



# wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 16, 2017 – 09:22 AM EST

PDB ID : 5MR3  
Title : Crystal structure of red abalone egg VERL repeat 2 with linker in complex with sperm lysin at 1.8 Å resolution  
Authors : Nishimura, K.; Raj, I.; Sadat Al-Hosseini, H.; De Sanctis, D.; Jovine, L.  
Deposited on : unknown  
Resolution : 1.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

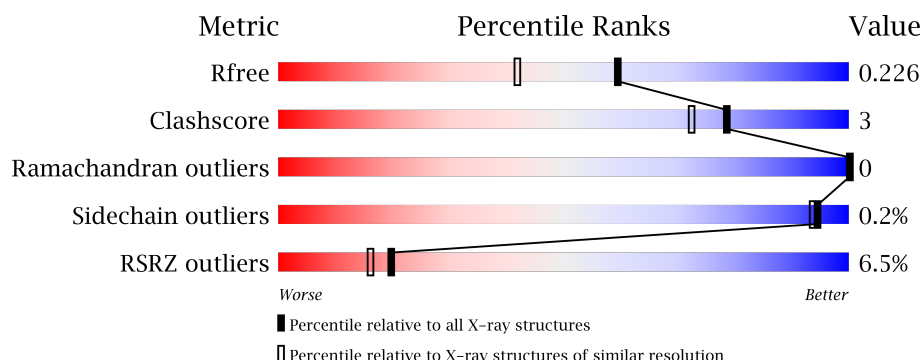
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	100719	4827 (1.80-1.80)
Clashscore	112137	5742 (1.80-1.80)
Ramachandran outliers	110173	5676 (1.80-1.80)
Sidechain outliers	110143	5675 (1.80-1.80)
RSRZ outliers	101464	4906 (1.80-1.80)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	136	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>7%</div> <div>8%</div> </div> </div>
1	C	136	<div> <div>4%</div> <div> <div></div> <div>89%</div> <div>• •</div> <div>7%</div> </div> </div>
1	E	136	<div> <div>3%</div> <div> <div></div> <div>96%</div> <div>• •</div> </div> </div>
1	G	136	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>8%</div> <div>5%</div> </div> </div>
2	B	134	<div> <div>7%</div> <div> <div></div> <div>78%</div> <div>10%</div> <div>13%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	D	134	
2	F	134	
2	H	134	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	201	-	-	-	X
3	GOL	B	901	-	-	-	X
3	GOL	B	902	-	-	-	X
3	GOL	B	904	-	-	-	X
3	GOL	D	902	-	-	-	X
3	GOL	D	903	-	-	-	X
3	GOL	E	900	-	-	-	X
3	GOL	E	901	-	-	-	X
3	GOL	H	902	-	-	-	X
4	CL	B	903	-	-	-	X
5	NAG	B	905	-	-	-	X
5	NAG	F	402	-	-	-	X
6	PGE	G	900	-	-	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 16585 atoms, of which 8042 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 Egg-lysin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	125	Total	C	H	N	O	S	0	0	0
			2129	687	1079	185	171	7			
1	C	127	Total	C	H	N	O	S	0	0	0
			2166	698	1100	188	173	7			
1	E	133	Total	C	H	N	O	S	0	1	0
			2293	743	1158	201	184	7			
1	G	129	Total	C	H	N	O	S	0	0	0
			2196	708	1114	190	177	7			

- Molecule 2 is a protein called Vitelline envelope sperm lysin receptor.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	117	Total	C	H	N	O	S	0	0	0
			1810	588	893	149	170	10			
2	D	110	Total	C	H	N	O	S	0	0	0
			1698	552	835	140	161	10			
2	F	107	Total	C	H	N	O	S	0	0	0
			1670	543	824	138	155	10			
2	H	111	Total	C	H	N	O	S	0	0	0
			1733	561	859	143	160	10			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	173	GLU	-	expression tag	UNP Q8WR62
B	174	THR	-	expression tag	UNP Q8WR62
B	175	GLY	-	expression tag	UNP Q8WR62
B	293	ALA	SER	engineered mutation	UNP Q8WR62
B	296	ALA	SER	engineered mutation	UNP Q8WR62
B	297	ALA	SER	engineered mutation	UNP Q8WR62
B	299	LEU	-	expression tag	UNP Q8WR62
B	300	GLU	-	expression tag	UNP Q8WR62

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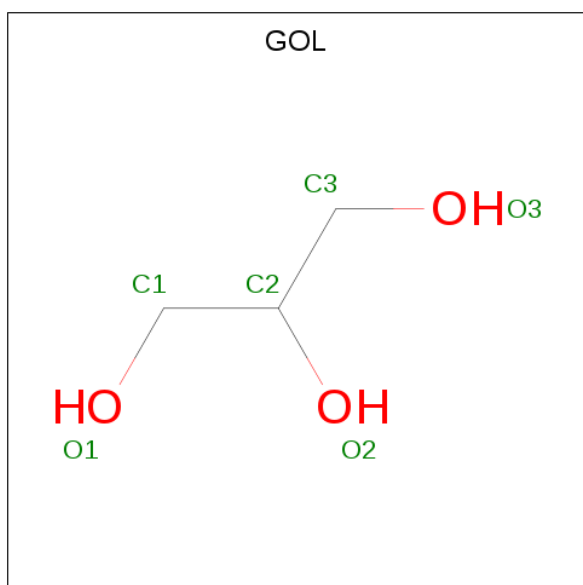
Chain	Residue	Modelled	Actual	Comment	Reference
B	301	HIS	-	expression tag	UNP Q8WR62
B	302	HIS	-	expression tag	UNP Q8WR62
B	303	HIS	-	expression tag	UNP Q8WR62
B	304	HIS	-	expression tag	UNP Q8WR62
B	305	HIS	-	expression tag	UNP Q8WR62
B	306	HIS	-	expression tag	UNP Q8WR62
D	173	GLU	-	expression tag	UNP Q8WR62
D	174	THR	-	expression tag	UNP Q8WR62
D	175	GLY	-	expression tag	UNP Q8WR62
D	293	ALA	SER	engineered mutation	UNP Q8WR62
D	296	ALA	SER	engineered mutation	UNP Q8WR62
D	297	ALA	SER	engineered mutation	UNP Q8WR62
D	299	LEU	-	expression tag	UNP Q8WR62
D	300	GLU	-	expression tag	UNP Q8WR62
D	301	HIS	-	expression tag	UNP Q8WR62
D	302	HIS	-	expression tag	UNP Q8WR62
D	303	HIS	-	expression tag	UNP Q8WR62
D	304	HIS	-	expression tag	UNP Q8WR62
D	305	HIS	-	expression tag	UNP Q8WR62
D	306	HIS	-	expression tag	UNP Q8WR62
F	173	GLU	-	expression tag	UNP Q8WR62
F	174	THR	-	expression tag	UNP Q8WR62
F	175	GLY	-	expression tag	UNP Q8WR62
F	293	ALA	SER	engineered mutation	UNP Q8WR62
F	296	ALA	SER	engineered mutation	UNP Q8WR62
F	297	ALA	SER	engineered mutation	UNP Q8WR62
F	299	LEU	-	expression tag	UNP Q8WR62
F	300	GLU	-	expression tag	UNP Q8WR62
F	301	HIS	-	expression tag	UNP Q8WR62
F	302	HIS	-	expression tag	UNP Q8WR62
F	303	HIS	-	expression tag	UNP Q8WR62
F	304	HIS	-	expression tag	UNP Q8WR62
F	305	HIS	-	expression tag	UNP Q8WR62
F	306	HIS	-	expression tag	UNP Q8WR62
H	173	GLU	-	expression tag	UNP Q8WR62
H	174	THR	-	expression tag	UNP Q8WR62
H	175	GLY	-	expression tag	UNP Q8WR62
H	293	ALA	SER	engineered mutation	UNP Q8WR62
H	296	ALA	SER	engineered mutation	UNP Q8WR62
H	297	ALA	SER	engineered mutation	UNP Q8WR62
H	299	LEU	-	expression tag	UNP Q8WR62
H	300	GLU	-	expression tag	UNP Q8WR62

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Chain	Residue	Modelled	Actual	Comment	Reference
H	301	HIS	-	expression tag	UNP Q8WR62
H	302	HIS	-	expression tag	UNP Q8WR62
H	303	HIS	-	expression tag	UNP Q8WR62
H	304	HIS	-	expression tag	UNP Q8WR62
H	305	HIS	-	expression tag	UNP Q8WR62
H	306	HIS	-	expression tag	UNP Q8WR62

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	D	1	Total	C	H	O	0	0
			14	3	8	3		
3	D	1	Total	C	H	O	0	0
			14	3	8	3		

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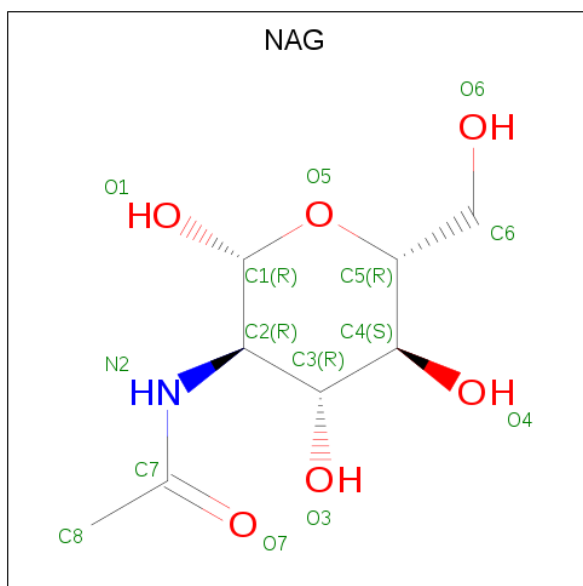
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	E	1	Total	C	H	O	0	0
			14	3	8	3		
3	E	1	Total	C	H	O	0	0
			14	3	8	3		
3	H	1	Total	C	H	O	0	0
			14	3	8	3		
3	H	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		
4	F	1	Total	Cl	0	0
			1	1		

- Molecule 5 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



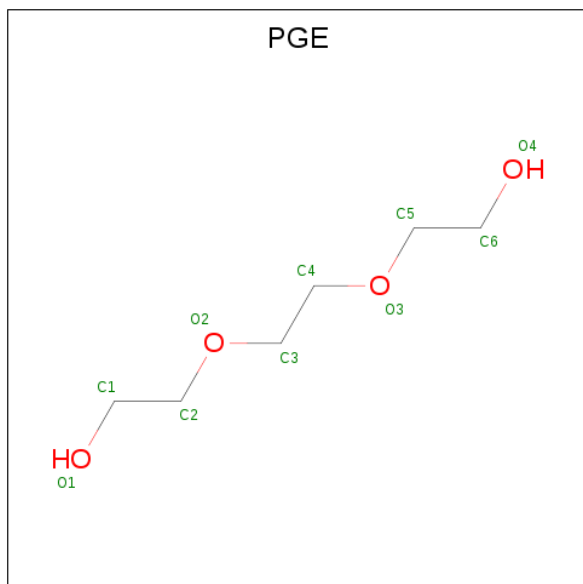
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
5	D	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
5	F	1	Total	C	H	N	O	0	0
			28	8	14	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	H	1	Total	C	H	N	O	
			28	8	14	1	5	
							0	0

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula:  $C_6H_{14}O_4$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	G	1	Total	C	H	O		
			24	6	14	4		
							0	0
6	G	1	Total	C	H	O		
			24	6	14	4		
							0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	70	Total	O		
			70	70		
					0	0
7	B	79	Total	O		
			79	79		
					0	0
7	C	57	Total	O		
			57	57		
					0	0
7	D	61	Total	O		
			61	61		
					0	0
7	E	94	Total	O		
			95	95		
					0	1
7	F	69	Total	O		
			69	69		
					0	0

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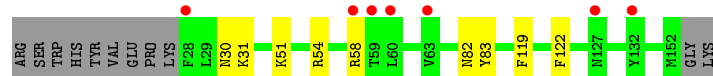
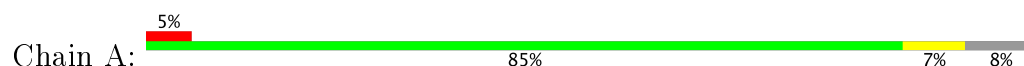
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	G	76	Total	O	0	0
			76	76		
7	H	53	Total	O	0	0
			53	53		

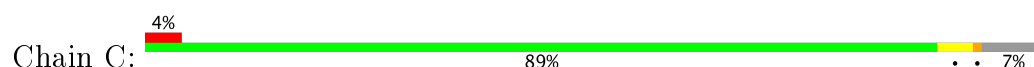
### 3 Residue-property plots [i](#)

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.

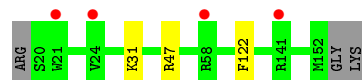
- Molecule 1: Egg-lysin



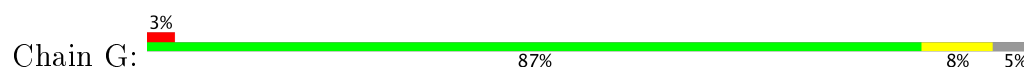
- Molecule 1: Egg-lysin



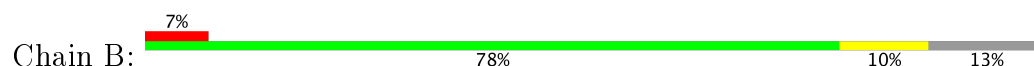
- Molecule 1: Egg-lysin



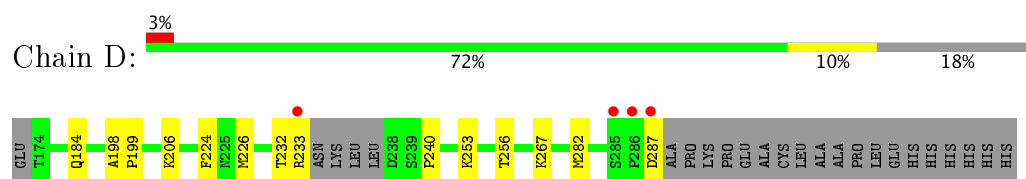
- Molecule 1: Egg-lysin



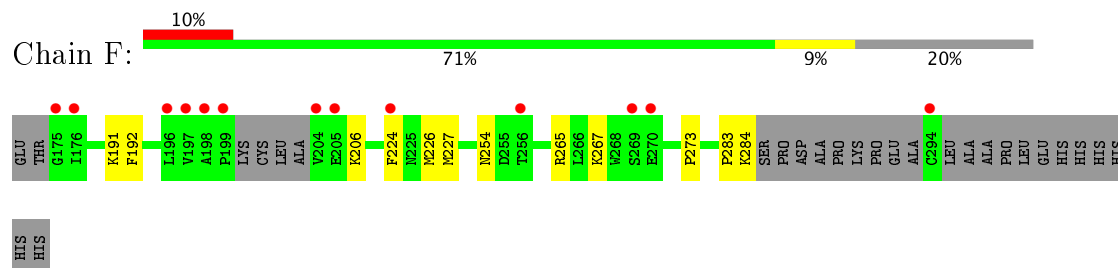
- Molecule 2: Vitelline envelope sperm lysin receptor



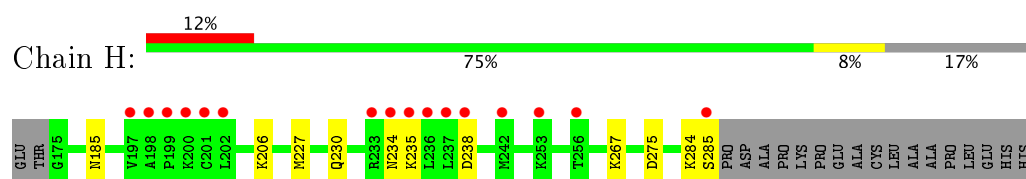
- Molecule 2: Vitelline envelope sperm lysin receptor



- Molecule 2: Vitelline envelope sperm lysin receptor



- Molecule 2: Vitelline envelope sperm lysin receptor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.87Å 87.86Å 92.55Å 90.00° 100.93° 90.00°	Depositor
Resolution (Å)	45.32 – 1.80 45.32 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (45.32-1.80) 89.8 (45.32-1.80)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.59 (at 1.79Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, $R_{free}$	0.201 , 0.227 0.200 , 0.226	Depositor DCC
$R_{free}$ test set	4971 reflections (5.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	30.0	Xtriage
Anisotropy	0.311	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 48.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	16585	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PGE, NAG, CL

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	0.29	0/1077	0.45	0/1451
1	C	0.29	0/1094	0.47	0/1473
1	E	0.29	0/1167	0.45	0/1574
1	G	0.34	0/1110	0.47	0/1496
2	B	0.31	0/943	0.55	0/1286
2	D	0.33	0/887	0.55	0/1209
2	F	0.30	0/868	0.53	0/1180
2	H	0.31	0/898	0.55	0/1223
All	All	0.31	0/8044	0.50	0/10892

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	1050	1079	1079	7	0
1	C	1066	1100	1100	3	1
1	E	1135	1158	1157	3	0
1	G	1082	1114	1114	9	1
2	B	917	893	893	10	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	863	835	835	10	1
2	F	846	824	823	7	0
2	H	874	859	859	12	1
3	A	6	8	8	0	0
3	B	18	24	24	0	0
3	C	6	8	8	0	0
3	D	18	24	24	1	0
3	E	12	16	16	0	0
3	H	12	16	16	1	0
4	B	1	0	0	0	0
4	F	1	0	0	0	0
5	B	14	14	13	0	0
5	D	14	14	13	0	0
5	F	14	14	13	1	0
5	H	14	14	13	0	0
6	G	20	28	28	0	0
7	A	70	0	0	0	0
7	B	79	0	0	2	0
7	C	57	0	0	0	0
7	D	61	0	0	2	0
7	E	95	0	0	2	0
7	F	69	0	0	0	0
7	G	76	0	0	1	0
7	H	53	0	0	5	0
All	All	8543	8042	8036	52	3

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.

The worst 5 of 52 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)
2:D:282:MET:HG2	7:D:1002:HOH:O	1.79	0.81
2:B:206:LYS:HB2	7:B:1041:HOH:O	1.85	0.75
2:H:238:ASP:OD1	2:H:238:ASP:O	2.06	0.74
1:C:55:VAL:O	1:C:58:ARG:NH2	2.22	0.72
1:E:47:ARG:NH2	7:E:1001:HOH:O	2.23	0.72

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:205:GLU:OE2	1:G:143:ARG:NH2[2_646]	1.97	0.23
2:B:184:GLN:OE1	1:C:143:ARG:HH21[2_555]	1.51	0.09
2:D:184:GLN:O	2:H:185:ASN:HD22[1_455]	1.58	0.02

## 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	123/136 (90%)	123 (100%)	0	0	100	100
1	C	125/136 (92%)	125 (100%)	0	0	100	100
1	E	132/136 (97%)	132 (100%)	0	0	100	100
1	G	127/136 (93%)	127 (100%)	0	0	100	100
2	B	113/134 (84%)	110 (97%)	3 (3%)	0	100	100
2	D	106/134 (79%)	105 (99%)	1 (1%)	0	100	100
2	F	102/134 (76%)	101 (99%)	1 (1%)	0	100	100
2	H	109/134 (81%)	108 (99%)	1 (1%)	0	100	100
All	All	937/1080 (87%)	931 (99%)	6 (1%)	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	109/119 (92%)	109 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	111/119 (93%)	110 (99%)	1 (1%)	82	78
1	E	118/119 (99%)	118 (100%)	0	100	100
1	G	113/119 (95%)	113 (100%)	0	100	100
2	B	106/120 (88%)	105 (99%)	1 (1%)	82	78
2	D	100/120 (83%)	100 (100%)	0	100	100
2	F	98/120 (82%)	98 (100%)	0	100	100
2	H	101/120 (84%)	101 (100%)	0	100	100
All	All	856/956 (90%)	854 (100%)	2 (0%)	94	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	220	HIS
1	C	58	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 20 ligands modelled in this entry, 2 are monoatomic - leaving 18 for Mogul analysis.

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	A	201	-	5,5,5	0.33	0	5,5,5	0.38	0
3	GOL	B	901	-	5,5,5	0.41	0	5,5,5	0.14	0
3	GOL	B	902	-	5,5,5	0.39	0	5,5,5	0.20	0
3	GOL	B	904	-	5,5,5	0.31	0	5,5,5	0.37	0
5	NAG	B	905	2	14,14,15	0.35	0	15,19,21	0.52	0
3	GOL	C	900	-	5,5,5	0.40	0	5,5,5	0.09	0
3	GOL	D	901	-	5,5,5	0.32	0	5,5,5	0.56	0
3	GOL	D	902	-	5,5,5	0.35	0	5,5,5	0.11	0
3	GOL	D	903	-	5,5,5	0.36	0	5,5,5	0.36	0
5	NAG	D	904	2	14,14,15	0.17	0	15,19,21	0.54	0
3	GOL	E	900	-	5,5,5	0.32	0	5,5,5	0.24	0
3	GOL	E	901	-	5,5,5	0.35	0	5,5,5	0.27	0
5	NAG	F	402	2	14,14,15	0.26	0	15,19,21	0.61	0
6	PGE	G	900	-	9,9,9	0.33	0	8,8,8	0.29	0
6	PGE	G	901	-	9,9,9	0.31	0	8,8,8	0.35	0
3	GOL	H	901	-	5,5,5	0.46	0	5,5,5	0.35	0
3	GOL	H	902	-	5,5,5	0.35	0	5,5,5	0.39	0
5	NAG	H	903	2	14,14,15	0.23	0	15,19,21	0.46	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	A	201	-	-	0/4/4/4	0/0/0/0
3	GOL	B	901	-	-	0/4/4/4	0/0/0/0
3	GOL	B	902	-	-	0/4/4/4	0/0/0/0
3	GOL	B	904	-	-	0/4/4/4	0/0/0/0
5	NAG	B	905	2	-	0/6/23/26	0/1/1/1
3	GOL	C	900	-	-	0/4/4/4	0/0/0/0
3	GOL	D	901	-	-	0/4/4/4	0/0/0/0
3	GOL	D	902	-	-	0/4/4/4	0/0/0/0
3	GOL	D	903	-	-	0/4/4/4	0/0/0/0
5	NAG	D	904	2	-	0/6/23/26	0/1/1/1
3	GOL	E	900	-	-	0/4/4/4	0/0/0/0
3	GOL	E	901	-	-	0/4/4/4	0/0/0/0
5	NAG	F	402	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PGE	G	900	-	-	0/7/7/7	0/0/0/0
6	PGE	G	901	-	-	0/7/7/7	0/0/0/0
3	GOL	H	901	-	-	0/4/4/4	0/0/0/0
3	GOL	H	902	-	-	0/4/4/4	0/0/0/0
5	NAG	H	903	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	902	GOL	1	0
5	F	402	NAG	1	0
3	H	902	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	125/136 (91%)	0.18	7 (5%) 25 21	30, 46, 74, 128	0
1	C	127/136 (93%)	0.03	5 (3%) 40 35	32, 49, 78, 111	0
1	E	133/136 (97%)	-0.08	4 (3%) 51 45	29, 42, 75, 107	0
1	G	129/136 (94%)	-0.09	4 (3%) 49 44	34, 46, 78, 101	0
2	B	117/134 (87%)	0.55	9 (7%) 14 11	28, 40, 101, 147	0
2	D	110/134 (82%)	0.33	4 (3%) 43 38	27, 42, 90, 131	0
2	F	107/134 (79%)	0.64	13 (12%) 5 4	28, 41, 104, 139	0
2	H	111/134 (82%)	0.65	16 (14%) 3 2	31, 43, 122, 139	0
All	All	959/1080 (88%)	0.25	62 (6%) 20 16	27, 45, 92, 147	0

The worst 5 of 62 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	237	LEU	11.0
2	H	201	CYS	8.7
2	B	197	VAL	8.0
2	F	198	ALA	7.8
2	H	235	LYS	6.6

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	A	201	6/6	0.84	0.55	19.99	109,130,134,137	0
3	GOL	H	902	6/6	0.68	0.52	19.16	117,141,148,151	0
3	GOL	D	902	6/6	0.76	0.47	11.72	91,110,116,116	0
3	GOL	E	901	6/6	0.54	0.38	11.41	121,146,147,147	0
3	GOL	E	900	6/6	0.44	0.38	10.64	134,161,167,168	0
3	GOL	B	904	6/6	0.72	0.22	8.70	75,91,95,95	0
3	GOL	D	903	6/6	0.82	0.41	6.63	148,178,182,185	0
5	NAG	B	905	14/15	0.79	0.23	4.47	76,91,103,103	0
5	NAG	F	402	14/15	0.58	0.30	3.73	98,104,123,125	0
6	PGE	G	900	10/10	0.78	0.19	2.91	70,84,96,97	0
3	GOL	B	901	6/6	0.86	0.20	2.39	110,132,144,148	0
4	CL	B	903	1/1	0.99	0.18	2.20	33,33,33,33	1
3	GOL	B	902	6/6	0.91	0.18	2.03	83,100,104,108	0
6	PGE	G	901	10/10	0.54	0.22	0.82	126,152,183,183	0
5	NAG	D	904	14/15	0.86	0.13	0.77	75,87,100,105	0
3	GOL	C	900	6/6	0.56	0.17	0.76	119,143,145,147	0
3	GOL	D	901	6/6	0.86	0.15	0.73	45,54,56,57	0
5	NAG	H	903	14/15	0.83	0.17	0.29	63,76,93,93	0
3	GOL	H	901	6/6	0.94	0.11	0.13	43,53,63,64	0
4	CL	F	401	1/1	0.98	0.09	-1.55	36,36,36,36	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.