



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:45 pm BST

PDB ID : 6IGQ
Title : Crystal structure of inactive state of S9 peptidase from *Deinococcus radiodurans* R1 (PMSF treated)
Authors : Yadav, P.; Goyal, V.D.; Kumar, A.; Makde, R.D.
Deposited on : 2018-09-25
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
Mogul	:	1.8.5 (274361), CSD as541be (2020)
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

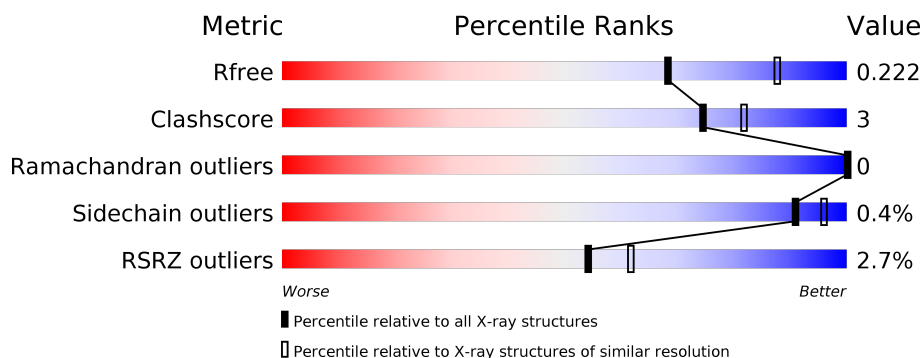
1 Overall quality at a glance

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 style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 6%, green 85%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 1% 85% 6% 9% </div> </div>
1	B	656	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 2%, yellow 5%, green 86%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 2% 86% 5% 9% </div> </div>
1	C	656	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 3%, yellow 6%, green 82%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 3% 82% 6% 11% </div> </div>
1	D	656	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 4%, orange 4%, yellow 9%, green 81%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 4% 81% 9% 11% </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	D	703	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 19250 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	599	Total	C	N	O	S	0	1	0
			4649	2950	837	853	9			
1	B	595	Total	C	N	O	S	0	3	0
			4610	2930	832	839	9			
1	C	583	Total	C	N	O	S	0	0	0
			4467	2843	796	819	9			
1	D	586	Total	C	N	O	S	0	0	0
			4446	2827	792	818	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 (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

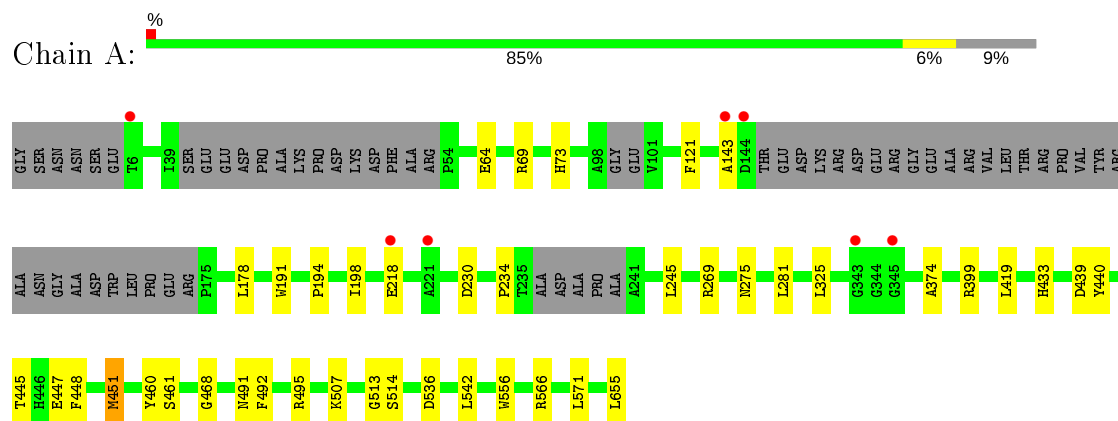
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	281	Total	O	0	0
			281	281		
4	B	277	Total	O	0	0
			277	277		
4	C	241	Total	O	0	0
			241	241		
4	D	193	Total	O	0	0
			193	193		

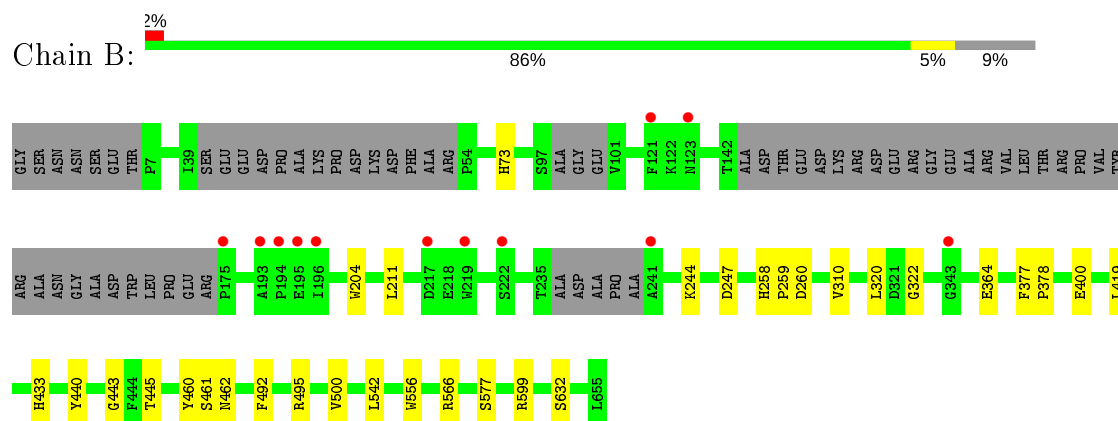
3 Residue-property plots [i](#)

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.

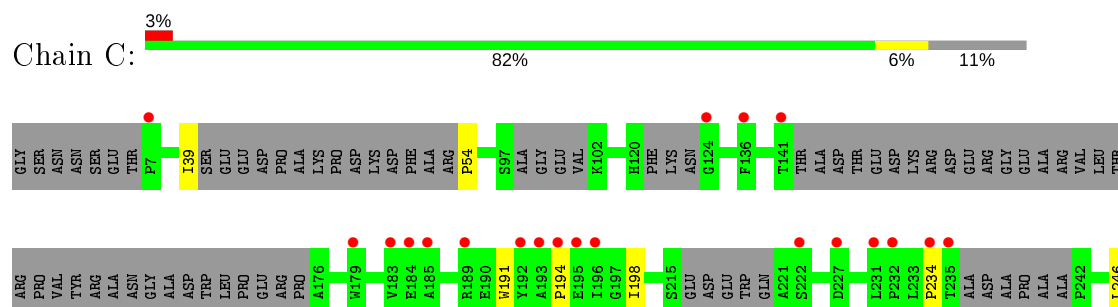
- Molecule 1: Acyl-peptide hydrolase, putative

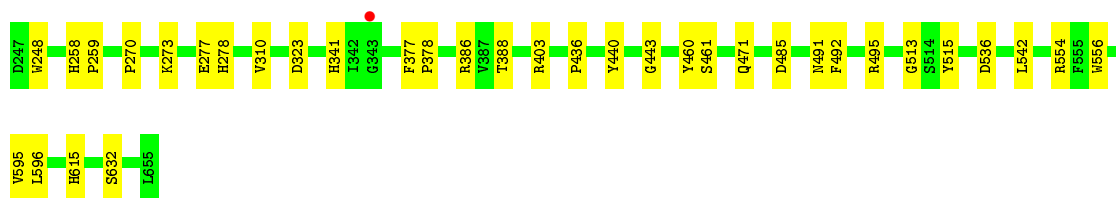


- Molecule 1: Acyl-peptide hydrolase, putative

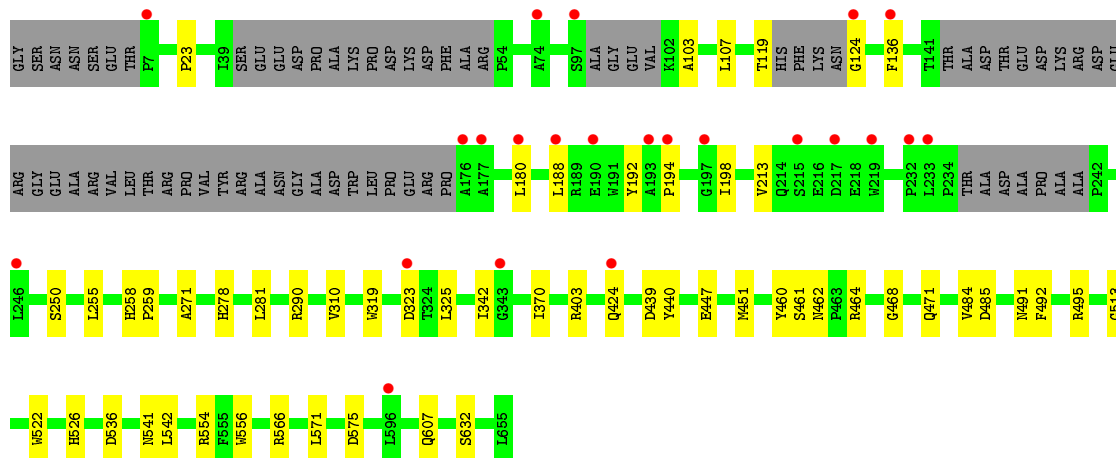
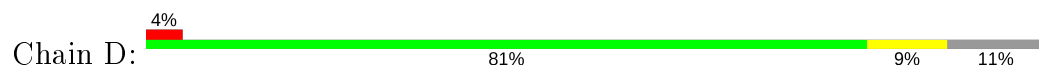


- Molecule 1: Acyl-peptide hydrolase, putative





- 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	119.66Å 129.99Å 194.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.87 – 2.30 39.87 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.87-2.30) 100.0 (39.87-2.30)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.29Å)	Xtrriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.192 , 0.224 0.192 , 0.222	Depositor DCC
R_{free} test set	6741 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtrriage
Anisotropy	0.269	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19250	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.61 % 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 1.9885e-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, SO4

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	0.36	0/4793	0.57	1/6539 (0.0%)
1	B	0.36	0/4761	0.56	0/6498
1	C	0.35	0/4600	0.55	0/6275
1	D	0.34	0/4578	0.55	0/6250
All	All	0.35	0/18732	0.56	1/25562 (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	A	451	MET	CG-SD-CE	5.17	108.48	100.20

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	4649	0	4385	27	0
1	B	4610	0	4340	24	0
1	C	4467	0	4195	24	0
1	D	4446	0	4137	38	0
2	A	18	0	24	3	0
2	B	18	0	24	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	12	0	16	1	0
2	D	18	0	24	4	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
4	A	281	0	0	0	0
4	B	277	0	0	0	0
4	C	241	0	0	1	0
4	D	193	0	0	3	0
All	All	19250	0	17145	112	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 (112) 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:440:TYR:H	1:B:462:ASN:HD21	1.11	0.98
1:B:258:HIS:HD2	1:B:260:ASP:H	1.04	0.92
1:D:440:TYR:H	1:D:462:ASN:HD21	1.12	0.92
1:B:258:HIS:CD2	1:B:260:ASP:H	1.93	0.87
1:C:278:HIS:H	1:C:471:GLN:HE22	1.24	0.86
1:D:571:LEU:HD11	2:D:703:GOL:H11	1.57	0.86
1:D:541:ASN:HD22	1:D:607:GLN:HE22	1.19	0.85
1:D:278:HIS:H	1:D:471:GLN:HE22	1.28	0.80
1:D:440:TYR:N	1:D:462:ASN:HD21	1.83	0.77
1:D:136:PHE:HB3	1:D:180:LEU:HD11	1.67	0.74
1:B:440:TYR:N	1:B:462:ASN:HD21	1.85	0.73
1:A:514:SER:HB3	2:A:703:GOL:H2	1.71	0.71
1:B:258:HIS:HD2	1:B:260:ASP:N	1.86	0.69
1:C:310:VAL:HG13	1:C:632:SER:HA	1.75	0.68
1:B:310:VAL:HG13	1:B:632:SER:HA	1.77	0.67
1:D:403:ARG:NH2	4:D:801:HOH:O	2.29	0.63
1:B:445:THR:HB	2:B:703:GOL:H11	1.81	0.63
1:D:522:TRP:O	1:D:526:HIS:HD2	1.83	0.60
1:A:64:GLU:OE2	1:A:69:ARG:NH2	2.33	0.59
1:D:310:VAL:HG13	1:D:632:SER:HA	1.84	0.59
1:A:194:PRO:HG3	1:A:198:ILE:HG13	1.84	0.59
1:A:447:GLU:HG2	1:A:451:MET:CE	2.34	0.57
1:A:447:GLU:HG2	1:A:451:MET:HE3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:278:HIS:H	1:C:471:GLN:NE2	1.98	0.57
1:D:440:TYR:H	1:D:462:ASN:ND2	1.93	0.57
1:B:443:GLY:O	2:B:703:GOL:O3	2.23	0.56
1:A:121:PHE:CD1	1:A:143:ALA:HA	2.41	0.56
1:D:103:ALA:HB3	1:D:124:GLY:HA3	1.87	0.56
1:C:194:PRO:HG3	1:C:198:ILE:HG13	1.88	0.56
1:A:571:LEU:HD13	2:D:703:GOL:H12	1.88	0.55
1:C:403:ARG:NH2	4:C:801:HOH:O	2.37	0.55
1:A:448:PHE:HD1	1:A:451:MET:HE3	1.71	0.55
1:A:399:ARG:HD3	1:A:419:LEU:O	2.06	0.55
1:A:447:GLU:O	1:A:451:MET:HG3	2.06	0.55
1:D:513:GLY:HA2	1:D:536:ASP:O	2.08	0.53
1:A:269:ARG:NH1	1:A:275:ASN:HD22	2.07	0.53
1:D:119:THR:HG22	1:D:188:LEU:HD21	1.91	0.53
1:C:386:ARG:NH1	1:C:388:THR:O	2.38	0.52
1:B:258:HIS:HE1	1:B:322:GLY:O	1.93	0.52
1:B:440:TYR:H	1:B:462:ASN:ND2	1.92	0.52
1:D:194:PRO:HG3	1:D:198:ILE:HG13	1.92	0.52
1:D:575:ASP:OD2	2:D:703:GOL:O3	2.26	0.52
1:B:440:TYR:CD2	1:B:460:TYR:HB2	2.45	0.52
1:C:191:TRP:NE1	1:C:234:PRO:HD3	2.26	0.50
1:D:250:SER:HB2	1:D:271:ALA:HA	1.93	0.50
1:D:290:ARG:NH2	1:D:342:ILE:HG23	2.26	0.50
1:A:440:TYR:CD2	1:A:460:TYR:HB2	2.47	0.49
1:C:270:PRO:HD2	1:C:273:LYS:HE3	1.93	0.49
1:D:278:HIS:H	1:D:471:GLN:NE2	2.04	0.49
1:A:439:ASP:OD2	1:A:468:GLY:HA3	2.13	0.49
1:D:571:LEU:HD21	2:D:703:GOL:H31	1.94	0.48
1:C:440:TYR:CD2	1:C:460:TYR:HB2	2.48	0.48
1:D:213:VAL:HG23	1:D:255:LEU:HD11	1.95	0.48
1:B:461:SER:HB3	1:B:492:PHE:CE2	2.49	0.48
1:C:595:VAL:HG23	1:C:596:LEU:HG	1.96	0.48
1:A:433:HIS:HB3	1:A:440:TYR:CZ	2.49	0.48
1:C:443:GLY:O	2:C:701:GOL:O1	2.33	0.47
1:A:491:ASN:O	1:A:495:ARG:HG2	2.14	0.47
1:A:445:THR:HB	2:A:702:GOL:H11	1.96	0.47
1:C:491:ASN:O	1:C:495:ARG:HG2	2.13	0.47
1:A:513:GLY:HA2	1:A:536:ASP:O	2.15	0.47
1:B:204:TRP:CZ2	1:B:211:LEU:HD21	2.50	0.47
1:C:273:LYS:HB3	1:C:277:GLU:HG3	1.97	0.46
1:C:542:LEU:HG	1:C:556:TRP:CZ2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:LEU:O	1:B:364:GLU:HG2	2.15	0.46
1:B:542:LEU:HD13	1:B:577:SER:HA	1.97	0.46
1:D:447:GLU:O	1:D:451:MET:HG3	2.15	0.46
1:A:230:ASP:HB2	1:A:245:LEU:HD13	1.98	0.46
1:A:461:SER:HB3	1:A:492:PHE:CE2	2.51	0.45
1:D:107:LEU:HA	1:D:107:LEU:HD12	1.85	0.45
1:C:513:GLY:HA2	1:C:536:ASP:O	2.16	0.45
1:D:542:LEU:HG	1:D:556:TRP:CZ2	2.51	0.45
1:D:440:TYR:CD2	1:D:460:TYR:HB2	2.51	0.45
1:D:461:SER:HB3	1:D:492:PHE:CE2	2.52	0.45
1:A:374:ALA:HB1	2:A:702:GOL:H2	1.99	0.44
1:A:507:LYS:HD3	1:A:655:LEU:OXT	2.18	0.44
1:C:258:HIS:CG	1:C:259:PRO:HD2	2.52	0.44
1:D:526:HIS:HE1	4:D:978:HOH:O	2.01	0.44
1:B:244:LYS:HD3	1:B:247:ASP:OD1	2.18	0.44
1:B:258:HIS:CG	1:B:259:PRO:HD2	2.53	0.44
1:B:542:LEU:HG	1:B:556:TRP:CZ2	2.53	0.43
1:B:461:SER:HB3	1:B:492:PHE:CD2	2.53	0.43
1:B:419:LEU:HD21	1:B:500:VAL:HG11	2.01	0.43
1:C:461:SER:HB3	1:C:492:PHE:CE2	2.54	0.42
1:A:178:LEU:HB2	1:A:198:ILE:HD12	2.01	0.42
1:A:542:LEU:HG	1:A:556:TRP:CZ2	2.54	0.42
1:B:433:HIS:HB3	1:B:440:TYR:CZ	2.54	0.42
1:D:323:ASP:OD1	1:D:342:ILE:HB	2.19	0.42
1:B:566:ARG:NH1	1:C:615:HIS:HE1	2.17	0.42
1:D:281:LEU:HD13	1:D:325:LEU:HD23	2.01	0.42
1:D:439:ASP:OD2	1:D:468:GLY:HA3	2.19	0.42
1:C:39:ILE:HD12	1:C:54:PRO:HB2	2.01	0.42
1:D:192:TYR:HD2	1:D:198:ILE:HD11	1.84	0.42
1:D:319:TRP:CE2	1:D:325:LEU:HD13	2.55	0.42
1:C:436:PRO:HB3	1:C:515:TYR:CD1	2.54	0.42
1:D:424:GLN:N	1:D:424:GLN:OE1	2.45	0.41
1:C:323:ASP:O	1:C:341:HIS:HA	2.21	0.41
1:D:23:PRO:HG3	1:D:370:ILE:HG13	2.01	0.41
1:D:485:ASP:N	1:D:485:ASP:OD1	2.53	0.41
1:D:491:ASN:O	1:D:495:ARG:HG2	2.21	0.41
1:C:377:PHE:CG	1:C:378:PRO:HD2	2.56	0.41
1:A:461:SER:HB3	1:A:492:PHE:CD2	2.56	0.41
1:D:258:HIS:CG	1:D:259:PRO:HD2	2.56	0.41
1:D:464:ARG:CZ	1:D:484:VAL:HB	2.51	0.41
1:C:485:ASP:OD1	1:C:485:ASP:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:566:ARG:NH2	4:D:803:HOH:O	2.42	0.41
1:A:191:TRP:NE1	1:A:234:PRO:HD3	2.36	0.40
1:A:281:LEU:HD13	1:A:325:LEU:HD23	2.02	0.40
1:C:246:LEU:HD21	1:C:248:TRP:CE3	2.56	0.40
1:A:269:ARG:HH11	1:A:275:ASN:HD22	1.67	0.40
1:B:377:PHE:CG	1:B:378:PRO:HD2	2.56	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	590/656 (90%)	576 (98%)	14 (2%)	0	100	100
1	B	588/656 (90%)	574 (98%)	14 (2%)	0	100	100
1	C	569/656 (87%)	559 (98%)	10 (2%)	0	100	100
1	D	574/656 (88%)	566 (99%)	8 (1%)	0	100	100
All	All	2321/2624 (88%)	2275 (98%)	46 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	469/526 (89%)	465 (99%)	4 (1%)	78	89
1	B	463/526 (88%)	459 (99%)	4 (1%)	78	89
1	C	445/526 (85%)	444 (100%)	1 (0%)	93	97
1	D	437/526 (83%)	436 (100%)	1 (0%)	93	97
All	All	1814/2104 (86%)	1804 (99%)	10 (1%)	91	94

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73[A]	HIS
1	A	73[B]	HIS
1	A	218	GLU
1	A	566	ARG
1	B	73[A]	HIS
1	B	73[B]	HIS
1	B	400	GLU
1	B	599	ARG
1	C	554	ARG
1	D	554	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	275	ASN
1	A	521	ASN
1	B	258	HIS
1	B	446	HIS
1	B	462	ASN
1	C	262	GLN
1	C	471	GLN
1	D	462	ASN
1	D	471	GLN
1	D	526	HIS
1	D	607	GLN

5.3.3 RNA ⓘ

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](#)

15 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	D	701	-	5,5,5	0.37	0	5,5,5	0.26	0
2	GOL	B	701	-	5,5,5	0.33	0	5,5,5	0.38	0
3	SO4	D	704	-	4,4,4	0.15	0	6,6,6	0.18	0
2	GOL	C	701	-	5,5,5	0.44	0	5,5,5	0.43	0
2	GOL	A	701	-	5,5,5	0.30	0	5,5,5	0.35	0
3	SO4	A	704	-	4,4,4	0.22	0	6,6,6	0.28	0
2	GOL	A	702	-	5,5,5	0.31	0	5,5,5	0.45	0
3	SO4	C	703	-	4,4,4	0.20	0	6,6,6	0.19	0
2	GOL	C	702	-	5,5,5	0.38	0	5,5,5	0.34	0
2	GOL	D	703	-	5,5,5	0.47	0	5,5,5	0.41	0
2	GOL	B	703	-	5,5,5	0.36	0	5,5,5	0.28	0
2	GOL	A	703	-	5,5,5	0.35	0	5,5,5	0.14	0
2	GOL	D	702	-	5,5,5	0.39	0	5,5,5	0.22	0
3	SO4	B	704	-	4,4,4	0.17	0	6,6,6	0.21	0
2	GOL	B	702	-	5,5,5	0.38	0	5,5,5	0.19	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	D	701	-	-	1/4/4/4	-
2	GOL	B	701	-	-	2/4/4/4	-
2	GOL	C	701	-	-	2/4/4/4	-
2	GOL	A	701	-	-	0/4/4/4	-
2	GOL	A	702	-	-	2/4/4/4	-
2	GOL	C	702	-	-	4/4/4/4	-
2	GOL	D	703	-	-	2/4/4/4	-
2	GOL	B	703	-	-	3/4/4/4	-
2	GOL	A	703	-	-	3/4/4/4	-
2	GOL	D	702	-	-	3/4/4/4	-
2	GOL	B	702	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	701	GOL	O1-C1-C2-C3
2	A	702	GOL	O1-C1-C2-C3
2	C	702	GOL	O1-C1-C2-C3
2	C	702	GOL	C1-C2-C3-O3
2	A	702	GOL	O1-C1-C2-O2
2	D	701	GOL	O1-C1-C2-C3
2	D	703	GOL	O1-C1-C2-C3
2	B	703	GOL	O1-C1-C2-C3
2	A	703	GOL	C1-C2-C3-O3
2	D	702	GOL	C1-C2-C3-O3
2	B	702	GOL	O1-C1-C2-C3
2	C	701	GOL	O1-C1-C2-O2
2	C	702	GOL	O1-C1-C2-O2
2	D	703	GOL	O1-C1-C2-O2
2	B	703	GOL	O1-C1-C2-O2
2	C	702	GOL	O2-C2-C3-O3
2	D	702	GOL	O2-C2-C3-O3
2	B	701	GOL	C1-C2-C3-O3
2	B	701	GOL	O2-C2-C3-O3
2	A	703	GOL	O1-C1-C2-C3
2	A	703	GOL	O2-C2-C3-O3
2	B	702	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	B	703	GOL	C1-C2-C3-O3
2	D	702	GOL	O1-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	701	GOL	1	0
2	A	702	GOL	2	0
2	D	703	GOL	4	0
2	B	703	GOL	2	0
2	A	703	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	599/656 (91%)	-0.32	7 (1%) 79 83	18, 26, 45, 69	0
1	B	595/656 (90%)	-0.26	12 (2%) 65 71	18, 27, 54, 67	0
1	C	583/656 (88%)	-0.13	21 (3%) 42 49	17, 30, 59, 75	0
1	D	586/656 (89%)	0.05	23 (3%) 39 46	18, 35, 61, 69	0
All	All	2363/2624 (90%)	-0.17	63 (2%) 54 62	17, 29, 57, 75	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	196	ILE	5.0
1	D	188	LEU	3.9
1	B	219	TRP	3.7
1	D	74	ALA	3.6
1	C	185	ALA	3.5
1	D	219	TRP	3.4
1	C	193	ALA	3.4
1	B	123	ASN	3.2
1	D	193	ALA	3.2
1	A	143	ALA	3.2
1	D	232	PRO	3.1
1	D	217	ASP	3.1
1	B	222	SER	3.0
1	D	197	GLY	3.0
1	C	179	TRP	3.0
1	B	196	ILE	3.0
1	A	144	ASP	2.9
1	D	233	LEU	2.9
1	B	193	ALA	2.9
1	B	195	GLU	2.8
1	D	177	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	7	PRO	2.8
1	B	343	GLY	2.8
1	C	222	SER	2.8
1	B	121	PHE	2.7
1	D	176	ALA	2.7
1	C	343	GLY	2.7
1	C	7	PRO	2.7
1	D	323	ASP	2.6
1	A	6	THR	2.5
1	C	194	PRO	2.5
1	C	189	ARG	2.5
1	D	424	GLN	2.5
1	C	195	GLU	2.5
1	D	180	LEU	2.5
1	A	343	GLY	2.5
1	B	175	PRO	2.4
1	D	136	PHE	2.4
1	A	345	GLY	2.4
1	C	184	GLU	2.4
1	D	596	LEU	2.4
1	D	97	SER	2.4
1	D	124	GLY	2.3
1	C	141	THR	2.3
1	C	235	THR	2.3
1	C	234	PRO	2.3
1	D	190	GLU	2.3
1	C	232	PRO	2.3
1	C	231	LEU	2.2
1	C	136	PHE	2.2
1	D	215	SER	2.2
1	C	183	VAL	2.2
1	A	218	GLU	2.2
1	B	194	PRO	2.2
1	B	241	ALA	2.1
1	D	194	PRO	2.1
1	A	221	ALA	2.1
1	D	246	LEU	2.1
1	B	217	ASP	2.1
1	C	192	TYR	2.1
1	D	343	GLY	2.0
1	C	227	ASP	2.0
1	C	124	GLY	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 carbohydrates 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(\AA^2)	Q<0.9
2	GOL	A	703	6/6	0.74	0.31	66,66,67,67	0
2	GOL	D	703	6/6	0.85	0.36	44,45,46,46	0
2	GOL	C	702	6/6	0.89	0.16	57,57,57,58	0
2	GOL	C	701	6/6	0.90	0.20	38,40,40,40	0
2	GOL	D	702	6/6	0.90	0.16	51,51,52,52	0
2	GOL	B	702	6/6	0.92	0.11	29,30,31,32	0
2	GOL	B	703	6/6	0.94	0.16	39,40,40,40	0
2	GOL	A	701	6/6	0.94	0.13	32,32,33,33	0
2	GOL	A	702	6/6	0.95	0.13	29,30,30,32	0
2	GOL	D	701	6/6	0.96	0.16	26,31,33,33	0
2	GOL	B	701	6/6	0.97	0.10	34,35,35,35	0
3	SO4	A	704	5/5	0.98	0.09	33,33,35,36	0
3	SO4	D	704	5/5	0.98	0.08	49,50,50,51	0
3	SO4	C	703	5/5	0.98	0.12	36,38,39,39	0
3	SO4	B	704	5/5	0.99	0.08	29,29,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.