



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:58 PM UTC

PDB ID : 5YZM / pdb_00005yzm
Title : Crystal structure of S9 peptidase (inactive 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 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
Mogul : 2022.3.0, CSD as543be (2022)
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

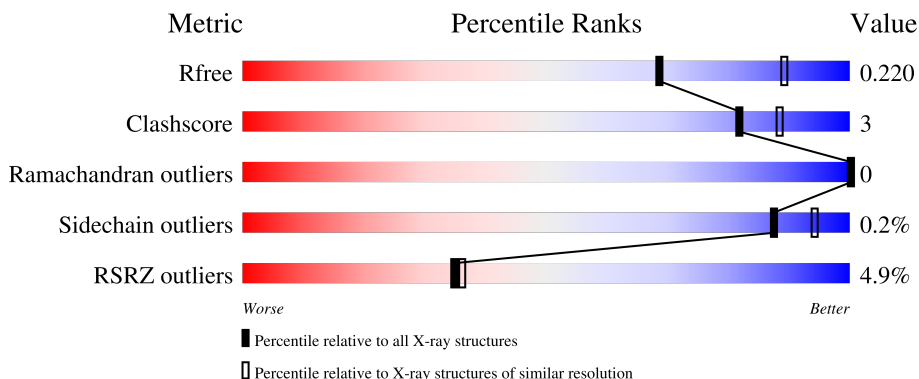
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 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	656	2% 87% 9%
1	B	656	4% 88% 8%
1	C	656	5% 85% 6% 10%
1	D	656	7% 81% 8% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19434 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	598	Total	C	N	O	S	0	0	0
			4633	2941	831	852	9			
1	B	601	Total	C	N	O	S	0	1	0
			4627	2935	829	854	9			
1	C	593	Total	C	N	O	S	0	0	0
			4560	2897	820	834	9			
1	D	590	Total	C	N	O	S	0	0	0
			4546	2891	815	831	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 ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

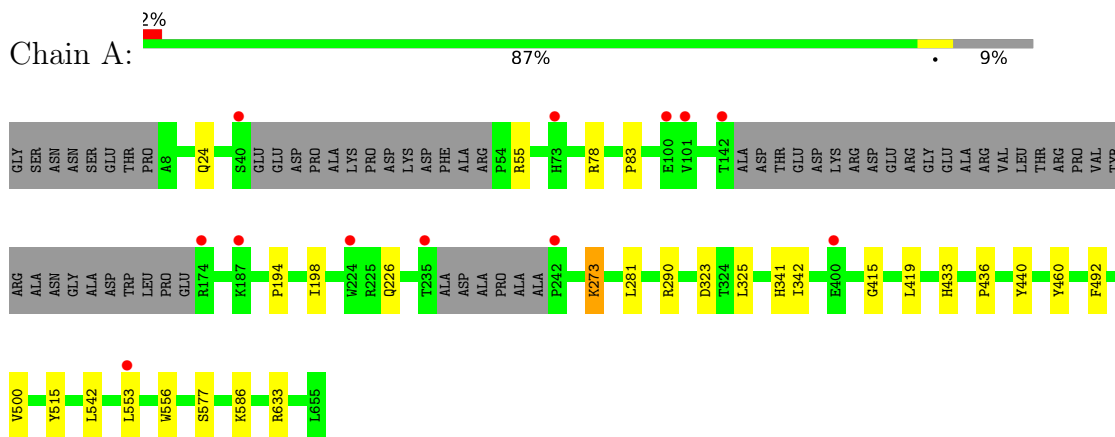
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	305	Total	O	0	0
			305	305		
3	B	327	Total	O	0	0
			327	327		
3	C	221	Total	O	0	0
			221	221		
3	D	199	Total	O	0	0
			199	199		

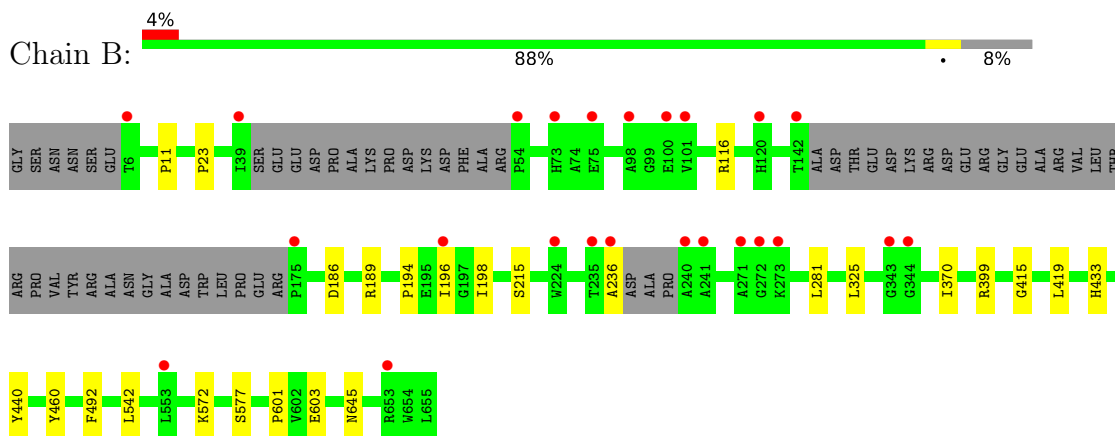
3 Residue-property plots [i](#)

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.

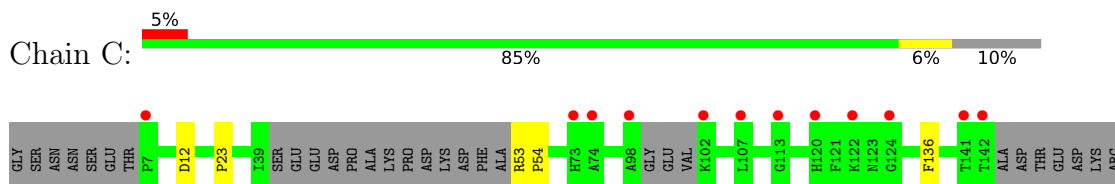
- Molecule 1: Acyl-peptide hydrolase, putative

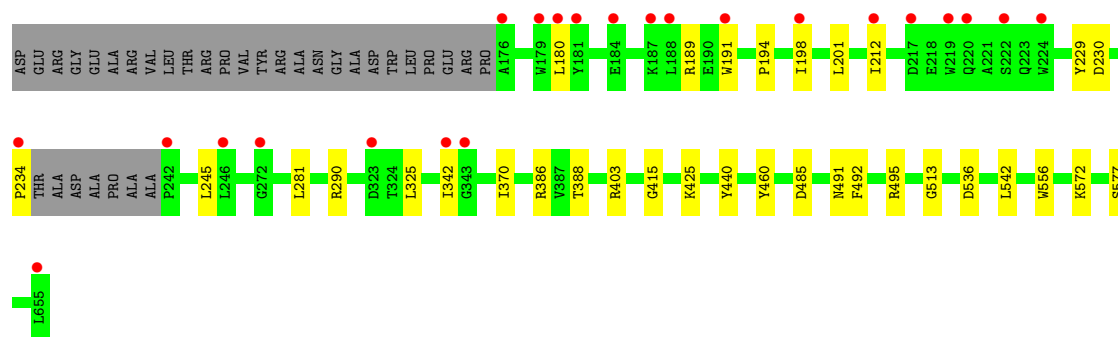


- Molecule 1: Acyl-peptide hydrolase, putative



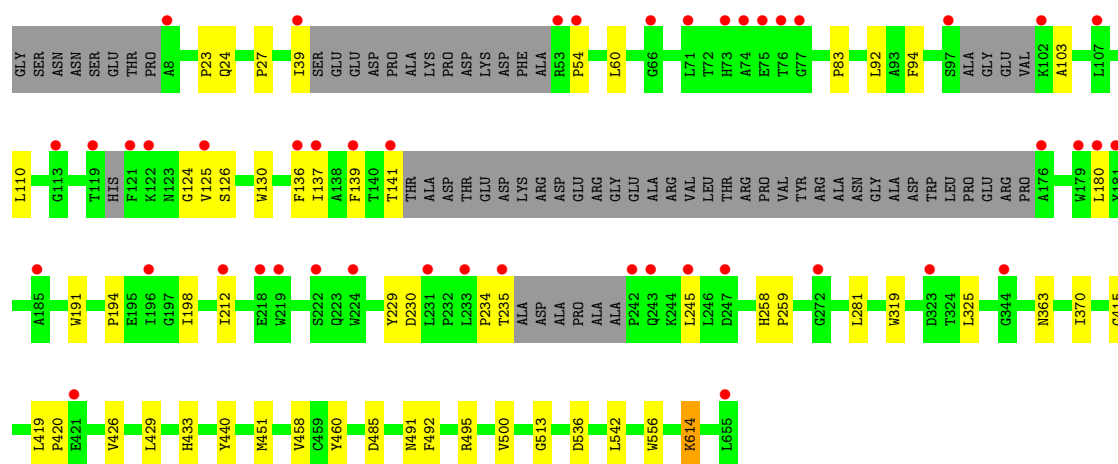
- Molecule 1: Acyl-peptide hydrolase, putative





- Molecule 1: Acyl-peptide hydrolase, putative

Chain D: 7% 81% 8% 10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.70Å 195.09Å 107.72Å 90.00° 99.97° 90.00°	Depositor
Resolution (Å)	46.87 – 2.30 46.87 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.87-2.30) 99.8 (46.87-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???), REFMAC	Depositor
R, R_{free}	0.189 , 0.219 0.190 , 0.220	Depositor DCC
R_{free} test set	6541 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.4	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	0.95	EDS
Total number of atoms	19434	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACT

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	0.13	0/4772	0.36	0/6507
1	B	0.14	0/4765	0.36	0/6504
1	C	0.12	0/4696	0.34	0/6406
1	D	0.12	0/4682	0.34	0/6387
All	All	0.13	0/18915	0.35	0/25804

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	4633	0	4372	18	0
1	B	4627	0	4351	15	0
1	C	4560	0	4289	24	0
1	D	4546	0	4273	34	0
2	A	4	0	3	0	0
2	B	4	0	3	0	0
2	C	4	0	3	0	0
2	D	4	0	3	0	0
3	A	305	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	327	0	0	1	0
3	C	221	0	0	3	0
3	D	199	0	0	0	0
All	All	19434	0	17297	90	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 (90) 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:403:ARG:NH2	3:C:801:HOH:O	2.23	0.71
1:A:226:GLN:NE2	3:A:801:HOH:O	2.21	0.71
1:C:290:ARG:NH2	1:C:342:ILE:O	2.28	0.67
1:C:491:ASN:OD1	1:C:495:ARG:NH1	2.30	0.65
1:D:429:LEU:HD12	1:D:451:MET:HE3	1.81	0.63
1:D:451:MET:HE2	1:D:458:VAL:HG22	1.79	0.63
1:B:194:PRO:HG3	1:B:198:ILE:HG13	1.82	0.61
1:C:194:PRO:HG3	1:C:198:ILE:HG13	1.84	0.60
1:A:586:LYS:NZ	3:A:803:HOH:O	2.35	0.60
1:D:194:PRO:HG3	1:D:198:ILE:HG13	1.84	0.58
1:C:53:ARG:HD2	1:C:54:PRO:HD2	1.85	0.58
1:B:572:LYS:NZ	3:B:806:HOH:O	2.36	0.57
1:D:27:PRO:O	1:D:363:ASN:ND2	2.38	0.57
1:C:189:ARG:HH21	1:C:234:PRO:HB2	1.70	0.55
1:D:103:ALA:HB3	1:D:125:VAL:H	1.71	0.55
1:B:189:ARG:HH22	1:B:236:ALA:HB3	1.71	0.54
1:C:191:TRP:CD1	1:C:234:PRO:HG3	2.43	0.54
1:A:194:PRO:HG3	1:A:198:ILE:HG13	1.89	0.54
1:A:290:ARG:NH2	1:A:342:ILE:O	2.42	0.53
1:A:55:ARG:NH2	3:A:809:HOH:O	2.42	0.53
1:C:281:LEU:HD13	1:C:325:LEU:HD23	1.91	0.52
1:A:553:LEU:HD22	1:D:614:LYS:HD3	1.90	0.52
1:D:136:PHE:HB3	1:D:180:LEU:HD11	1.91	0.52
1:B:542:LEU:HD13	1:B:577:SER:HA	1.92	0.52
1:D:542:LEU:HG	1:D:556:TRP:CZ2	2.45	0.52
1:B:433:HIS:HB3	1:B:440:TYR:CZ	2.45	0.51
1:D:420:PRO:HG2	1:D:426:VAL:HG11	1.91	0.51
1:A:281:LEU:HD13	1:A:325:LEU:HD23	1.92	0.50
1:B:116:ARG:HH22	1:B:186:ASP:CG	2.20	0.50
1:C:386:ARG:NH1	1:C:388:THR:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:HIS:HB3	1:A:440:TYR:CZ	2.46	0.49
1:D:92:LEU:HG	1:D:110:LEU:HD11	1.95	0.49
1:B:281:LEU:HD13	1:B:325:LEU:HD23	1.93	0.49
1:C:230:ASP:HB2	1:C:245:LEU:HD12	1.94	0.49
1:D:440:TYR:CD2	1:D:460:TYR:HB2	2.48	0.48
1:D:39:ILE:HD12	1:D:54:PRO:HB2	1.94	0.48
1:C:542:LEU:HG	1:C:556:TRP:CZ2	2.48	0.48
1:A:542:LEU:HD13	1:A:577:SER:HA	1.96	0.47
1:C:201:LEU:HD23	1:C:212:ILE:HG22	1.96	0.47
1:C:415:GLY:HA3	1:C:492:PHE:CE1	2.50	0.47
1:D:103:ALA:O	1:D:124:GLY:HA2	2.15	0.47
1:B:440:TYR:CD2	1:B:460:TYR:HB2	2.49	0.47
1:D:230:ASP:HB2	1:D:245:LEU:HG	1.97	0.46
1:D:281:LEU:HD13	1:D:325:LEU:HD23	1.98	0.46
1:C:491:ASN:O	1:C:495:ARG:HG2	2.16	0.46
1:D:124:GLY:O	1:D:141:THR:OG1	2.30	0.46
1:C:23:PRO:HG3	1:C:370:ILE:HG13	1.96	0.46
1:C:440:TYR:CD2	1:C:460:TYR:HB2	2.50	0.46
1:B:601:PRO:HB2	1:B:603:GLU:HG3	1.97	0.46
1:D:103:ALA:HB3	1:D:124:GLY:HA3	1.98	0.45
1:B:116:ARG:NH2	1:B:186:ASP:OD2	2.50	0.45
1:D:258:HIS:CG	1:D:259:PRO:HD2	2.51	0.45
1:A:542:LEU:HG	1:A:556:TRP:CZ2	2.52	0.45
1:C:425:LYS:NZ	3:C:813:HOH:O	2.48	0.45
1:D:23:PRO:HG3	1:D:370:ILE:HG13	1.98	0.45
1:C:513:GLY:HA2	1:C:536:ASP:O	2.17	0.45
1:D:433:HIS:HB3	1:D:440:TYR:CZ	2.52	0.45
1:C:212:ILE:HG13	1:C:229:TYR:HB2	1.98	0.45
1:A:440:TYR:CD2	1:A:460:TYR:HB2	2.53	0.44
1:B:415:GLY:HA3	1:B:492:PHE:CE1	2.53	0.44
1:D:485:ASP:OD1	1:D:485:ASP:N	2.51	0.44
1:B:23:PRO:HG3	1:B:370:ILE:HG13	2.00	0.44
1:D:125:VAL:HG23	1:D:139:PHE:CD1	2.53	0.44
1:C:542:LEU:HD13	1:C:577:SER:HA	2.00	0.43
1:D:191:TRP:NE1	1:D:234:PRO:HD3	2.33	0.43
1:A:24:GLN:HG3	1:A:83:PRO:O	2.19	0.43
1:D:513:GLY:HA2	1:D:536:ASP:O	2.18	0.43
1:D:419:LEU:HD21	1:D:500:VAL:HG11	2.01	0.43
1:A:273:LYS:H	1:A:273:LYS:HG2	1.65	0.43
1:C:572:LYS:NZ	3:C:805:HOH:O	2.42	0.43
1:C:485:ASP:OD1	1:C:485:ASP:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:319:TRP:CE2	1:D:325:LEU:HD13	2.54	0.42
1:C:136:PHE:HB3	1:C:180:LEU:HD11	2.01	0.42
1:A:78:ARG:HD3	1:A:633:ARG:O	2.19	0.42
1:A:323:ASP:O	1:A:341:HIS:HA	2.20	0.42
1:D:130:TRP:CH2	1:D:137:ILE:HD11	2.55	0.41
1:D:415:GLY:HA3	1:D:492:PHE:CE1	2.55	0.41
1:B:11:PRO:HG2	1:B:645:ASN:OD1	2.21	0.41
1:C:12:ASP:OD1	1:C:53:ARG:NH2	2.53	0.41
1:D:24:GLN:HG3	1:D:83:PRO:O	2.21	0.41
1:D:60:LEU:HD11	1:D:94:PHE:CG	2.56	0.41
1:A:436:PRO:HB3	1:A:515:TYR:CD2	2.56	0.41
1:D:491:ASN:O	1:D:495:ARG:HG2	2.21	0.41
1:A:419:LEU:HD21	1:A:500:VAL:HG11	2.03	0.41
1:B:196:ILE:HG21	1:B:215:SER:O	2.21	0.40
1:B:399:ARG:HD3	1:B:419:LEU:O	2.21	0.40
1:D:92:LEU:HD23	1:D:92:LEU:HA	1.94	0.40
1:A:415:GLY:HA3	1:A:492:PHE:CE1	2.57	0.40
1:D:125:VAL:HG22	1:D:126:SER:N	2.36	0.40
1:D:212:ILE:HG13	1:D:229:TYR:HB2	2.04	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	590/656 (90%)	578 (98%)	12 (2%)	0	100	100
1	B	594/656 (90%)	582 (98%)	12 (2%)	0	100	100
1	C	583/656 (89%)	572 (98%)	11 (2%)	0	100	100
1	D	578/656 (88%)	566 (98%)	12 (2%)	0	100	100
All	All	2345/2624 (89%)	2298 (98%)	47 (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	466/526 (89%)	465 (100%)	1 (0%)	87	94
1	B	463/526 (88%)	463 (100%)	0	100	100
1	C	454/526 (86%)	454 (100%)	0	100	100
1	D	454/526 (86%)	452 (100%)	2 (0%)	84	92
All	All	1837/2104 (87%)	1834 (100%)	3 (0%)	87	94

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

Mol	Chain	Res	Type
1	A	273	LYS
1	D	235	THR
1	D	614	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	286	ASN
1	A	431	ASN
1	B	223	GLN
1	B	275	ASN
1	B	431	ASN
1	B	616	GLN
1	B	628	ASN
1	C	129	GLN
1	C	286	ASN
1	C	289	HIS
1	C	628	ASN
1	D	90	GLN
1	D	286	ASN
1	D	442	HIS

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Mol	Chain	Res	Type
1	D	628	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ACT	C	701	-	3,3,3	1.37	0	3,3,3	1.46	0
2	ACT	B	701	-	3,3,3	1.35	0	3,3,3	1.51	0
2	ACT	A	701	-	3,3,3	1.43	1 (33%)	3,3,3	1.31	0
2	ACT	D	701	-	3,3,3	1.45	1 (33%)	3,3,3	1.45	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	701	ACT	CH3-C	2.09	1.57	1.49
2	A	701	ACT	CH3-C	2.05	1.57	1.49

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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	598/656 (91%)	-0.03	12 (2%)	65	66	23, 34, 59, 88	0
1	B	601/656 (91%)	0.04	24 (3%)	42	44	17, 33, 61, 88	1 (0%)
1	C	593/656 (90%)	0.28	35 (5%)	28	30	23, 41, 75, 96	0
1	D	590/656 (89%)	0.39	46 (7%)	19	21	23, 44, 77, 95	0
All	All	2382/2624 (90%)	0.17	117 (4%)	35	36	17, 37, 71, 96	1 (0%)

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	240	ALA	5.5
1	B	236	ALA	5.3
1	D	74	ALA	4.7
1	C	120	HIS	4.6
1	C	219	TRP	4.6
1	C	242	PRO	4.5
1	C	655	LEU	4.5
1	A	242	PRO	4.1
1	A	73	HIS	4.1
1	D	655	LEU	4.1
1	B	142	THR	3.9
1	C	98	ALA	3.9
1	A	553	LEU	3.9
1	D	219	TRP	3.9
1	D	113	GLY	3.8
1	C	74	ALA	3.8
1	D	102	LYS	3.8
1	C	142	THR	3.7
1	A	235	THR	3.7
1	C	224	TRP	3.6
1	D	272	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	113	GLY	3.5
1	D	176	ALA	3.3
1	D	235	THR	3.3
1	A	174	ARG	3.2
1	C	176	ALA	3.1
1	D	121	PHE	3.1
1	A	142	THR	3.1
1	B	224	TRP	3.0
1	D	141	THR	2.9
1	B	73	HIS	2.9
1	D	53	ARG	2.9
1	B	273	LYS	2.9
1	D	323	ASP	2.9
1	D	179	TRP	2.8
1	D	122	LYS	2.8
1	D	242	PRO	2.8
1	D	180	LEU	2.8
1	D	224	TRP	2.8
1	C	217	ASP	2.8
1	A	400	GLU	2.8
1	B	553	LEU	2.8
1	D	107	LEU	2.8
1	D	181	TYR	2.8
1	B	98	ALA	2.8
1	D	73	HIS	2.7
1	B	75	GLU	2.7
1	B	343	GLY	2.7
1	D	76	THR	2.7
1	C	7	PRO	2.7
1	D	71	LEU	2.7
1	B	100	GLU	2.6
1	B	241	ALA	2.6
1	B	344	GLY	2.6
1	C	107	LEU	2.6
1	A	100	GLU	2.6
1	D	344	GLY	2.6
1	D	245	LEU	2.6
1	B	196	ILE	2.6
1	D	97	SER	2.6
1	C	102	LYS	2.5
1	D	243	GLN	2.5
1	C	180	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	222	SER	2.5
1	B	120	HIS	2.5
1	C	220	GLN	2.5
1	B	272	GLY	2.4
1	D	39	ILE	2.4
1	C	234	PRO	2.4
1	D	125	VAL	2.4
1	C	191	TRP	2.4
1	B	175	PRO	2.3
1	D	218	GLU	2.3
1	B	6	THR	2.3
1	C	141	THR	2.3
1	C	222	SER	2.3
1	C	342	ILE	2.3
1	B	101	VAL	2.3
1	D	231	LEU	2.3
1	D	54	PRO	2.3
1	D	8	ALA	2.3
1	C	73	HIS	2.3
1	C	122	LYS	2.3
1	B	271	ALA	2.3
1	C	323	ASP	2.2
1	C	198	ILE	2.2
1	D	196	ILE	2.2
1	C	188	LEU	2.2
1	D	233	LEU	2.2
1	D	139	PHE	2.2
1	D	75	GLU	2.2
1	D	119	THR	2.2
1	C	179	TRP	2.2
1	D	247	ASP	2.2
1	C	184	GLU	2.2
1	C	181	TYR	2.2
1	A	187	LYS	2.2
1	B	39	ILE	2.1
1	C	212	ILE	2.1
1	C	272	GLY	2.1
1	B	235	THR	2.1
1	A	101	VAL	2.1
1	A	224	TRP	2.1
1	D	77	GLY	2.1
1	B	653	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	187	LYS	2.1
1	D	137	ILE	2.1
1	B	54	PRO	2.1
1	C	124	GLY	2.1
1	C	343	GLY	2.1
1	D	185	ALA	2.1
1	D	136	PHE	2.1
1	D	421	GLU	2.0
1	C	246	LEU	2.0
1	D	212	ILE	2.0
1	D	66	GLY	2.0
1	A	40	SER	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ACT	B	701	4/4	0.90	0.11	33,35,36,39	0
2	ACT	A	701	4/4	0.91	0.12	33,39,42,46	0
2	ACT	D	701	4/4	0.92	0.14	27,31,33,38	0
2	ACT	C	701	4/4	0.96	0.09	29,29,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.