Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:49 PM GMT

PDB ID	:	4Q20
Title	:	Crystal structure of a C-terminal part of tyrosine kinase (DivL) from Caulobacter crescentus CB15 at 2.50 A resolution (PSI Community Target, Shapiro)
Authors	:	Joint Center for Structural Genomics (JCSG)
Deposited on	:	2014-04-04
Resolution	:	2.50 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
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

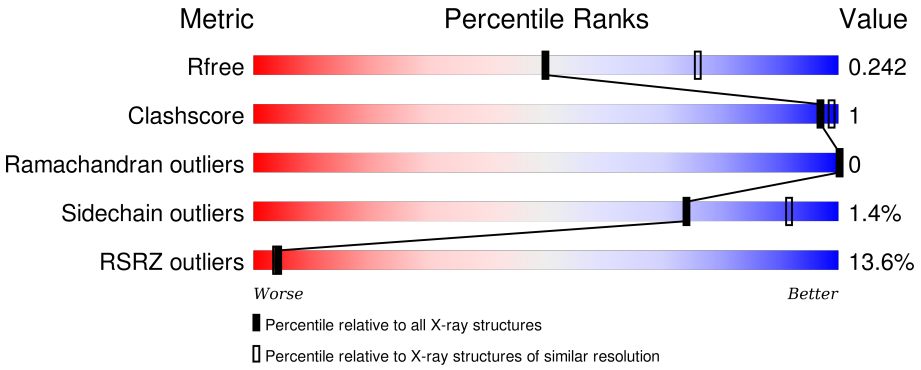
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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. 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 <=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	268	9% 82% · 14%
1	B	268	15% 80% · 16%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3430 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 Sensor protein DivL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	0	0
			1720	1064	318	334	4			
1	B	225	Total	C	N	O	S	0	1	0
			1691	1046	311	330	4			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	EXPRESSION TAG	UNP Q9RQQ9
A	-19	GLY	-	EXPRESSION TAG	UNP Q9RQQ9
A	-18	SER	-	EXPRESSION TAG	UNP Q9RQQ9
A	-17	SER	-	EXPRESSION TAG	UNP Q9RQQ9
A	-16	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
A	-15	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
A	-14	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
A	-13	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
A	-12	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
A	-11	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
A	-10	SER	-	EXPRESSION TAG	UNP Q9RQQ9
A	-9	SER	-	EXPRESSION TAG	UNP Q9RQQ9
A	-8	GLY	-	EXPRESSION TAG	UNP Q9RQQ9
A	-7	LEU	-	EXPRESSION TAG	UNP Q9RQQ9
A	-6	VAL	-	EXPRESSION TAG	UNP Q9RQQ9
A	-5	PRO	-	EXPRESSION TAG	UNP Q9RQQ9
A	-4	ARG	-	EXPRESSION TAG	UNP Q9RQQ9
A	-3	GLY	-	EXPRESSION TAG	UNP Q9RQQ9
A	-2	SER	-	EXPRESSION TAG	UNP Q9RQQ9
A	-1	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
A	0	MET	-	EXPRESSION TAG	UNP Q9RQQ9
B	-20	MET	-	EXPRESSION TAG	UNP Q9RQQ9
B	-19	GLY	-	EXPRESSION TAG	UNP Q9RQQ9
B	-18	SER	-	EXPRESSION TAG	UNP Q9RQQ9
B	-17	SER	-	EXPRESSION TAG	UNP Q9RQQ9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
B	-15	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
B	-14	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
B	-13	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
B	-12	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
B	-11	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
B	-10	SER	-	EXPRESSION TAG	UNP Q9RQQ9
B	-9	SER	-	EXPRESSION TAG	UNP Q9RQQ9
B	-8	GLY	-	EXPRESSION TAG	UNP Q9RQQ9
B	-7	LEU	-	EXPRESSION TAG	UNP Q9RQQ9
B	-6	VAL	-	EXPRESSION TAG	UNP Q9RQQ9
B	-5	PRO	-	EXPRESSION TAG	UNP Q9RQQ9
B	-4	ARG	-	EXPRESSION TAG	UNP Q9RQQ9
B	-3	GLY	-	EXPRESSION TAG	UNP Q9RQQ9
B	-2	SER	-	EXPRESSION TAG	UNP Q9RQQ9
B	-1	HIS	-	EXPRESSION TAG	UNP Q9RQQ9
B	0	MET	-	EXPRESSION TAG	UNP Q9RQQ9

- Molecule 2 is water.

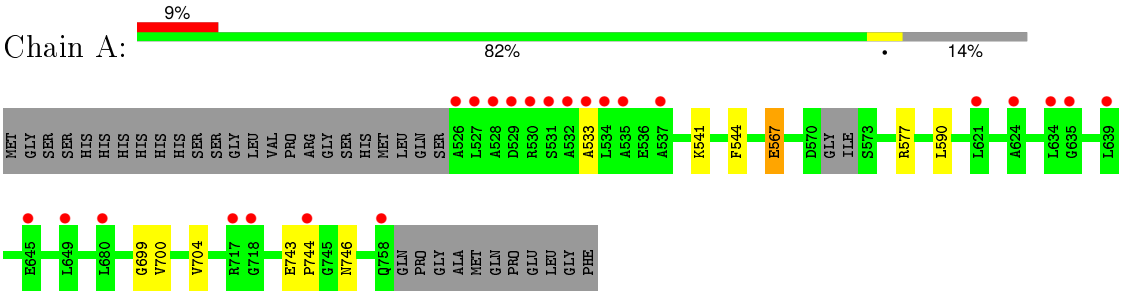
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	12	Total O 12 12	0	0
2	B	7	Total O 7 7	0	0

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 errors 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.

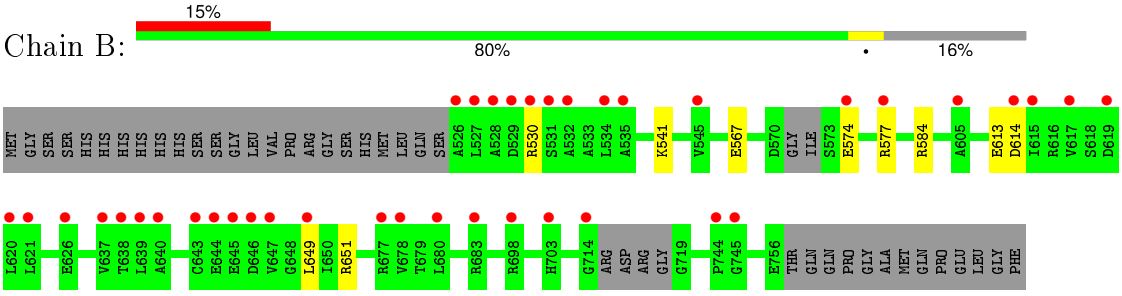
- Molecule 1: Sensor protein DivL

Chain A:



- Molecule 1: Sensor protein DivL

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	69.53Å 69.53Å 194.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.11 – 2.50 44.11 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.1 (44.11-2.50) 99.1 (44.11-2.50)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.07 (at 2.51Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.202 , 0.232 0.213 , 0.242	Depositor DCC
R_{free} test set	965 reflections (5.22%)	DCC
Wilson B-factor (Å^2)	80.2	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 77.0	EDS
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 19453 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3430	wwPDB-VP
Average B, all atoms (Å^2)	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.51	0/1740	0.64	0/2357
1	B	0.49	0/1713	0.65	0/2318
All	All	0.50	0/3453	0.64	0/4675

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

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	1720	0	1706	7	0
1	B	1691	0	1680	6	0
2	A	12	0	0	0	0
2	B	7	0	0	0	0
All	All	3430	0	3386	9	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 (9) 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:574:GLU:HG2	1:B:577:ARG:HH21	1.64	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:567:GLU:CD	1:B:584:ARG:HE	2.15	0.50
1:A:743:GLU:HB3	1:A:746:ASN:HB2	1.94	0.49
1:A:699:GLY:HA3	1:A:744:PRO:HA	1.95	0.47
1:A:533:ALA:HB1	1:B:530:ARG:HA	1.96	0.47
1:B:614:ASP:HB3	1:B:649:LEU:HD21	2.02	0.42
1:A:544:PHE:CE2	1:B:541:LYS:HG3	2.56	0.40
1:A:577:ARG:NH1	1:B:567:GLU:O	2.54	0.40
1:A:700:VAL:CG1	1:A:704:VAL:HG23	2.51	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	227/268 (85%)	222 (98%)	5 (2%)	0	100	100
1	B	220/268 (82%)	214 (97%)	6 (3%)	0	100	100
All	All	447/536 (83%)	436 (98%)	11 (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	174/210 (83%)	171 (98%)	3 (2%)	68	89

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	173/210 (82%)	171 (99%)	2 (1%)	78 93
All	All	347/420 (83%)	342 (99%)	5 (1%)	74 91

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	541	LYS
1	A	567	GLU
1	A	590	LEU
1	B	613	GLU
1	B	651	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

There are no ligands in this entry.

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^{th} 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^2)	Q<0.9
1	A	231/268 (86%)	0.66	23 (9%)	9	10	60, 98, 145, 181	0
1	B	225/268 (83%)	1.02	39 (17%)	2	2	62, 112, 160, 182	0
All	All	456/536 (85%)	0.84	62 (13%)	4	4	60, 103, 157, 182	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	526	ALA	8.9
1	B	528	ALA	7.5
1	B	534	LEU	7.0
1	B	531	SER	6.5
1	B	527	LEU	6.0
1	A	635	GLY	5.5
1	A	744	PRO	5.5
1	B	744	PRO	5.0
1	B	530	ARG	4.8
1	A	717	ARG	4.3
1	B	647	VAL	4.3
1	B	645	GLU	4.2
1	A	526	ALA	4.2
1	B	619	ASP	4.1
1	A	718	GLY	4.0
1	B	529	ASP	4.0
1	A	531	SER	3.7
1	A	527	LEU	3.7
1	A	639	LEU	3.7
1	B	620	LEU	3.7
1	A	758	GLN	3.6
1	B	617	VAL	3.6
1	A	532	ALA	3.4
1	B	643	CYS	3.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	698	ARG	3.3
1	B	535	ALA	3.2
1	B	644	GLU	3.2
1	B	677	ARG	3.1
1	B	640	ALA	3.0
1	B	626	GLU	3.0
1	B	680	LEU	2.9
1	B	574	GLU	2.8
1	A	534	LEU	2.8
1	A	645	GLU	2.8
1	A	621	LEU	2.7
1	B	683	ARG	2.7
1	B	646	ASP	2.7
1	A	649	LEU	2.7
1	A	530	ARG	2.7
1	A	533	ALA	2.6
1	B	703	HIS	2.6
1	B	714	GLY	2.6
1	B	614	ASP	2.5
1	B	605	ALA	2.5
1	A	528	ALA	2.5
1	B	638	THR	2.4
1	B	649	LEU	2.4
1	B	621	LEU	2.3
1	A	529	ASP	2.3
1	B	545	VAL	2.3
1	A	535	ALA	2.3
1	B	615	ILE	2.3
1	B	745	GLY	2.2
1	A	634	LEU	2.2
1	A	680	LEU	2.2
1	B	639	LEU	2.2
1	A	537	ALA	2.2
1	B	577	ARG	2.2
1	B	532	ALA	2.1
1	A	624	ALA	2.1
1	B	637	VAL	2.1
1	B	678	VAL	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

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.