



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 9, 2017 – 05:28 PM EST

PDB ID : 4NAU  
Title : S. aureus CoaD with Inhibitor  
Authors : Lahiri, S.D.  
Deposited on : unknown  
Resolution : 2.33 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  
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

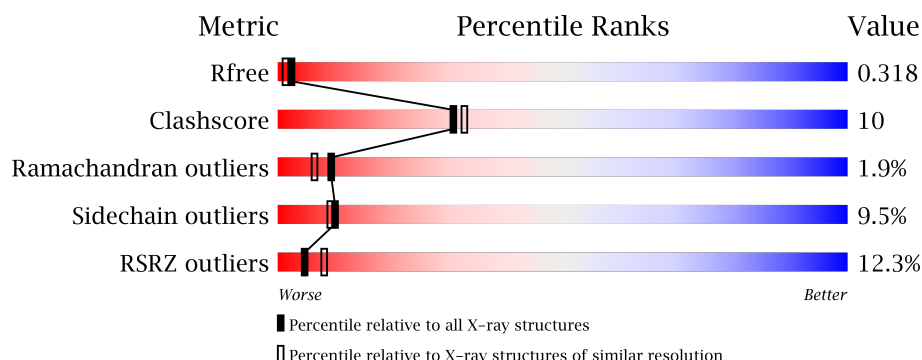
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.33 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	1570 (2.36-2.32)
Clashscore	112137	1673 (2.36-2.32)
Ramachandran outliers	110173	1654 (2.36-2.32)
Sidechain outliers	110143	1655 (2.36-2.32)
RSRZ outliers	101464	1576 (2.36-2.32)

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	160	<div> <div>14%</div> <div>73%</div> <div>21%</div> <div>• •</div> </div>
1	B	160	<div> <div>12%</div> <div>81%</div> <div>14%</div> <div>• •</div> </div>
1	C	160	<div> <div>11%</div> <div>77%</div> <div>16%</div> <div>7%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	AGS	A	201	-	-	-	X
2	AGS	C	201	-	-	-	X

## 2 Entry composition [i](#)

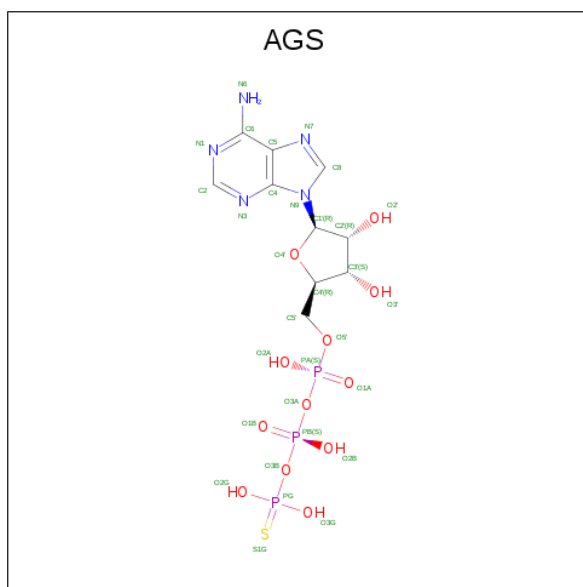
There are 4 unique types of molecules in this entry. The entry contains 4062 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Phosphopantetheine adenylyltransferase.

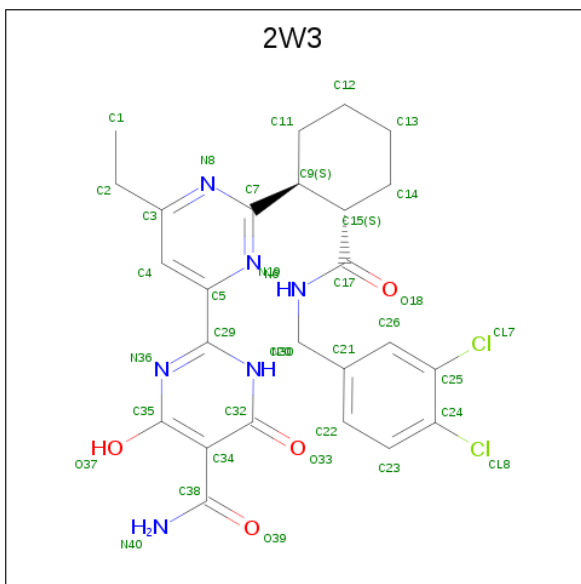
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	0	0
			1282	819	213	244	6			
1	B	160	Total	C	N	O	S	0	0	0
			1290	824	214	245	7			
1	C	160	Total	C	N	O	S	0	0	0
			1290	824	214	245	7			

- Molecule 2 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (three-letter code: AGS) (formula:  $C_{10}H_{16}N_5O_{12}P_3S$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
2	C	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		

- Molecule 3 is 2-[2-[(1S,2S)-2-[(3,4-dichlorophenyl)methylcarbamoyl]cyclohexyl]-6-ethyl-pyrimidin-4-yl]-4-oxidanyl-6-oxidanylidene-1H-pyrimidine-5-carboxamide (three-letter code: 2W3) (formula:  $C_{25}H_{26}Cl_2N_6O_4$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			37	25	2	6	4		

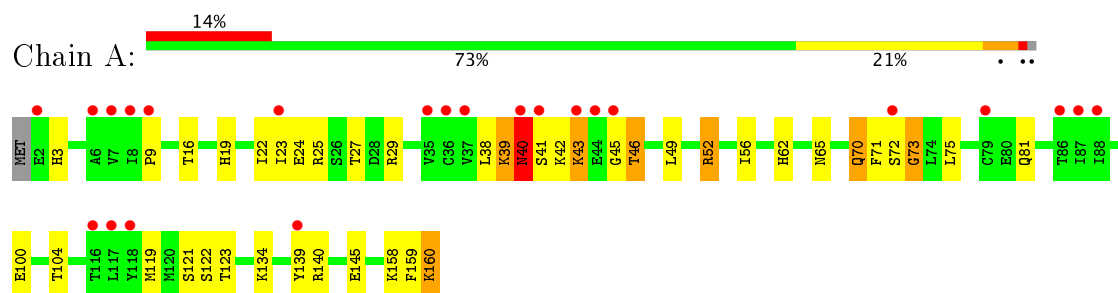
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	28	Total	O	0	0
			28	28		
4	B	20	Total	O	0	0
			20	20		
4	C	22	Total	O	0	0
			22	22		

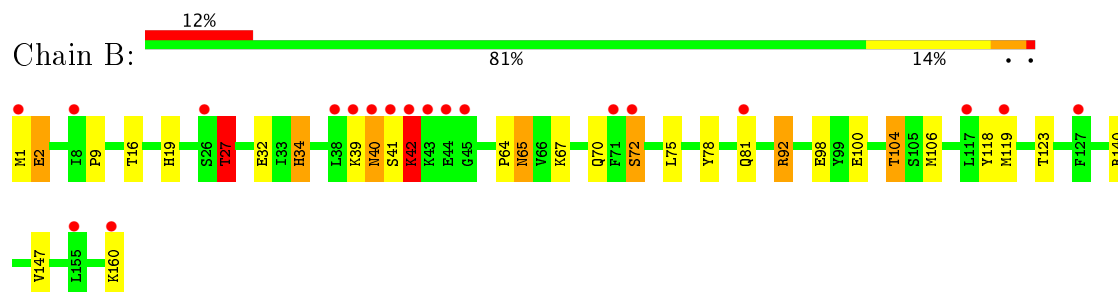
### 3 Residue-property plots [i](#)

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.

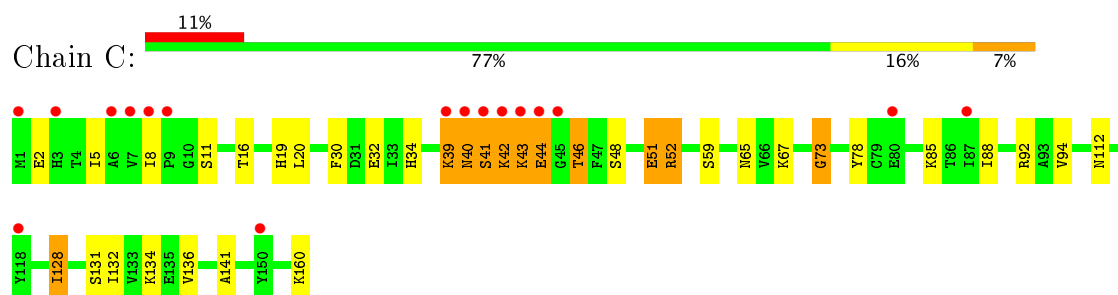
- Molecule 1: Phosphopantetheine adenylyltransferase



- Molecule 1: Phosphopantetheine adenylyltransferase



- Molecule 1: Phosphopantetheine adenylyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.09Å 127.78Å 127.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.94 – 2.33 31.78 – 2.33	Depositor EDS
% Data completeness (in resolution range)	98.8 (31.94-2.33) 98.8 (31.78-2.33)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.68 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.4.0066	Depositor
R, $R_{free}$	0.256 , 0.320 0.260 , 0.318	Depositor DCC
$R_{free}$ test set	1353 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.8	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 34.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.034 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	4062	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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: 2W3, AGS

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	0.81	0/1306	0.86	1/1759 (0.1%)
1	B	0.74	0/1314	0.83	2/1769 (0.1%)
1	C	0.69	0/1314	0.84	2/1769 (0.1%)
All	All	0.75	0/3934	0.84	5/5297 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	52	ARG	NE-CZ-NH1	-7.54	116.53	120.30
1	C	52	ARG	NE-CZ-NH2	5.31	122.96	120.30
1	B	27	THR	CB-CA-C	-5.25	97.41	111.60
1	A	52	ARG	NE-CZ-NH2	5.02	122.81	120.30
1	B	92	ARG	NE-CZ-NH1	5.00	122.80	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1282	0	1292	30	0
1	B	1290	0	1304	24	0
1	C	1290	0	1304	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	31	0	12	3	0
2	B	31	0	12	1	0
2	C	31	0	12	3	0
3	A	37	0	26	5	0
4	A	28	0	0	2	0
4	B	20	0	0	0	0
4	C	22	0	0	0	0
All	All	4062	0	3962	83	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 10.

All (83) 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:32:GLU:OE1	1:B:67:LYS:NZ	1.97	0.97
1:A:71:PHE:HB3	3:A:202:2W3:CL7	2.04	0.94
1:C:39:LYS:HG2	1:C:52:ARG:HH12	1.33	0.94
1:C:39:LYS:CG	1:C:52:ARG:HH12	1.82	0.92
1:A:40:ASN:OD1	1:A:42:LYS:HG2	1.74	0.88
1:A:3:HIS:CE1	4:A:308:HOH:O	2.29	0.84
1:B:27:THR:HG22	1:B:65:ASN:HD21	1.43	0.83
1:B:27:THR:CG2	1:B:65:ASN:HD21	1.93	0.81
1:C:32:GLU:OE1	1:C:67:LYS:NZ	2.18	0.76
1:B:27:THR:HG22	1:B:65:ASN:ND2	2.02	0.75
1:B:16:THR:H	1:B:19:HIS:HD2	1.38	0.71
1:A:40:ASN:HB3	1:A:41:SER:HA	1.72	0.70
1:C:39:LYS:HG3	1:C:52:ARG:HH12	1.60	0.66
1:C:39:LYS:HG3	1:C:52:ARG:NH1	2.11	0.66
1:C:42:LYS:HA	1:C:44:GLU:N	2.10	0.66
2:A:201:AGS:O2A	4:A:327:HOH:O	2.13	0.66
1:A:39:LYS:HE2	1:A:70:GLN:OE1	1.96	0.65
1:A:100:GLU:O	1:A:104:THR:HG22	1.98	0.64
1:C:39:LYS:CG	1:C:52:ARG:NH1	2.58	0.63
1:A:16:THR:H	1:A:19:HIS:HD2	1.50	0.59
1:B:104:THR:CG2	1:B:118:TYR:OH	2.50	0.59
3:A:202:2W3:N36	3:A:202:2W3:H18	2.18	0.58
2:A:201:AGS:H8	2:A:201:AGS:O2B	2.04	0.58
1:A:40:ASN:CB	1:A:41:SER:HA	2.35	0.57
1:B:41:SER:O	1:B:42:LYS:C	2.45	0.55
1:C:34:HIS:HD2	1:C:78:TYR:OH	1.90	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:PHE:O	1:A:160:LYS:HB2	2.07	0.55
1:B:65:ASN:HD22	1:B:65:ASN:H	1.55	0.55
1:A:16:THR:H	1:A:19:HIS:CD2	2.25	0.54
1:B:39:LYS:HD3	1:B:72:SER:HB3	1.90	0.53
1:B:27:THR:CG2	1:B:65:ASN:ND2	2.63	0.53
1:B:39:LYS:O	1:B:40:ASN:C	2.46	0.53
2:C:201:AGS:H8	2:C:201:AGS:O2B	2.09	0.52
1:C:46:THR:HG21	1:C:134:LYS:HD2	1.92	0.52
1:A:25:ARG:CZ	1:A:119:MET:HE3	2.40	0.51
1:B:65:ASN:HD22	1:B:65:ASN:N	2.06	0.51
1:A:72:SER:O	1:A:73:GLY:O	2.28	0.51
1:B:34:HIS:HD2	1:B:78:TYR:OH	1.94	0.51
1:A:42:LYS:O	1:A:43:LYS:O	2.29	0.50
1:A:25:ARG:HB2	1:A:119:MET:HE1	1.94	0.50
1:B:16:THR:N	1:B:19:HIS:HD2	2.06	0.50
1:A:73:GLY:O	3:A:202:2W3:H23	2.12	0.49
1:A:40:ASN:OD1	1:A:42:LYS:CG	2.53	0.49
1:C:40:ASN:O	1:C:41:SER:C	2.50	0.49
1:B:9:PRO:HG3	1:B:75:LEU:HD21	1.95	0.49
1:B:100:GLU:O	1:B:104:THR:HG23	2.13	0.49
1:C:128:ILE:HA	1:C:132:ILE:HD11	1.95	0.49
1:B:27:THR:HG21	1:B:64:PRO:HD2	1.95	0.48
1:C:39:LYS:O	1:C:41:SER:N	2.47	0.48
1:B:27:THR:HG23	1:B:65:ASN:HD21	1.75	0.48
1:A:49:LEU:HD21	1:A:70:GLN:OE1	2.14	0.48
1:C:48:SER:OG	1:C:51:GLU:HG3	2.14	0.47
1:A:38:LEU:HD21	3:A:202:2W3:H25	1.79	0.47
1:A:46:THR:HG21	1:A:134:LYS:HD2	1.97	0.47
1:C:136:VAL:HG13	1:C:141:ALA:HB3	1.97	0.46
1:A:41:SER:N	1:A:42:LYS:HA	2.30	0.46
1:A:52:ARG:O	1:A:56:ILE:HG13	2.16	0.46
1:C:42:LYS:N	1:C:43:LYS:HB2	2.31	0.46
1:B:100:GLU:HB3	1:B:118:TYR:OH	2.15	0.45
1:C:16:THR:CB	1:C:128:ILE:HD11	2.46	0.45
1:A:9:PRO:HG3	1:A:75:LEU:HD21	1.99	0.45
1:C:16:THR:OG1	1:C:128:ILE:HD11	2.17	0.44
1:A:45:GLY:HA2	1:A:46:THR:HA	1.81	0.44
1:A:23:ILE:O	1:A:27:THR:HG23	2.16	0.44
1:B:98:GLU:OE2	1:C:94:VAL:HB	2.17	0.44
3:A:202:2W3:C5	3:A:202:2W3:H18	2.48	0.43
1:C:30:PHE:CE1	1:C:88:ILE:HD11	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:ILE:HG12	1:C:32:GLU:HB2	2.01	0.42
1:C:92:ARG:HD3	2:C:201:AGS:C5	2.48	0.42
1:A:22:ILE:HD12	1:A:119:MET:HE2	2.02	0.42
1:C:8:ILE:HG12	1:C:88:ILE:HB	2.00	0.42
1:B:39:LYS:HD3	1:B:40:ASN:OD1	2.19	0.42
1:C:39:LYS:HD2	1:C:40:ASN:N	2.34	0.42
1:B:92:ARG:HD3	2:B:201:AGS:C6	2.50	0.41
1:C:16:THR:H	1:C:19:HIS:HD2	1.67	0.41
1:A:139:TYR:CD1	1:C:73:GLY:HA2	2.55	0.41
1:A:41:SER:N	1:A:42:LYS:CA	2.82	0.41
1:C:92:ARG:HD3	2:C:201:AGS:C6	2.51	0.41
1:B:106:MET:HE2	1:C:132:ILE:HD13	2.03	0.40
1:A:24:GLU:OE2	1:A:62:HIS:HE1	2.03	0.40
1:B:16:THR:HA	1:B:147:VAL:HG12	2.02	0.40
1:A:29:ARG:HG3	1:A:29:ARG:HH11	1.85	0.40
1:A:121:SER:OG	2:A:201:AGS:C2	2.70	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	157/160 (98%)	149 (95%)	4 (2%)	4 (2%)	6	4
1	B	158/160 (99%)	151 (96%)	5 (3%)	2 (1%)	14	12
1	C	158/160 (99%)	145 (92%)	10 (6%)	3 (2%)	9	6
All	All	473/480 (98%)	445 (94%)	19 (4%)	9 (2%)	9	6

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	39	LYS

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Mol	Chain	Res	Type
1	A	43	LYS
1	A	73	GLY
1	C	40	ASN
1	B	2	GLU
1	C	41	SER
1	C	73	GLY
1	B	42	LYS
1	A	40	ASN

### 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	146/147 (99%)	135 (92%)	11 (8%)	16	16
1	B	147/147 (100%)	132 (90%)	15 (10%)	8	8
1	C	147/147 (100%)	131 (89%)	16 (11%)	7	7
All	All	440/441 (100%)	398 (90%)	42 (10%)	10	9

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	46	THR
1	A	65	ASN
1	A	70	GLN
1	A	81	GLN
1	A	122	SER
1	A	123	THR
1	A	140	ARG
1	A	145	GLU
1	A	158	LYS
1	A	160	LYS
1	B	1	MET
1	B	2	GLU
1	B	27	THR

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Mol	Chain	Res	Type
1	B	34	HIS
1	B	40	ASN
1	B	42	LYS
1	B	65	ASN
1	B	70	GLN
1	B	72	SER
1	B	81	GLN
1	B	104	THR
1	B	119	MET
1	B	123	THR
1	B	140	ARG
1	B	160	LYS
1	C	2	GLU
1	C	11	SER
1	C	20	LEU
1	C	39	LYS
1	C	42	LYS
1	C	43	LYS
1	C	44	GLU
1	C	46	THR
1	C	51	GLU
1	C	59	SER
1	C	65	ASN
1	C	85	LYS
1	C	112	ASN
1	C	128	ILE
1	C	131	SER
1	C	160	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	HIS
1	A	34	HIS
1	A	62	HIS
1	A	65	ASN
1	A	69	HIS
1	A	107	ASN
1	B	19	HIS
1	B	34	HIS
1	B	65	ASN
1	B	107	ASN

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Mol	Chain	Res	Type
1	C	19	HIS
1	C	34	HIS
1	C	65	ASN
1	C	81	GLN
1	C	107	ASN
1	C	112	ASN

### 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 ⓘ

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	AGS	A	201	-	26,33,33	1.27	2 (7%)	22,52,52	2.08	2 (9%)
3	2W3	A	202	-	39,40,40	1.10	2 (5%)	49,57,57	2.61	21 (42%)
2	AGS	B	201	-	26,33,33	1.53	2 (7%)	22,52,52	2.28	4 (18%)
2	AGS	C	201	-	26,33,33	1.19	2 (7%)	22,52,52	2.33	2 (9%)

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	AGS	A	201	-	-	0/17/38/38	0/3/3/3
3	2W3	A	202	-	-	0/21/34/34	0/4/4/4
2	AGS	B	201	-	-	0/17/38/38	0/3/3/3
2	AGS	C	201	-	-	0/17/38/38	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	201	AGS	O4'-C1'	2.87	1.45	1.41
3	A	202	2W3	C38-N40	2.99	1.38	1.33
3	A	202	2W3	C32-N30	3.20	1.38	1.33
2	B	201	AGS	O4'-C1'	3.24	1.45	1.41
2	C	201	AGS	PG-S1G	3.24	1.96	1.90
2	A	201	AGS	O4'-C1'	3.61	1.46	1.41
2	A	201	AGS	PG-S1G	3.93	1.98	1.90
2	B	201	AGS	PG-S1G	5.62	2.01	1.90

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	201	AGS	N3-C2-N1	-10.17	120.00	128.86
2	B	201	AGS	N3-C2-N1	-8.81	121.19	128.86
2	A	201	AGS	N3-C2-N1	-8.49	121.47	128.86
3	A	202	2W3	C34-C32-N30	-4.94	119.37	124.12
3	A	202	2W3	C25-C24-CL8	-4.36	109.46	120.85
3	A	202	2W3	N8-C7-N6	-4.11	119.97	126.08
3	A	202	2W3	N30-C29-N36	-3.90	119.15	125.31
3	A	202	2W3	C4-C5-N6	-3.83	117.78	122.38
3	A	202	2W3	C20-C21-C26	-3.65	112.41	120.59
3	A	202	2W3	C13-C14-C15	-3.22	106.16	111.79
3	A	202	2W3	O39-C38-N40	-2.87	118.50	122.58
2	B	201	AGS	C1'-N9-C4	-2.80	121.80	126.64
3	A	202	2W3	C25-C26-C21	-2.37	118.89	120.46
2	A	201	AGS	O5'-PA-O1A	-2.04	101.00	109.25
3	A	202	2W3	C34-C38-N40	2.12	121.99	117.12
3	A	202	2W3	C22-C21-C26	2.13	121.58	118.53
2	C	201	AGS	O2A-PA-O1A	2.19	123.62	112.28
2	B	201	AGS	O2A-PA-O1A	2.25	123.94	112.28
2	B	201	AGS	O2'-C2'-C3'	2.39	119.49	111.83
3	A	202	2W3	C20-C21-C22	2.40	126.01	120.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	202	2W3	C26-C25-CL7	2.87	123.03	118.50
3	A	202	2W3	C14-C15-C9	2.90	112.76	109.68
3	A	202	2W3	C5-C29-N30	3.22	121.36	117.41
3	A	202	2W3	C32-N30-C29	3.67	121.63	116.17
3	A	202	2W3	C9-C7-N6	4.02	121.05	116.97
3	A	202	2W3	C29-N36-C35	4.80	123.31	115.93
3	A	202	2W3	C5-N6-C7	4.81	123.17	116.66
3	A	202	2W3	C4-C5-C29	4.83	125.85	120.17
3	A	202	2W3	C23-C24-CL8	4.91	128.57	118.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	AGS	3	0
3	A	202	2W3	5	0
2	B	201	AGS	1	0
2	C	201	AGS	3	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	159/160 (99%)	0.78	23 (14%) <b>3</b> <b>5</b>	31, 42, 50, 50	0
1	B	160/160 (100%)	0.74	19 (11%) <b>5</b> <b>9</b>	33, 46, 50, 50	1 (0%)
1	C	160/160 (100%)	0.79	17 (10%) <b>7</b> <b>11</b>	33, 48, 50, 50	0
All	All	479/480 (99%)	0.77	59 (12%) <b>5</b> <b>8</b>	31, 46, 50, 50	1 (0%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	40	ASN	10.6
1	C	41	SER	7.3
1	C	39	LYS	6.8
1	C	1	MET	5.4
1	B	40	ASN	5.4
1	B	39	LYS	5.3
1	A	40	ASN	5.2
1	A	41	SER	4.8
1	C	42	LYS	4.7
1	A	44	GLU	4.7
1	B	1	MET	4.7
1	B	41	SER	4.7
1	C	44	GLU	4.5
1	B	38	LEU	4.5
1	B	42	LYS	4.4
1	B	43	LYS	4.4
1	B	26	SER	4.4
1	B	44	GLU	4.1
1	A	88	ILE	4.0
1	A	7	VAL	3.9
1	C	3	HIS	3.8
1	B	81	GLN	3.8
1	A	8	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	43	LYS	3.6
1	C	118	TYR	3.5
1	A	36	CYS	3.5
1	A	2	GLU	3.5
1	C	8	ILE	3.1
1	A	117	LEU	3.1
1	B	71	PHE	3.1
1	A	45	GLY	3.1
1	A	87	ILE	2.8
1	A	6	ALA	2.8
1	A	86	THR	2.8
1	A	35	VAL	2.7
1	C	80	GLU	2.6
1	A	43	LYS	2.6
1	A	116	THR	2.6
1	B	155	LEU	2.6
1	A	139	TYR	2.6
1	A	37	VAL	2.5
1	C	7	VAL	2.5
1	C	6	ALA	2.4
1	B	45	GLY	2.4
1	B	8	ILE	2.4
1	C	150	TYR	2.4
1	C	87	ILE	2.4
1	A	72	SER	2.4
1	A	23	ILE	2.2
1	A	79	CYS	2.2
1	A	118	TYR	2.2
1	B	72	SER	2.2
1	B	117	LEU	2.2
1	B	119	MET	2.2
1	C	9	PRO	2.2
1	B	160	LYS	2.1
1	C	45	GLY	2.1
1	A	9	PRO	2.1
1	B	127	PHE	2.1

## 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 [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( $\text{\AA}^2$ )	Q<0.9
2	AGS	C	201	31/31	0.72	0.30	2.77	47,50,50,50	0
2	AGS	A	201	31/31	0.69	0.29	2.47	46,50,50,50	0
3	2W3	A	202	37/37	0.65	0.36	1.90	46,50,50,50	0
2	AGS	B	201	31/31	0.81	0.22	1.60	50,50,50,50	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.