



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:13 PM GMT

PDB ID : 4KWB
Title : Structure of signal peptide peptidase A with C-termini bound in the active sites: insights into specificity, self-processing and regulation
Authors : Nam, S.E.; Paetzel, M.
Deposited on : 2013-05-23
Resolution : 2.39 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

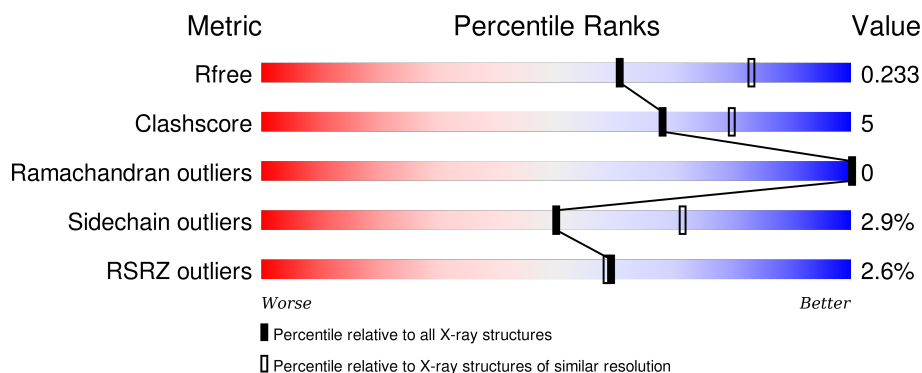
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	273	<div> <div></div> <div>79% 7% 13%</div> </div>
1	B	273	<div> <div>3%</div> <div>78% 7% 15%</div> </div>
1	C	273	<div> <div></div> <div>77% 8% 13%</div> </div>
1	D	273	<div> <div>3%</div> <div>72% 10% 16%</div> </div>
1	E	273	<div> <div>5%</div> <div>74% 11% 15%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	273	<div><div>2%</div><div><div></div><div>75%</div><div>11%</div><div>14%</div></div></div>
1	G	273	<div><div>5%</div><div><div></div><div>77%</div><div>7%</div><div>15%</div></div></div>
1	H	273	<div><div>%</div><div><div></div><div>74%</div><div>11%</div><div>• 15%</div></div></div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 14324 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 Signal peptide peptidase SppA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1797	1133	297	356	11			
1	B	233	Total	C	N	O	S	0	0	0
			1754	1103	292	348	11			
1	C	237	Total	C	N	O	S	0	0	0
			1808	1138	300	359	11			
1	D	228	Total	C	N	O	S	0	0	0
			1686	1063	280	332	11			
1	E	232	Total	C	N	O	S	0	0	0
			1747	1097	293	346	11			
1	F	234	Total	C	N	O	S	0	0	0
			1779	1122	294	352	11			
1	G	232	Total	C	N	O	S	0	0	0
			1756	1102	294	349	11			
1	H	232	Total	C	N	O	S	0	0	0
			1776	1119	297	349	11			

There are 104 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
A	?	-	LEU	SEE REMARK 999	UNP O34525
A	?	-	GLY	SEE REMARK 999	UNP O34525
A	?	-	SER	SEE REMARK 999	UNP O34525
A	?	-	LEU	SEE REMARK 999	UNP O34525
A	?	-	PHE	SEE REMARK 999	UNP O34525
A	?	-	SER	SEE REMARK 999	UNP O34525
A	?	-	MET	SEE REMARK 999	UNP O34525
A	?	-	GLY	SEE REMARK 999	UNP O34525
A	?	-	ALA	SEE REMARK 999	UNP O34525
A	?	-	ASN	SEE REMARK 999	UNP O34525
A	?	-	LYS	SEE REMARK 999	UNP O34525
A	?	-	MET	SEE REMARK 999	UNP O34525

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Chain	Residue	Modelled	Actual	Comment	Reference
B	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
B	?	-	LEU	SEE REMARK 999	UNP O34525
B	?	-	GLY	SEE REMARK 999	UNP O34525
B	?	-	SER	SEE REMARK 999	UNP O34525
B	?	-	LEU	SEE REMARK 999	UNP O34525
B	?	-	PHE	SEE REMARK 999	UNP O34525
B	?	-	SER	SEE REMARK 999	UNP O34525
B	?	-	MET	SEE REMARK 999	UNP O34525
B	?	-	GLY	SEE REMARK 999	UNP O34525
B	?	-	ALA	SEE REMARK 999	UNP O34525
B	?	-	ASN	SEE REMARK 999	UNP O34525
B	?	-	LYS	SEE REMARK 999	UNP O34525
B	?	-	MET	SEE REMARK 999	UNP O34525
C	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
C	?	-	LEU	SEE REMARK 999	UNP O34525
C	?	-	GLY	SEE REMARK 999	UNP O34525
C	?	-	SER	SEE REMARK 999	UNP O34525
C	?	-	LEU	SEE REMARK 999	UNP O34525
C	?	-	PHE	SEE REMARK 999	UNP O34525
C	?	-	SER	SEE REMARK 999	UNP O34525
C	?	-	MET	SEE REMARK 999	UNP O34525
C	?	-	GLY	SEE REMARK 999	UNP O34525
C	?	-	ALA	SEE REMARK 999	UNP O34525
C	?	-	ASN	SEE REMARK 999	UNP O34525
C	?	-	LYS	SEE REMARK 999	UNP O34525
C	?	-	MET	SEE REMARK 999	UNP O34525
D	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
D	?	-	LEU	SEE REMARK 999	UNP O34525
D	?	-	GLY	SEE REMARK 999	UNP O34525
D	?	-	SER	SEE REMARK 999	UNP O34525
D	?	-	LEU	SEE REMARK 999	UNP O34525
D	?	-	PHE	SEE REMARK 999	UNP O34525
D	?	-	SER	SEE REMARK 999	UNP O34525
D	?	-	MET	SEE REMARK 999	UNP O34525
D	?	-	GLY	SEE REMARK 999	UNP O34525
D	?	-	ALA	SEE REMARK 999	UNP O34525
D	?	-	ASN	SEE REMARK 999	UNP O34525
D	?	-	LYS	SEE REMARK 999	UNP O34525
D	?	-	MET	SEE REMARK 999	UNP O34525
E	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
E	?	-	LEU	SEE REMARK 999	UNP O34525
E	?	-	GLY	SEE REMARK 999	UNP O34525

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Chain	Residue	Modelled	Actual	Comment	Reference
E	?	-	SER	SEE REMARK 999	UNP O34525
E	?	-	LEU	SEE REMARK 999	UNP O34525
E	?	-	PHE	SEE REMARK 999	UNP O34525
E	?	-	SER	SEE REMARK 999	UNP O34525
E	?	-	MET	SEE REMARK 999	UNP O34525
E	?	-	GLY	SEE REMARK 999	UNP O34525
E	?	-	ALA	SEE REMARK 999	UNP O34525
E	?	-	ASN	SEE REMARK 999	UNP O34525
E	?	-	LYS	SEE REMARK 999	UNP O34525
E	?	-	MET	SEE REMARK 999	UNP O34525
F	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
F	?	-	LEU	SEE REMARK 999	UNP O34525
F	?	-	GLY	SEE REMARK 999	UNP O34525
F	?	-	SER	SEE REMARK 999	UNP O34525
F	?	-	LEU	SEE REMARK 999	UNP O34525
F	?	-	PHE	SEE REMARK 999	UNP O34525
F	?	-	SER	SEE REMARK 999	UNP O34525
F	?	-	MET	SEE REMARK 999	UNP O34525
F	?	-	GLY	SEE REMARK 999	UNP O34525
F	?	-	ALA	SEE REMARK 999	UNP O34525
F	?	-	ASN	SEE REMARK 999	UNP O34525
F	?	-	LYS	SEE REMARK 999	UNP O34525
F	?	-	MET	SEE REMARK 999	UNP O34525
G	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
G	?	-	LEU	SEE REMARK 999	UNP O34525
G	?	-	GLY	SEE REMARK 999	UNP O34525
G	?	-	SER	SEE REMARK 999	UNP O34525
G	?	-	LEU	SEE REMARK 999	UNP O34525
G	?	-	PHE	SEE REMARK 999	UNP O34525
G	?	-	SER	SEE REMARK 999	UNP O34525
G	?	-	MET	SEE REMARK 999	UNP O34525
G	?	-	GLY	SEE REMARK 999	UNP O34525
G	?	-	ALA	SEE REMARK 999	UNP O34525
G	?	-	ASN	SEE REMARK 999	UNP O34525
G	?	-	LYS	SEE REMARK 999	UNP O34525
G	?	-	MET	SEE REMARK 999	UNP O34525
H	199	ALA	LYS	ENGINEERED MUTATION	UNP O34525
H	?	-	LEU	SEE REMARK 999	UNP O34525
H	?	-	GLY	SEE REMARK 999	UNP O34525
H	?	-	SER	SEE REMARK 999	UNP O34525
H	?	-	LEU	SEE REMARK 999	UNP O34525
H	?	-	PHE	SEE REMARK 999	UNP O34525

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Chain	Residue	Modelled	Actual	Comment	Reference
H	?	-	SER	SEE REMARK 999	UNP O34525
H	?	-	MET	SEE REMARK 999	UNP O34525
H	?	-	GLY	SEE REMARK 999	UNP O34525
H	?	-	ALA	SEE REMARK 999	UNP O34525
H	?	-	ASN	SEE REMARK 999	UNP O34525
H	?	-	LYS	SEE REMARK 999	UNP O34525
H	?	-	MET	SEE REMARK 999	UNP O34525

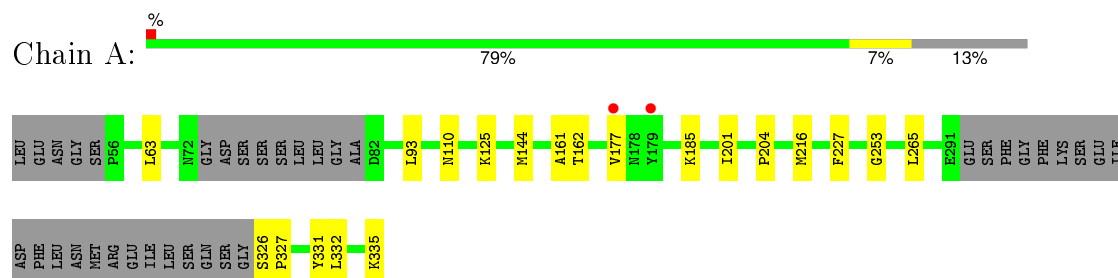
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	36	Total O 36 36	0	0
2	B	22	Total O 22 22	0	0
2	C	31	Total O 31 31	0	0
2	D	19	Total O 19 19	0	0
2	E	21	Total O 21 21	0	0
2	F	35	Total O 35 35	0	0
2	G	24	Total O 24 24	0	0
2	H	33	Total O 33 33	0	0

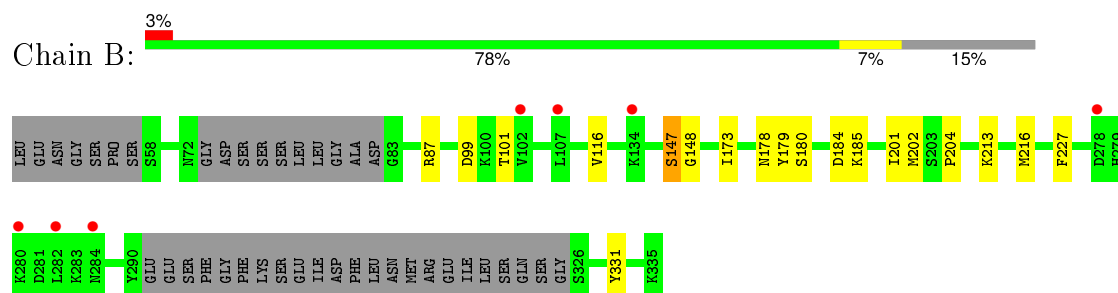
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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.

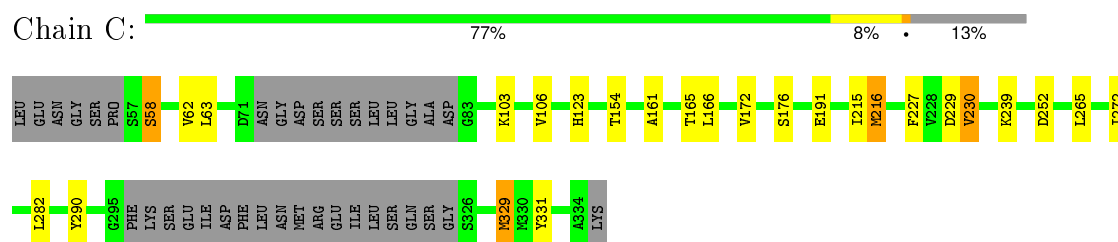
- Molecule 1: Signal peptide peptidase SppA



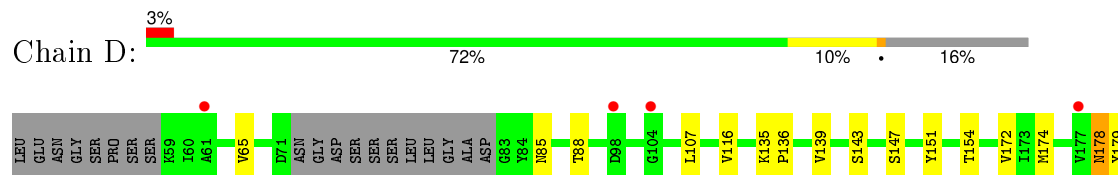
- Molecule 1: Signal peptide peptidase SppA

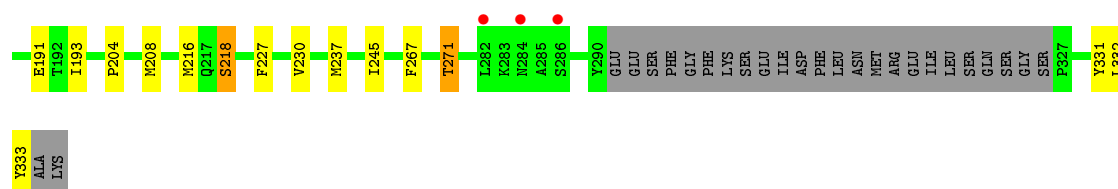


- Molecule 1: Signal peptide peptidase SppA

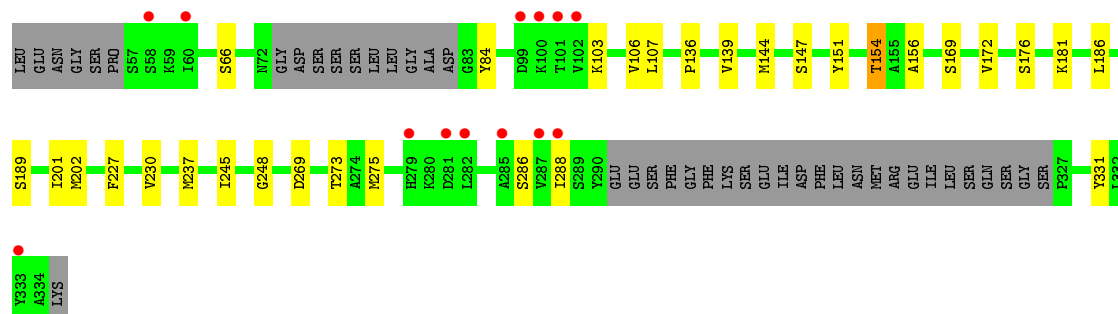


- Molecule 1: Signal peptide peptidase SppA

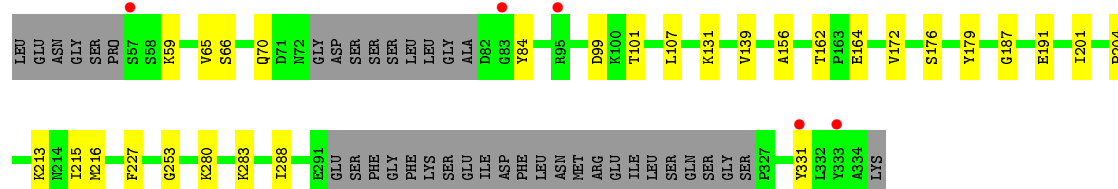




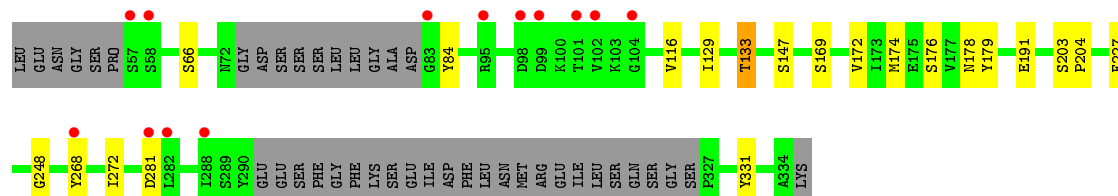
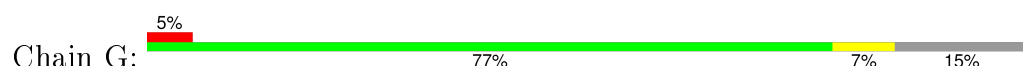
- Molecule 1: Signal peptide peptidase SppA



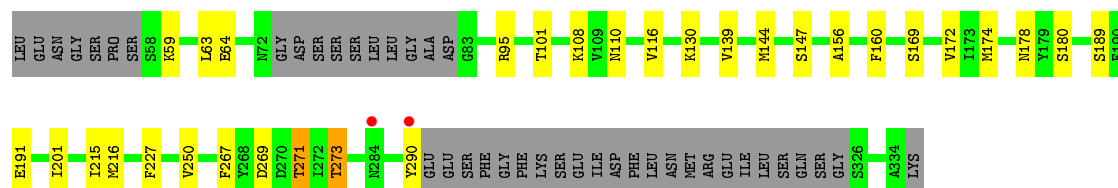
- Molecule 1: Signal peptide peptidase SppA



- Molecule 1: Signal peptide peptidase SppA



- Molecule 1: Signal peptide peptidase SppA



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.80Å 130.98Å 207.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.39 44.60 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.7 (50.00-2.39) 98.7 (44.60-2.39)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.13 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.206 , 0.240 0.202 , 0.233	Depositor DCC
R_{free} test set	4716 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 93821 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14324	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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 [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	0.45	0/1821	0.57	0/2441
1	B	0.47	0/1778	0.52	0/2390
1	C	0.46	0/1833	0.56	0/2458
1	D	0.43	0/1708	0.54	0/2298
1	E	0.45	0/1769	0.54	0/2374
1	F	0.45	0/1803	0.55	0/2417
1	G	0.42	0/1778	0.54	0/2386
1	H	0.44	0/1800	0.56	0/2413
All	All	0.45	0/14290	0.55	0/19177

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	1797	0	1786	13	0
1	B	1754	0	1721	15	0
1	C	1808	0	1799	31	0
1	D	1686	0	1629	26	0
1	E	1747	0	1727	26	0
1	F	1779	0	1781	21	0
1	G	1756	0	1742	14	0
1	H	1776	0	1784	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	36	0	0	0	0
2	B	22	0	0	0	0
2	C	31	0	0	0	0
2	D	19	0	0	1	0
2	E	21	0	0	0	0
2	F	35	0	0	0	0
2	G	24	0	0	0	0
2	H	33	0	0	1	0
All	All	14324	0	13969	132	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 (132) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:SER:HB2	1:A:327:PRO:HD3	1.52	0.89
1:D:218:SER:HB3	2:D:409:HOH:O	1.71	0.88
1:E:176:SER:OG	1:F:191:GLU:HB3	1.87	0.74
1:B:213:LYS:HE2	1:C:191:GLU:OE1	1.87	0.73
1:D:333:TYR:CD2	1:E:144:MET:HE3	2.25	0.71
1:C:216:MET:HE2	1:C:216:MET:HA	1.72	0.71
1:F:59:LYS:HD3	1:F:288:ILE:HG12	1.73	0.71
1:F:176:SER:OG	1:G:191:GLU:HB3	1.91	0.71
1:C:62:VAL:HG22	1:C:106:VAL:CG1	2.22	0.70
1:F:201:ILE:HD13	1:F:216:MET:HE3	1.73	0.69
1:G:176:SER:OG	1:H:191:GLU:HB3	1.93	0.68
1:C:215:ILE:HG22	1:C:216:MET:CE	2.26	0.66
1:D:333:TYR:CD2	1:E:144:MET:CE	2.79	0.66
1:C:106:VAL:HG11	1:C:272:ILE:HD11	1.79	0.65
1:D:151:TYR:O	1:D:154:THR:HG22	1.96	0.65
1:F:99:ASP:OD1	1:F:101:THR:HG22	1.97	0.64
1:B:179:TYR:CE2	1:B:204:PRO:HB2	2.33	0.64
1:D:333:TYR:HD2	1:E:144:MET:CE	2.13	0.62
1:C:216:MET:HB3	1:D:193:ILE:HD13	1.82	0.61
1:C:154:THR:OG1	1:C:230:VAL:HG13	2.01	0.61
1:D:331:TYR:CE1	1:E:172:VAL:HG22	2.35	0.61
1:F:65:VAL:HG23	1:F:107:LEU:HD11	1.82	0.61
1:C:329:MET:CE	1:C:329:MET:H	2.13	0.61
1:C:62:VAL:HG22	1:C:106:VAL:HG13	1.82	0.60
1:A:201:ILE:HG12	1:A:216:MET:HE2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:333:TYR:HD2	1:E:144:MET:HE1	1.66	0.60
1:F:215:ILE:HG22	1:F:216:MET:HE1	1.84	0.60
1:B:99:ASP:OD1	1:B:101:THR:HG22	2.02	0.60
1:H:201:ILE:HD13	1:H:216:MET:HE3	1.83	0.60
1:E:331:TYR:CE1	1:F:172:VAL:HG22	2.37	0.59
1:D:65:VAL:HG23	1:D:107:LEU:HD11	1.84	0.59
1:G:129:ILE:O	1:G:133:THR:HG22	2.03	0.59
1:H:215:ILE:HG22	1:H:216:MET:CE	2.32	0.59
1:C:216:MET:CE	1:C:216:MET:CA	2.80	0.59
1:D:331:TYR:HE1	1:E:172:VAL:HG22	1.67	0.58
1:A:201:ILE:HD13	1:A:216:MET:HE1	1.85	0.58
1:E:237:MET:HE3	1:E:245:ILE:HD12	1.86	0.58
1:C:216:MET:HE2	1:C:216:MET:CA	2.34	0.57
1:B:213:LYS:CE	1:C:191:GLU:OE1	2.52	0.57
1:D:107:LEU:HB3	1:D:139:VAL:HG12	1.85	0.57
1:H:269:ASP:O	1:H:273:THR:HG23	2.06	0.56
1:F:331:TYR:CE1	1:G:172:VAL:HG22	2.41	0.56
1:C:216:MET:N	1:C:216:MET:HE3	2.21	0.55
1:H:130:LYS:HG2	2:H:414:HOH:O	2.05	0.55
1:H:267:PHE:O	1:H:271:THR:HG23	2.07	0.55
1:B:201:ILE:HD13	1:B:216:MET:HE1	1.89	0.55
1:B:178:ASN:OD1	1:B:180:SER:HB3	2.06	0.54
1:C:329:MET:HE2	1:C:329:MET:H	1.72	0.54
1:H:178:ASN:HD22	1:H:180:SER:H	1.55	0.54
1:A:326:SER:HB2	1:A:327:PRO:CD	2.31	0.53
1:H:215:ILE:HG22	1:H:216:MET:HE1	1.89	0.53
1:B:331:TYR:CE1	1:C:172:VAL:HG22	2.44	0.53
1:A:332:LEU:HD23	1:A:335:LYS:HD2	1.92	0.52
1:G:268:TYR:CZ	1:G:272:ILE:HD11	2.45	0.52
1:B:201:ILE:HD13	1:B:216:MET:CE	2.40	0.51
1:C:106:VAL:HG11	1:C:272:ILE:CD1	2.40	0.51
1:C:331:TYR:CE1	1:D:172:VAL:HG22	2.45	0.51
1:C:331:TYR:CD2	1:D:116:VAL:HA	2.47	0.50
1:E:331:TYR:HE1	1:F:172:VAL:HG22	1.75	0.50
1:C:215:ILE:HG22	1:C:216:MET:HE1	1.93	0.50
1:F:139:VAL:HG23	1:F:156:ALA:HB2	1.95	0.49
1:F:164:GLU:HB3	1:G:116:VAL:HG23	1.95	0.49
1:E:269:ASP:O	1:E:273:THR:HG23	2.13	0.49
1:D:237:MET:HE1	1:D:245:ILE:HD12	1.95	0.49
1:D:267:PHE:O	1:D:271:THR:HG23	2.12	0.48
1:C:58:SER:HB3	1:C:103:LYS:HD2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:66:SER:O	1:G:84:TYR:HB2	2.13	0.48
1:C:331:TYR:HE1	1:D:172:VAL:HG22	1.78	0.48
1:F:66:SER:O	1:F:84:TYR:HB2	2.13	0.48
1:C:216:MET:CE	1:C:216:MET:N	2.77	0.48
1:H:63:LEU:HD21	1:H:290:TYR:HD2	1.77	0.48
1:E:106:VAL:HG23	1:E:275:MET:CE	2.43	0.48
1:D:178:ASN:ND2	1:E:189:SER:HB3	2.29	0.48
1:E:154:THR:HG21	1:E:230:VAL:HB	1.95	0.48
1:H:169:SER:HB2	1:H:250:VAL:HG12	1.96	0.47
1:G:331:TYR:CE1	1:H:172:VAL:HG22	2.49	0.47
1:A:177:VAL:O	1:A:204:PRO:HA	2.14	0.47
1:E:107:LEU:HB3	1:E:139:VAL:HG22	1.97	0.47
1:F:162:THR:O	1:F:253:GLY:HA3	2.15	0.47
1:E:103:LYS:O	1:E:136:PRO:HD2	2.14	0.47
1:F:331:TYR:HE1	1:G:172:VAL:HG22	1.81	0.46
1:F:179:TYR:CE2	1:F:204:PRO:HB2	2.51	0.46
1:C:229:ASP:OD1	1:C:239:LYS:NZ	2.41	0.46
1:F:215:ILE:HG22	1:F:216:MET:CE	2.45	0.46
1:A:161:ALA:O	1:A:265:LEU:HA	2.15	0.46
1:C:123:HIS:CE1	1:C:230:VAL:HG22	2.51	0.45
1:A:331:TYR:CD2	1:B:116:VAL:HA	2.51	0.45
1:C:166:LEU:HB2	1:C:329:MET:HE1	1.99	0.45
1:G:176:SER:HG	1:H:191:GLU:HB3	1.80	0.45
1:G:169:SER:HA	1:G:248:GLY:O	2.15	0.45
1:E:169:SER:HA	1:E:248:GLY:O	2.17	0.45
1:D:85:ASN:HB3	1:D:88:THR:HB	1.99	0.44
1:C:176:SER:HB3	1:D:191:GLU:O	2.16	0.44
1:E:151:TYR:O	1:E:154:THR:HB	2.18	0.44
1:H:215:ILE:HG22	1:H:216:MET:HE2	1.99	0.44
1:C:63:LEU:HD21	1:C:290:TYR:HD2	1.83	0.44
1:A:162:THR:O	1:A:253:GLY:HA3	2.17	0.44
1:D:179:TYR:CE2	1:D:204:PRO:HB2	2.52	0.44
1:C:216:MET:CE	1:C:216:MET:HA	2.42	0.44
1:C:161:ALA:O	1:C:265:LEU:HA	2.18	0.44
1:H:59:LYS:N	1:H:101:THR:O	2.51	0.44
1:C:165:THR:O	1:C:252:ASP:HA	2.18	0.44
1:F:201:ILE:CD1	1:F:216:MET:HE3	2.47	0.43
1:H:160:PHE:CG	1:H:271:THR:HG22	2.53	0.43
1:H:110:ASN:HA	1:H:144:MET:O	2.18	0.43
1:E:237:MET:HB3	1:E:237:MET:HE2	1.77	0.43
1:B:201:ILE:CD1	1:B:216:MET:CE	2.97	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:TYR:O	1:B:147:SER:HB3	2.19	0.43
1:D:208:MET:HE1	1:D:216:MET:HG3	2.01	0.42
1:H:64:GLU:HG2	1:H:108:LYS:HB2	2.01	0.42
1:B:213:LYS:NZ	1:C:191:GLU:OE1	2.52	0.42
1:B:201:ILE:O	1:B:202:MET:HB2	2.20	0.42
1:A:63:LEU:HD13	1:A:93:LEU:HD13	2.01	0.42
1:E:66:SER:O	1:E:84:TYR:HB2	2.20	0.42
1:E:201:ILE:O	1:E:202:MET:HB2	2.19	0.42
1:D:267:PHE:O	1:D:271:THR:CG2	2.67	0.41
1:D:178:ASN:HD21	1:E:189:SER:HB3	1.85	0.41
1:F:213:LYS:HD3	1:G:191:GLU:OE2	2.21	0.41
1:G:331:TYR:CD2	1:H:116:VAL:HA	2.56	0.41
1:A:110:ASN:HA	1:A:144:MET:O	2.21	0.41
1:D:331:TYR:C	1:E:147:SER:OG	2.58	0.41
1:F:280:LYS:O	1:F:283:LYS:HB2	2.21	0.41
1:E:106:VAL:HG23	1:E:275:MET:HE2	2.01	0.41
1:H:139:VAL:HG23	1:H:156:ALA:HB2	2.02	0.41
1:A:125:LYS:HA	1:A:125:LYS:HD3	1.97	0.41
1:E:139:VAL:HG23	1:E:156:ALA:HB2	2.02	0.41
1:D:135:LYS:HA	1:D:136:PRO:HD3	1.96	0.41
1:B:87:ARG:HA	1:B:87:ARG:HD3	1.84	0.41
1:B:147:SER:HB3	1:B:148:GLY:H	1.66	0.40
1:E:181:LYS:HB2	1:F:187:GLY:HA3	2.04	0.40
1:D:154:THR:HG21	1:D:230:VAL:HB	2.04	0.40
1:G:203:SER:HA	1:G:204:PRO:HD3	1.95	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	231/273 (85%)	227 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	227/273 (83%)	220 (97%)	7 (3%)	0	100	100
1	C	231/273 (85%)	228 (99%)	3 (1%)	0	100	100
1	D	222/273 (81%)	216 (97%)	6 (3%)	0	100	100
1	E	226/273 (83%)	220 (97%)	6 (3%)	0	100	100
1	F	228/273 (84%)	225 (99%)	3 (1%)	0	100	100
1	G	226/273 (83%)	219 (97%)	7 (3%)	0	100	100
1	H	226/273 (83%)	220 (97%)	6 (3%)	0	100	100
All	All	1817/2184 (83%)	1775 (98%)	42 (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	190/231 (82%)	188 (99%)	2 (1%)	80	92
1	B	184/231 (80%)	179 (97%)	5 (3%)	52	73
1	C	194/231 (84%)	188 (97%)	6 (3%)	47	69
1	D	170/231 (74%)	162 (95%)	8 (5%)	32	50
1	E	184/231 (80%)	179 (97%)	5 (3%)	52	73
1	F	191/231 (83%)	188 (98%)	3 (2%)	70	86
1	G	187/231 (81%)	180 (96%)	7 (4%)	41	62
1	H	191/231 (83%)	184 (96%)	7 (4%)	41	62
All	All	1491/1848 (81%)	1448 (97%)	43 (3%)	50	71

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	LYS
1	A	227	PHE

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Mol	Chain	Res	Type
1	B	147	SER
1	B	173	ILE
1	B	184	ASP
1	B	185	LYS
1	B	227	PHE
1	C	58	SER
1	C	216	MET
1	C	227	PHE
1	C	230	VAL
1	C	282	LEU
1	C	329	MET
1	D	143	SER
1	D	147	SER
1	D	174	MET
1	D	178	ASN
1	D	218	SER
1	D	227	PHE
1	D	271	THR
1	D	332	LEU
1	E	154	THR
1	E	186	LEU
1	E	227	PHE
1	E	286	SER
1	E	288	ILE
1	F	70	GLN
1	F	131	LYS
1	F	227	PHE
1	G	133	THR
1	G	147	SER
1	G	174	MET
1	G	178	ASN
1	G	179	TYR
1	G	227	PHE
1	G	281	ASP
1	H	95	ARG
1	H	147	SER
1	H	174	MET
1	H	189	SER
1	H	227	PHE
1	H	271	THR
1	H	273	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	178	ASN
1	A	217	GLN
1	B	110	ASN
1	B	198	HIS
1	C	70	GLN
1	C	214	ASN
1	D	222	ASN
1	E	110	ASN
1	E	214	ASN
1	F	110	ASN
1	G	70	GLN
1	G	110	ASN
1	G	178	ASN
1	G	214	ASN
1	G	222	ASN
1	H	70	GLN
1	H	178	ASN

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

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	237/273 (86%)	-0.02	2 (0%) 87 87	26, 40, 61, 70	0
1	B	233/273 (85%)	0.17	7 (3%) 54 53	33, 50, 92, 106	0
1	C	237/273 (86%)	-0.04	0 100 100	30, 44, 64, 74	0
1	D	228/273 (83%)	0.25	7 (3%) 52 52	36, 55, 106, 122	0
1	E	232/273 (84%)	0.25	13 (5%) 28 28	32, 52, 101, 122	0
1	F	234/273 (85%)	0.03	5 (2%) 67 66	27, 42, 68, 82	0
1	G	232/273 (84%)	0.23	13 (5%) 28 28	30, 49, 93, 114	0
1	H	232/273 (84%)	0.09	2 (0%) 85 85	30, 46, 74, 90	0
All	All	1865/2184 (85%)	0.12	49 (2%) 59 58	26, 46, 86, 122	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	282	LEU	5.3
1	B	282	LEU	5.0
1	G	288	ILE	3.7
1	G	282	LEU	3.6
1	E	287	VAL	3.4
1	E	100	LYS	3.3
1	E	58	SER	3.2
1	G	95	ARG	3.2
1	G	102	VAL	3.2
1	B	280	LYS	3.1
1	E	285	ALA	3.0
1	A	177	VAL	2.9
1	G	58	SER	2.9
1	E	101	THR	2.9
1	H	290	TYR	2.9
1	H	284	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	G	83	GLY	2.8
1	F	331	TYR	2.8
1	D	284	ASN	2.7
1	G	57	SER	2.6
1	G	281	ASP	2.6
1	D	177	VAL	2.6
1	F	83	GLY	2.6
1	G	101	THR	2.5
1	E	333	TYR	2.4
1	D	98	ASP	2.4
1	B	284	ASN	2.3
1	B	102	VAL	2.3
1	G	99	ASP	2.3
1	G	268	TYR	2.3
1	D	282	LEU	2.3
1	E	288	ILE	2.3
1	B	134	LYS	2.3
1	G	104	GLY	2.2
1	E	102	VAL	2.2
1	E	281	ASP	2.2
1	F	95	ARG	2.2
1	D	61	ALA	2.2
1	A	179	TYR	2.2
1	F	57	SER	2.1
1	D	286	SER	2.1
1	B	278	ASP	2.1
1	F	333	TYR	2.1
1	E	60	ILE	2.1
1	G	98	ASP	2.1
1	D	104	GLY	2.0
1	E	279	HIS	2.0
1	E	99	ASP	2.0
1	B	107	LEU	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 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.