



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:15 AM UTC

PDB ID : 6IGP / pdb_00006igp
Title : Crystal structure of S9 peptidase (inactive state) from *Deinococcus radiodurans* R1 in P212121
Authors : Yadav, P.; Goyal, V.D.; Kumar, A.; Makde, R.D.
Deposited on : 2018-09-25
Resolution : 2.40 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

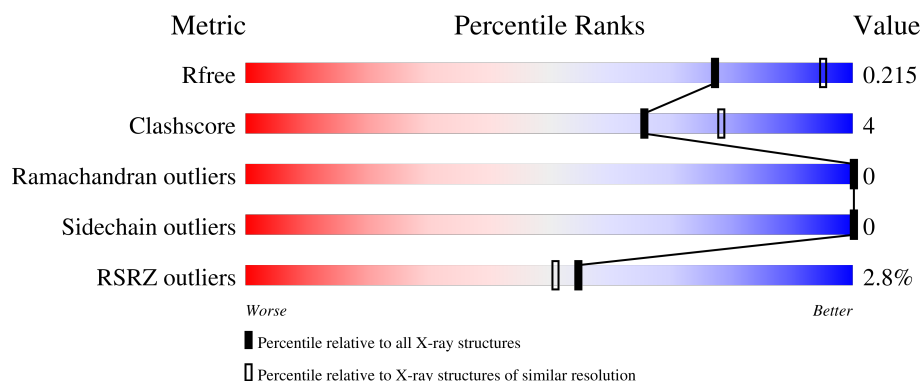
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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	0.2% 85% 5% 9%
1	B	656	2% 81% 8% 11%
1	C	656	3% 81% 8% 11%
1	D	656	4% 80% 8% 12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 18910 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	594	Total	C	N	O	S	0	0	0
			4612	2929	830	844	9			
1	B	587	Total	C	N	O	S	0	0	0
			4510	2870	808	823	9			
1	C	582	Total	C	N	O	S	0	0	0
			4423	2823	780	811	9			
1	D	579	Total	C	N	O	S	0	0	0
			4381	2792	777	803	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 GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

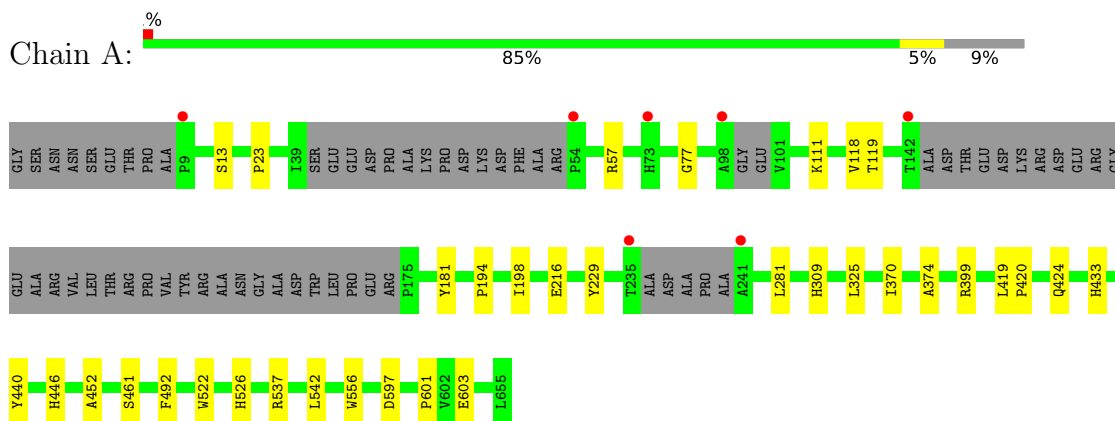
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	274	Total	O	0	0
			274	274		
3	B	279	Total	O	0	0
			279	279		
3	C	216	Total	O	0	0
			216	216		
3	D	191	Total	O	0	0
			191	191		

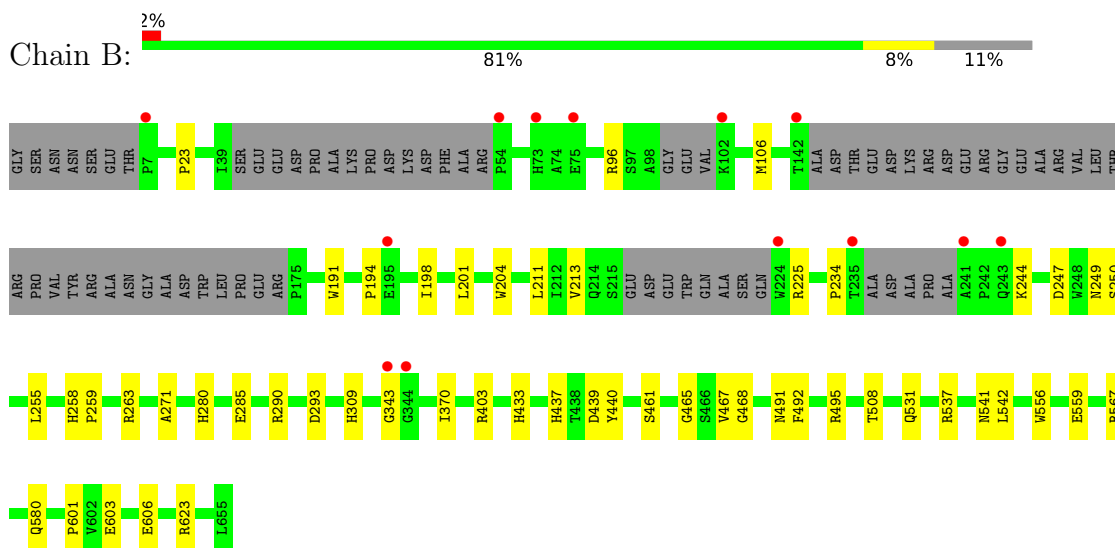
3 Residue-property plots

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.

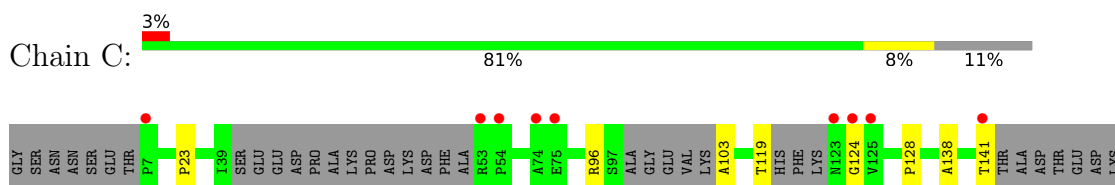
• Molecule 1: Acyl-peptide hydrolase, putative

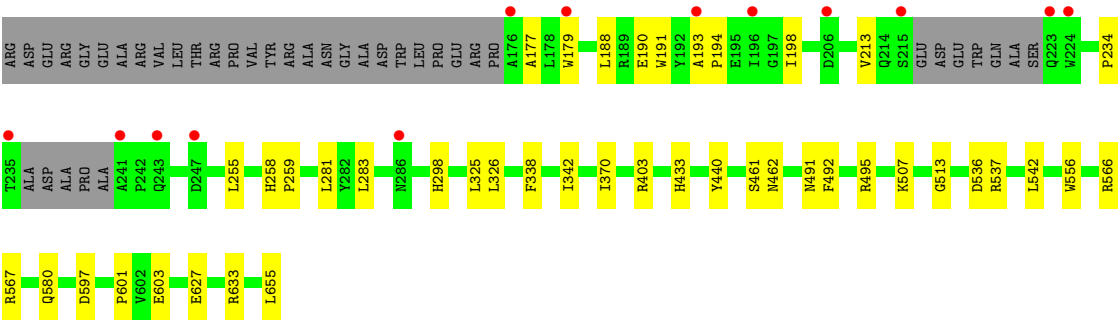


• Molecule 1: Acyl-peptide hydrolase, putative

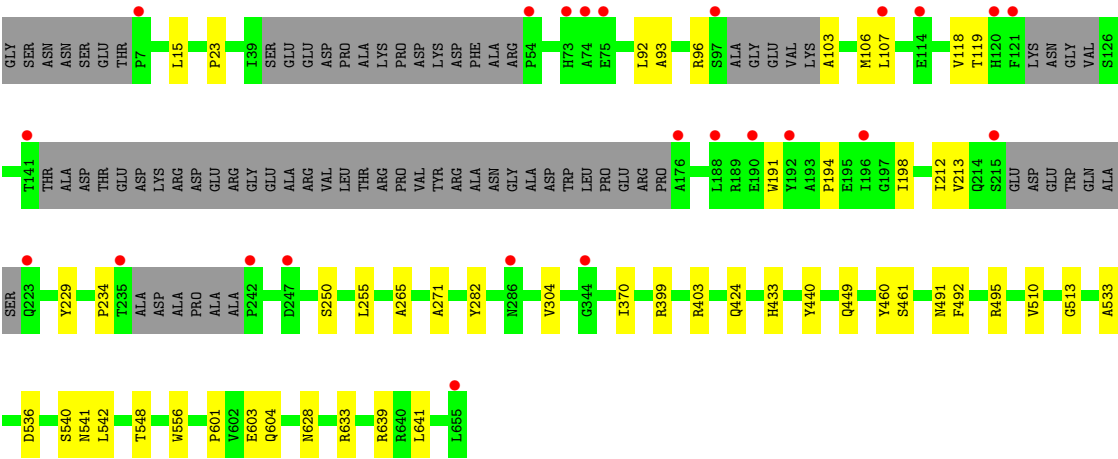
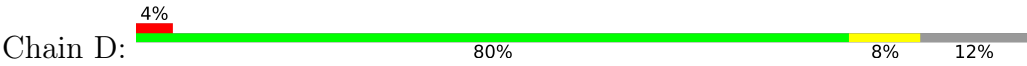


• Molecule 1: Acyl-peptide hydrolase, putative





● Molecule 1: Acyl-peptide hydrolase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	119.86Å 130.49Å 194.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.74 – 2.40 48.74 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.74-2.40) 100.0 (48.74-2.40)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.185 , 0.217 0.185 , 0.215	Depositor DCC
R_{free} test set	5991 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.6	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.1	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	0.95	EDS
Total number of atoms	18910	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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 44.73 % 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 1.4625e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.32	0/4751	0.53	0/6479
1	B	0.32	0/4646	0.53	0/6341
1	C	0.28	0/4558	0.50	0/6228
1	D	0.28	0/4512	0.51	1/6162 (0.0%)
All	All	0.30	0/18467	0.52	1/25210 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	304	VAL	CG1-CB-CG2	-5.55	98.58	110.80

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	4612	0	4361	24	0
1	B	4510	0	4240	38	0
1	C	4423	0	4106	41	0
1	D	4381	0	4066	36	0
2	A	6	0	8	1	0
2	B	6	0	8	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	12	0	16	2	0
3	A	274	0	0	3	0
3	B	279	0	0	1	0
3	C	216	0	0	2	0
3	D	191	0	0	2	0
All	All	18910	0	16805	133	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 4.

All (133) 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:298:HIS:HE1	2:C:702:GOL:H12	1.34	0.88
1:B:508:THR:H	1:B:531:GLN:HE21	1.22	0.88
1:B:580:GLN:HE21	1:C:567:ARG:HE	1.22	0.86
1:D:399:ARG:H	1:D:449:GLN:HE22	1.22	0.85
1:D:541:ASN:H	1:D:604:GLN:HE22	1.20	0.85
1:B:96:ARG:HD2	1:B:106:MET:HE3	1.63	0.79
1:B:567:ARG:HE	1:C:580:GLN:HE21	1.34	0.74
1:A:522:TRP:O	1:A:526:HIS:HD2	1.72	0.73
1:D:540:SER:H	1:D:604:GLN:HE21	1.36	0.72
1:B:194:PRO:HG3	1:B:198:ILE:HG13	1.73	0.70
1:C:298:HIS:CE1	2:C:702:GOL:H12	2.23	0.69
1:C:194:PRO:HG3	1:C:198:ILE:HG13	1.74	0.69
1:D:540:SER:H	1:D:604:GLN:NE2	1.91	0.69
1:B:508:THR:H	1:B:531:GLN:NE2	1.92	0.67
1:C:566:ARG:HH11	1:C:566:ARG:HG2	1.61	0.66
1:B:567:ARG:NE	1:C:580:GLN:HE21	1.93	0.66
1:D:118:VAL:HG23	1:D:119:THR:HG23	1.78	0.65
1:A:216:GLU:HG3	1:A:229:TYR:OH	2.01	0.61
1:B:437:HIS:HE1	1:B:559:GLU:OE1	1.85	0.60
1:B:204:TRP:CZ2	1:B:211:LEU:HD21	2.36	0.60
1:B:606:GLU:OE2	1:B:623:ARG:NH2	2.33	0.60
1:B:280:HIS:HD2	1:B:293:ASP:OD2	1.84	0.60
1:C:403:ARG:NH2	3:C:801:HOH:O	2.29	0.59
1:D:194:PRO:HG3	1:D:198:ILE:HG13	1.84	0.59
1:D:23:PRO:HG3	1:D:370:ILE:HG13	1.85	0.59
1:D:191:TRP:NE1	1:D:234:PRO:HD3	2.22	0.55
1:D:541:ASN:H	1:D:604:GLN:NE2	1.98	0.54
1:B:437:HIS:HD2	1:B:465:GLY:O	1.91	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:THR:HG22	1:C:188:LEU:HD11	1.90	0.53
1:D:250:SER:HB2	1:D:271:ALA:HA	1.90	0.52
1:C:191:TRP:NE1	1:C:234:PRO:HD3	2.25	0.52
1:B:541:ASN:HD22	1:B:541:ASN:C	2.16	0.52
1:C:23:PRO:HG3	1:C:370:ILE:HG13	1.91	0.52
1:C:177:ALA:HB3	1:C:179:TRP:HE1	1.74	0.52
1:B:437:HIS:CE1	1:B:559:GLU:OE1	2.64	0.51
1:A:420:PRO:HB3	1:A:452:ALA:O	2.10	0.51
1:C:513:GLY:HA2	1:C:536:ASP:O	2.11	0.50
1:A:57:ARG:NE	1:A:77:GLY:O	2.44	0.50
1:B:250:SER:HB2	1:B:271:ALA:HA	1.93	0.50
1:C:627:GLU:HB2	1:C:633:ARG:NH2	2.27	0.49
1:D:15:LEU:HD21	1:D:641:LEU:HD11	1.95	0.49
1:D:628:ASN:O	1:D:633:ARG:NH2	2.43	0.49
1:B:433:HIS:HB3	1:B:440:TYR:CZ	2.48	0.49
1:A:433:HIS:HB3	1:A:440:TYR:CZ	2.48	0.48
1:B:309:HIS:HE1	3:B:801:HOH:O	1.94	0.48
1:D:107:LEU:HD13	1:D:118:VAL:HG13	1.96	0.48
1:D:403:ARG:NH2	3:D:705:HOH:O	2.40	0.48
1:D:399:ARG:H	1:D:449:GLN:NE2	2.02	0.48
1:D:461:SER:HB3	1:D:492:PHE:CD2	2.49	0.48
1:B:461:SER:HB3	1:B:492:PHE:CD2	2.49	0.48
1:C:141:THR:HG22	1:C:179:TRP:CD1	2.49	0.48
1:B:461:SER:HB3	1:B:492:PHE:CE2	2.48	0.48
1:A:461:SER:HB3	1:A:492:PHE:CE2	2.49	0.47
1:B:201:LEU:HA	1:B:211:LEU:O	2.15	0.47
1:C:281:LEU:HD13	1:C:325:LEU:HD23	1.95	0.47
1:C:433:HIS:HB3	1:C:440:TYR:CZ	2.50	0.47
1:C:601:PRO:HB2	1:C:603:GLU:HG3	1.96	0.47
1:A:281:LEU:HD13	1:A:325:LEU:HD23	1.97	0.47
1:D:107:LEU:N	1:D:107:LEU:HD12	2.30	0.47
1:A:526:HIS:HE1	3:A:1027:HOH:O	1.97	0.47
1:C:507:LYS:NZ	1:C:655:LEU:HD22	2.30	0.46
1:C:461:SER:HB3	1:C:492:PHE:CE2	2.51	0.46
1:C:542:LEU:HG	1:C:556:TRP:CZ2	2.50	0.46
1:B:263:ARG:HG2	1:B:285:GLU:HB2	1.96	0.46
1:B:491:ASN:O	1:B:495:ARG:HG2	2.16	0.46
1:D:440:TYR:CD2	1:D:460:TYR:HB2	2.50	0.46
1:A:601:PRO:HB2	1:A:603:GLU:HG3	1.98	0.46
1:C:298:HIS:HD2	3:C:803:HOH:O	1.98	0.46
1:D:213:VAL:HG23	1:D:255:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:542:LEU:HG	1:B:556:TRP:CZ2	2.51	0.46
1:B:580:GLN:HE21	1:C:567:ARG:NE	2.02	0.46
1:D:96:ARG:O	1:D:103:ALA:HA	2.16	0.46
1:B:244:LYS:HD3	1:B:247:ASP:OD1	2.16	0.46
1:A:194:PRO:HG3	1:A:198:ILE:HG13	1.98	0.45
1:A:537:ARG:NE	1:A:597:ASP:OD2	2.45	0.45
1:C:566:ARG:HG2	1:C:566:ARG:NH1	2.31	0.45
1:D:542:LEU:HG	1:D:556:TRP:CZ2	2.52	0.45
1:B:403:ARG:HH11	1:B:403:ARG:HD3	1.64	0.45
1:D:433:HIS:HB3	1:D:440:TYR:CZ	2.51	0.45
1:D:424:GLN:OE1	1:D:424:GLN:N	2.27	0.45
1:B:580:GLN:HG3	1:C:567:ARG:HG3	1.99	0.44
1:B:258:HIS:CG	1:B:259:PRO:HD2	2.52	0.44
1:C:128:PRO:HA	1:C:138:ALA:O	2.18	0.44
1:C:179:TRP:CZ2	1:C:190:GLU:HB2	2.53	0.44
1:D:107:LEU:HD13	1:D:118:VAL:CG1	2.47	0.44
1:A:111:LYS:HD3	1:A:111:LYS:HA	1.58	0.44
1:C:213:VAL:HG23	1:C:255:LEU:HD11	1.99	0.44
1:C:433:HIS:HB3	1:C:440:TYR:CE2	2.52	0.44
1:B:437:HIS:C	1:B:467:VAL:HB	2.42	0.44
1:A:424:GLN:OE1	1:A:424:GLN:N	2.38	0.44
1:A:118:VAL:HG21	1:A:181:TYR:CZ	2.53	0.44
1:D:513:GLY:HA2	1:D:536:ASP:O	2.17	0.44
1:C:326:LEU:HD12	1:C:338:PHE:O	2.17	0.43
1:D:106:MET:C	1:D:107:LEU:HD12	2.43	0.43
1:C:440:TYR:N	1:C:462:ASN:OD1	2.41	0.43
1:C:491:ASN:OD1	1:C:495:ARG:NH1	2.51	0.43
1:B:213:VAL:HG23	1:B:255:LEU:HD11	2.00	0.43
1:C:258:HIS:CG	1:C:259:PRO:HD2	2.54	0.43
1:A:399:ARG:HD3	1:A:419:LEU:O	2.19	0.43
1:D:548:THR:OG1	1:D:601:PRO:HG2	2.19	0.43
1:B:601:PRO:HB2	1:B:603:GLU:HG3	2.01	0.43
1:A:309:HIS:HE1	3:A:839:HOH:O	2.02	0.43
1:B:290:ARG:CZ	1:B:343:GLY:HA2	2.49	0.43
1:B:567:ARG:HE	1:C:580:GLN:NE2	2.10	0.43
1:B:225:ARG:HG2	1:B:249:ASN:O	2.19	0.42
1:B:439:ASP:OD2	1:B:468:GLY:HA3	2.19	0.42
1:C:103:ALA:O	1:C:124:GLY:HA2	2.19	0.42
1:C:491:ASN:O	1:C:495:ARG:HG2	2.19	0.42
1:C:537:ARG:NE	1:C:597:ASP:OD2	2.44	0.42
1:D:93:ALA:HA	1:D:106:MET:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:446:HIS:HD2	3:A:957:HOH:O	2.02	0.42
1:B:537:ARG:HH22	2:B:701:GOL:H31	1.85	0.42
1:D:461:SER:HB3	1:D:492:PHE:CE2	2.54	0.42
1:C:190:GLU:OE2	1:C:193:ALA:HB2	2.19	0.42
1:C:179:TRP:CE2	1:C:190:GLU:HB2	2.53	0.41
1:D:212:ILE:HG13	1:D:229:TYR:HB2	2.02	0.41
1:B:23:PRO:HG3	1:B:370:ILE:HG13	2.03	0.41
1:D:601:PRO:HB2	1:D:603:GLU:HG3	2.02	0.41
1:D:92:LEU:HD12	1:D:92:LEU:HA	1.94	0.41
1:A:374:ALA:HB1	2:A:701:GOL:H2	2.02	0.41
1:A:461:SER:HB3	1:A:492:PHE:CD2	2.56	0.41
1:C:283:LEU:HD22	1:C:342:ILE:HG12	2.03	0.41
1:D:510:VAL:O	1:D:533:ALA:HA	2.20	0.41
1:D:639:ARG:NH2	3:D:703:HOH:O	2.39	0.41
1:A:13:SER:OG	1:A:446:HIS:HE1	2.03	0.41
1:A:118:VAL:HG13	1:A:119:THR:HG23	2.03	0.41
1:B:191:TRP:NE1	1:B:234:PRO:HD3	2.36	0.41
1:A:542:LEU:HG	1:A:556:TRP:CZ2	2.56	0.40
1:C:96:ARG:O	1:C:103:ALA:HA	2.22	0.40
1:A:522:TRP:O	1:A:526:HIS:CD2	2.62	0.40
1:D:491:ASN:O	1:D:495:ARG:HG2	2.21	0.40
1:A:23:PRO:HG3	1:A:370:ILE:HG13	2.03	0.40
1:D:265:ALA:HA	1:D:282:TYR:O	2.22	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	584/656 (89%)	569 (97%)	15 (3%)	0	100	100
1	B	575/656 (88%)	562 (98%)	13 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	568/656 (87%)	557 (98%)	11 (2%)	0	100	100
1	D	565/656 (86%)	556 (98%)	9 (2%)	0	100	100
All	All	2292/2624 (87%)	2244 (98%)	48 (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%)	466 (100%)	0	100	100
1	B	450/526 (86%)	450 (100%)	0	100	100
1	C	433/526 (82%)	433 (100%)	0	100	100
1	D	428/526 (81%)	428 (100%)	0	100	100
All	All	1777/2104 (84%)	1777 (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 (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	38	GLN
1	A	129	GLN
1	A	214	GLN
1	A	253	HIS
1	A	298	HIS
1	A	365	HIS
1	A	431	ASN
1	A	446	HIS
1	A	526	HIS
1	B	24	GLN
1	B	120	HIS

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Mol	Chain	Res	Type
1	B	226	GLN
1	B	280	HIS
1	B	354	GLN
1	B	437	HIS
1	B	446	HIS
1	B	531	GLN
1	B	541	ASN
1	B	580	GLN
1	B	598	HIS
1	C	24	GLN
1	C	298	HIS
1	C	385	GLN
1	C	580	GLN
1	D	38	GLN
1	D	91	ASN
1	D	262	GLN
1	D	354	GLN
1	D	402	GLN
1	D	446	HIS
1	D	449	GLN
1	D	604	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	$\# Z > 2$
2	GOL	C	702	-	5,5,5	0.34	0	5,5,5	0.54	0
2	GOL	B	701	-	5,5,5	0.41	0	5,5,5	0.20	0
2	GOL	A	701	-	5,5,5	0.34	0	5,5,5	0.51	0
2	GOL	C	701	-	5,5,5	0.34	0	5,5,5	0.41	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
2	GOL	C	702	-	-	0/4/4/4	-
2	GOL	B	701	-	-	4/4/4/4	-
2	GOL	A	701	-	-	3/4/4/4	-
2	GOL	C	701	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	701	GOL	O1-C1-C2-O2
2	B	701	GOL	O1-C1-C2-C3
2	B	701	GOL	C1-C2-C3-O3
2	B	701	GOL	O2-C2-C3-O3
2	A	701	GOL	C1-C2-C3-O3
2	A	701	GOL	O1-C1-C2-C3
2	C	701	GOL	C1-C2-C3-O3
2	A	701	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	702	GOL	2	0
2	B	701	GOL	1	0
2	A	701	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	594/656 (90%)	-0.43	7 (1%) 76 73	15, 27, 49, 64	0
1	B	587/656 (89%)	-0.35	13 (2%) 62 58	14, 26, 54, 73	0
1	C	582/656 (88%)	-0.18	22 (3%) 44 40	17, 32, 63, 76	0
1	D	579/656 (88%)	-0.07	24 (4%) 41 37	16, 38, 66, 81	0
All	All	2342/2624 (89%)	-0.26	66 (2%) 55 51	14, 30, 60, 81	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	241	ALA	4.9
1	D	54	PRO	4.8
1	C	241	ALA	4.4
1	B	142	THR	4.3
1	D	121	PHE	4.2
1	C	224	TRP	4.0
1	D	7	PRO	3.9
1	D	235	THR	3.9
1	C	123	ASN	3.8
1	D	141	THR	3.8
1	C	223	GLN	3.6
1	C	75	GLU	3.6
1	C	235	THR	3.6
1	D	75	GLU	3.6
1	D	120	HIS	3.4
1	D	74	ALA	3.3
1	C	74	ALA	3.2
1	C	176	ALA	3.2
1	D	223	GLN	3.2
1	D	215	SER	3.2
1	A	142	THR	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	235	THR	3.1
1	D	188	LEU	3.0
1	D	176	ALA	3.0
1	B	7	PRO	2.9
1	B	54	PRO	2.9
1	D	97	SER	2.9
1	C	286	ASN	2.9
1	D	242	PRO	2.9
1	C	53	ARG	2.8
1	B	224	TRP	2.8
1	B	73	HIS	2.8
1	B	343	GLY	2.8
1	D	247	ASP	2.7
1	C	141	THR	2.7
1	B	195	GLU	2.7
1	D	196	ILE	2.7
1	D	286	ASN	2.7
1	D	107	LEU	2.6
1	A	54	PRO	2.6
1	C	7	PRO	2.6
1	B	241	ALA	2.6
1	A	235	THR	2.6
1	C	193	ALA	2.5
1	C	179	TRP	2.5
1	B	75	GLU	2.5
1	C	125	VAL	2.5
1	A	98	ALA	2.5
1	B	243	GLN	2.5
1	C	215	SER	2.5
1	A	9	PRO	2.4
1	D	73	HIS	2.3
1	C	54	PRO	2.3
1	B	344	GLY	2.3
1	D	114	GLU	2.3
1	C	124	GLY	2.3
1	D	192	TYR	2.3
1	A	73	HIS	2.2
1	D	344	GLY	2.2
1	B	102	LYS	2.2
1	C	243	GLN	2.2
1	D	655	LEU	2.1
1	C	206	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	190	GLU	2.1
1	C	196	ILE	2.0
1	C	247	ASP	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	GOL	B	701	6/6	0.64	0.21	58,60,61,61	0
2	GOL	C	702	6/6	0.74	0.20	59,60,60,61	0
2	GOL	C	701	6/6	0.92	0.10	28,33,34,36	0
2	GOL	A	701	6/6	0.94	0.17	26,32,33,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.