



Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2020 – 01:33 pm BST

PDB ID : 5YZN
Title : Crystal structure of S9 peptidase (active form) from *Deinococcus radiodurans* R1
Authors : Yadav, P.; Jamdar, S.N.; Kumar, A.; Ghosh, B.; Makde, R.D.
Deposited on : 2017-12-15
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

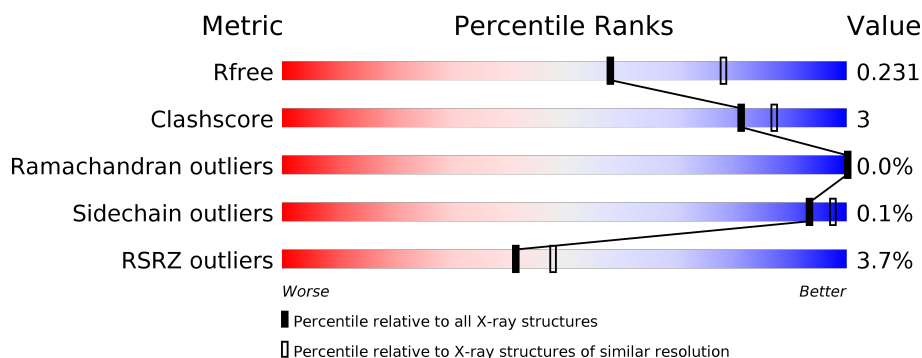
1 Overall quality at a glance

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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 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	656	<div> <div>3%</div> <div> <div></div> <div>93%</div> <div>6%</div> </div> </div>
1	B	656	<div> <div>3%</div> <div> <div></div> <div>92%</div> <div>7%</div> </div> </div>
1	C	656	<div> <div>4%</div> <div> <div></div> <div>93%</div> <div>6%</div> </div> </div>
1	D	656	<div> <div>4%</div> <div> <div></div> <div>90%</div> <div>8%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 21431 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 Acyl-peptide hydrolase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	649	Total	C	N	O	S	0	0	0
			5054	3197	912	936	9			
1	B	649	Total	C	N	O	S	0	0	0
			5038	3186	909	934	9			
1	C	646	Total	C	N	O	S	0	0	0
			5002	3164	903	926	9			
1	D	644	Total	C	N	O	S	0	0	0
			5006	3166	905	926	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q9RXY9
A	1	SER	-	expression tag	UNP Q9RXY9
B	0	GLY	-	expression tag	UNP Q9RXY9
B	1	SER	-	expression tag	UNP Q9RXY9
C	0	GLY	-	expression tag	UNP Q9RXY9
C	1	SER	-	expression tag	UNP Q9RXY9
D	0	GLY	-	expression tag	UNP Q9RXY9
D	1	SER	-	expression tag	UNP Q9RXY9

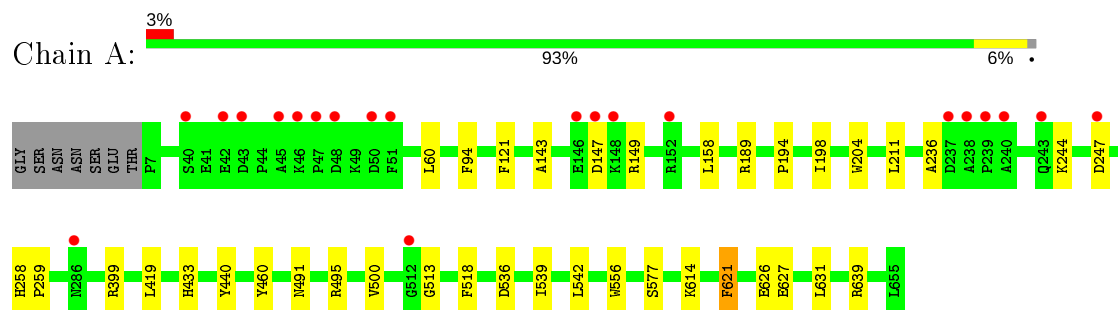
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	331	Total	O	0	0
			331	331		
2	B	373	Total	O	0	0
			373	373		
2	C	292	Total	O	0	0
			292	292		
2	D	335	Total	O	0	0
			335	335		

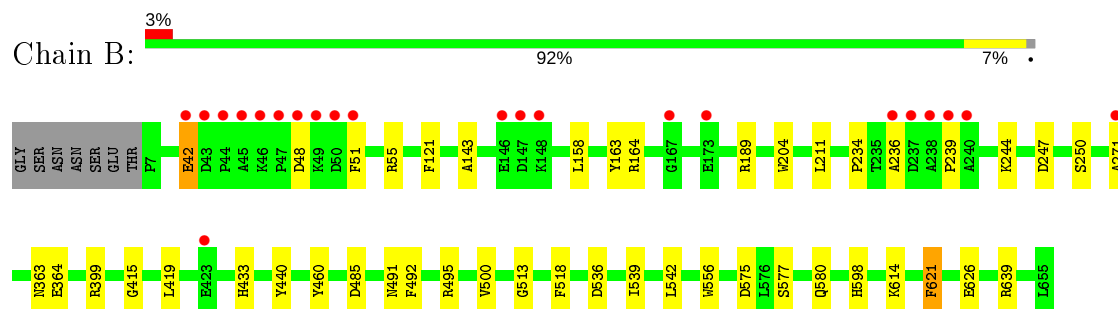
3 Residue-property plots [i](#)

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

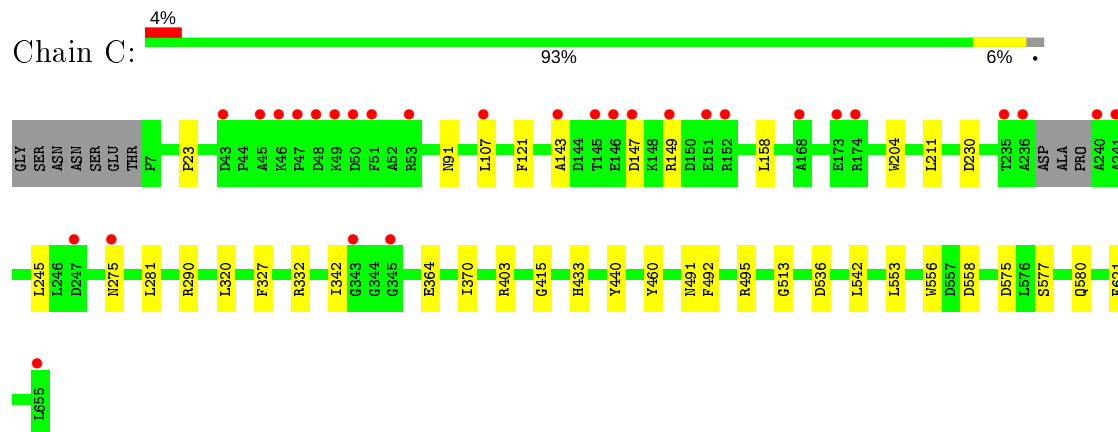
- Molecule 1: Acyl-peptide hydrolase, putative



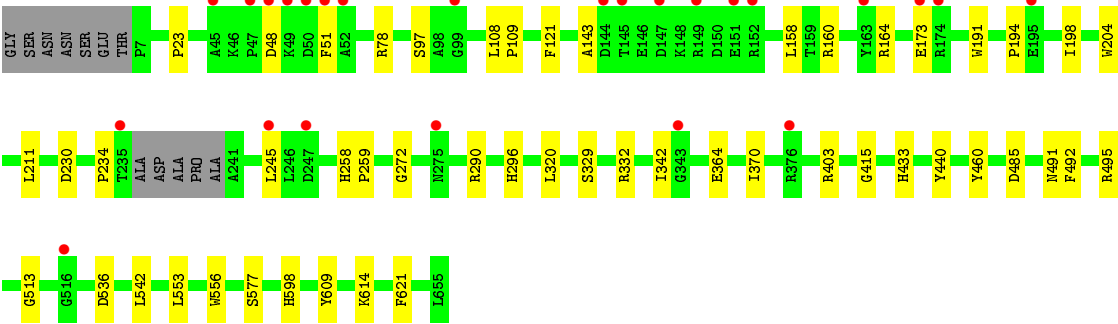
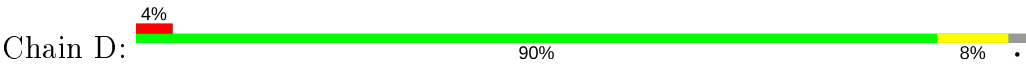
- Molecule 1: Acyl-peptide hydrolase, putative



- Molecule 1: Acyl-peptide hydrolase, putative



- Molecule 1: Acyl-peptide hydrolase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	120.43Å 148.22Å 190.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.28 – 2.30 39.28 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (39.28-2.30) 99.8 (39.28-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???), REFMAC	Depositor
R, R_{free}	0.204 , 0.231 0.205 , 0.231	Depositor DCC
R_{free} test set	7518 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.843	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.2	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	0.96	EDS
Total number of atoms	21431	wwPDB-VP
Average B, all atoms (Å ²)	35.0	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 49.92 % 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 6.9602e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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 [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	0.26	0/5209	0.48	0/7111
1	B	0.26	0/5193	0.48	0/7094
1	C	0.25	0/5155	0.49	0/7043
1	D	0.25	0/5159	0.49	0/7045
All	All	0.25	0/20716	0.49	0/28293

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	5054	0	4771	22	0
1	B	5038	0	4734	29	0
1	C	5002	0	4685	25	0
1	D	5006	0	4701	30	0
2	A	331	0	0	1	0
2	B	373	0	0	2	0
2	C	292	0	0	3	0
2	D	335	0	0	2	0
All	All	21431	0	18891	100	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 3.

All (100) 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:B:42:GLU:HG2	1:B:55:ARG:HG3	1.61	0.80
1:D:230:ASP:HB2	1:D:245:LEU:HD23	1.72	0.70
1:C:230:ASP:HB2	1:C:245:LEU:HD23	1.75	0.67
1:C:91:ASN:ND2	1:C:107:LEU:HD21	2.11	0.65
1:A:626:GLU:OE1	1:A:639:ARG:NH2	2.30	0.64
1:B:189:ARG:NH2	1:B:236:ALA:O	2.30	0.62
1:A:244:LYS:NZ	1:A:247:ASP:OD1	2.23	0.61
1:B:626:GLU:OE1	1:B:639:ARG:NH2	2.33	0.61
1:D:204:TRP:CZ2	1:D:211:LEU:HD21	2.38	0.59
1:C:332:ARG:NH2	2:C:702:HOH:O	2.37	0.58
1:B:158:LEU:HB2	1:C:621:PHE:HB3	1.88	0.56
1:A:204:TRP:CZ2	1:A:211:LEU:HD21	2.42	0.55
1:C:204:TRP:CZ2	1:C:211:LEU:HD21	2.42	0.54
1:B:234:PRO:HB3	1:B:239:PRO:HB3	1.90	0.54
1:B:621:PHE:HB3	1:C:158:LEU:HB2	1.92	0.52
1:C:440:TYR:CD2	1:C:460:TYR:HB2	2.45	0.52
1:A:491:ASN:O	1:A:495:ARG:HG2	2.10	0.51
1:D:403:ARG:NH2	2:D:706:HOH:O	2.42	0.51
1:D:440:TYR:CD2	1:D:460:TYR:HB2	2.46	0.51
1:A:147:ASP:OD1	1:A:149:ARG:HD3	2.10	0.51
1:D:78:ARG:HB3	1:D:97:SER:OG	2.11	0.51
1:A:614:LYS:HG3	1:D:553:LEU:HD22	1.94	0.50
1:C:147:ASP:OD1	1:C:149:ARG:HD3	2.12	0.50
1:A:158:LEU:HB2	1:D:621:PHE:HB3	1.94	0.50
1:C:491:ASN:OD1	1:C:495:ARG:NH1	2.45	0.49
1:A:440:TYR:CD2	1:A:460:TYR:HB2	2.47	0.49
1:B:204:TRP:CZ2	1:B:211:LEU:HD21	2.48	0.49
1:B:491:ASN:O	1:B:495:ARG:HG2	2.13	0.49
1:B:244:LYS:NZ	1:B:247:ASP:OD1	2.43	0.49
1:D:290:ARG:NH2	1:D:342:ILE:HG23	2.28	0.49
1:A:621:PHE:HB3	1:D:158:LEU:HB2	1.96	0.48
1:D:191:TRP:NE1	1:D:234:PRO:HD3	2.29	0.48
1:D:23:PRO:HG3	1:D:370:ILE:HG13	1.96	0.48
1:B:364:GLU:HG2	2:B:994:HOH:O	2.14	0.47
1:A:542:LEU:HG	1:A:556:TRP:CZ2	2.49	0.47
1:B:542:LEU:HG	1:B:556:TRP:CZ2	2.49	0.47
1:A:194:PRO:HG3	1:A:198:ILE:HG13	1.97	0.47
1:A:189:ARG:NH2	1:A:236:ALA:O	2.38	0.47
1:B:419:LEU:HD21	1:B:500:VAL:HG11	1.96	0.47
1:B:250:SER:HB2	1:B:271:ALA:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:440:TYR:CD2	1:B:460:TYR:HB2	2.49	0.47
1:A:542:LEU:HD13	1:A:577:SER:HA	1.95	0.47
1:A:399:ARG:NE	1:A:419:LEU:O	2.45	0.46
1:A:419:LEU:HD21	1:A:500:VAL:HG11	1.97	0.46
1:C:320:LEU:O	1:C:364:GLU:HG2	2.16	0.46
1:C:491:ASN:O	1:C:495:ARG:HG2	2.16	0.46
1:B:164:ARG:HD2	1:B:598:HIS:CE1	2.51	0.46
1:D:542:LEU:HG	1:D:556:TRP:CZ2	2.49	0.46
1:B:48:ASP:HB3	1:B:51:PHE:O	2.16	0.46
1:D:415:GLY:HA3	1:D:492:PHE:CE1	2.51	0.46
1:C:433:HIS:HB3	1:C:440:TYR:CZ	2.51	0.46
1:A:518:PHE:HB2	1:A:539:ILE:HB	1.98	0.45
1:D:160:ARG:NH2	1:D:173:GLU:OE1	2.49	0.45
1:C:403:ARG:NH2	2:C:715:HOH:O	2.49	0.45
1:D:320:LEU:O	1:D:364:GLU:HG2	2.17	0.45
1:C:290:ARG:NH2	1:C:342:ILE:HG23	2.33	0.45
1:C:415:GLY:HA3	1:C:492:PHE:CE1	2.52	0.44
1:A:258:HIS:CG	1:A:259:PRO:HD2	2.53	0.44
1:A:433:HIS:HB3	1:A:440:TYR:CZ	2.53	0.44
1:D:491:ASN:O	1:D:495:ARG:HG2	2.18	0.44
1:C:23:PRO:HG3	1:C:370:ILE:HG13	2.00	0.44
1:B:415:GLY:HA3	1:B:492:PHE:CE1	2.53	0.44
1:C:542:LEU:HD13	1:C:577:SER:HA	2.00	0.44
1:B:163:TYR:CD2	1:B:164:ARG:HG3	2.53	0.43
1:B:121:PHE:CD1	1:B:143:ALA:HA	2.53	0.43
1:C:275:ASN:ND2	1:C:558:ASP:OD2	2.37	0.43
1:D:296:HIS:HE1	1:D:329:SER:OG	2.01	0.43
1:A:121:PHE:CD1	1:A:143:ALA:HA	2.53	0.43
1:B:399:ARG:NE	1:B:419:LEU:O	2.50	0.43
1:B:433:HIS:HB3	1:B:440:TYR:CZ	2.54	0.43
1:D:433:HIS:HB3	1:D:440:TYR:CZ	2.53	0.43
1:D:121:PHE:CD1	1:D:143:ALA:HA	2.54	0.43
1:B:363:ASN:HB2	2:B:994:HOH:O	2.19	0.42
1:B:513:GLY:HA2	1:B:536:ASP:O	2.19	0.42
1:C:121:PHE:CD1	1:C:143:ALA:HA	2.54	0.42
1:B:542:LEU:HD13	1:B:577:SER:HA	2.00	0.42
1:C:281:LEU:HD11	1:C:327:PHE:CG	2.54	0.42
1:A:627:GLU:HG3	1:A:631:LEU:HD13	2.01	0.42
1:C:575:ASP:HA	1:C:580:GLN:NE2	2.34	0.42
1:B:575:ASP:HA	1:B:580:GLN:NE2	2.34	0.42
1:C:542:LEU:HG	1:C:556:TRP:CZ2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:332:ARG:NH2	2:D:719:HOH:O	2.51	0.42
1:B:485:ASP:OD1	1:B:485:ASP:N	2.53	0.42
1:B:614:LYS:HG2	1:C:553:LEU:HD22	2.01	0.42
1:D:513:GLY:HA2	1:D:536:ASP:O	2.20	0.42
1:D:48:ASP:HB3	1:D:51:PHE:O	2.19	0.42
1:D:542:LEU:HD13	1:D:577:SER:HA	2.02	0.42
1:C:149:ARG:NH2	2:C:718:HOH:O	2.50	0.42
2:A:742:HOH:O	1:D:614:LYS:HE2	2.20	0.41
1:D:485:ASP:OD1	1:D:485:ASP:N	2.52	0.41
1:D:258:HIS:CG	1:D:259:PRO:HD2	2.56	0.41
1:C:513:GLY:HA2	1:C:536:ASP:O	2.20	0.41
1:D:164:ARG:HD2	1:D:598:HIS:CE1	2.55	0.41
1:D:194:PRO:HG3	1:D:198:ILE:HG13	2.03	0.41
1:D:108:LEU:HD12	1:D:109:PRO:HD2	2.02	0.41
1:B:575:ASP:HA	1:B:580:GLN:HE22	1.86	0.41
1:A:513:GLY:HA2	1:A:536:ASP:O	2.21	0.41
1:D:609:TYR:CD1	1:D:621:PHE:HB2	2.56	0.40
1:B:518:PHE:HB2	1:B:539:ILE:HB	2.03	0.40
1:A:60:LEU:HD11	1:A:94:PHE:CG	2.56	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	647/656 (99%)	626 (97%)	21 (3%)	0	100	100
1	B	647/656 (99%)	626 (97%)	21 (3%)	0	100	100
1	C	642/656 (98%)	619 (96%)	23 (4%)	0	100	100
1	D	640/656 (98%)	620 (97%)	19 (3%)	1 (0%)	47	58
All	All	2576/2624 (98%)	2491 (97%)	84 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	272	GLY

5.3.2 Protein sidechains [i](#)

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	510/526 (97%)	509 (100%)	1 (0%)	93	97
1	B	506/526 (96%)	504 (100%)	2 (0%)	91	96
1	C	500/526 (95%)	500 (100%)	0	100	100
1	D	503/526 (96%)	503 (100%)	0	100	100
All	All	2019/2104 (96%)	2016 (100%)	3 (0%)	93	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	621	PHE
1	B	42	GLU
1	B	621	PHE

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

Mol	Chain	Res	Type
1	D	296	HIS

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	649/656 (98%)	0.02	21 (3%)	47 54	22, 32, 57, 95	0
1	B	649/656 (98%)	-0.09	22 (3%)	45 52	20, 31, 53, 84	0
1	C	646/656 (98%)	0.07	29 (4%)	33 40	22, 35, 58, 85	0
1	D	644/656 (98%)	0.01	25 (3%)	39 46	20, 33, 56, 73	0
All	All	2588/2624 (98%)	0.00	97 (3%)	41 48	20, 33, 57, 95	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	51	PHE	4.9
1	B	45	ALA	4.8
1	C	47	PRO	4.5
1	C	240	ALA	4.5
1	A	237	ASP	4.5
1	C	45	ALA	4.3
1	A	238	ALA	4.3
1	B	50	ASP	4.3
1	B	51	PHE	4.3
1	B	237	ASP	4.3
1	C	50	ASP	4.0
1	A	45	ALA	3.9
1	C	236	ALA	3.9
1	D	50	ASP	3.8
1	C	235	THR	3.8
1	D	47	PRO	3.8
1	C	51	PHE	3.8
1	B	238	ALA	3.8
1	B	239	PRO	3.7
1	C	168	ALA	3.5
1	D	48	ASP	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	50	ASP	3.5
1	C	48	ASP	3.5
1	B	173	GLU	3.4
1	A	148	LYS	3.4
1	C	46	LYS	3.4
1	D	235	THR	3.4
1	B	240	ALA	3.3
1	B	48	ASP	3.3
1	D	51	PHE	3.3
1	C	241	ALA	3.2
1	D	152	ARG	3.2
1	C	146	GLU	3.1
1	C	343	GLY	3.1
1	D	145	THR	3.0
1	D	173	GLU	3.0
1	D	343	GLY	3.0
1	A	240	ALA	2.9
1	D	247	ASP	2.8
1	B	236	ALA	2.7
1	C	345	GLY	2.7
1	A	147	ASP	2.7
1	C	43	ASP	2.7
1	A	46	LYS	2.6
1	B	146	GLU	2.6
1	C	173	GLU	2.6
1	C	149	ARG	2.6
1	D	45	ALA	2.6
1	A	48	ASP	2.6
1	D	195	GLU	2.6
1	B	49	LYS	2.5
1	A	152	ARG	2.5
1	A	47	PRO	2.5
1	D	52	ALA	2.5
1	A	239	PRO	2.5
1	B	46	LYS	2.5
1	D	245	LEU	2.5
1	B	271	ALA	2.5
1	C	152	ARG	2.5
1	D	376	ARG	2.4
1	B	44	PRO	2.4
1	D	144	ASP	2.4
1	D	99	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	243	GLN	2.4
1	B	42	GLU	2.4
1	D	174	ARG	2.3
1	B	47	PRO	2.3
1	A	43	ASP	2.3
1	D	147	ASP	2.3
1	C	53	ARG	2.3
1	C	145	THR	2.3
1	B	43	ASP	2.3
1	A	146	GLU	2.3
1	D	49	LYS	2.3
1	A	40	SER	2.2
1	B	148	LYS	2.2
1	C	151	GLU	2.2
1	D	151	GLU	2.2
1	C	107	LEU	2.1
1	C	147	ASP	2.1
1	B	423	GLU	2.1
1	C	275	ASN	2.1
1	C	655	LEU	2.1
1	B	167	GLY	2.1
1	A	42	GLU	2.1
1	A	286	ASN	2.1
1	C	49	LYS	2.1
1	D	516	GLY	2.1
1	B	147	ASP	2.0
1	D	149	ARG	2.0
1	C	143	ALA	2.0
1	C	174	ARG	2.0
1	D	163	TYR	2.0
1	A	247	ASP	2.0
1	D	275	ASN	2.0
1	A	512	GLY	2.0
1	C	247	ASP	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 [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.