



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:37 AM UTC

PDB ID : 6KJ9 / pdb_00006kj9
Title : E. coli ATCase catalytic subunit mutant - G128/130A
Authors : Lei, Z.; Zheng, J.; Jia, Z.C.
Deposited on : 2019-07-22
Resolution : 2.50 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

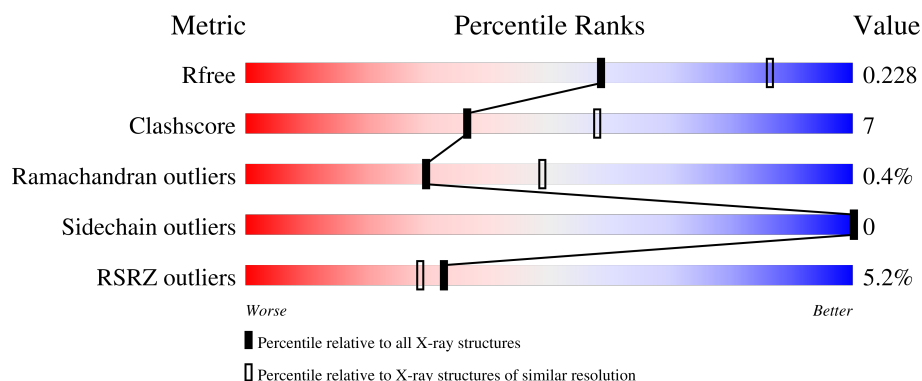
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

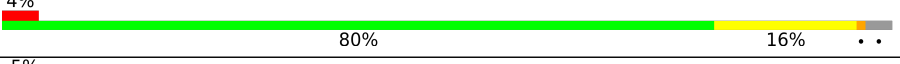
The reported resolution of this entry is 2.50 Å.

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	310	
1	B	310	
1	C	310	
1	D	310	
1	E	310	

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Mol	Chain	Length	Quality of chain
1	F	310	4% 83% 13%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14188 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	301	Total	C	N	O	S	0	0	0
			2339	1482	406	442	9			
1	B	301	Total	C	N	O	S	0	0	0
			2331	1480	405	437	9			
1	C	291	Total	C	N	O	S	0	0	0
			2243	1423	388	423	9			
1	D	284	Total	C	N	O	S	0	0	0
			2216	1407	386	414	9			
1	E	283	Total	C	N	O	S	0	0	0
			2163	1377	370	408	8			
1	F	298	Total	C	N	O	S	0	0	0
			2313	1469	396	439	9			

There are 12 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
B	128	ALA	GLY	engineered mutation	UNP P0A786
B	130	ALA	GLY	engineered mutation	UNP P0A786
C	128	ALA	GLY	engineered mutation	UNP P0A786
C	130	ALA	GLY	engineered mutation	UNP P0A786
D	128	ALA	GLY	engineered mutation	UNP P0A786
D	130	ALA	GLY	engineered mutation	UNP P0A786
E	128	ALA	GLY	engineered mutation	UNP P0A786
E	130	ALA	GLY	engineered mutation	UNP P0A786
F	128	ALA	GLY	engineered mutation	UNP P0A786
F	130	ALA	GLY	engineered mutation	UNP P0A786

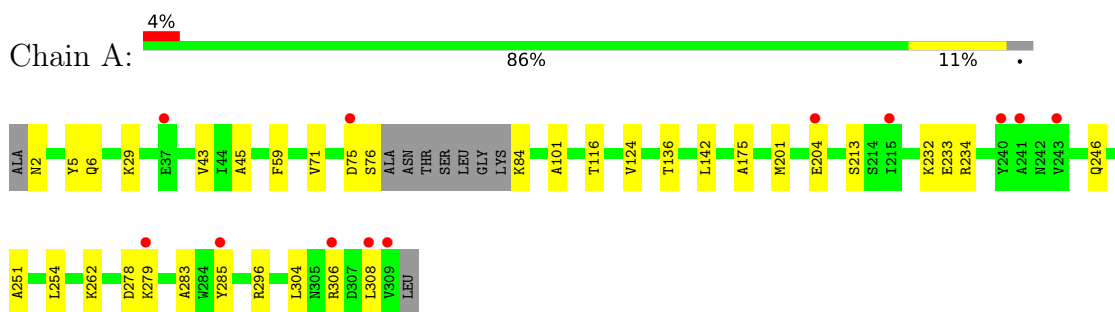
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	69	Total 69	O 69	0	0
2	B	120	Total 120	O 120	0	0
2	C	89	Total 89	O 89	0	0
2	D	120	Total 120	O 120	0	0
2	E	47	Total 47	O 47	0	0
2	F	138	Total 138	O 138	0	0

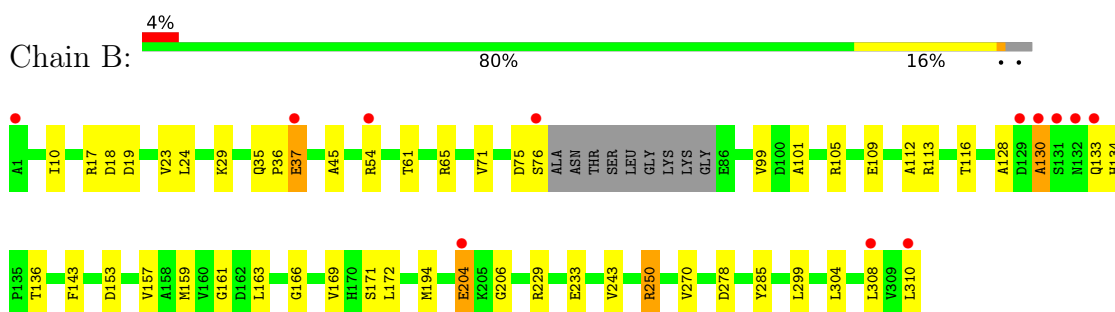
3 Residue-property plots [i](#)

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.

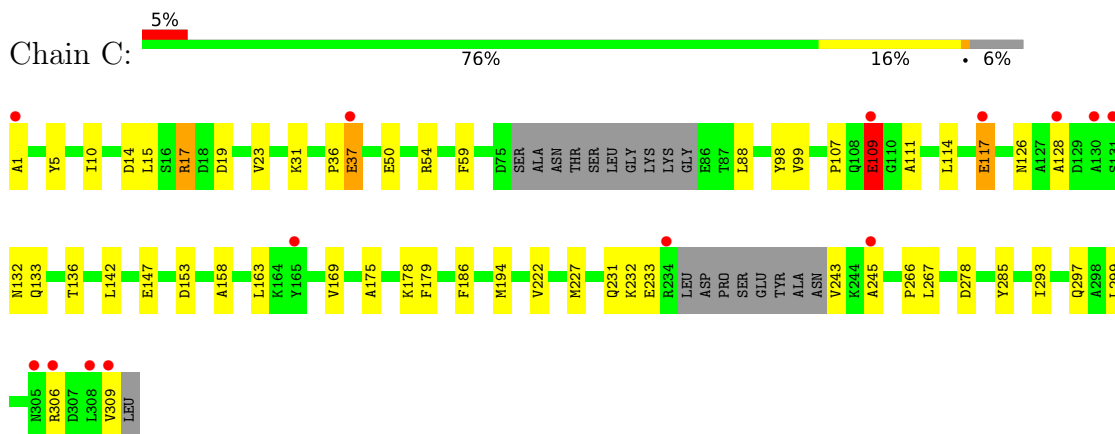
- Molecule 1: Aspartate carbamoyltransferase catalytic subunit



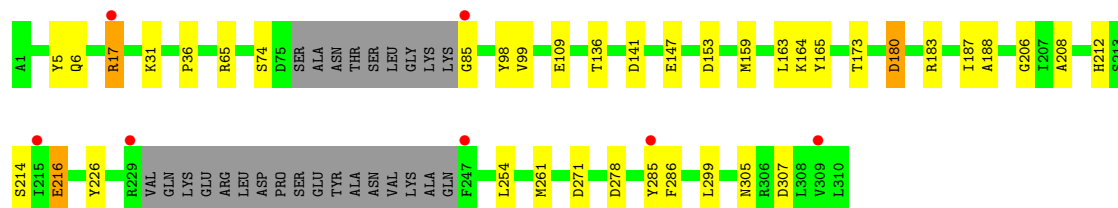
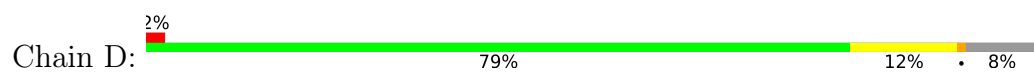
- Molecule 1: Aspartate carbamoyltransferase catalytic subunit



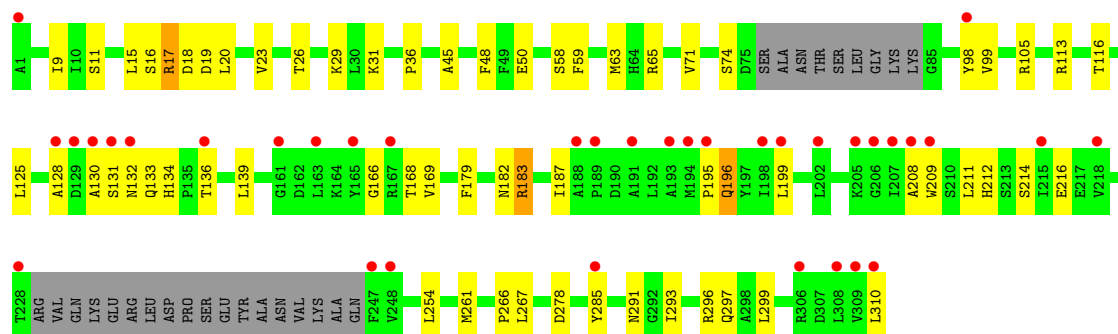
- Molecule 1: Aspartate carbamoyltransferase catalytic subunit



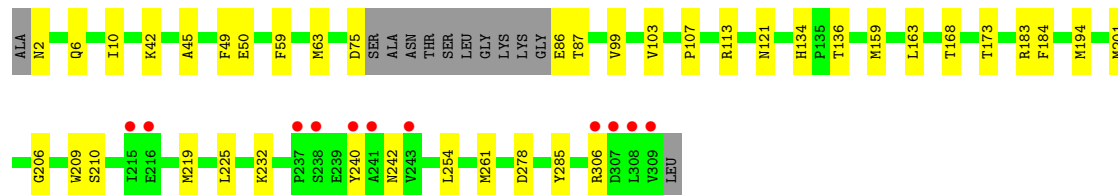
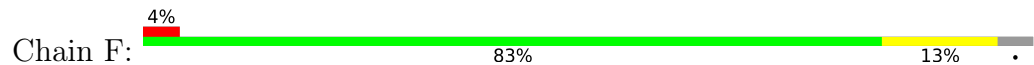
- Molecule 1: Aspartate carbamoyltransferase catalytic subunit



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- Molecule 1: Aspartate carbamoyltransferase catalytic subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.76Å 96.73Å 121.69Å 90.00° 94.16° 90.00°	Depositor
Resolution (Å)	48.36 – 2.50 48.36 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.36-2.50) 93.4 (48.36-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.14-3260	Depositor
R, R_{free}	0.192 , 0.228 0.192 , 0.228	Depositor DCC
R_{free} test set	2005 reflections (3.05%)	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.4	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	0.94	EDS
Total number of atoms	14188	wwPDB-VP
Average B, all atoms (Å ²)	45.0	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 40.52 % 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 2.7037e-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)

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	0.31	0/2384	0.56	0/3240
1	B	0.45	1/2376 (0.0%)	0.74	8/3231 (0.2%)
1	C	0.41	1/2285 (0.0%)	0.66	4/3109 (0.1%)
1	D	0.37	0/2258	0.66	5/3066 (0.2%)
1	E	0.34	0/2205	0.68	9/3004 (0.3%)
1	F	0.35	0/2358	0.58	1/3207 (0.0%)
All	All	0.37	2/13866 (0.0%)	0.65	27/18857 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	161	GLY	C-O	6.13	1.27	1.24
1	C	17	ARG	CZ-NH1	-5.63	1.24	1.32

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	37	GLU	CA-CB-CG	-13.69	86.72	114.10
1	D	17	ARG	CB-CG-CD	-9.77	88.83	111.30
1	B	37	GLU	CB-CG-CD	8.87	127.67	112.60
1	E	183	ARG	CB-CG-CD	-8.71	91.28	111.30
1	B	204	GLU	CA-CB-CG	-7.86	98.39	114.10
1	C	37	GLU	CA-CB-CG	-7.43	99.25	114.10
1	C	117	GLU	CA-CB-CG	7.26	128.62	114.10
1	E	196	GLN	CA-CB-CG	-7.04	100.03	114.10
1	B	54	ARG	CG-CD-NE	-6.82	97.01	112.00
1	E	17	ARG	CA-CB-CG	-6.82	100.47	114.10
1	C	17	ARG	CB-CG-CD	6.78	126.88	111.30
1	D	180	ASP	N-CA-CB	-6.20	100.33	110.43
1	B	29	LYS	CA-CB-CG	5.99	126.08	114.10
1	F	306	ARG	CD-NE-CZ	5.87	132.61	124.40
1	D	17	ARG	CA-CB-CG	5.67	125.44	114.10
1	B	54	ARG	CA-CB-CG	-5.65	102.80	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	17	ARG	NE-CZ-NH2	-5.62	114.14	119.20
1	D	216	GLU	N-CA-CB	-5.60	101.59	110.28
1	D	216	GLU	CB-CG-CD	-5.55	103.16	112.60
1	E	196	GLN	N-CA-CB	5.41	118.55	110.22
1	B	250	ARG	NE-CZ-NH2	-5.41	114.33	119.20
1	C	109	GLU	CB-CG-CD	5.15	121.35	112.60
1	B	250	ARG	CB-CG-CD	-5.11	99.54	111.30
1	E	195	PRO	CA-C-N	-5.04	113.03	120.28
1	E	195	PRO	C-N-CA	-5.04	113.03	120.28
1	E	16	SER	CA-C-N	-5.00	113.94	120.44
1	E	16	SER	C-N-CA	-5.00	113.94	120.44

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	2339	0	2319	27	0
1	B	2331	0	2319	37	0
1	C	2243	0	2211	48	0
1	D	2216	0	2219	22	0
1	E	2163	0	2124	40	0
1	F	2313	0	2291	27	0
2	A	69	0	0	7	0
2	B	120	0	0	7	0
2	C	89	0	0	4	1
2	D	120	0	0	4	0
2	E	47	0	0	5	1
2	F	138	0	0	4	1
All	All	14188	0	13483	196	2

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 7.

All (196) 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:C:17:ARG:NH2	1:C:179:PHE:HA	1.50	1.25
1:C:17:ARG:HH22	1:C:179:PHE:HA	0.87	0.99
1:C:17:ARG:HH22	1:C:179:PHE:CA	1.80	0.90
1:F:75:ASP:OD2	2:F:401:HOH:O	1.91	0.87
1:E:182:ASN:O	2:E:401:HOH:O	1.93	0.87
1:A:2:ASN:HB3	1:A:308:LEU:HD21	1.57	0.86
1:E:50:GLU:OE1	2:E:402:HOH:O	1.99	0.81
1:F:183:ARG:NH2	1:F:210:SER:OG	2.12	0.81
1:C:36:PRO:C	1:C:37:GLU:HG2	2.06	0.78
1:C:1:ALA:N	2:C:402:HOH:O	2.16	0.77
1:B:153:ASP:OD1	2:B:402:HOH:O	2.03	0.76
1:B:36:PRO:O	1:B:37:GLU:HG2	1.85	0.75
1:B:130:ALA:HB3	1:B:133:GLN:HB2	1.68	0.74
1:C:17:ARG:NH1	1:C:179:PHE:CD2	2.53	0.74
1:B:229:ARG:HH11	1:B:270:VAL:HG21	1.54	0.72
1:C:14:ASP:OD2	2:C:401:HOH:O	2.07	0.72
1:C:153:ASP:OD1	1:C:179:PHE:HB3	1.89	0.71
1:E:58:SER:OG	1:E:296:ARG:NH1	2.24	0.71
1:E:196:GLN:O	1:E:196:GLN:HG3	1.90	0.71
1:C:31:LYS:NZ	1:C:147:GLU:OE2	2.18	0.70
1:A:75:ASP:OD1	1:A:76:SER:N	2.23	0.70
1:C:278:ASP:OD1	1:C:285:TYR:OH	2.10	0.68
1:A:29:LYS:NZ	2:A:401:HOH:O	2.08	0.68
1:C:114:LEU:HA	1:C:117:GLU:OE1	1.93	0.68
1:C:50:GLU:HB2	1:C:107:PRO:HD3	1.77	0.67
1:C:17:ARG:NH2	1:C:178:LYS:O	2.28	0.66
1:B:243:VAL:O	2:B:404:HOH:O	2.14	0.66
1:A:234:ARG:NH1	2:A:409:HOH:O	2.26	0.66
1:C:243:VAL:N	2:C:407:HOH:O	2.28	0.66
1:C:17:ARG:NH2	1:C:179:PHE:CA	2.44	0.65
1:D:305:ASN:ND2	1:D:307:ASP:O	2.28	0.65
1:F:2:ASN:N	2:F:404:HOH:O	2.30	0.64
1:E:15:LEU:HD22	1:E:19:ASP:HB3	1.80	0.64
1:A:59:PHE:HZ	1:A:136:THR:HG21	1.63	0.63
1:E:17:ARG:O	1:E:20:LEU:N	2.32	0.63
1:A:75:ASP:OD2	2:A:402:HOH:O	2.16	0.61
1:E:9:ILE:HB	1:E:125:LEU:HD23	1.83	0.61
1:E:296:ARG:NH2	2:E:408:HOH:O	2.27	0.61
1:B:134:HIS:NE2	1:B:171:SER:OG	2.31	0.60
1:B:36:PRO:C	1:B:37:GLU:HG2	2.23	0.60
1:C:109:GLU:OE2	1:C:132:ASN:HB2	2.01	0.60
1:F:45:ALA:HB2	1:F:99:VAL:HG11	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:GLN:HG3	1:B:310:LEU:HD22	1.82	0.60
1:C:88:LEU:HD23	1:C:114:LEU:HD22	1.84	0.59
1:C:163:LEU:HD22	1:C:169:VAL:HG11	1.83	0.59
1:C:54:ARG:NH2	1:C:267:LEU:HB3	2.17	0.59
1:F:50:GLU:HB3	1:F:107:PRO:HD3	1.85	0.59
1:C:17:ARG:NH1	1:C:179:PHE:CE2	2.71	0.58
1:B:105:ARG:HG3	1:B:128:ALA:HA	1.83	0.58
1:E:179:PHE:O	2:E:403:HOH:O	2.17	0.58
1:C:36:PRO:O	1:C:37:GLU:HG2	2.04	0.57
1:C:231:GLN:O	1:C:233:GLU:N	2.37	0.57
1:A:279:LYS:O	2:A:403:HOH:O	2.18	0.57
1:B:109:GLU:O	1:B:133:GLN:NE2	2.32	0.57
1:F:219:MET:HE1	1:F:225:LEU:HD22	1.87	0.57
1:B:308:LEU:C	1:B:308:LEU:HD23	2.30	0.56
1:A:59:PHE:CE1	1:A:296:ARG:HG2	2.39	0.56
1:C:309:VAL:O	2:C:403:HOH:O	2.17	0.56
1:D:214:SER:OG	1:D:216:GLU:OE1	2.08	0.56
1:A:59:PHE:CZ	1:A:136:THR:HG21	2.41	0.56
1:B:250:ARG:NH1	2:B:412:HOH:O	2.37	0.56
1:C:98:TYR:HB2	1:C:99:VAL:HG13	1.87	0.56
1:E:278:ASP:OD1	1:E:285:TYR:OH	2.23	0.56
1:D:85:GLY:N	2:D:409:HOH:O	2.39	0.55
1:B:278:ASP:OD1	1:B:285:TYR:OH	2.21	0.55
1:E:296:ARG:NE	2:E:408:HOH:O	2.39	0.55
1:E:11:SER:HA	1:E:133:GLN:CB	2.37	0.55
1:A:84:LYS:N	2:A:413:HOH:O	2.39	0.54
1:A:278:ASP:OD1	1:A:285:TYR:OH	2.24	0.54
1:B:233:GLU:OE2	1:D:17:ARG:NH2	2.40	0.54
1:A:201:MET:HA	1:A:204:GLU:OE1	2.08	0.54
1:B:250:ARG:NH1	2:B:413:HOH:O	2.40	0.54
1:C:136:THR:HG22	1:C:299:LEU:HD12	1.90	0.54
1:C:306:ARG:HH21	1:C:306:ARG:HG3	1.73	0.53
1:E:17:ARG:O	1:E:17:ARG:HG2	2.07	0.53
1:E:26:THR:HG23	1:E:310:LEU:HD13	1.90	0.53
1:F:6:GLN:NE2	1:F:121:ASN:O	2.41	0.53
1:B:204:GLU:O	1:B:204:GLU:HG3	2.08	0.53
1:B:166:GLY:O	1:B:169:VAL:HG22	2.09	0.53
1:C:54:ARG:NH1	1:F:86:GLU:OE1	2.41	0.53
1:C:5:TYR:CD2	1:C:306:ARG:NH2	2.77	0.53
1:C:54:ARG:HH12	1:F:86:GLU:CD	2.16	0.53
1:B:18:ASP:OD1	2:B:405:HOH:O	2.19	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:158:ALA:HB2	1:C:222:VAL:HG11	1.91	0.52
1:A:142:LEU:HD11	1:A:175:ALA:HB1	1.91	0.52
1:D:278:ASP:OD1	1:D:285:TYR:OH	2.27	0.51
1:C:186:PHE:CD1	1:C:194:MET:HE3	2.46	0.51
1:E:187:ILE:HG13	1:E:212:HIS:HB2	1.93	0.51
1:F:63:MET:HE2	1:F:103:VAL:HG21	1.94	0.50
1:F:75:ASP:CG	2:F:401:HOH:O	2.44	0.50
1:E:113:ARG:O	1:E:116:THR:OG1	2.23	0.50
1:B:45:ALA:HB2	1:B:99:VAL:HG11	1.94	0.50
1:E:196:GLN:HA	1:E:199:LEU:HD12	1.93	0.50
1:B:113:ARG:O	1:B:116:THR:OG1	2.27	0.50
1:C:54:ARG:NH1	1:F:86:GLU:CD	2.70	0.50
1:E:105:ARG:HG3	1:E:128:ALA:HA	1.94	0.49
1:B:159:MET:HE2	1:B:172:LEU:HD23	1.93	0.49
1:D:163:LEU:HG	1:D:188:ALA:HB2	1.94	0.49
1:A:232:LYS:HG2	1:A:233:GLU:OE1	2.13	0.49
1:E:166:GLY:O	1:E:169:VAL:HG22	2.12	0.49
1:A:45:ALA:HA	1:A:71:VAL:O	2.13	0.49
1:F:278:ASP:OD1	1:F:285:TYR:OH	2.29	0.49
1:D:271:ASP:OD1	2:D:401:HOH:O	2.20	0.49
1:B:10:ILE:HD12	1:B:112:ALA:HB1	1.95	0.48
1:B:35:GLN:HG3	1:B:310:LEU:CD2	2.43	0.48
1:E:136:THR:CG2	1:E:299:LEU:HD12	2.43	0.48
1:D:164:LYS:HD3	1:D:165:TYR:CE2	2.48	0.48
1:C:163:LEU:HB3	1:C:194:MET:HE2	1.96	0.47
1:F:163:LEU:HD23	1:F:194:MET:HG2	1.97	0.47
1:D:285:TYR:HE2	1:D:286:PHE:CZ	2.33	0.47
1:A:283:ALA:O	2:A:404:HOH:O	2.20	0.47
1:B:24:LEU:HD22	1:B:143:PHE:HB2	1.96	0.47
1:A:5:TYR:CE1	1:A:306:ARG:HG3	2.50	0.47
1:B:250:ARG:NE	2:B:409:HOH:O	2.27	0.47
1:C:111:ALA:O	1:C:114:LEU:HB3	2.15	0.46
1:F:59:PHE:CZ	1:F:136:THR:HG21	2.50	0.46
1:D:109:GLU:HA	2:D:455:HOH:O	2.15	0.46
1:D:183:ARG:HG3	1:D:208:ALA:HB3	1.97	0.46
1:E:214:SER:OG	1:E:216:GLU:OE1	2.33	0.46
1:A:251:ALA:HA	1:A:254:LEU:HD12	1.98	0.46
1:C:54:ARG:HH21	1:C:267:LEU:HB3	1.79	0.46
1:D:153:ASP:OD1	1:D:180:ASP:HB2	2.16	0.46
1:E:17:ARG:O	1:E:18:ASP:C	2.59	0.46
1:E:36:PRO:HA	1:E:65:ARG:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:136:THR:HG22	1:E:299:LEU:HD12	1.98	0.45
1:C:306:ARG:HG3	1:C:306:ARG:NH2	2.29	0.45
1:D:187:ILE:HG13	1:D:212:HIS:HB2	1.99	0.45
1:C:59:PHE:CZ	1:C:136:THR:HG21	2.51	0.45
1:D:36:PRO:HA	1:D:65:ARG:O	2.15	0.45
1:D:98:TYR:HB2	1:D:99:VAL:HG13	1.98	0.45
1:E:209:TRP:HZ3	1:E:211:LEU:CD2	2.30	0.45
1:E:31:LYS:NZ	1:E:291:ASN:OD1	2.49	0.45
1:B:35:GLN:CG	1:B:310:LEU:HD22	2.46	0.45
1:B:136:THR:HG22	1:B:299:LEU:HD12	1.99	0.45
1:C:299:LEU:HD23	1:C:299:LEU:HA	1.90	0.45
1:D:5:TYR:CE2	1:D:6:GLN:HG3	2.52	0.45
1:E:183:ARG:HG3	1:E:208:ALA:HB3	1.98	0.45
1:F:159:MET:HE3	1:F:173:THR:OG1	2.17	0.45
1:C:227:MET:HE3	1:C:227:MET:HB3	1.82	0.44
1:C:293:ILE:HG22	1:C:297:GLN:HE21	1.82	0.44
1:E:23:VAL:HG11	1:E:139:LEU:HD22	1.99	0.44
1:C:15:LEU:HD23	1:C:19:ASP:HB3	1.99	0.44
1:C:243:VAL:HG12	1:C:245:ALA:HB3	1.98	0.44
1:D:254:LEU:HD22	1:D:261:MET:HE1	1.99	0.44
1:E:134:HIS:CD2	1:E:168:THR:HG22	2.52	0.44
1:F:254:LEU:HD22	1:F:261:MET:HE1	2.00	0.44
1:A:262:LYS:NZ	2:A:416:HOH:O	2.50	0.44
1:B:45:ALA:HA	1:B:71:VAL:O	2.18	0.44
1:A:213:SER:O	1:A:246:GLN:NE2	2.24	0.44
1:C:10:ILE:HA	1:C:126:ASN:HB2	2.00	0.44
1:B:163:LEU:HB3	1:B:194:MET:HE2	1.98	0.43
1:E:45:ALA:HA	1:E:71:VAL:O	2.17	0.43
1:E:254:LEU:HD22	1:E:261:MET:HE1	2.00	0.43
1:F:49:PHE:O	1:F:75:ASP:HB3	2.17	0.43
1:B:101:ALA:HB2	1:B:304:LEU:HD21	2.00	0.43
1:A:5:TYR:CD1	1:A:306:ARG:HG3	2.54	0.42
1:B:36:PRO:HA	1:B:65:ARG:O	2.18	0.42
1:F:242:ASN:ND2	2:F:413:HOH:O	2.52	0.42
1:B:17:ARG:NH1	2:B:402:HOH:O	2.53	0.42
1:F:86:GLU:HG2	1:F:87:THR:N	2.33	0.42
1:E:29:LYS:HE2	1:E:29:LYS:HB2	1.83	0.42
1:E:293:ILE:HG22	1:E:297:GLN:HE21	1.84	0.42
1:A:5:TYR:CE2	1:A:6:GLN:HG3	2.55	0.42
1:A:234:ARG:HD3	1:A:234:ARG:HA	1.75	0.42
1:B:36:PRO:C	1:B:37:GLU:CG	2.90	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:LEU:O	1:C:117:GLU:HB2	2.20	0.42
1:E:48:PHE:O	1:E:74:SER:HA	2.20	0.42
1:F:42:LYS:HA	1:F:42:LYS:HD3	1.87	0.42
1:A:101:ALA:HB2	1:A:304:LEU:HD21	2.01	0.41
1:C:19:ASP:O	1:C:23:VAL:HG23	2.20	0.41
1:F:10:ILE:HG21	1:F:113:ARG:HB2	2.02	0.41
1:F:134:HIS:NE2	1:F:168:THR:HA	2.35	0.41
1:B:157:VAL:CG1	1:B:159:MET:HE3	2.51	0.41
1:C:266:PRO:O	1:C:267:LEU:HB2	2.19	0.41
1:E:17:ARG:HE	1:E:17:ARG:HB3	1.50	0.41
1:E:131:SER:O	1:E:132:ASN:C	2.64	0.41
1:E:266:PRO:O	1:E:267:LEU:HB2	2.20	0.41
1:A:116:THR:HG22	1:A:124:VAL:HB	2.03	0.41
1:D:31:LYS:NZ	1:D:147:GLU:OE1	2.43	0.41
1:D:141:ASP:OD2	1:D:226:TYR:OH	2.31	0.41
1:E:59:PHE:O	1:E:63:MET:HG3	2.21	0.41
1:F:232:LYS:HD2	1:F:240:TYR:CD2	2.55	0.41
1:D:136:THR:HG22	1:D:299:LEU:HD12	2.03	0.41
1:D:159:MET:HE3	1:D:173:THR:OG1	2.21	0.41
1:C:142:LEU:HD11	1:C:175:ALA:HB1	2.01	0.41
1:F:201:MET:HE3	1:F:201:MET:HB3	1.95	0.41
1:A:59:PHE:CZ	1:A:296:ARG:HG2	2.56	0.41
1:B:75:ASP:O	1:B:76:SER:OG	2.33	0.41
1:C:128:ALA:HB3	1:C:133:GLN:O	2.20	0.40
1:B:19:ASP:O	1:B:23:VAL:HG23	2.21	0.40
1:E:299:LEU:HD23	1:E:299:LEU:HA	1.80	0.40
1:F:184:PHE:O	1:F:209:TRP:HA	2.21	0.40
1:A:43:VAL:HG11	1:B:61:THR:HG23	2.04	0.40
1:E:98:TYR:HB2	1:E:99:VAL:HG13	2.03	0.40
1:F:194:MET:HE3	1:F:194:MET:HB3	1.97	0.40
1:D:74:SER:HB3	2:D:493:HOH:O	2.21	0.40

All (2) 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)
2:E:439:HOH:O	2:F:474:HOH:O[1_565]	1.90	0.30
2:C:481:HOH:O	2:C:484:HOH:O[2_557]	2.11	0.09

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	297/310 (96%)	286 (96%)	11 (4%)	0	100	100
1	B	297/310 (96%)	283 (95%)	12 (4%)	2 (1%)	18	34
1	C	285/310 (92%)	274 (96%)	9 (3%)	2 (1%)	18	34
1	D	278/310 (90%)	265 (95%)	12 (4%)	1 (0%)	30	49
1	E	277/310 (89%)	267 (96%)	9 (3%)	1 (0%)	30	49
1	F	294/310 (95%)	281 (96%)	12 (4%)	1 (0%)	36	55
All	All	1728/1860 (93%)	1656 (96%)	65 (4%)	7 (0%)	30	49

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	232	LYS
1	E	130	ALA
1	B	206	GLY
1	C	109	GLU
1	D	206	GLY
1	F	206	GLY
1	B	130	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	251/261 (96%)	251 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	249/261 (95%)	249 (100%)	0	100	100
1	C	238/261 (91%)	238 (100%)	0	100	100
1	D	239/261 (92%)	239 (100%)	0	100	100
1	E	228/261 (87%)	228 (100%)	0	100	100
1	F	249/261 (95%)	249 (100%)	0	100	100
All	All	1454/1566 (93%)	1454 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	297	GLN
1	A	305	ASN
1	B	132	ASN
1	B	255	HIS
1	B	260	ASN
1	C	21	ASN
1	E	35	GLN
1	E	64	HIS
1	E	146	GLN
1	E	196	GLN
1	E	297	GLN
1	F	41	HIS
1	F	297	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	301/310 (97%)	0.13	12 (3%) 42 38	27, 49, 83, 100	0
1	B	301/310 (97%)	-0.09	12 (3%) 42 38	24, 38, 66, 92	0
1	C	291/310 (93%)	0.07	14 (4%) 35 31	23, 42, 68, 100	0
1	D	284/310 (91%)	-0.31	7 (2%) 58 54	19, 32, 60, 78	0
1	E	283/310 (91%)	0.66	36 (12%) 8 6	31, 55, 101, 114	0
1	F	298/310 (96%)	-0.13	11 (3%) 45 40	23, 35, 73, 110	0
All	All	1758/1860 (94%)	0.05	92 (5%) 33 29	19, 41, 83, 114	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	309	VAL	5.2
1	B	1	ALA	4.5
1	A	309	VAL	4.4
1	E	1	ALA	4.4
1	C	130	ALA	4.0
1	C	1	ALA	3.5
1	E	161	GLY	3.5
1	C	309	VAL	3.4
1	F	241	ALA	3.4
1	B	37	GLU	3.4
1	E	228	THR	3.3
1	B	54	ARG	3.2
1	E	132	ASN	3.2
1	E	202	LEU	3.1
1	F	308	LEU	3.1
1	E	130	ALA	3.1
1	E	310	LEU	3.1
1	C	234	ARG	3.1
1	C	308	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	76	SER	3.0
1	B	310	LEU	3.0
1	F	306	ARG	3.0
1	E	247	PHE	3.0
1	E	206	GLY	2.9
1	E	128	ALA	2.8
1	E	163	LEU	2.8
1	A	285	TYR	2.8
1	E	198	ILE	2.8
1	B	308	LEU	2.8
1	D	309	VAL	2.8
1	A	241	ALA	2.8
1	E	308	LEU	2.7
1	B	130	ALA	2.7
1	E	194	MET	2.7
1	E	215	ILE	2.7
1	F	243	VAL	2.7
1	C	128	ALA	2.7
1	C	109	GLU	2.6
1	E	189	PRO	2.6
1	B	131	SER	2.6
1	E	195	PRO	2.5
1	A	37	GLU	2.5
1	E	191	ALA	2.5
1	B	204	GLU	2.5
1	B	132	ASN	2.5
1	E	193	ALA	2.5
1	D	85	GLY	2.5
1	C	131	SER	2.5
1	E	131	SER	2.5
1	E	209	TRP	2.5
1	B	133	GLN	2.4
1	C	306	ARG	2.4
1	C	305	ASN	2.4
1	E	205	LYS	2.4
1	E	165	TYR	2.4
1	D	215	ILE	2.4
1	A	308	LEU	2.4
1	A	215	ILE	2.4
1	A	75	ASP	2.4
1	E	218	VAL	2.4
1	E	129	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	279	LYS	2.3
1	F	215	ILE	2.3
1	D	247	PHE	2.3
1	A	306	ARG	2.3
1	D	229	ARG	2.3
1	E	167	ARG	2.3
1	E	285	TYR	2.3
1	E	306	ARG	2.3
1	F	216	GLU	2.2
1	E	309	VAL	2.2
1	E	188	ALA	2.2
1	E	199	LEU	2.2
1	E	207	ILE	2.2
1	C	117	GLU	2.2
1	E	136	THR	2.2
1	C	37	GLU	2.2
1	D	17	ARG	2.1
1	E	248	VAL	2.1
1	A	204	GLU	2.1
1	C	165	TYR	2.1
1	D	285	TYR	2.1
1	F	238	SER	2.1
1	E	208	ALA	2.1
1	F	240	TYR	2.1
1	A	240	TYR	2.1
1	B	129	ASP	2.1
1	F	237	PRO	2.1
1	A	243	VAL	2.0
1	C	245	ALA	2.0
1	E	98	TYR	2.0
1	F	307	ASP	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.