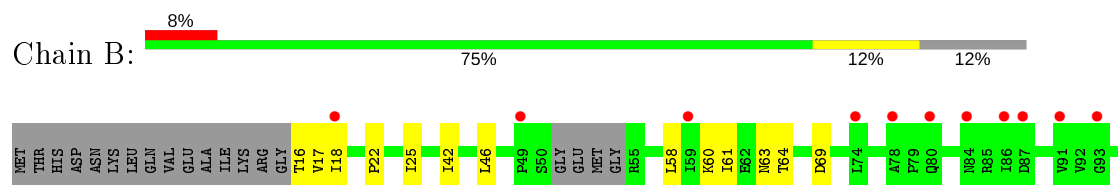
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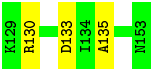
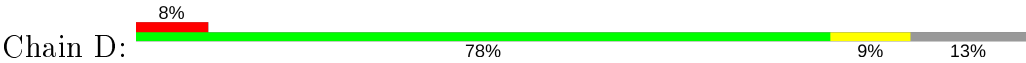


- Molecule 2: Aspartate carbamoyltransferase regulatory chain

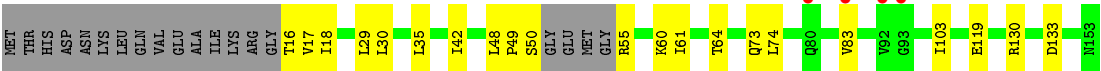




• Molecule 2: Aspartate carbamoyltransferase regulatory chain



• Molecule 2: Aspartate carbamoyltransferase regulatory chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.41Å 127.41Å 197.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.29 – 3.06 42.29 – 3.06	Depositor EDS
% Data completeness (in resolution range)	98.6 (42.29-3.06) 98.6 (42.29-3.06)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.02 (at 3.06Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, $R_{free}$	0.249 , 0.283 0.250 , 0.282	Depositor DCC
$R_{free}$ test set	1966 reflections (5.62%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	68.7	Xtriage
Anisotropy	0.235	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 43.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.55$ , $\langle L^2 \rangle = 0.40$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10019	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	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 48.63 % 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 8.3693e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: ZN

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	0/2399	0.45	0/3260
1	C	0.24	0/2400	0.44	0/3263
1	E	0.24	0/2401	0.46	1/3261 (0.0%)
2	B	0.25	0/989	0.48	0/1348
2	D	0.25	0/968	0.48	0/1323
2	F	0.26	0/1024	0.47	0/1393
All	All	0.24	0/10181	0.46	1/13848 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	140	LEU	CA-CB-CG	5.70	128.41	115.30

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	2354	0	2344	17	0
1	C	2355	0	2341	22	0
1	E	2356	0	2352	30	0
2	B	976	0	944	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	955	0	897	9	0
2	F	1009	0	995	11	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
3	F	1	0	0	0	0
4	A	4	0	0	2	0
4	B	2	0	0	1	0
4	C	2	0	0	0	0
4	D	2	0	0	1	0
4	E	1	0	0	0	0
All	All	10019	0	9873	94	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 (94) 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)
2:D:130:ARG:NH2	2:D:133:ASP:OD2	1.90	1.04
2:F:130:ARG:NH2	2:F:133:ASP:OD2	2.06	0.87
2:D:122:SER:O	4:D:301:HOH:O	1.97	0.82
1:C:11:SER:HA	1:C:133:GLN:HG2	1.66	0.76
1:A:18:ASP:OD2	4:A:401:HOH:O	2.04	0.75
1:A:141:ASP:OD1	1:A:288:GLN:NE2	2.23	0.71
1:C:243:VAL:HG22	1:C:247:PHE:HE2	1.59	0.68
1:E:9:ILE:HG21	1:E:299:LEU:HD21	1.76	0.68
1:C:141:ASP:OD1	1:C:288:GLN:NE2	2.28	0.67
1:E:243:VAL:HG22	1:E:247:PHE:HE2	1.61	0.65
2:F:42:ILE:HG12	2:F:61:ILE:HG22	1.79	0.65
2:B:17:VAL:HG13	2:B:60:LYS:HG2	1.81	0.62
2:B:42:ILE:HG12	2:B:61:ILE:HG22	1.83	0.60
2:F:29:LEU:HD22	2:F:74:LEU:HD22	1.83	0.59
1:C:9:ILE:HG21	1:C:299:LEU:HD21	1.84	0.59
1:A:14:ASP:OD2	1:A:113:ARG:NH2	2.36	0.59
1:C:87:THR:OG1	2:D:119:GLU:OE2	2.19	0.59
1:C:45:ALA:HB2	1:C:99:VAL:HG11	1.84	0.57
1:E:293:ILE:O	1:E:297:GLN:HB2	2.05	0.57
2:D:61:ILE:HD11	2:D:64:THR:CB	2.35	0.57
1:E:141:ASP:OD1	1:E:288:GLN:NE2	2.37	0.57
1:E:29:LYS:HD3	1:E:310:LEU:HB3	1.87	0.56
1:A:2:ASN:HB2	1:A:22:LEU:HD21	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:18:ILE:HG23	2:F:83:VAL:HG22	1.87	0.56
1:C:160:VAL:HG11	1:C:215:ILE:HD11	1.87	0.55
2:F:16:THR:OG1	2:F:64:THR:O	2.19	0.54
1:A:9:ILE:HG21	1:A:299:LEU:HD21	1.89	0.54
1:E:59:PHE:CZ	1:E:136:THR:HG21	2.43	0.54
1:E:229:ARG:HA	1:E:272:GLU:OE2	2.07	0.54
1:C:101:ALA:HB2	1:C:304:LEU:HD21	1.89	0.53
1:E:218:VAL:HG22	1:E:222:VAL:HG13	1.90	0.53
1:E:130:ALA:O	1:E:167:ARG:NH1	2.36	0.52
1:A:59:PHE:CZ	1:A:136:THR:HG21	2.44	0.52
1:C:187:ILE:HG12	1:C:212:HIS:HB2	1.91	0.52
2:D:19:ASP:OD1	2:D:20:HIS:N	2.43	0.52
2:B:16:THR:OG1	2:B:64:THR:O	2.21	0.51
1:A:129:ASP:OD1	1:A:129:ASP:N	2.40	0.51
1:A:136:THR:HB	1:A:296:ARG:HE	1.76	0.51
1:C:189:PRO:HB3	1:C:243:VAL:HG21	1.93	0.51
2:F:30:LEU:HD23	2:F:35:LEU:HD12	1.92	0.50
1:E:59:PHE:HZ	1:E:136:THR:HG21	1.77	0.50
1:E:35:GLN:HG2	1:E:38:LEU:HD12	1.92	0.50
1:E:185:TYR:CG	1:E:218:VAL:HG21	2.48	0.49
2:D:42:ILE:HG23	2:D:61:ILE:HG22	1.95	0.49
2:F:17:VAL:HG13	2:F:60:LYS:HG2	1.94	0.49
1:C:88:LEU:N	2:D:119:GLU:OE2	2.46	0.48
1:C:290:GLY:O	1:C:293:ILE:HG12	2.14	0.48
2:F:50:SER:HG	2:F:55:ARG:N	2.11	0.48
1:C:269:ARG:NH2	1:C:278:ASP:OD2	2.43	0.47
1:A:59:PHE:HZ	1:A:136:THR:HG21	1.79	0.47
1:A:185:TYR:CG	1:A:218:VAL:HG21	2.50	0.47
2:F:73:GLN:HE22	2:F:103:ILE:HA	1.78	0.47
1:C:154:ASN:HA	1:C:181:GLY:O	2.15	0.47
1:C:128:ALA:HA	1:C:133:GLN:O	2.16	0.46
1:C:137:GLN:O	1:C:140:LEU:HG	2.16	0.46
1:A:244:LYS:NZ	4:A:402:HOH:O	2.33	0.46
1:E:189:PRO:HB3	1:E:243:VAL:HG21	1.97	0.46
1:A:293:ILE:HG22	1:A:297:GLN:HE21	1.81	0.45
1:C:100:ASP:OD2	1:E:65:ARG:NH2	2.48	0.45
1:C:219:MET:HE3	1:C:254:LEU:HD23	1.99	0.45
1:E:156:HIS:ND1	1:E:185:TYR:OH	2.34	0.45
1:E:168:THR:HG21	1:E:266:PRO:HG2	1.98	0.45
1:E:43:VAL:HG22	1:E:69:SER:HB3	2.00	0.44
1:E:223:ASP:O	1:E:224:ILE:HD12	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:135:ALA:HB1	2:B:144:GLU:HG2	1.99	0.44
2:B:128:ARG:HG3	2:B:135:ALA:HB3	1.99	0.44
2:D:48:LEU:N	2:D:56:LYS:O	2.47	0.44
1:E:92:ILE:HD13	1:E:115:ALA:CB	2.47	0.44
1:A:218:VAL:HG22	1:A:222:VAL:HG13	1.99	0.44
1:E:156:HIS:HB3	1:E:185:TYR:HE1	1.83	0.43
1:E:145:ILE:HG12	1:E:224:ILE:HG12	2.01	0.43
1:E:128:ALA:HB2	1:E:134:HIS:CG	2.53	0.43
1:E:64:HIS:CE1	1:E:70:VAL:HG21	2.54	0.43
1:C:133:GLN:HG3	1:C:134:HIS:N	2.33	0.43
1:A:278:ASP:OD1	1:A:278:ASP:N	2.52	0.43
2:B:18:ILE:O	2:B:58:LEU:HD23	2.18	0.43
1:C:278:ASP:OD1	1:C:278:ASP:N	2.51	0.43
1:A:137:GLN:O	1:A:140:LEU:HG	2.20	0.42
1:E:92:ILE:HD13	1:E:115:ALA:HB2	2.02	0.42
1:A:256:ASN:N	1:A:256:ASN:OD1	2.52	0.41
2:F:48:LEU:HB3	2:F:49:PRO:HD2	2.01	0.41
2:B:22:PRO:HB2	2:B:25:ILE:HG13	2.01	0.41
1:E:128:ALA:HB2	1:E:134:HIS:ND1	2.35	0.41
1:E:88:LEU:N	2:F:119:GLU:OE1	2.52	0.41
1:A:127:ALA:O	1:A:128:ALA:HB3	2.21	0.41
1:C:136:THR:OG1	1:C:296:ARG:NE	2.49	0.41
1:E:47:CYS:O	1:E:104:MET:HA	2.21	0.41
1:E:278:ASP:OD1	1:E:278:ASP:N	2.54	0.41
2:B:140:TYR:HB2	4:B:301:HOH:O	2.20	0.41
2:B:102:ARG:NH2	2:B:124:SER:OG	2.54	0.40
2:D:128:ARG:HG2	2:D:135:ALA:HB3	2.03	0.40
1:E:244:LYS:O	1:E:248:VAL:HG13	2.22	0.40
1:E:35:GLN:HG2	1:E:38:LEU:HB2	2.03	0.40
1:C:174:GLN:NE2	1:C:201:MET:HG3	2.37	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	299/310 (96%)	288 (96%)	10 (3%)	1 (0%)	41	70
1	C	299/310 (96%)	290 (97%)	8 (3%)	1 (0%)	41	70
1	E	300/310 (97%)	289 (96%)	10 (3%)	1 (0%)	41	70
2	B	130/153 (85%)	118 (91%)	12 (9%)	0	100	100
2	D	129/153 (84%)	119 (92%)	10 (8%)	0	100	100
2	F	130/153 (85%)	119 (92%)	11 (8%)	0	100	100
All	All	1287/1389 (93%)	1223 (95%)	61 (5%)	3 (0%)	47	77

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	128	ALA
1	A	128	ALA
1	E	128	ALA

### 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	253/261 (97%)	248 (98%)	5 (2%)	55	78
1	C	252/261 (97%)	248 (98%)	4 (2%)	62	83
1	E	252/261 (97%)	248 (98%)	4 (2%)	62	83
2	B	104/137 (76%)	101 (97%)	3 (3%)	42	70
2	D	99/137 (72%)	98 (99%)	1 (1%)	76	89
2	F	112/137 (82%)	112 (100%)	0	100	100
All	All	1072/1194 (90%)	1055 (98%)	17 (2%)	62	83

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	PHE
1	A	134	HIS
1	A	146	GLN
1	A	244	LYS
1	A	261	MET
2	B	46	LEU
2	B	63	ASN
2	B	69	ASP
1	C	59	PHE
1	C	261	MET
1	C	278	ASP
1	C	285	TYR
2	D	104	ASP
1	E	59	PHE
1	E	65	ARG
1	E	261	MET
1	E	285	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	133	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](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	303/310 (97%)	-0.21	3 (0%) 82 63	35, 62, 97, 120	0
1	C	303/310 (97%)	-0.27	1 (0%) 94 85	35, 52, 85, 123	0
1	E	304/310 (98%)	-0.23	2 (0%) 87 72	32, 55, 96, 141	0
2	B	134/153 (87%)	0.36	13 (9%) 7 2	49, 96, 128, 152	0
2	D	133/153 (86%)	0.44	13 (9%) 7 2	54, 98, 142, 164	0
2	F	134/153 (87%)	0.21	4 (2%) 50 25	44, 81, 118, 135	0
All	All	1311/1389 (94%)	-0.06	36 (2%) 54 28	32, 63, 118, 164	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	93	GLY	3.9
1	C	77	ALA	3.9
2	B	93	GLY	3.5
1	E	1	ALA	3.4
1	A	309	VAL	3.4
2	D	18	ILE	3.4
2	B	84	ASN	3.3
2	F	92	VAL	3.2
2	B	18	ILE	3.2
1	A	77	ALA	3.1
2	D	87	ASP	3.0
2	B	59	ILE	2.9
2	B	87	ASP	2.8
2	D	83	VAL	2.8
2	B	91	VAL	2.8
2	D	59	ILE	2.7
2	D	91	VAL	2.7
2	F	80	GLN	2.6
2	D	84	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
2	F	93	GLY	2.5
2	B	80	GLN	2.5
2	D	48	LEU	2.4
2	B	49	PRO	2.3
1	A	78	ASN	2.3
2	B	74	LEU	2.2
1	E	84	LYS	2.2
2	B	78	ALA	2.2
2	F	83	VAL	2.2
2	D	44	ILE	2.1
2	B	99	LEU	2.1
2	D	49	PRO	2.0
2	D	17	VAL	2.0
2	B	86	ILE	2.0
2	D	79	PRO	2.0
2	D	85	ARG	2.0
2	B	153	ASN	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	ZN	D	201	1/1	0.98	0.12	66,66,66,66	0
3	ZN	B	201	1/1	0.99	0.08	58,58,58,58	0
3	ZN	F	201	1/1	0.99	0.13	48,48,48,48	0



## 6.5 Other polymers

There are no such residues in this entry.