



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 05:41 pm BST

PDB ID : 1WBQ  
Title : ZnMg substituted aminopeptidase P from E. coli  
Authors : Graham, S.C.; Bond, C.S.; Freeman, H.C.; Guss, J.M.  
Deposited on : 2004-11-05  
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](mailto: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.

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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 : 1.8.5 (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

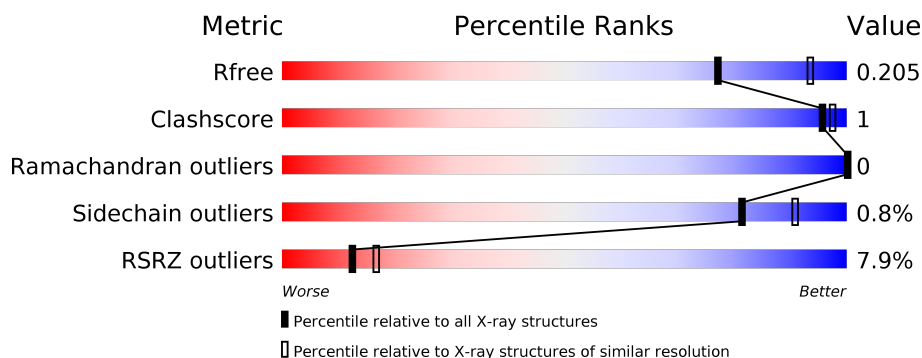
# 1 Overall quality at a glance

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  $\geq 3$ , 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	440	<div> <div>7%</div> <div>98%</div> <div>.</div> </div>
1	B	440	<div> <div>4%</div> <div>95%</div> <div>.</div> </div>
1	C	440	<div> <div>15%</div> <div>97%</div> <div>.</div> </div>
1	D	440	<div> <div>6%</div> <div>96%</div> <div>.</div> </div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 14814 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 XAA-PRO AMINOPEPTIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	0	1
			3494	2198	617	664	15			
1	B	439	Total	C	N	O	S	0	1	0
			3494	2199	615	664	16			
1	C	439	Total	C	N	O	S	0	0	0
			3490	2196	615	664	15			
1	D	439	Total	C	N	O	S	0	0	0
			3490	2196	615	664	15			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Cl 1	0	0
4	A	1	Total 1	Cl 1	0	0
4	D	1	Total 1	Cl 1	0	0
4	C	1	Total 1	Cl 1	0	0

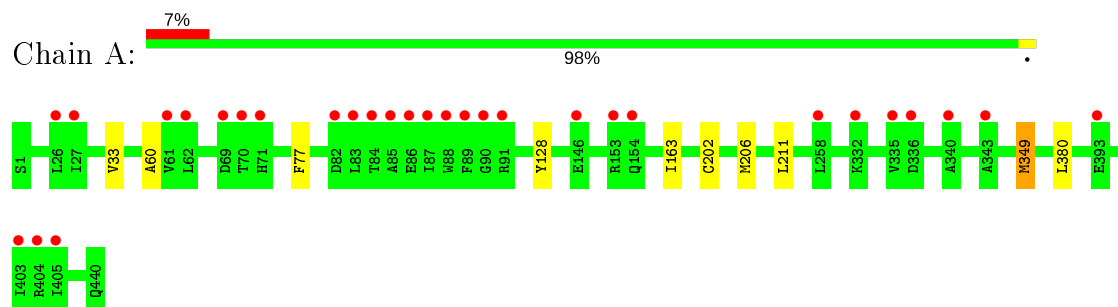
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	191	Total 191	O 191	0	0
5	B	285	Total 285	O 285	0	0
5	C	100	Total 100	O 100	0	0
5	D	256	Total 256	O 256	0	0

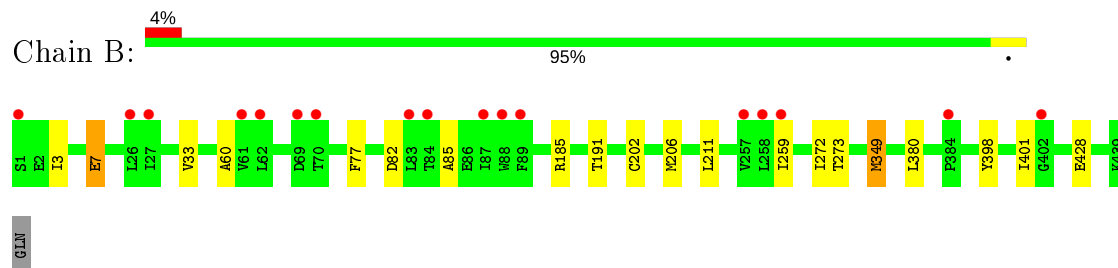
### 3 Residue-property plots [i](#)

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.

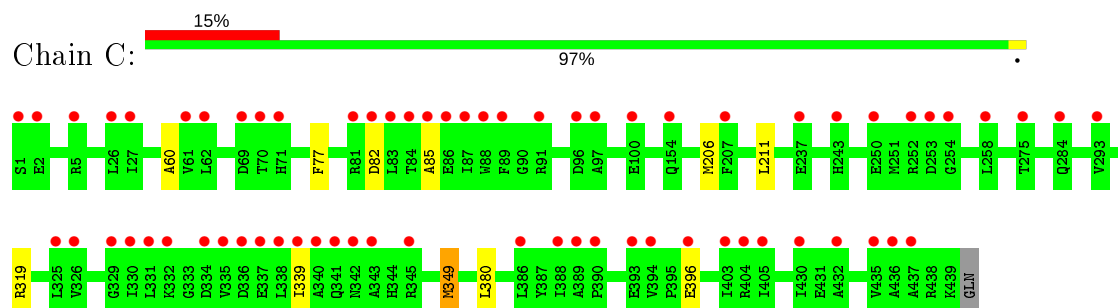
#### • Molecule 1: XAA-PRO AMINOPEPTIDASE



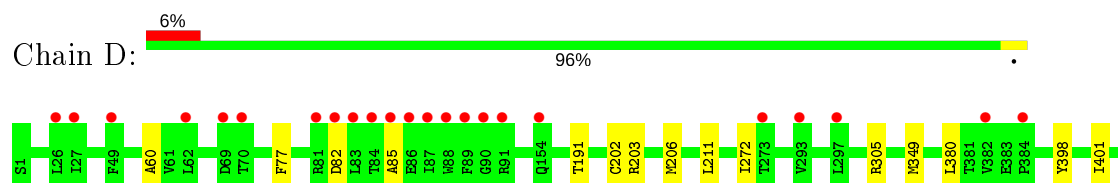
#### • Molecule 1: XAA-PRO AMINOPEPTIDASE



#### • Molecule 1: XAA-PRO AMINOPEPTIDASE



#### • Molecule 1: XAA-PRO AMINOPEPTIDASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.97Å 236.70Å 137.64Å 90.00° 106.14° 90.00°	Depositor
Resolution (Å)	29.97 – 2.30 29.97 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.97-2.30) 98.9 (29.97-2.30)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.38 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.167 , 0.197 0.177 , 0.205	Depositor DCC
$R_{free}$ test set	7512 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.9	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 40.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	14814	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% 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: MG, ZN, OCS, 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	0.45	0/3556	0.58	0/4815
1	B	0.48	0/3560	0.61	1/4820 (0.0%)
1	C	0.43	0/3552	0.55	0/4810
1	D	0.48	0/3552	0.59	0/4810
All	All	0.46	0/14220	0.58	1/19255 (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	B	185	ARG	NE-CZ-NH1	5.90	123.25	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	3494	0	3420	8	0
1	B	3494	0	3417	12	0
1	C	3490	0	3412	6	0
1	D	3490	0	3412	8	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	2	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	191	0	0	0	0
5	B	285	0	0	0	0
5	C	100	0	0	0	0
5	D	256	0	0	0	0
All	All	14814	0	13661	31	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 1.

All (31) 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:D:202:CYS:HA	1:D:206:MET:HE1	1.75	0.68
1:A:202:CYS:HA	1:A:206:MET:HE1	1.78	0.65
1:C:206:MET:HE1	1:C:211:LEU:HD21	1.82	0.61
1:B:202:CYS:HA	1:B:206:MET:HE1	1.85	0.58
1:A:33:VAL:CG1	1:B:33:VAL:CG1	2.82	0.57
1:A:206:MET:HE1	1:A:211:LEU:HD21	1.87	0.55
1:B:206:MET:HE1	1:B:211:LEU:HD21	1.89	0.55
1:A:33:VAL:HG11	1:B:33:VAL:CG1	2.37	0.54
1:A:33:VAL:CG1	1:B:33:VAL:HG11	2.39	0.53
1:D:206:MET:HE1	1:D:211:LEU:HD21	1.96	0.46
1:C:206:MET:HE1	1:C:211:LEU:CD2	2.46	0.46
1:B:398:TYR:HA	1:B:401:ILE:HD12	1.98	0.44
1:B:60:ALA:HA	1:B:77:PHE:O	2.17	0.44
1:D:82:ASP:HB3	1:D:85:ALA:HB3	1.99	0.44
1:D:202:CYS:HA	1:D:206:MET:CE	2.46	0.43
1:D:398:TYR:HA	1:D:401:ILE:HD12	2.00	0.43
1:C:60:ALA:HA	1:C:77:PHE:O	2.19	0.43
1:D:191:THR:HG22	1:D:272:ILE:HD12	2.00	0.43
1:A:349:MET:SD	1:A:349:MET:N	2.92	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:349:MET:SD	1:C:349:MET:N	2.89	0.43
1:D:60:ALA:HA	1:D:77:PHE:O	2.18	0.43
1:A:60:ALA:HA	1:A:77:PHE:O	2.19	0.42
1:B:3:ILE:HG23	1:B:7:GLU:HG2	1.99	0.42
1:C:82:ASP:HB3	1:C:85:ALA:HB3	2.01	0.42
1:D:203:ARG:H	1:D:206:MET:HE2	1.85	0.42
1:C:319:ARG:HA	1:C:339:ILE:HD11	2.02	0.42
1:A:128:TYR:HA	1:A:163:ILE:O	2.21	0.41
1:B:259:ILE:O	1:B:273:THR:HA	2.21	0.41
1:B:191:THR:HG22	1:B:272:ILE:HD12	2.03	0.40
1:B:82:ASP:HB3	1:B:85:ALA:HB3	2.03	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	437/440 (99%)	426 (98%)	11 (2%)	0	100	100
1	B	437/440 (99%)	427 (98%)	10 (2%)	0	100	100
1	C	436/440 (99%)	426 (98%)	10 (2%)	0	100	100
1	D	436/440 (99%)	425 (98%)	11 (2%)	0	100	100
All	All	1746/1760 (99%)	1704 (98%)	42 (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	369/370 (100%)	367 (100%)	2 (0%)	88	95
1	B	369/370 (100%)	364 (99%)	5 (1%)	67	81
1	C	368/370 (100%)	365 (99%)	3 (1%)	81	91
1	D	368/370 (100%)	365 (99%)	3 (1%)	81	91
All	All	1474/1480 (100%)	1461 (99%)	13 (1%)	81	89

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	349	MET
1	A	380	LEU
1	B	7	GLU
1	B	349[A]	MET
1	B	349[B]	MET
1	B	380	LEU
1	B	428	GLU
1	C	349	MET
1	C	380	LEU
1	C	396	GLU
1	D	305	ARG
1	D	349	MET
1	D	380	LEU

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 ⓘ

4 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	OCS	B	249	1	7,8,9	1.08	0	6,11,13	1.37	1 (16%)
1	OCS	A	249	1	7,8,9	0.91	0	6,11,13	1.57	1 (16%)
1	OCS	D	249	1	7,8,9	1.28	1 (14%)	6,11,13	1.65	1 (16%)
1	OCS	C	249	1	7,8,9	1.01	0	6,11,13	1.81	2 (33%)

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
1	OCS	B	249	1	-	0/4/7/9	-
1	OCS	A	249	1	-	0/4/7/9	-
1	OCS	D	249	1	-	0/4/7/9	-
1	OCS	C	249	1	-	0/4/7/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	249	OCS	CB-CA	-2.17	1.51	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	249	OCS	OD2-SG-CB	3.39	111.15	105.74
1	A	249	OCS	OD3-SG-CB	2.97	110.47	106.94
1	C	249	OCS	OD3-SG-CB	2.90	110.38	106.94
1	C	249	OCS	OD2-SG-CB	2.64	109.95	105.74
1	B	249	OCS	OD3-SG-CB	2.50	109.91	106.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 14 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 [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	439/440 (99%)	0.32	30 (6%) 17 22	29, 34, 57, 90	0
1	B	438/440 (99%)	0.15	17 (3%) 39 46	29, 34, 57, 90	0
1	C	438/440 (99%)	0.60	67 (15%) 2 3	29, 35, 57, 90	0
1	D	438/440 (99%)	0.25	25 (5%) 23 30	29, 34, 56, 90	0
All	All	1753/1760 (99%)	0.33	139 (7%) 12 17	29, 34, 57, 90	0

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	89	PHE	8.1
1	A	85	ALA	7.7
1	D	83	LEU	7.6
1	C	89	PHE	7.4
1	C	394	VAL	7.0
1	C	83	LEU	6.4
1	A	84	THR	6.4
1	D	89	PHE	6.1
1	C	84	THR	5.7
1	A	83	LEU	5.6
1	C	437	ALA	5.1
1	B	83	LEU	5.0
1	C	403	ILE	4.7
1	B	89	PHE	4.7
1	D	84	THR	4.6
1	C	1	SER	4.6
1	C	88	TRP	4.6
1	C	70	THR	4.3
1	D	85	ALA	4.3
1	B	70	THR	4.2
1	C	85	ALA	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	335	VAL	4.2
1	A	87	ILE	4.1
1	C	338	LEU	4.1
1	D	91	ARG	4.0
1	D	70	THR	4.0
1	D	87	ILE	3.9
1	A	27	ILE	3.8
1	C	87	ILE	3.8
1	C	430	ILE	3.8
1	C	97	ALA	3.7
1	C	82	ASP	3.7
1	C	336	ASP	3.6
1	A	62	LEU	3.6
1	C	252	ARG	3.5
1	B	27	ILE	3.5
1	C	334	ASP	3.5
1	A	82	ASP	3.5
1	C	340	ALA	3.4
1	C	26	LEU	3.4
1	C	284	GLN	3.4
1	C	331	LEU	3.3
1	C	62	LEU	3.3
1	C	254	GLY	3.3
1	A	26	LEU	3.3
1	B	87	ILE	3.3
1	C	405	ILE	3.3
1	A	70	THR	3.3
1	C	343	ALA	3.2
1	A	71	HIS	3.2
1	A	61	VAL	3.1
1	D	26	LEU	3.1
1	D	382	VAL	3.1
1	C	96	ASP	3.1
1	A	332	LYS	3.1
1	B	69	ASP	3.0
1	A	91	ARG	3.0
1	A	340	ALA	3.0
1	D	88	TRP	3.0
1	C	337	GLU	3.0
1	A	86	GLU	3.0
1	C	237	GLU	2.9
1	C	329	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	153	ARG	2.9
1	A	405	ILE	2.9
1	C	275	THR	2.9
1	D	86	GLU	2.9
1	C	325	LEU	2.9
1	C	332	LYS	2.9
1	D	90	GLY	2.8
1	B	26	LEU	2.8
1	A	90	GLY	2.8
1	D	82	ASP	2.8
1	C	393	GLU	2.7
1	D	49	PHE	2.7
1	B	1	SER	2.7
1	C	341	GLN	2.7
1	A	403	ILE	2.7
1	C	27	ILE	2.7
1	B	84	THR	2.7
1	D	62	LEU	2.7
1	B	259	ILE	2.6
1	C	339	ILE	2.6
1	D	27	ILE	2.6
1	C	2	GLU	2.6
1	C	69	ASP	2.6
1	C	432	ALA	2.6
1	B	402	GLY	2.6
1	C	5	ARG	2.6
1	C	386	LEU	2.6
1	A	69	ASP	2.5
1	D	384	PRO	2.5
1	B	88	TRP	2.5
1	C	243	HIS	2.5
1	A	88	TRP	2.5
1	C	207	PHE	2.5
1	A	335	VAL	2.5
1	C	86	GLU	2.4
1	C	396	GLU	2.4
1	D	69	ASP	2.4
1	B	257	VAL	2.4
1	A	146	GLU	2.4
1	B	61	VAL	2.4
1	A	393	GLU	2.3
1	C	436	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	273	THR	2.3
1	C	253	ASP	2.3
1	B	62	LEU	2.3
1	C	91	ARG	2.3
1	C	389	ALA	2.3
1	A	336	ASP	2.3
1	C	330	ILE	2.3
1	D	154	GLN	2.3
1	A	404	ARG	2.3
1	D	405	ILE	2.3
1	D	81	ARG	2.2
1	B	384	PRO	2.2
1	C	250	GLU	2.2
1	C	258	LEU	2.2
1	C	293	VAL	2.2
1	D	293	VAL	2.2
1	D	439	LYS	2.2
1	C	388	ILE	2.1
1	B	258	LEU	2.1
1	C	404	ARG	2.1
1	A	343	ALA	2.1
1	C	154	GLN	2.1
1	A	258	LEU	2.1
1	C	390	PRO	2.1
1	C	61	VAL	2.1
1	A	154	GLN	2.1
1	C	342	ASN	2.1
1	C	71	HIS	2.1
1	C	100	GLU	2.1
1	C	345	ARG	2.1
1	C	326	VAL	2.1
1	C	435	VAL	2.1
1	D	297	LEU	2.0
1	C	81	ARG	2.0

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

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, 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	B-factors( $\text{\AA}^2$ )	Q<0.9
1	OCS	C	249	9/10	0.94	0.13	26,31,35,42	3
1	OCS	A	249	9/10	0.96	0.14	26,31,36,44	3
1	OCS	B	249	9/10	0.98	0.09	26,31,36,43	3
1	OCS	D	249	9/10	0.99	0.09	24,31,35,44	3

### 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, 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	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MG	C	1441	1/1	0.90	0.14	60,60,60,60	0
3	MG	A	1441	1/1	0.91	0.06	31,31,31,31	0
3	MG	B	1441	1/1	0.93	0.05	26,26,26,26	0
3	MG	D	1442	1/1	0.94	0.04	38,38,38,38	0
4	CL	C	1442	1/1	0.95	0.09	48,48,48,48	0
3	MG	B	1442	1/1	0.96	0.03	37,37,37,37	0
4	CL	D	1443	1/1	0.98	0.05	45,45,45,45	0
2	ZN	C	1440	1/1	0.98	0.04	67,67,67,67	0
3	MG	D	1441	1/1	0.98	0.03	29,29,29,29	0
4	CL	A	1442	1/1	0.98	0.08	49,49,49,49	0
4	CL	B	1443	1/1	0.99	0.06	38,38,38,38	0
2	ZN	D	1440	1/1	0.99	0.08	46,46,46,46	0
2	ZN	A	1440	1/1	0.99	0.05	53,53,53,53	0
2	ZN	B	1440	1/1	1.00	0.04	44,44,44,44	0

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

There are no such residues in this entry.