



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 16, 2018 – 12:27 PM EST

PDB ID : 2RCV  
Title : Crystal structure of the Bacillus subtilis superoxide dismutase  
Authors : Liu, P.; Ewis, H.E.; Huang, Y.J.; Lu, C.D.; Tai, P.C.; Weber, I.T.  
Deposited on : 2007-09-20  
Resolution : 1.60 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](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  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030736  
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-20030736

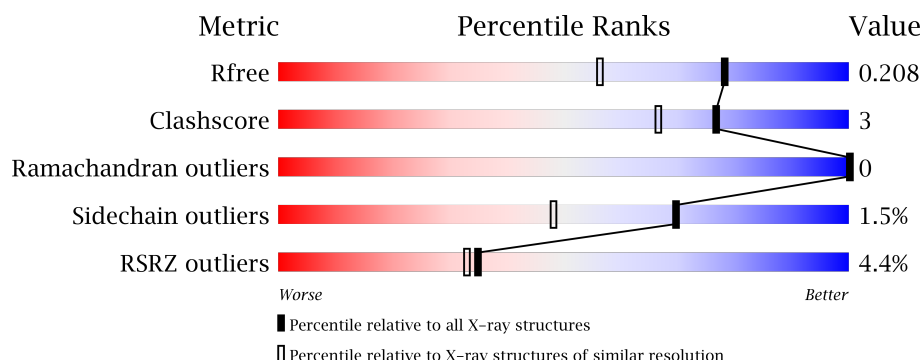
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 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	2696 (1.60-1.60)
Clashscore	112137	2967 (1.60-1.60)
Ramachandran outliers	110173	2887 (1.60-1.60)
Sidechain outliers	110143	2886 (1.60-1.60)
RSRZ outliers	101464	2714 (1.60-1.60)

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	202	<div> <div>4%</div> <div>94%</div> <div>6%</div> <div>..</div> </div>
1	B	202	<div> <div>5%</div> <div>91%</div> <div>6%</div> <div>..</div> </div>
1	C	202	<div> <div>4%</div> <div>89%</div> <div>9%</div> <div>.</div> </div>
1	D	202	<div> <div>6%</div> <div>87%</div> <div>10%</div> <div>..</div> </div>
1	E	202	<div> <div>3%</div> <div>92%</div> <div>6%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	202	
1	G	202	
1	H	202	

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
2	MN	A	203	-	-	-	X
2	MN	C	203	-	-	-	X
2	MN	F	203	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13464 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 Superoxide dismutase [Mn].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1576	1000	270	305	1			
1	B	197	Total	C	N	O	S	0	0	0
			1555	989	266	299	1			
1	C	200	Total	C	N	O	S	0	0	0
			1576	1000	270	305	1			
1	D	197	Total	C	N	O	S	0	0	0
			1556	987	267	301	1			
1	E	200	Total	C	N	O	S	0	0	0
			1576	1000	270	305	1			
1	F	200	Total	C	N	O	S	0	0	0
			1576	1000	270	305	1			
1	G	200	Total	C	N	O	S	0	0	0
			1576	1000	270	305	1			
1	H	197	Total	C	N	O	S	0	0	0
			1555	989	266	299	1			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		
2	E	1	Total	Mn	0	0
			1	1		
2	H	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Mn 1	0	0
2	F	1	Total 1	Mn 1	0	0

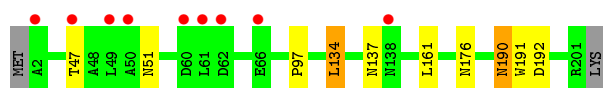
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	115	Total 115	O 115	0	0
3	B	114	Total 114	O 114	0	0
3	C	98	Total 98	O 98	0	0
3	D	103	Total 103	O 103	0	0
3	E	123	Total 123	O 123	0	0
3	F	117	Total 117	O 117	0	0
3	G	111	Total 111	O 111	0	0
3	H	129	Total 129	O 129	0	0

### 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: Superoxide dismutase [Mn]



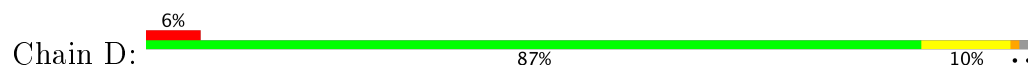
- Molecule 1: Superoxide dismutase [Mn]



- Molecule 1: Superoxide dismutase [Mn]



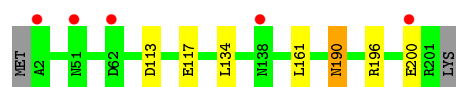
- Molecule 1: Superoxide dismutase [Mn]



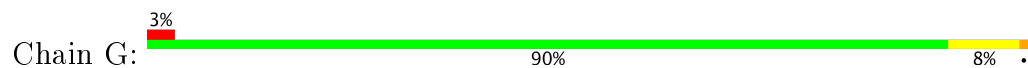
- Molecule 1: Superoxide dismutase [Mn]



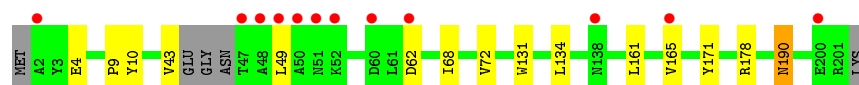
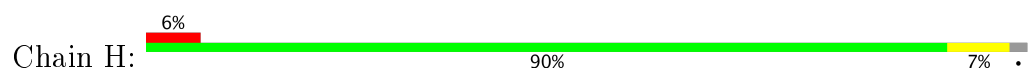
- Molecule 1: Superoxide dismutase [Mn]



- Molecule 1: Superoxide dismutase [Mn]



- Molecule 1: Superoxide dismutase [Mn]



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	68.37Å 84.03Å 91.95Å 99.13° 105.98° 105.58°	Depositor
Resolution (Å)	45.90 – 1.60 45.90 – 1.59	Depositor EDS
% Data completeness (in resolution range)	89.9 (45.90-1.60) 84.8 (45.90-1.59)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.13 (at 1.59Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.211 , 0.230 0.212 , 0.208	Depositor DCC
$R_{free}$ test set	10942 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	12.1	Xtriage
Anisotropy	0.834	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 45.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13464	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.86% 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: MN

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	0/1621	0.54	1/2214 (0.0%)
1	B	0.32	0/1599	0.53	1/2183 (0.0%)
1	C	0.30	0/1621	0.51	0/2214
1	D	0.32	0/1600	0.54	0/2183
1	E	0.32	0/1621	0.54	0/2214
1	F	0.33	0/1621	0.55	0/2214
1	G	0.32	0/1621	0.54	0/2214
1	H	0.33	0/1599	0.55	0/2183
All	All	0.32	0/12903	0.54	2/17619 (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	B	134	LEU	N-CA-C	-5.35	96.56	111.00
1	A	134	LEU	N-CA-C	-5.17	97.04	111.00

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	1576	0	1481	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1555	0	1465	11	0
1	C	1576	0	1481	12	0
1	D	1556	0	1457	14	0
1	E	1576	0	1481	9	0
1	F	1576	0	1481	4	0
1	G	1576	0	1481	18	0
1	H	1555	0	1465	10	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	115	0	0	0	0
3	B	114	0	0	1	0
3	C	98	0	0	0	0
3	D	103	0	0	0	0
3	E	123	0	0	1	0
3	F	117	0	0	0	0
3	G	111	0	0	0	0
3	H	129	0	0	1	0
All	All	13464	0	11792	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 3.

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:B:6:PRO:HG3	1:B:83:LYS:HD2	1.74	0.69
1:G:101:LEU:HA	1:G:104:GLU:OE1	1.94	0.67
1:C:134:LEU:HB3	1:C:161:LEU:HB3	1.77	0.66
1:B:96:GLU:HG3	1:B:112:PHE:CD2	2.31	0.65
1:E:134:LEU:HB3	1:E:161:LEU:HB3	1.79	0.65
1:F:113:ASP:O	1:F:117:GLU:HG3	1.99	0.63
1:B:113:ASP:O	1:B:117:GLU:HG3	1.99	0.63
1:C:131:TRP:HA	1:C:165:VAL:HG23	1.81	0.62
1:A:134:LEU:HB3	1:A:161:LEU:HB3	1.80	0.62
1:D:61:LEU:HD12	1:D:69:ARG:HG3	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:97:PRO:HG3	1:G:191:TRP:CE3	2.37	0.60
1:H:131:TRP:HA	1:H:165:VAL:HG23	1.84	0.60
1:G:135:VAL:HG12	1:G:159:PRO:HA	1.84	0.59
1:G:190:ASN:C	1:G:190:ASN:HD22	2.06	0.58
1:G:130:GLY:O	1:G:165:VAL:HG22	2.04	0.57
1:E:131:TRP:HA	1:E:165:VAL:HG23	1.87	0.56
1:D:190:ASN:C	1:D:190:ASN:HD22	2.10	0.56
1:G:134:LEU:HB3	1:G:161:LEU:HB3	1.87	0.55
1:G:100:ALA:O	1:G:104:GLU:CD	2.45	0.54
1:H:4:GLU:HG3	3:H:286:HOH:O	2.07	0.54
1:B:68:ILE:O	1:B:72:VAL:HG23	2.08	0.53
1:D:134:LEU:HB3	1:D:161:LEU:HB3	1.90	0.53
1:C:114:LYS:O	1:C:117:GLU:HG2	2.09	0.53
1:G:9:PRO:HG2	1:G:10:TYR:CD2	2.44	0.53
1:C:97:PRO:HG3	1:C:191:TRP:CE3	2.45	0.52
1:B:190:ASN:C	1:B:190:ASN:HD22	2.13	0.51
1:E:41:LYS:HG3	3:E:248:HOH:O	2.10	0.51
1:A:190:ASN:C	1:A:190:ASN:HD22	2.14	0.51
1:A:47:THR:HG22	1:A:51:ASN:HD21	1.76	0.51
1:G:9:PRO:HG2	1:G:10:TYR:CE2	2.47	0.51
1:G:100:ALA:O	1:G:104:GLU:OE2	2.28	0.50
1:D:96:GLU:HG3	1:D:112:PHE:CD2	2.45	0.50
1:G:97:PRO:HG3	1:G:191:TRP:CD2	2.46	0.50
1:C:97:PRO:HG3	1:C:191:TRP:CD2	2.46	0.49
1:A:176:ASN:HD22	1:A:176:ASN:H	1.60	0.49
1:A:97:PRO:HG3	1:A:191:TRP:CE3	2.48	0.49
1:D:43:VAL:HG12	1:D:43:VAL:O	2.13	0.49
1:H:190:ASN:HD22	1:H:190:ASN:C	2.15	0.48
1:H:134:LEU:HB3	1:H:161:LEU:HB3	1.95	0.48
1:E:47:THR:HG22	1:E:51:ASN:ND2	2.28	0.48
1:G:21:LYS:O	1:G:25:THR:HG23	2.13	0.48
1:H:68:ILE:O	1:H:72:VAL:HG23	2.12	0.48
1:F:134:LEU:HB3	1:F:161:LEU:HB3	1.95	0.48
1:G:104:GLU:CD	1:G:198:TYR:CZ	2.87	0.48
1:G:176:ASN:H	1:G:176:ASN:HD22	1.62	0.48
1:E:190:ASN:C	1:E:190:ASN:HD22	2.17	0.47
1:B:6:PRO:HG3	1:B:83:LYS:CD	2.44	0.47
1:D:97:PRO:HG3	1:D:191:TRP:CE3	2.49	0.47
1:F:190:ASN:C	1:F:190:ASN:HD22	2.15	0.47
1:A:190:ASN:HD21	1:A:192:ASP:HB2	1.80	0.47
1:C:53:SER:OG	1:C:56:GLU:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:165:VAL:HG12	1:H:178:ARG:CZ	2.45	0.47
1:E:165:VAL:HG12	1:E:178:ARG:CZ	2.44	0.46
1:G:130:GLY:C	1:G:165:VAL:HG22	2.35	0.46
1:C:176:ASN:HD22	1:C:176:ASN:H	1.64	0.46
1:B:97:PRO:HG3	1:B:191:TRP:CD2	2.50	0.46
1:F:196:ARG:NH2	1:F:200:GLU:OE2	2.50	0.45
1:A:47:THR:HG22	1:A:51:ASN:ND2	2.31	0.45
1:C:190:ASN:C	1:C:190:ASN:HD22	2.19	0.45
1:H:43:VAL:CG1	1:H:49:LEU:HB2	2.46	0.44
1:D:65:PRO:HD2	1:D:68:ILE:HD11	2.00	0.44
1:D:113:ASP:O	1:D:117:GLU:HG3	2.17	0.44
1:H:43:VAL:HG13	1:H:49:LEU:HB2	2.00	0.43
1:D:54:VAL:O	1:D:58:VAL:HG23	2.19	0.43
1:G:104:GLU:CD	1:G:198:TYR:OH	2.56	0.43
1:C:167:GLU:HB2	1:D:167:GLU:HB2	2.01	0.43
1:H:9:PRO:HG2	1:H:10:TYR:CD2	2.54	0.43
1:E:47:THR:HG22	1:E:51:ASN:HD21	1.84	0.43
1:C:104:GLU:O	1:C:108:VAL:HG23	2.19	0.43
1:G:165:VAL:O	1:G:165:VAL:HG23	2.18	0.43
1:B:97:PRO:HG3	1:B:191:TRP:CE3	2.53	0.42
1:H:9:PRO:HG2	1:H:10:TYR:CE2	2.55	0.42
1:E:176:ASN:HD22	1:E:176:ASN:H	1.68	0.42
1:B:6:PRO:HB3	1:B:83:LYS:HE2	2.02	0.42
1:C:165:VAL:HG12	1:C:178:ARG:CZ	2.49	0.41
1:D:37:THR:HG22	1:D:41:LYS:HE3	2.02	0.41
1:B:134:LEU:HB3	1:B:161:LEU:HB3	2.01	0.41
1:C:8:LEU:HA	1:C:9:PRO:HD3	1.85	0.41
1:D:97:PRO:HG3	1:D:191:TRP:CD2	2.55	0.41
1:E:135:VAL:HG12	1:E:159:PRO:HA	2.02	0.41
1:B:96:GLU:HG2	3:B:266:HOH:O	2.20	0.41
1:D:67:ASN:ND2	1:D:68:ILE:HG23	2.36	0.41
1:A:97:PRO:HG3	1:A:191:TRP:CD2	2.56	0.40
1:G:190:ASN:C	1:G:190:ASN:ND2	2.72	0.40
1:D:193:GLU:OE2	1:D:196:ARG:NH1	2.54	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	198/202 (98%)	193 (98%)	5 (2%)	0	100	100
1	B	193/202 (96%)	186 (96%)	7 (4%)	0	100	100
1	C	198/202 (98%)	193 (98%)	5 (2%)	0	100	100
1	D	193/202 (96%)	187 (97%)	6 (3%)	0	100	100
1	E	198/202 (98%)	193 (98%)	5 (2%)	0	100	100
1	F	198/202 (98%)	192 (97%)	6 (3%)	0	100	100
1	G	198/202 (98%)	191 (96%)	7 (4%)	0	100	100
1	H	193/202 (96%)	187 (97%)	6 (3%)	0	100	100
All	All	1569/1616 (97%)	1522 (97%)	47 (3%)	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	164/167 (98%)	162 (99%)	2 (1%)	75	58
1	B	162/167 (97%)	160 (99%)	2 (1%)	75	58
1	C	164/167 (98%)	161 (98%)	3 (2%)	64	40
1	D	162/167 (97%)	160 (99%)	2 (1%)	75	58
1	E	164/167 (98%)	160 (98%)	4 (2%)	54	26
1	F	164/167 (98%)	163 (99%)	1 (1%)	89	80

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	164/167 (98%)	161 (98%)	3 (2%)	64	40
1	H	162/167 (97%)	159 (98%)	3 (2%)	62	37
All	All	1306/1336 (98%)	1286 (98%)	20 (2%)	70	48

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	190	ASN
1	B	171	TYR
1	B	190	ASN
1	C	113	ASP
1	C	137	ASN
1	C	190	ASN
1	D	61	LEU
1	D	190	ASN
1	E	113	ASP
1	E	137	ASN
1	E	165	VAL
1	E	190	ASN
1	F	190	ASN
1	G	113	ASP
1	G	165	VAL
1	G	190	ASN
1	H	62	ASP
1	H	171	TYR
1	H	190	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	92	ASN
1	A	137	ASN
1	A	148	ASN
1	A	176	ASN
1	A	190	ASN
1	B	33	ASN
1	B	148	ASN
1	B	190	ASN

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Mol	Chain	Res	Type
1	C	51	ASN
1	C	137	ASN
1	C	148	ASN
1	C	176	ASN
1	C	190	ASN
1	D	190	ASN
1	E	33	ASN
1	E	51	ASN
1	E	92	ASN
1	E	137	ASN
1	E	148	ASN
1	E	176	ASN
1	E	190	ASN
1	F	148	ASN
1	F	190	ASN
1	G	137	ASN
1	G	148	ASN
1	G	176	ASN
1	G	190	ASN
1	H	148	ASN
1	H	190	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	200/202 (99%)	0.22	9 (4%) 34 32	8, 16, 34, 45	0
1	B	197/202 (97%)	0.33	10 (5%) 29 28	8, 16, 41, 49	0
1	C	200/202 (99%)	0.23	9 (4%) 34 32	9, 19, 36, 44	0
1	D	197/202 (97%)	0.27	12 (6%) 22 21	9, 18, 41, 48	0
1	E	200/202 (99%)	0.26	6 (3%) 51 49	8, 16, 35, 46	0
1	F	200/202 (99%)	0.15	5 (2%) 58 57	8, 15, 34, 43	0
1	G	200/202 (99%)	0.22	7 (3%) 44 43	8, 17, 35, 44	0
1	H	197/202 (97%)	0.31	12 (6%) 22 21	8, 16, 35, 44	0
All	All	1591/1616 (98%)	0.25	70 (4%) 35 33	8, 17, 36, 49	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	50	ALA	13.1
1	B	47	THR	6.2
1	H	51	ASN	5.6
1	G	47	THR	4.8
1	E	2	ALA	4.7
1	F	51	ASN	4.6
1	C	2	ALA	4.5
1	A	47	THR	4.3
1	B	138	ASN	4.1
1	C	47	THR	4.0
1	A	2	ALA	3.9
1	H	138	ASN	3.9
1	D	44	GLU	3.9
1	H	62	ASP	3.7
1	H	48	ALA	3.7
1	B	48	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	H	49	LEU	3.7
1	B	60	ASP	3.6
1	B	50	ALA	3.6
1	E	50	ALA	3.6
1	C	165	VAL	3.5
1	A	61	LEU	3.5
1	G	104	GLU	3.5
1	G	62	ASP	3.5
1	D	61	LEU	3.4
1	G	2	ALA	3.4
1	C	50	ALA	3.3
1	B	66	GLU	3.3
1	G	50	ALA	3.2
1	D	60	ASP	3.2
1	B	56	GLU	3.2
1	F	2	ALA	3.2
1	A	60	ASP	3.1
1	H	2	ALA	3.1
1	G	155	GLU	3.0
1	A	62	ASP	3.0
1	E	47	THR	3.0
1	H	47	THR	2.9
1	D	50	ALA	2.9
1	D	63	SER	2.8
1	H	165	VAL	2.8
1	B	62	ASP	2.8
1	H	200	GLU	2.7
1	C	138	ASN	2.7
1	B	49	LEU	2.6
1	C	155	GLU	2.6
1	D	138	ASN	2.6
1	G	138	ASN	2.6
1	H	52	LYS	2.5
1	C	66	GLU	2.5
1	B	63	SER	2.4
1	A	138	ASN	2.4
1	D	51	ASN	2.3
1	F	138	ASN	2.3
1	D	43	VAL	2.3
1	D	2	ALA	2.2
1	D	62	ASP	2.2
1	C	60	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	66	GLU	2.2
1	D	45	GLY	2.2
1	D	200	GLU	2.2
1	E	155	GLU	2.2
1	E	165	VAL	2.1
1	F	62	ASP	2.1
1	A	50	ALA	2.1
1	E	66	GLU	2.1
1	C	62	ASP	2.1
1	H	60	ASP	2.1
1	A	49	LEU	2.0
1	F	200	GLU	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. 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
2	MN	C	203	1/1	1.00	0.10	2.95	10,10,10,10	0
2	MN	F	203	1/1	1.00	0.11	2.46	9,9,9,9	0
2	MN	A	203	1/1	1.00	0.10	2.35	9,9,9,9	0
2	MN	G	203	1/1	1.00	0.10	1.56	8,8,8,8	0
2	MN	H	203	1/1	1.00	0.10	1.06	10,10,10,10	0
2	MN	B	203	1/1	1.00	0.09	-0.80	10,10,10,10	0
2	MN	D	203	1/1	1.00	0.07	-2.01	9,9,9,9	0
2	MN	E	203	1/1	1.00	0.08	-2.23	8,8,8,8	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.