



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:49 AM GMT

PDB ID : 2JGO
Title : Structure of the arsenated de novo designed peptide Coil Ser L9C
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Deposited on : 2007-02-13
Resolution : 1.81 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

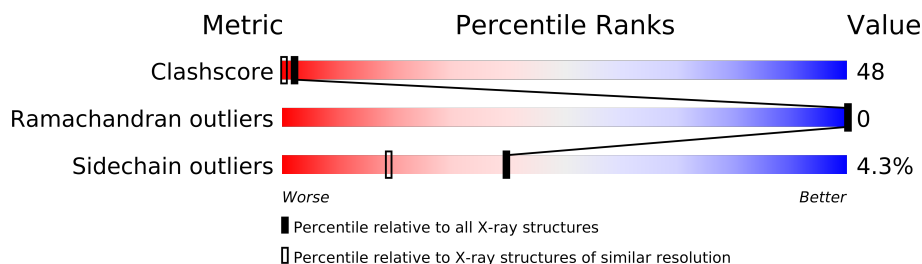
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.81 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))
Clashscore	79885	5140 (1.84-1.80)
Ramachandran outliers	78287	5077 (1.84-1.80)
Sidechain outliers	78261	5077 (1.84-1.80)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	31	
1	B	31	
1	C	31	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 805 atoms, of which 0 are hydrogen and 0 are deuterium.

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 COIL SER L9C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	31	Total	C	N	O	S	5	4	1
			261	173	41	46	1			
1	B	31	Total	C	N	O	S	2	2	1
			246	158	41	46	1			
1	C	31	Total	C	N	O	S	0	1	1
			240	153	40	46	1			

- Molecule 2 is ARSENIC (three-letter code: ARS) (formula: As).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	As	0	0
			1	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Zn	0	0
			2	2		
3	A	2	Total	Zn	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	19	Total	O	0	0
			19	19		
4	B	16	Total	O	0	0
			16	16		
4	C	18	Total	O	0	0
			18	18		

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Note EDS was not executed.

- Chain A:



Chain B:



Chain C:



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	77.28Å 29.42Å 44.20Å 90.00° 119.49° 90.00°	Depositor
Resolution (Å)	13.74 – 1.81	Depositor
% Data completeness (in resolution range)	99.2 (13.74-1.81)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 6.0	Depositor
R, R_{free}	0.198 , 0.257	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	805	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, ARS, ACE, NH2

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	1.28	2/274 (0.7%)	1.22	2/366 (0.5%)
1	B	0.43	0/251	1.08	1/332 (0.3%)
1	C	0.42	0/242	1.22	2/321 (0.6%)
All	All	0.84	2/767 (0.3%)	1.18	5/1019 (0.5%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	0	ACE	C-N	-17.13	0.94	1.34
1	A	1	GLU	C-N	-8.16	1.15	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	0	ACE	O-C-N	9.23	137.47	122.70
1	A	1	GLU	O-C-N	-6.43	112.41	122.70
1	C	20	GLU	CG-CD-OE2	-6.18	105.94	118.30
1	B	1	GLU	CA-CB-CG	-5.80	100.65	113.40
1	C	23	LEU	CB-CG-CD1	5.34	120.08	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	261	0	280	43	1
1	B	246	0	264	38	1
1	C	240	0	251	40	0
2	A	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	19	0	0	3	0
4	B	16	0	0	2	0
4	C	18	0	0	8	0
All	All	805	0	795	73	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 48.

All (73) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:2[A]:TRP:CZ3	1:B:5:LEU:HD12	1.67	1.28
1:C:22[B]:LYS:NZ	4:C:2015:HOH:O	1.73	1.19
1:B:2:TRP:CZ3	1:C:2:TRP:CZ3	2.32	1.17
1:B:2:TRP:HZ3	1:C:2:TRP:CZ3	1.65	1.12
1:B:2:TRP:CZ3	1:C:2:TRP:HZ3	1.65	1.11
1:A:26[A]:LEU:HD12	1:C:23:LEU:HD21	1.16	1.09
1:B:2:TRP:CH2	1:C:2:TRP:CE3	2.44	1.06
1:A:26[A]:LEU:CD1	1:C:23:LEU:HD21	1.90	1.01
1:A:26[A]:LEU:CD1	1:C:23:LEU:CD2	2.38	1.00
1:A:18:ALA:O	1:A:22[B]:LYS:HD3	1.63	0.99
1:B:24:GLU:OE1	4:C:2015:HOH:O	1.80	0.97
1:B:2:TRP:CH2	1:C:2:TRP:HE3	1.83	0.96
1:A:2[A]:TRP:CH2	1:B:5:LEU:HD12	2.00	0.95
1:A:26[A]:LEU:HD12	1:C:23:LEU:CD2	1.94	0.95
1:A:19:LEU:CD2	1:C:23:LEU:HD12	1.99	0.92
1:B:2:TRP:HZ3	1:C:2:TRP:HZ3	0.97	0.91
1:C:6:GLU:OE1	4:C:2006:HOH:O	1.93	0.86
1:B:2:TRP:CH2	1:C:2:TRP:CZ3	2.61	0.86
1:B:2:TRP:HH2	1:C:2:TRP:CE3	1.91	0.86
1:B:27:GLU:OE2	4:C:2015:HOH:O	1.94	0.84
1:B:20:GLU:HG2	1:C:19:LEU:HD11	1.60	0.83
1:A:2[A]:TRP:HZ2	1:B:1:GLU:HG2	1.45	0.82
1:B:2:TRP:CZ3	1:C:2:TRP:CE3	2.66	0.81
1:A:19:LEU:HD21	1:C:23:LEU:HD12	1.60	0.80

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:20:GLU:CG	1:C:19:LEU:HD11	2.12	0.80
1:A:20:GLU:HG2	1:B:19:LEU:HD21	1.64	0.79
1:A:20:GLU:CG	1:B:19:LEU:HD11	2.13	0.78
1:A:2[A]:TRP:HZ3	1:B:5:LEU:HD12	1.40	0.77
1:A:2[A]:TRP:CZ2	1:B:1:GLU:HG2	2.20	0.76
1:A:19:LEU:HD23	1:C:23:LEU:HD12	1.69	0.74
1:C:7:LYS:NZ	4:C:2008:HOH:O	2.14	0.74
1:A:2[A]:TRP:CZ3	1:B:5:LEU:CD1	2.62	0.71
1:A:20:GLU:HG3	1:B:19:LEU:HD11	1.71	0.71
1:A:1:GLU:OE2	4:C:2006:HOH:O	2.11	0.69
1:A:20:GLU:OE1	4:A:2013:HOH:O	2.11	0.67
1:A:20:GLU:HG2	1:B:19:LEU:HD11	1.77	0.65
1:B:17:GLN:NE2	4:B:2011:HOH:O	2.30	0.64
1:A:16[A]:LEU:HD11	1:B:19:LEU:HD12	1.82	0.62
1:A:22[B]:LYS:HD3	1:A:22[B]:LYS:H	1.67	0.59
1:B:20:GLU:HG2	1:C:19:LEU:HD21	1.84	0.58
1:A:26[A]:LEU:CD1	1:C:23:LEU:HD22	2.30	0.58
1:A:16[A]:LEU:CD1	1:B:19:LEU:HD12	2.35	0.56
1:A:26[A]:LEU:HD11	1:C:23:LEU:CD2	2.36	0.55
1:C:15:LYS:HG3	4:C:2009:HOH:O	2.07	0.55
1:B:20:GLU:HG2	1:C:19:LEU:CD1	2.34	0.54
1:A:26[A]:LEU:HD11	1:C:23:LEU:HD22	1.89	0.54
1:A:21:LYS:HE3	4:A:2014:HOH:O	2.07	0.53
1:C:2:TRP:HA	1:C:2:TRP:CE3	2.42	0.53
1:A:22[B]:LYS:CD	1:A:22[B]:LYS:H	2.23	0.50
1:B:23:LEU:HD21	1:C:22[B]:LYS:HB3	1.94	0.50
1:B:20:GLU:HG3	1:C:19:LEU:HD11	1.92	0.49
1:C:1:GLU:OE1	4:C:2003:HOH:O	2.19	0.49
1:A:2[B]:TRP:CZ2	1:B:1:GLU:HG2	2.48	0.48
1:A:19:LEU:HD23	1:C:23:LEU:CD1	2.42	0.48
1:A:5:LEU:HD23	1:C:2:TRP:CH2	2.50	0.47
1:A:15:LYS:HB3	1:C:16:LEU:HD11	1.97	0.47
1:A:18:ALA:O	1:A:22[B]:LYS:CD	2.50	0.47
1:B:24:GLU:OE1	1:C:22[B]:LYS:NZ	2.37	0.46
1:A:18:ALA:O	1:A:22[A]:LYS:HG3	2.17	0.45
1:A:20:GLU:HG2	1:B:19:LEU:CD2	2.43	0.45
1:A:2[A]:TRP:HZ3	1:B:5:LEU:CD1	2.18	0.44
1:B:20:GLU:CD	4:B:2013:HOH:O	2.46	0.44
1:B:16:LEU:O	1:B:20:GLU:HG3	2.18	0.44
1:B:5:LEU:HD13	1:C:5:LEU:HD21	2.00	0.43
1:C:16:LEU:O	1:C:16:LEU:HD23	2.19	0.43
1:A:30:NH2:N	4:A:2018:HOH:O	2.51	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:15:LYS:CB	1:C:16:LEU:HD11	2.50	0.42
1:C:19:LEU:HD23	1:C:19:LEU:HA	1.86	0.42
1:A:5:LEU:CD2	1:C:2:TRP:CH2	3.02	0.42
1:A:2[A]:TRP:HZ2	1:B:1:GLU:CG	2.25	0.40

All (1) 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	Distance(Å)	Clash(Å)
1:A:22[B]:LYS:NZ	1:B:30:NH2:N[4_454]	1.75	0.45

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	33/31 (106%)	33 (100%)	0	0	100	100
1	B	31/31 (100%)	31 (100%)	0	0	100	100
1	C	30/31 (97%)	30 (100%)	0	0	100	100
All	All	94/93 (101%)	94 (100%)	0	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	27/23 (117%)	26 (96%)	1 (4%)	45	25
1	B	25/23 (109%)	24 (96%)	1 (4%)	42	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	24/23 (104%)	23 (96%)	1 (4%)	40	19
All	All	76/69 (110%)	73 (96%)	3 (4%)	40	22

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	B	5	LEU
1	C	23	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	17	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 5 ligands modelled in this entry, 5 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.

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.