



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:52 AM GMT

PDB ID : 3ND5
Title : Crystal structure of phosphopantetheine adenylyltransferase (PPAT) from *Enterococcus faecalis*
Authors : Yoon, H.J.; Lee, H.H.; Suh, S.W.
Deposited on : 2010-06-07
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

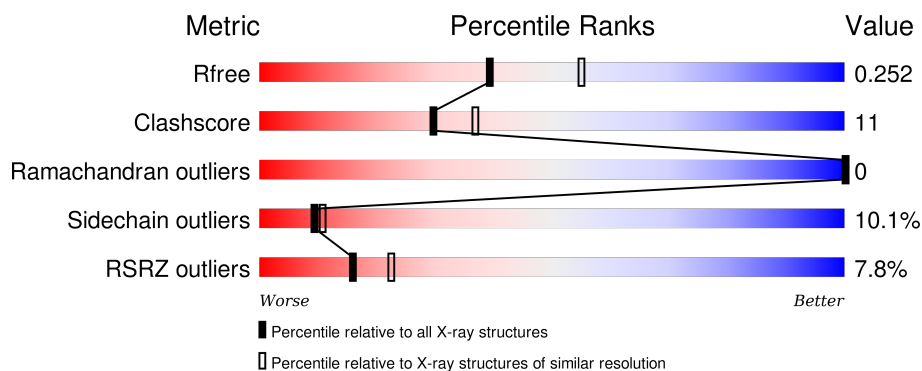
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	171	<div> <div>3%</div> <div>73%</div> <div>14%</div> <div>• •</div> <div>11%</div> </div>
1	B	171	<div> <div>9%</div> <div>68%</div> <div>17%</div> <div>•</div> <div>11%</div> </div>
1	C	171	<div> <div>9%</div> <div>62%</div> <div>20%</div> <div>6%</div> <div>•</div> <div>11%</div> </div>
1	D	171	<div> <div>10%</div> <div>63%</div> <div>22%</div> <div>• •</div> <div>11%</div> </div>
1	E	171	<div> <div>4%</div> <div>71%</div> <div>11%</div> <div>5%</div> <div>•</div> <div>11%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	171	<div><div></div><div>7%</div><div>59%</div><div>23%</div><div>6%</div><div>•</div><div>11%</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7878 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 Phosphopantetheine adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	152	Total	C	N	O	S	0	0	0
			1224	791	201	227	5			
1	B	152	Total	C	N	O	S	0	0	0
			1224	791	201	227	5			
1	C	152	Total	C	N	O	S	0	0	0
			1224	791	201	227	5			
1	D	152	Total	C	N	O	S	0	0	0
			1224	791	201	227	5			
1	E	152	Total	C	N	O	S	0	0	0
			1224	791	201	227	5			
1	F	152	Total	C	N	O	S	0	0	0
			1224	791	201	227	5			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	164	LEU	-	EXPRESSION TAG	UNP Q831P9
A	165	GLU	-	EXPRESSION TAG	UNP Q831P9
A	166	HIS	-	EXPRESSION TAG	UNP Q831P9
A	167	HIS	-	EXPRESSION TAG	UNP Q831P9
A	168	HIS	-	EXPRESSION TAG	UNP Q831P9
A	169	HIS	-	EXPRESSION TAG	UNP Q831P9
A	170	HIS	-	EXPRESSION TAG	UNP Q831P9
A	171	HIS	-	EXPRESSION TAG	UNP Q831P9
B	164	LEU	-	EXPRESSION TAG	UNP Q831P9
B	165	GLU	-	EXPRESSION TAG	UNP Q831P9
B	166	HIS	-	EXPRESSION TAG	UNP Q831P9
B	167	HIS	-	EXPRESSION TAG	UNP Q831P9
B	168	HIS	-	EXPRESSION TAG	UNP Q831P9
B	169	HIS	-	EXPRESSION TAG	UNP Q831P9
B	170	HIS	-	EXPRESSION TAG	UNP Q831P9
B	171	HIS	-	EXPRESSION TAG	UNP Q831P9
C	164	LEU	-	EXPRESSION TAG	UNP Q831P9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	165	GLU	-	EXPRESSION TAG	UNP Q831P9
C	166	HIS	-	EXPRESSION TAG	UNP Q831P9
C	167	HIS	-	EXPRESSION TAG	UNP Q831P9
C	168	HIS	-	EXPRESSION TAG	UNP Q831P9
C	169	HIS	-	EXPRESSION TAG	UNP Q831P9
C	170	HIS	-	EXPRESSION TAG	UNP Q831P9
C	171	HIS	-	EXPRESSION TAG	UNP Q831P9
D	164	LEU	-	EXPRESSION TAG	UNP Q831P9
D	165	GLU	-	EXPRESSION TAG	UNP Q831P9
D	166	HIS	-	EXPRESSION TAG	UNP Q831P9
D	167	HIS	-	EXPRESSION TAG	UNP Q831P9
D	168	HIS	-	EXPRESSION TAG	UNP Q831P9
D	169	HIS	-	EXPRESSION TAG	UNP Q831P9
D	170	HIS	-	EXPRESSION TAG	UNP Q831P9
D	171	HIS	-	EXPRESSION TAG	UNP Q831P9
E	164	LEU	-	EXPRESSION TAG	UNP Q831P9
E	165	GLU	-	EXPRESSION TAG	UNP Q831P9
E	166	HIS	-	EXPRESSION TAG	UNP Q831P9
E	167	HIS	-	EXPRESSION TAG	UNP Q831P9
E	168	HIS	-	EXPRESSION TAG	UNP Q831P9
E	169	HIS	-	EXPRESSION TAG	UNP Q831P9
E	170	HIS	-	EXPRESSION TAG	UNP Q831P9
E	171	HIS	-	EXPRESSION TAG	UNP Q831P9
F	164	LEU	-	EXPRESSION TAG	UNP Q831P9
F	165	GLU	-	EXPRESSION TAG	UNP Q831P9
F	166	HIS	-	EXPRESSION TAG	UNP Q831P9
F	167	HIS	-	EXPRESSION TAG	UNP Q831P9
F	168	HIS	-	EXPRESSION TAG	UNP Q831P9
F	169	HIS	-	EXPRESSION TAG	UNP Q831P9
F	170	HIS	-	EXPRESSION TAG	UNP Q831P9
F	171	HIS	-	EXPRESSION TAG	UNP Q831P9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	102	Total O 102 102	0	0
2	B	82	Total O 82 82	0	0
2	C	74	Total O 74 74	0	0
2	D	82	Total O 82 82	0	0

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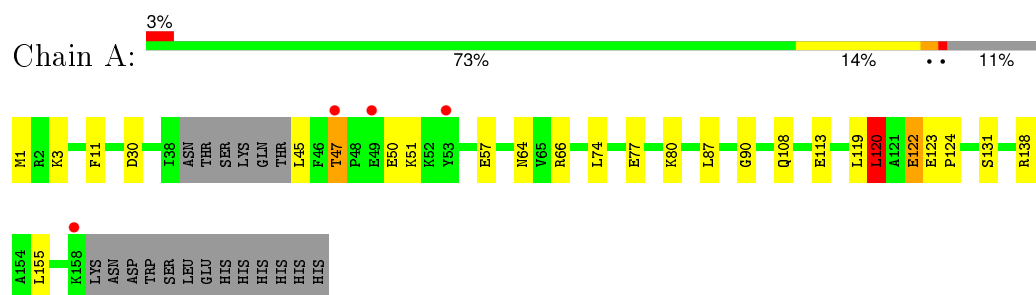
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	100	Total	O	0	0
			100	100		
2	F	94	Total	O	0	0
			94	94		

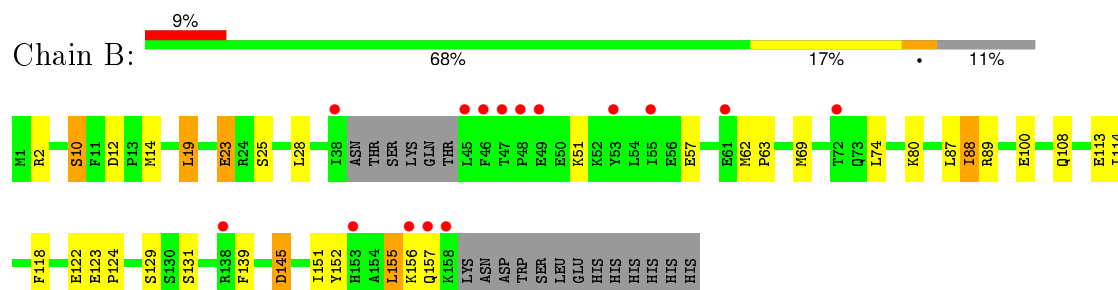
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

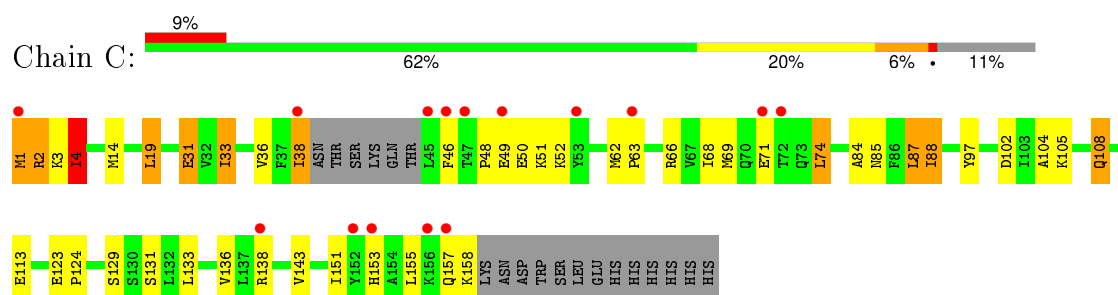
• Molecule 1: Phosphopantetheine adenylyltransferase



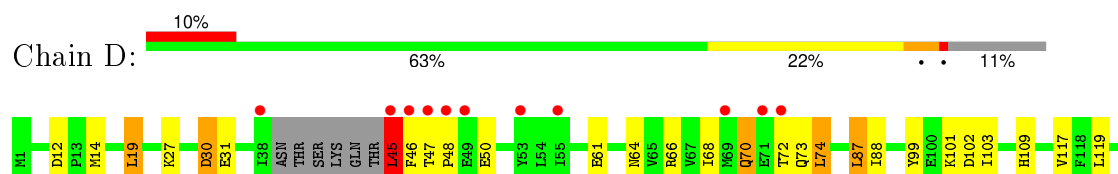
• Molecule 1: Phosphopantetheine adenylyltransferase

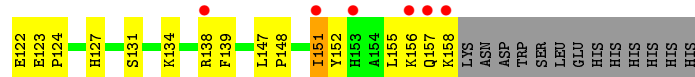


• Molecule 1: Phosphopantetheine adenylyltransferase

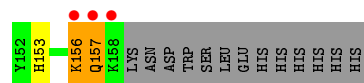


• Molecule 1: Phosphopantetheine adenylyltransferase

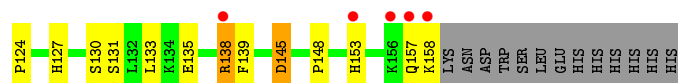




• Molecule 1: Phosphopantetheine adenylyltransferase



• Molecule 1: Phosphopantetheine adenylyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	110.20Å 125.68Å 125.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.89 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.3 (20.00-2.30) 98.1 (19.89-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.53 (at 2.30Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.196 , 0.249 0.202 , 0.252	Depositor DCC
R_{free} test set	3845 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.390	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.4	EDS
Estimated twinning fraction	0.026 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 76152 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7878	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	1.20	3/1248 (0.2%)	0.97	3/1683 (0.2%)
1	B	1.17	3/1248 (0.2%)	1.02	2/1683 (0.1%)
1	C	1.24	6/1248 (0.5%)	1.01	2/1683 (0.1%)
1	D	1.18	2/1248 (0.2%)	0.98	5/1683 (0.3%)
1	E	1.24	4/1248 (0.3%)	1.06	6/1683 (0.4%)
1	F	1.09	1/1248 (0.1%)	1.00	3/1683 (0.2%)
All	All	1.19	19/7488 (0.3%)	1.01	21/10098 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	77	GLU	CB-CG	-8.84	1.35	1.52
1	B	57	GLU	CG-CD	7.48	1.63	1.51
1	C	113	GLU	CB-CG	7.43	1.66	1.52
1	F	88	ILE	CB-CG2	7.10	1.74	1.52
1	C	113	GLU	CG-CD	6.30	1.61	1.51
1	C	31	GLU	CB-CG	6.06	1.63	1.52
1	E	113	GLU	CG-CD	5.80	1.60	1.51
1	A	113	GLU	CG-CD	5.68	1.60	1.51
1	B	113	GLU	CG-CD	5.62	1.60	1.51
1	A	113	GLU	CB-CG	5.54	1.62	1.52
1	E	122	GLU	CD-OE2	5.49	1.31	1.25
1	C	97	TYR	CE1-CZ	5.46	1.45	1.38
1	B	23	GLU	CG-CD	5.32	1.59	1.51
1	D	27	LYS	CD-CE	5.30	1.64	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	94	VAL	CB-CG1	5.20	1.63	1.52
1	A	122	GLU	CD-OE2	5.14	1.31	1.25
1	D	117	VAL	CA-CB	5.13	1.65	1.54
1	C	4	ILE	CB-CG2	5.05	1.68	1.52
1	C	19	LEU	CG-CD2	5.01	1.70	1.51

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2	ARG	NE-CZ-NH1	-10.28	115.16	120.30
1	B	2	ARG	NE-CZ-NH2	9.69	125.14	120.30
1	C	102	ASP	CB-CG-OD1	7.59	125.13	118.30
1	D	30	ASP	CB-CG-OD2	-6.61	112.35	118.30
1	E	102	ASP	CB-CG-OD2	6.19	123.87	118.30
1	F	145	ASP	CB-CG-OD1	-6.01	112.89	118.30
1	D	45	LEU	CA-CB-CG	5.92	128.91	115.30
1	E	89	ARG	CG-CD-NE	-5.78	99.66	111.80
1	E	2	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	E	102	ASP	CB-CG-OD1	-5.66	113.21	118.30
1	F	133	LEU	CB-CG-CD2	-5.64	101.41	111.00
1	A	1	MET	CG-SD-CE	5.50	109.00	100.20
1	A	147	LEU	CA-CB-CG	-5.39	102.90	115.30
1	F	19	LEU	CB-CG-CD2	5.35	120.09	111.00
1	A	120	LEU	CB-CG-CD1	5.22	119.88	111.00
1	C	87	LEU	CA-CB-CG	5.18	127.21	115.30
1	D	102	ASP	CB-CG-OD2	-5.13	113.68	118.30
1	D	87	LEU	CB-CG-CD1	5.09	119.65	111.00
1	E	2	ARG	CD-NE-CZ	5.08	130.72	123.60
1	E	120	LEU	CB-CG-CD1	5.05	119.58	111.00
1	D	87	LEU	CA-CB-CG	5.02	126.84	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1	MET	Peptide

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	1224	0	1245	21	0
1	B	1224	0	1246	19	0
1	C	1224	0	1246	39	0
1	D	1224	0	1246	27	0
1	E	1224	0	1246	21	0
1	F	1224	0	1246	53	0
2	A	102	0	0	7	0
2	B	82	0	0	2	0
2	C	74	0	0	3	0
2	D	82	0	0	5	0
2	E	100	0	0	7	0
2	F	94	0	0	10	0
All	All	7878	0	7475	166	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 11.

All (166) 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:F:88:ILE:CG2	1:F:88:ILE:CB	1.74	1.61
1:F:68:ILE:CD1	1:F:70:GLN:HG2	1.99	0.91
1:A:66:ARG:CZ	2:A:244:HOH:O	2.18	0.90
1:E:105:LYS:HE3	1:F:123:GLU:OE2	1.72	0.90
1:A:77:GLU:HG2	2:A:298:HOH:O	1.76	0.84
1:F:3:LYS:HG3	2:F:306:HOH:O	1.79	0.83
1:A:123:GLU:HG3	1:F:123:GLU:HG2	1.63	0.79
1:C:38:ILE:HD11	1:C:69:MET:CE	2.13	0.78
1:D:123:GLU:HG3	1:D:124:PRO:HD3	1.66	0.78
1:F:2:ARG:HH11	1:F:2:ARG:CG	1.96	0.77
1:F:68:ILE:HD12	1:F:68:ILE:O	1.86	0.76
1:F:130:SER:HB3	2:F:238:HOH:O	1.86	0.76
1:C:123:GLU:HA	2:C:229:HOH:O	1.84	0.76
1:D:147:LEU:HD12	1:D:152:TYR:HD1	1.52	0.75
1:F:47:THR:HG21	1:F:49:GLU:OE1	1.89	0.72
1:F:12:ASP:HB3	2:F:266:HOH:O	1.88	0.72
1:F:2:ARG:HG3	1:F:2:ARG:HH11	1.53	0.72
1:E:57:GLU:HG2	2:E:223:HOH:O	1.92	0.69
1:A:123:GLU:CG	1:F:123:GLU:HG2	2.23	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:68:ILE:O	1:D:68:ILE:HD12	1.92	0.68
1:F:123:GLU:N	1:F:124:PRO:HD2	2.08	0.68
1:F:68:ILE:HD13	1:F:70:GLN:HG2	1.73	0.68
1:C:153:HIS:HB3	2:C:349:HOH:O	1.93	0.68
1:F:74:LEU:HB2	1:F:77:GLU:HG3	1.75	0.67
1:D:147:LEU:HD12	1:D:152:TYR:CD1	2.29	0.67
1:E:89:ARG:HD3	2:E:227:HOH:O	1.93	0.67
1:F:52:LYS:HD2	1:F:69:MET:HG2	1.76	0.67
1:B:19:LEU:O	1:B:23:GLU:HG3	1.94	0.66
1:D:109:HIS:ND1	2:D:569:HOH:O	2.29	0.66
1:F:153:HIS:HB3	2:F:438:HOH:O	1.96	0.65
1:C:4:ILE:HD11	1:C:33:ILE:HD11	1.76	0.65
1:B:89:ARG:HD3	1:B:100:GLU:OE1	1.97	0.65
1:C:123:GLU:N	1:C:124:PRO:HD2	2.12	0.64
1:F:88:ILE:CG2	1:F:88:ILE:HB	2.15	0.64
1:C:52:LYS:HB2	1:C:69:MET:SD	2.38	0.64
1:B:123:GLU:OE1	2:B:296:HOH:O	2.15	0.63
1:E:148:PRO:HB2	1:E:151:ILE:HG13	1.79	0.63
1:C:1:MET:HB3	1:C:85:ASN:HD21	1.64	0.63
1:D:139:PHE:HB2	1:F:74:LEU:HD13	1.79	0.63
1:C:4:ILE:HG13	1:C:84:ALA:HA	1.81	0.63
1:A:123:GLU:N	1:A:124:PRO:HD2	2.13	0.62
1:C:31:GLU:OE2	1:C:66:ARG:HD3	1.99	0.62
1:D:12:ASP:OD1	1:D:45:LEU:HD11	2.00	0.62
1:C:31:GLU:OE2	1:C:66:ARG:NH1	2.28	0.62
1:A:66:ARG:NH1	2:A:244:HOH:O	2.28	0.61
1:F:1:MET:O	1:F:3:LYS:HD3	2.00	0.61
1:D:12:ASP:O	1:D:134:LYS:HE2	2.00	0.61
1:A:120:LEU:HD22	1:B:118:PHE:CE2	2.35	0.60
1:C:38:ILE:HD11	1:C:69:MET:HE3	1.83	0.59
1:F:47:THR:CG2	1:F:49:GLU:OE1	2.51	0.59
1:C:38:ILE:HD11	1:C:69:MET:HE2	1.85	0.59
1:E:2:ARG:NH1	2:E:245:HOH:O	2.35	0.58
1:F:88:ILE:CG2	1:F:88:ILE:CG1	2.77	0.57
1:A:66:ARG:NE	2:A:244:HOH:O	2.30	0.57
1:B:139:PHE:HB2	1:D:74:LEU:HD13	1.87	0.56
1:D:73:GLN:HG3	1:D:74:LEU:N	2.20	0.55
1:C:74:LEU:HD13	1:E:139:PHE:HB2	1.87	0.55
1:F:12:ASP:CB	2:F:266:HOH:O	2.48	0.55
1:B:145:ASP:OD1	1:B:145:ASP:N	2.37	0.55
1:C:4:ILE:HD11	1:C:33:ILE:CD1	2.36	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:3:LYS:HG3	2:F:306:HOH:O	2.06	0.54
1:A:57:GLU:HG3	1:A:57:GLU:O	2.08	0.54
1:C:4:ILE:CD1	1:C:33:ILE:HD11	2.39	0.53
1:F:38:ILE:O	1:F:72:THR:HG23	2.08	0.53
1:D:73:GLN:CG	1:D:74:LEU:N	2.72	0.53
1:F:88:ILE:CG2	1:F:88:ILE:CA	2.82	0.53
1:E:123:GLU:N	1:E:124:PRO:HD2	2.24	0.53
1:D:122:GLU:HG3	2:D:371:HOH:O	2.09	0.52
1:B:123:GLU:N	1:B:124:PRO:HD2	2.24	0.51
1:E:149:PRO:HD2	2:E:394:HOH:O	2.10	0.51
1:C:138:ARG:O	1:C:138:ARG:HG2	2.09	0.51
1:C:46:PHE:HB3	1:C:50:GLU:HB3	1.93	0.50
1:C:14:MET:HB3	1:C:151:ILE:HD13	1.94	0.50
1:A:57:GLU:HB2	2:A:316:HOH:O	2.10	0.50
1:A:138:ARG:HB3	1:A:138:ARG:CZ	2.42	0.50
1:C:136:VAL:HG11	1:C:143:VAL:HG13	1.94	0.50
1:D:127:HIS:HB3	2:D:291:HOH:O	2.11	0.50
1:C:123:GLU:N	1:C:124:PRO:CD	2.75	0.49
1:F:74:LEU:HB2	1:F:77:GLU:CG	2.41	0.49
1:E:89:ARG:CD	2:E:227:HOH:O	2.58	0.49
1:A:3:LYS:HE3	1:B:28:LEU:O	2.12	0.49
1:F:105:LYS:NZ	2:F:277:HOH:O	2.41	0.49
1:E:113:GLU:HB2	2:E:201:HOH:O	2.13	0.48
1:A:66:ARG:HD3	2:A:807:HOH:O	2.12	0.48
1:B:152:TYR:CE2	1:B:156:LYS:HD2	2.49	0.48
1:C:66:ARG:HE	1:C:68:ILE:HD11	1.79	0.48
1:F:14:MET:HE3	1:F:148:PRO:HG2	1.96	0.48
1:E:153:HIS:CE1	1:E:157:GLN:OE1	2.67	0.48
1:C:2:ARG:HG3	1:C:3:LYS:N	2.29	0.47
1:C:71:GLU:H	1:C:71:GLU:CD	2.18	0.47
1:D:151:ILE:O	1:D:155:LEU:HB2	2.13	0.47
1:F:68:ILE:HD13	1:F:70:GLN:CG	2.43	0.47
1:F:90:GLY:HA2	1:F:119:LEU:O	2.14	0.47
1:F:68:ILE:HD12	1:F:68:ILE:C	2.34	0.47
1:F:49:GLU:HG2	1:F:50:GLU:N	2.29	0.47
1:E:2:ARG:HB2	2:E:321:HOH:O	2.15	0.47
1:F:89:ARG:HB2	1:F:118:PHE:CD1	2.50	0.47
1:D:30:ASP:O	1:D:64:ASN:HB2	2.15	0.47
1:F:127:HIS:HB3	2:F:264:HOH:O	2.15	0.47
1:E:2:ARG:HG2	1:E:2:ARG:HH11	1.80	0.46
1:F:52:LYS:HB2	1:F:69:MET:SD	2.56	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:SER:O	1:C:133:LEU:HB2	2.16	0.46
1:D:139:PHE:CB	1:F:74:LEU:HD13	2.45	0.46
1:C:31:GLU:CD	1:C:66:ARG:HH11	2.16	0.46
1:A:30:ASP:O	1:A:64:ASN:HB2	2.16	0.45
1:B:157:GLN:HG3	1:B:157:GLN:O	2.16	0.45
1:C:1:MET:HB3	1:C:85:ASN:ND2	2.31	0.45
1:B:14:MET:HE2	1:B:14:MET:HB3	1.73	0.45
1:F:71:GLU:H	1:F:71:GLU:CD	2.20	0.45
1:F:2:ARG:HG3	1:F:2:ARG:NH1	2.26	0.45
1:B:69:MET:CA	1:B:69:MET:HE2	2.46	0.45
1:E:156:LYS:HE3	1:E:156:LYS:HB2	1.61	0.45
1:C:123:GLU:HB3	1:C:124:PRO:HD3	1.98	0.45
1:A:90:GLY:HA2	1:A:119:LEU:O	2.16	0.45
1:E:24:ARG:NE	1:F:115:GLU:OE1	2.33	0.45
1:F:74:LEU:HB3	1:F:77:GLU:HG2	1.99	0.45
1:F:12:ASP:HA	1:F:13:PRO:HA	1.78	0.45
1:F:6:LEU:HB3	1:F:87:LEU:HD12	1.99	0.44
1:F:10:SER:O	1:F:11:PHE:C	2.55	0.44
1:E:135:GLU:O	1:E:138:ARG:HG2	2.18	0.44
1:C:31:GLU:CD	1:C:66:ARG:NH1	2.71	0.44
1:F:108:GLN:HG3	2:F:208:HOH:O	2.18	0.44
1:B:122:GLU:HG3	2:B:287:HOH:O	2.18	0.43
1:B:25:SER:OG	1:B:88:ILE:HD12	2.19	0.43
1:A:123:GLU:CG	1:F:123:GLU:CG	2.94	0.43
1:D:68:ILE:HD13	1:D:70:GLN:NE2	2.33	0.43
1:C:104:ALA:O	1:C:108:GLN:HB3	2.17	0.43
1:E:123:GLU:N	1:E:124:PRO:CD	2.81	0.43
1:C:36:VAL:CG1	1:C:51:LYS:HE3	2.48	0.43
1:C:62:MET:HA	1:C:63:PRO:HD2	1.92	0.42
1:D:19:LEU:HD12	1:D:148:PRO:HG3	2.01	0.42
1:C:46:PHE:HB3	1:C:50:GLU:CB	2.48	0.42
1:A:66:ARG:CD	2:A:244:HOH:O	2.65	0.42
1:C:38:ILE:HD12	1:C:48:PRO:HB3	2.02	0.42
1:F:77:GLU:HG2	1:F:77:GLU:H	1.59	0.42
1:D:46:PHE:HB3	1:D:50:GLU:HB2	2.01	0.42
1:B:10:SER:HB3	1:B:12:ASP:OD2	2.18	0.42
1:F:61:GLU:HB2	2:F:271:HOH:O	2.20	0.42
1:D:99:TYR:CZ	1:D:103:ILE:HD11	2.55	0.42
1:B:151:ILE:O	1:B:155:LEU:HB2	2.20	0.42
1:A:123:GLU:HG2	1:F:123:GLU:CG	2.50	0.42
1:D:47:THR:O	1:D:48:PRO:C	2.57	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:MET:HA	1:B:63:PRO:HD3	1.83	0.42
1:A:139:PHE:HB2	1:E:74:LEU:HD13	2.02	0.42
1:D:73:GLN:CD	1:D:74:LEU:H	2.24	0.41
1:C:105:LYS:O	2:C:172:HOH:O	2.22	0.41
1:F:45:LEU:HG	1:F:158:LYS:HD2	2.02	0.41
1:C:88:ILE:HD13	1:C:88:ILE:C	2.41	0.41
1:D:147:LEU:CD1	1:D:152:TYR:HD1	2.29	0.41
1:F:14:MET:CE	1:F:148:PRO:HG2	2.51	0.41
1:D:101:LYS:NZ	2:D:788:HOH:O	2.42	0.41
1:D:88:ILE:HD11	1:D:119:LEU:HG	2.02	0.41
1:F:130:SER:O	1:F:131:SER:C	2.57	0.41
1:C:4:ILE:CD1	1:C:33:ILE:HG13	2.51	0.41
1:C:74:LEU:HD13	1:E:139:PHE:CB	2.51	0.41
1:C:136:VAL:HG11	1:C:143:VAL:CG1	2.51	0.41
1:B:69:MET:HE3	1:B:69:MET:HB2	1.87	0.41
1:D:72:THR:HG23	2:D:294:HOH:O	2.20	0.41
1:E:122:GLU:C	1:E:124:PRO:HD2	2.42	0.41
1:D:14:MET:CE	1:D:148:PRO:HG2	2.51	0.41
1:A:47:THR:HB	1:A:50:GLU:H	1.86	0.41
1:C:38:ILE:HD12	1:C:48:PRO:HG3	2.03	0.40
1:F:38:ILE:HG12	1:F:38:ILE:H	1.61	0.40
1:A:11:PHE:O	1:A:51:LYS:HE2	2.21	0.40
1:F:138:ARG:HG3	1:F:139:PHE:N	2.37	0.40
1:B:80:LYS:HG3	1:B:114:ILE:HD11	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	148/171 (86%)	144 (97%)	4 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	148/171 (86%)	146 (99%)	2 (1%)	0	100	100
1	C	148/171 (86%)	144 (97%)	4 (3%)	0	100	100
1	D	148/171 (86%)	146 (99%)	2 (1%)	0	100	100
1	E	148/171 (86%)	144 (97%)	4 (3%)	0	100	100
1	F	148/171 (86%)	145 (98%)	3 (2%)	0	100	100
All	All	888/1026 (86%)	869 (98%)	19 (2%)	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	135/154 (88%)	124 (92%)	11 (8%)	15	18
1	B	135/154 (88%)	124 (92%)	11 (8%)	15	18
1	C	135/154 (88%)	121 (90%)	14 (10%)	9	10
1	D	135/154 (88%)	121 (90%)	14 (10%)	9	10
1	E	135/154 (88%)	120 (89%)	15 (11%)	8	8
1	F	135/154 (88%)	118 (87%)	17 (13%)	5	5
All	All	810/924 (88%)	728 (90%)	82 (10%)	9	11

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LEU
1	A	47	THR
1	A	74	LEU
1	A	80	LYS
1	A	87	LEU
1	A	108	GLN
1	A	120	LEU
1	A	122	GLU

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Mol	Chain	Res	Type
1	A	131	SER
1	A	147	LEU
1	A	155	LEU
1	B	10	SER
1	B	19	LEU
1	B	51	LYS
1	B	74	LEU
1	B	87	LEU
1	B	88	ILE
1	B	108	GLN
1	B	129	SER
1	B	131	SER
1	B	145	ASP
1	B	155	LEU
1	C	2	ARG
1	C	4	ILE
1	C	19	LEU
1	C	33	ILE
1	C	38	ILE
1	C	49	GLU
1	C	74	LEU
1	C	87	LEU
1	C	88	ILE
1	C	108	GLN
1	C	131	SER
1	C	155	LEU
1	C	157	GLN
1	C	158	LYS
1	D	19	LEU
1	D	31	GLU
1	D	45	LEU
1	D	61	GLU
1	D	66	ARG
1	D	70	GLN
1	D	74	LEU
1	D	87	LEU
1	D	131	SER
1	D	138	ARG
1	D	151	ILE
1	D	156	LYS
1	D	157	GLN
1	D	158	LYS

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Mol	Chain	Res	Type
1	E	2	ARG
1	E	19	LEU
1	E	45	LEU
1	E	61	GLU
1	E	74	LEU
1	E	77	GLU
1	E	80	LYS
1	E	87	LEU
1	E	105	LYS
1	E	120	LEU
1	E	122	GLU
1	E	131	SER
1	E	151	ILE
1	E	156	LYS
1	E	157	GLN
1	F	2	ARG
1	F	4	ILE
1	F	19	LEU
1	F	38	ILE
1	F	46	PHE
1	F	60	LYS
1	F	74	LEU
1	F	77	GLU
1	F	87	LEU
1	F	88	ILE
1	F	105	LYS
1	F	122	GLU
1	F	123	GLU
1	F	135	GLU
1	F	138	ARG
1	F	145	ASP
1	F	157	GLN

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

Mol	Chain	Res	Type
1	A	153	HIS
1	D	157	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	152/171 (88%)	-0.19	5 (3%)	50	59	18, 34, 67, 81	0
1	B	152/171 (88%)	-0.06	15 (9%)	9	14	18, 37, 75, 90	0
1	C	152/171 (88%)	0.16	15 (9%)	9	14	18, 42, 83, 98	0
1	D	152/171 (88%)	0.10	17 (11%)	7	10	18, 42, 90, 101	0
1	E	152/171 (88%)	-0.06	7 (4%)	36	45	18, 34, 68, 84	0
1	F	152/171 (88%)	0.10	12 (7%)	15	22	20, 41, 84, 95	0
All	All	912/1026 (88%)	0.01	71 (7%)	16	22	18, 39, 79, 101	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	45	LEU	6.5
1	C	1	MET	4.5
1	D	47	THR	4.0
1	F	49	GLU	4.0
1	F	72	THR	3.9
1	C	47	THR	3.8
1	C	53	TYR	3.7
1	B	157	GLN	3.5
1	C	157	GLN	3.4
1	D	38	ILE	3.4
1	D	48	PRO	3.4
1	C	72	THR	3.4
1	D	45	LEU	3.4
1	E	156	LYS	3.4
1	D	72	THR	3.4
1	B	45	LEU	3.3
1	F	156	LYS	3.3
1	D	138	ARG	3.2
1	D	153	HIS	3.2

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Mol	Chain	Res	Type	RSRZ
1	F	71	GLU	3.2
1	D	157	GLN	3.1
1	F	157	GLN	3.1
1	D	156	LYS	3.0
1	C	156	LYS	3.0
1	D	158	LYS	3.0
1	B	48	PRO	3.0
1	D	53	TYR	2.9
1	D	46	PHE	2.9
1	B	49	GLU	2.8
1	B	153	HIS	2.8
1	F	138	ARG	2.8
1	C	138	ARG	2.8
1	C	38	ILE	2.7
1	F	158	LYS	2.7
1	B	47	THR	2.7
1	D	71	GLU	2.7
1	F	153	HIS	2.7
1	B	158	LYS	2.7
1	A	153	HIS	2.7
1	A	47	THR	2.6
1	B	53	TYR	2.6
1	F	61	GLU	2.6
1	B	55	ILE	2.6
1	E	157	GLN	2.6
1	D	151	ILE	2.6
1	C	49	GLU	2.5
1	C	46	PHE	2.5
1	B	138	ARG	2.5
1	F	38	ILE	2.4
1	B	46	PHE	2.4
1	E	158	LYS	2.4
1	C	71	GLU	2.4
1	F	45	LEU	2.4
1	C	63	PRO	2.4
1	F	60	LYS	2.4
1	E	53	TYR	2.3
1	B	72	THR	2.3
1	B	38	ILE	2.3
1	A	158	LYS	2.2
1	A	49	GLU	2.2
1	D	49	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	156	LYS	2.2
1	C	152	TYR	2.2
1	E	138	ARG	2.2
1	C	153	HIS	2.1
1	E	47	THR	2.1
1	B	61	GLU	2.1
1	A	53	TYR	2.1
1	E	45	LEU	2.1
1	D	55	ILE	2.1
1	D	69	MET	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.