



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:37 PM UTC

PDB ID : 6IKG / pdb_00006ikg
Title : Crystal structure of substrate-bound S9 peptidase (S514A mutant) from *Deinococcus radiodurans*
Authors : Yadav, P.; Kumar, A.; Goyal, V.D.; Makde, R.D.
Deposited on : 2018-10-16
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

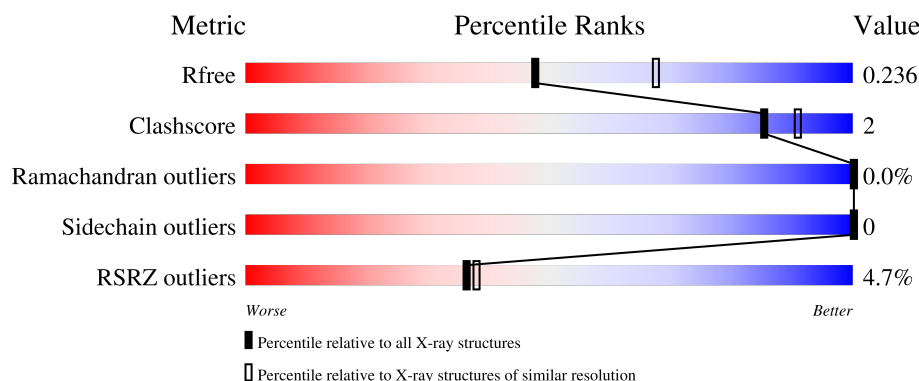
1 Overall quality at a glance

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	4% 93% 5%
1	B	656	2% 91% 6%
1	C	656	3% 93% 5%
1	D	656	9% 92% 6%
2	E	3	100% 100%

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Mol	Chain	Length	Quality of chain
2	F	3	100%
			100%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20572 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	644	Total	C	N	O	S	0	0	0
			4982	3156	898	919	9			
1	B	640	Total	C	N	O	S	0	0	0
			4945	3135	891	910	9			
1	C	641	Total	C	N	O	S	0	0	0
			4947	3138	887	913	9			
1	D	640	Total	C	N	O	S	0	0	0
			4878	3097	874	898	9			

There are 12 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
A	514	ALA	SER	engineered mutation	UNP Q9RXY9
B	0	GLY	-	expression tag	UNP Q9RXY9
B	1	SER	-	expression tag	UNP Q9RXY9
B	514	ALA	SER	engineered mutation	UNP Q9RXY9
C	0	GLY	-	expression tag	UNP Q9RXY9
C	1	SER	-	expression tag	UNP Q9RXY9
C	514	ALA	SER	engineered mutation	UNP Q9RXY9
D	0	GLY	-	expression tag	UNP Q9RXY9
D	1	SER	-	expression tag	UNP Q9RXY9
D	514	ALA	SER	engineered mutation	UNP Q9RXY9

- Molecule 2 is a protein called MET-ALA-ALA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	S	0	0	0
			19	11	3	4	1			
2	F	3	Total	C	N	O	S	0	0	0
			19	11	3	4	1			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		

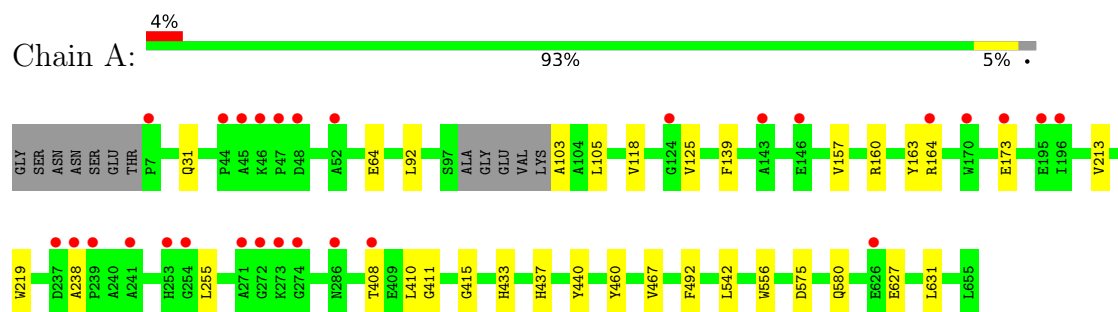
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	239	Total	O	0	0
			239	239		
4	B	224	Total	O	0	0
			224	224		
4	C	181	Total	O	0	0
			181	181		
4	D	114	Total	O	0	0
			114	114		

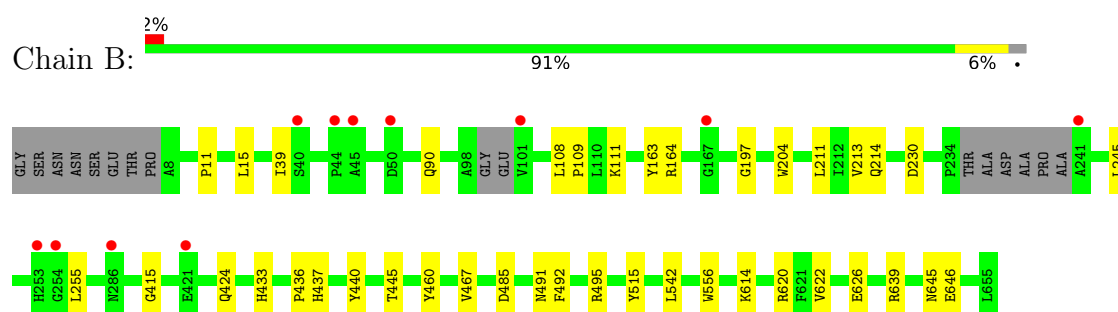
3 Residue-property plots [i](#)

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.

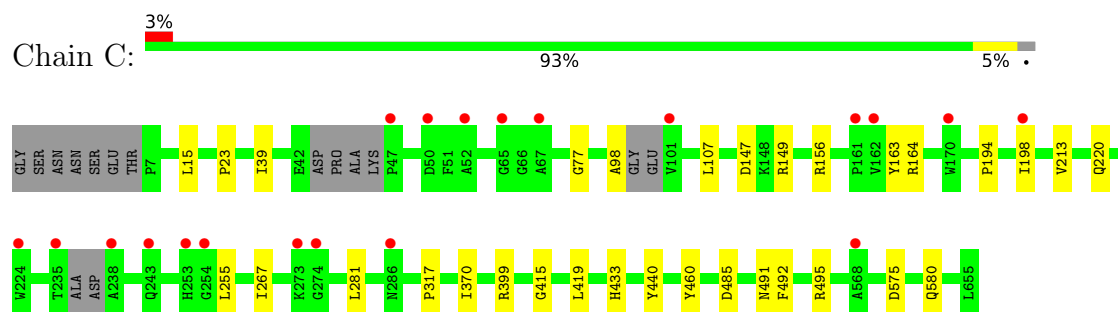
- Molecule 1: Acyl-peptide hydrolase, putative



- Molecule 1: Acyl-peptide hydrolase, putative

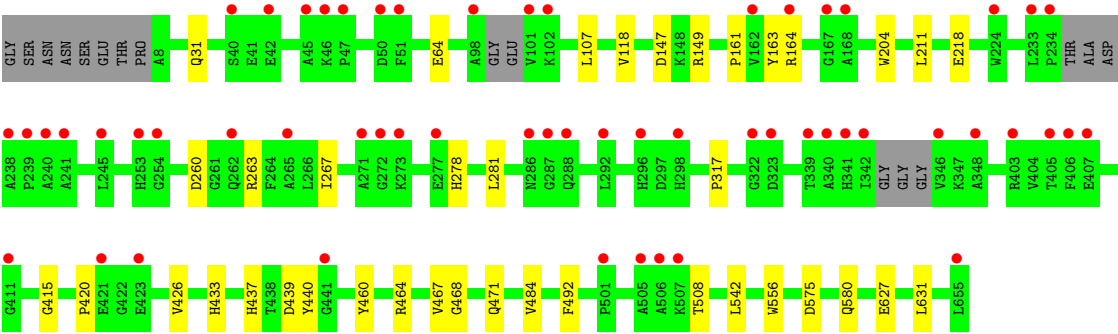


- Molecule 1: Acyl-peptide hydrolase, putative



- Molecule 1: Acyl-peptide hydrolase, putative





• Molecule 2: MET-ALA-ALA



• Molecule 2: MET-ALA-ALA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.65Å 214.64Å 92.02Å 90.00° 102.98° 90.00°	Depositor
Resolution (Å)	46.98 – 2.30 46.98 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.98-2.30) 99.9 (46.98-2.30)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.207 , 0.234 0.209 , 0.236	Depositor DCC
R_{free} test set	5981 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.367	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20572	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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: GOL

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	0.14	0/5135	0.39	0/7014
1	B	0.13	0/5095	0.37	0/6958
1	C	0.14	0/5098	0.39	1/6963 (0.0%)
1	D	0.15	0/5028	0.38	1/6880 (0.0%)
2	E	0.12	0/18	0.32	0/21
2	F	0.23	0/18	0.60	0/21
All	All	0.14	0/20392	0.38	2/27857 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	220	GLN	CA-CB-CG	-7.45	99.21	114.10
1	D	508	THR	OG1-CB-CG2	5.13	119.56	109.30

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	4982	0	4681	21	0
1	B	4945	0	4640	25	0
1	C	4947	0	4631	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	4878	0	4509	22	0
2	E	19	0	21	0	0
2	F	19	0	21	0	0
3	A	6	0	8	0	0
3	B	12	0	16	1	0
3	C	6	0	8	0	0
4	A	239	0	0	0	0
4	B	224	0	0	0	0
4	C	181	0	0	0	0
4	D	114	0	0	0	0
All	All	20572	0	18535	85	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 2.

All (85) 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:D:278:HIS:H	1:D:471:GLN:HE22	1.17	0.92
1:A:408:THR:HG22	1:A:410:LEU:H	1.38	0.88
1:A:157:VAL:HG22	1:B:622:VAL:HG22	1.70	0.74
1:C:399:ARG:NH1	1:C:419:LEU:O	2.25	0.68
1:C:267:ILE:HD11	1:C:317:PRO:HG2	1.77	0.66
1:D:278:HIS:H	1:D:471:GLN:NE2	1.92	0.63
1:D:267:ILE:HD11	1:D:317:PRO:HG2	1.81	0.63
1:D:278:HIS:N	1:D:471:GLN:HE22	1.97	0.56
1:D:440:TYR:CD2	1:D:460:TYR:HB2	2.41	0.55
1:B:445:THR:HB	3:B:702:GOL:H11	1.88	0.55
1:D:433:HIS:HB3	1:D:440:TYR:CZ	2.41	0.55
1:C:163:TYR:CD1	1:C:164:ARG:HG3	2.41	0.55
1:D:204:TRP:CZ2	1:D:211:LEU:HD21	2.42	0.55
1:D:420:PRO:HG2	1:D:426:VAL:HG11	1.91	0.53
1:C:433:HIS:HB3	1:C:440:TYR:CZ	2.45	0.52
1:B:230:ASP:HB2	1:B:245:LEU:HD12	1.92	0.51
1:D:163:TYR:CD1	1:D:164:ARG:HG3	2.45	0.51
1:D:260:ASP:OD2	1:D:263:ARG:HD3	2.10	0.51
1:A:163:TYR:CD1	1:A:164:ARG:HG3	2.46	0.51
1:B:163:TYR:CD1	1:B:164:ARG:HG3	2.46	0.50
1:A:125:VAL:HG23	1:A:139:PHE:CD2	2.46	0.50
1:D:107:LEU:HD13	1:D:118:VAL:HG21	1.93	0.49
1:C:267:ILE:HG13	1:C:281:LEU:HD23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:440:TYR:CD2	1:A:460:TYR:HB2	2.49	0.48
1:B:197:GLY:O	1:B:214:GLN:HG2	2.13	0.48
1:D:147:ASP:OD1	1:D:149:ARG:HD3	2.13	0.47
1:B:542:LEU:HG	1:B:556:TRP:CZ2	2.50	0.47
1:B:90:GLN:OE1	1:B:111:LYS:NZ	2.48	0.47
1:C:440:TYR:CD2	1:C:460:TYR:HB2	2.49	0.47
1:B:433:HIS:HB3	1:B:440:TYR:CZ	2.49	0.47
1:C:23:PRO:HG3	1:C:370:ILE:HG13	1.97	0.47
1:D:464:ARG:CZ	1:D:484:VAL:HB	2.46	0.46
1:B:415:GLY:HA3	1:B:492:PHE:CE1	2.51	0.46
1:C:194:PRO:HG3	1:C:198:ILE:HG12	1.97	0.46
1:B:424:GLN:H	1:B:424:GLN:CD	2.21	0.46
1:B:204:TRP:CZ2	1:B:211:LEU:HD21	2.51	0.45
1:C:15:LEU:HD22	1:C:39:ILE:HG21	1.98	0.45
1:A:433:HIS:HB3	1:A:440:TYR:CZ	2.52	0.45
1:C:491:ASN:O	1:C:495:ARG:HG2	2.17	0.45
1:D:161:PRO:HD2	1:D:218:GLU:HB3	1.99	0.45
1:B:15:LEU:HD22	1:B:39:ILE:HG21	1.99	0.45
1:A:408:THR:HB	1:A:411:GLY:O	2.17	0.45
1:B:440:TYR:CD2	1:B:460:TYR:HB2	2.52	0.44
1:A:433:HIS:HB3	1:A:440:TYR:CE2	2.53	0.44
1:A:103:ALA:O	1:A:125:VAL:HG12	2.17	0.43
1:A:105:LEU:O	1:A:118:VAL:HG22	2.18	0.43
1:B:108:LEU:HD12	1:B:109:PRO:HD2	1.99	0.43
1:B:626:GLU:OE1	1:B:639:ARG:NH2	2.50	0.43
1:D:542:LEU:HG	1:D:556:TRP:CZ2	2.54	0.43
1:A:213:VAL:HG23	1:A:255:LEU:HD11	2.00	0.43
1:A:31:GLN:HG2	1:A:64:GLU:HG3	2.01	0.42
1:B:213:VAL:HG23	1:B:255:LEU:HD11	2.00	0.42
1:C:149:ARG:NH2	1:C:156:ARG:HG2	2.34	0.42
1:D:415:GLY:HA3	1:D:492:PHE:CE1	2.54	0.42
1:A:157:VAL:HG11	1:B:620:ARG:CZ	2.50	0.42
1:A:415:GLY:HA3	1:A:492:PHE:CE2	2.54	0.42
1:A:542:LEU:HG	1:A:556:TRP:CZ2	2.55	0.42
1:D:267:ILE:HG13	1:D:281:LEU:HD23	2.01	0.42
1:D:439:ASP:OD2	1:D:468:GLY:HA3	2.19	0.42
1:D:575:ASP:HA	1:D:580:GLN:HE22	1.85	0.42
1:D:627:GLU:HG3	1:D:631:LEU:HD13	2.01	0.42
1:A:627:GLU:HG3	1:A:631:LEU:HD13	2.01	0.42
1:B:491:ASN:O	1:B:495:ARG:HG2	2.19	0.42
1:B:622:VAL:HG21	1:B:646:GLU:HG3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:415:GLY:HA3	1:C:492:PHE:CE1	2.55	0.42
1:C:213:VAL:HG23	1:C:255:LEU:HD11	2.01	0.42
1:C:485:ASP:OD1	1:C:485:ASP:N	2.52	0.42
1:C:575:ASP:HA	1:C:580:GLN:HE22	1.84	0.42
1:A:160:ARG:NH2	1:A:173:GLU:O	2.46	0.41
1:B:436:PRO:HB3	1:B:515:TYR:CE2	2.55	0.41
1:C:440:TYR:CE2	1:C:460:TYR:HB2	2.56	0.41
1:D:31:GLN:HG2	1:D:64:GLU:HG2	2.00	0.41
1:B:437:HIS:C	1:B:467:VAL:HB	2.45	0.41
1:B:11:PRO:HG2	1:B:645:ASN:OD1	2.19	0.41
1:B:485:ASP:OD1	1:B:485:ASP:N	2.52	0.41
1:A:219:TRP:CH2	1:B:614:LYS:HA	2.56	0.41
1:C:107:LEU:HD12	1:C:107:LEU:HA	1.87	0.41
1:D:437:HIS:C	1:D:467:VAL:HB	2.45	0.41
1:B:436:PRO:HB3	1:B:515:TYR:CD2	2.56	0.40
1:C:147:ASP:OD1	1:C:149:ARG:HD3	2.20	0.40
1:A:92:LEU:HD23	1:A:92:LEU:HA	1.98	0.40
1:C:77:GLY:HA3	1:C:98:ALA:HA	2.03	0.40
1:C:491:ASN:OD1	1:C:495:ARG:NH1	2.53	0.40
1:A:575:ASP:HA	1:A:580:GLN:HE22	1.86	0.40
1:A:437:HIS:C	1:A:467:VAL:HB	2.46	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	640/656 (98%)	617 (96%)	22 (3%)	1 (0%)	43	55
1	B	634/656 (97%)	616 (97%)	18 (3%)	0	100	100
1	C	633/656 (96%)	615 (97%)	18 (3%)	0	100	100
1	D	632/656 (96%)	614 (97%)	18 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	1/3 (33%)	1 (100%)	0	0	100	100
2	F	1/3 (33%)	1 (100%)	0	0	100	100
All	All	2541/2630 (97%)	2464 (97%)	76 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	238	ALA

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	497/525 (95%)	497 (100%)	0	100	100
1	B	492/525 (94%)	492 (100%)	0	100	100
1	C	492/525 (94%)	492 (100%)	0	100	100
1	D	475/525 (90%)	475 (100%)	0	100	100
2	E	1/1 (100%)	1 (100%)	0	100	100
2	F	1/1 (100%)	1 (100%)	0	100	100
All	All	1958/2102 (93%)	1958 (100%)	0	100	100

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

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

Mol	Chain	Res	Type
1	A	31	GLN
1	B	73	HIS
1	B	91	ASN
1	B	120	HIS
1	B	289	HIS
1	B	431	ASN

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Mol	Chain	Res	Type
1	C	253	HIS
1	C	275	ASN
1	C	341	HIS
1	D	90	GLN
1	D	91	ASN
1	D	249	ASN
1	D	341	HIS
1	D	471	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 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	$\# Z > 2$
3	GOL	B	701	-	5,5,5	0.41	0	5,5,5	0.29	0
3	GOL	A	701	-	5,5,5	0.40	0	5,5,5	0.34	0
3	GOL	B	702	-	5,5,5	0.36	0	5,5,5	0.22	0
3	GOL	C	701	-	5,5,5	0.40	0	5,5,5	0.29	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	701	-	-	0/4/4/4	-
3	GOL	A	701	-	-	0/4/4/4	-
3	GOL	B	702	-	-	2/4/4/4	-
3	GOL	C	701	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	702	GOL	O1-C1-C2-C3
3	B	702	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	702	GOL	1	0

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	644/656 (98%)	0.12	28 (4%)	40	41	16, 27, 45, 66	0
1	B	640/656 (97%)	0.13	11 (1%)	69	70	17, 28, 45, 68	0
1	C	641/656 (97%)	0.26	20 (3%)	51	53	16, 31, 50, 64	0
1	D	640/656 (97%)	0.79	57 (8%)	15	17	24, 40, 58, 75	0
2	E	3/3 (100%)	4.58	3 (100%)	0	0	52, 52, 58, 68	0
2	F	3/3 (100%)	5.44	3 (100%)	0	0	59, 59, 62, 64	0
All	All	2571/2630 (97%)	0.34	122 (4%)	36	38	16, 32, 52, 75	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	2	ALA	6.7
2	F	3	ALA	6.1
1	D	238	ALA	5.6
1	A	237	ASP	5.4
1	D	253	HIS	5.1
2	E	2	ALA	5.1
1	A	254	GLY	5.0
1	D	342	ILE	4.7
1	D	346	VAL	4.7
2	E	1	MET	4.4
2	E	3	ALA	4.3
1	A	253	HIS	4.3
1	A	271	ALA	4.3
1	D	272	GLY	4.1
1	A	273	LYS	4.0
1	D	45	ALA	4.0
1	B	241	ALA	4.0
1	D	98	ALA	3.9
2	F	1	MET	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	348	ALA	3.6
1	D	239	PRO	3.5
1	D	505	ALA	3.5
1	B	50	ASP	3.5
1	D	286	ASN	3.5
1	D	50	ASP	3.4
1	D	655	LEU	3.3
1	C	273	LYS	3.3
1	B	254	GLY	3.3
1	C	50	ASP	3.3
1	A	7	PRO	3.3
1	A	143	ALA	3.2
1	D	501	PRO	3.1
1	A	274	GLY	3.1
1	D	421	GLU	3.1
1	B	286	ASN	3.0
1	D	101	VAL	3.0
1	C	238	ALA	2.9
1	D	46	LYS	2.9
1	A	44	PRO	2.9
1	D	506	ALA	2.9
1	D	403	ARG	2.9
1	A	45	ALA	2.9
1	A	241	ALA	2.8
1	C	161	PRO	2.8
1	C	274	GLY	2.8
1	C	253	HIS	2.8
1	D	241	ALA	2.8
1	C	254	GLY	2.7
1	D	287	GLY	2.7
1	A	239	PRO	2.7
1	A	238	ALA	2.7
1	B	45	ALA	2.7
1	D	254	GLY	2.7
1	D	341	HIS	2.6
1	B	167	GLY	2.6
1	C	65	GLY	2.6
1	D	262	GLN	2.6
1	D	507	LYS	2.6
1	D	233	LEU	2.6
1	D	273	LYS	2.6
1	A	52	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	626	GLU	2.5
1	A	170	TRP	2.5
1	D	406	PHE	2.5
1	D	288	GLN	2.5
1	D	224	TRP	2.5
1	C	286	ASN	2.5
1	D	168	ALA	2.5
1	D	47	PRO	2.5
1	D	322	GLY	2.5
1	D	245	LEU	2.4
1	A	146	GLU	2.4
1	C	47	PRO	2.4
1	A	408	THR	2.4
1	C	235	THR	2.4
1	D	296	HIS	2.4
1	A	286	ASN	2.4
1	D	277	GLU	2.4
1	D	423	GLU	2.4
1	C	101	VAL	2.3
1	C	162	VAL	2.3
1	A	47	PRO	2.3
1	C	224	TRP	2.3
1	D	40	SER	2.3
1	D	240	ALA	2.3
1	A	173	GLU	2.3
1	A	195	GLU	2.3
1	B	421	GLU	2.3
1	B	253	HIS	2.2
1	C	52	ALA	2.2
1	C	67	ALA	2.2
1	B	40	SER	2.2
1	B	44	PRO	2.2
1	D	167	GLY	2.2
1	D	42	GLU	2.2
1	D	405	THR	2.2
1	D	292	LEU	2.2
1	D	323	ASP	2.2
1	D	234	PRO	2.2
1	D	441	GLY	2.2
1	C	568	ALA	2.2
1	C	170	TRP	2.1
1	D	51	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	339	THR	2.1
1	D	340	ALA	2.1
1	D	407	GLU	2.1
1	D	162	VAL	2.1
1	D	411	GLY	2.1
1	D	265	ALA	2.1
1	D	298	HIS	2.1
1	A	196	ILE	2.1
1	A	124	GLY	2.1
1	A	272	GLY	2.1
1	D	271	ALA	2.0
1	C	243	GLN	2.0
1	B	101	VAL	2.0
1	D	164	ARG	2.0
1	A	46	LYS	2.0
1	A	48	ASP	2.0
1	A	164	ARG	2.0
1	C	198	ILE	2.0
1	D	102	LYS	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
3	GOL	C	701	6/6	0.88	0.12	45,45,45,46	0
3	GOL	B	701	6/6	0.90	0.12	43,47,47,48	0
3	GOL	B	702	6/6	0.93	0.13	35,35,36,36	0
3	GOL	A	701	6/6	0.94	0.08	26,28,30,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.