



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2020 – 11:21 AM EDT

PDB ID : 6E6U
Title : Variant C89S of Dieckmann cyclase, NcmC
Authors : Cogan, D.P.; Nair, S.K.
Deposited on : 2018-07-25
Resolution : 1.55 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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 : 2.14.6
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) : 2.14.6

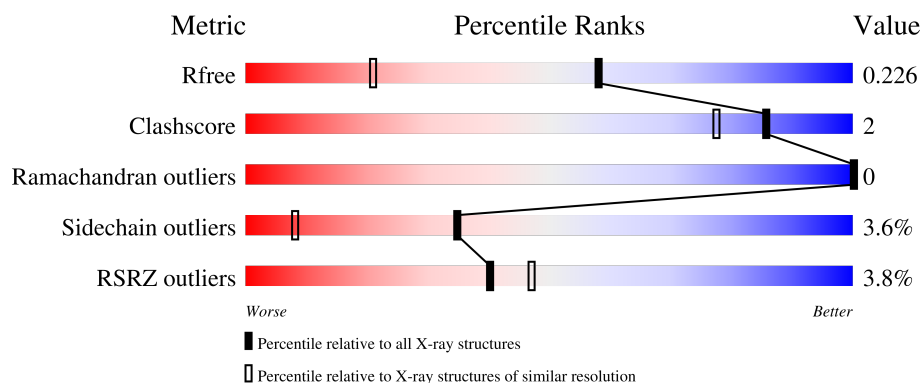
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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 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	275	<div> <div>6%</div> <div> <div></div> <div>87%</div> <div>8%</div> <div>..</div> </div> </div>
1	B	275	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div>..</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4465 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 Dieckmann cyclase, NcmC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	0	0
			1972	1238	352	379	3			
1	B	266	Total	C	N	O	S	0	0	0
			1964	1234	352	375	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP A0A1X9WEN9
A	-1	ASN	-	expression tag	UNP A0A1X9WEN9
A	0	ALA	-	expression tag	UNP A0A1X9WEN9
A	89	SER	CYS	engineered mutation	UNP A0A1X9WEN9
B	-2	SER	-	expression tag	UNP A0A1X9WEN9
B	-1	ASN	-	expression tag	UNP A0A1X9WEN9
B	0	ALA	-	expression tag	UNP A0A1X9WEN9
B	89	SER	CYS	engineered mutation	UNP A0A1X9WEN9

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

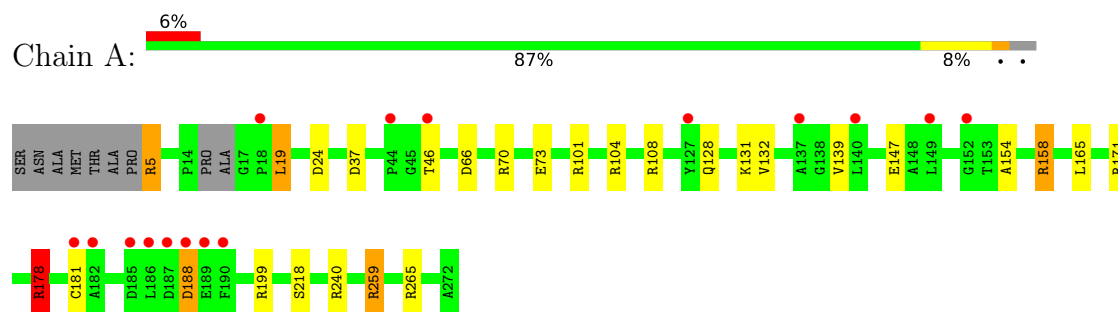
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	245	Total	O	0	0
			245	245		
3	B	274	Total	O	0	0
			274	274		

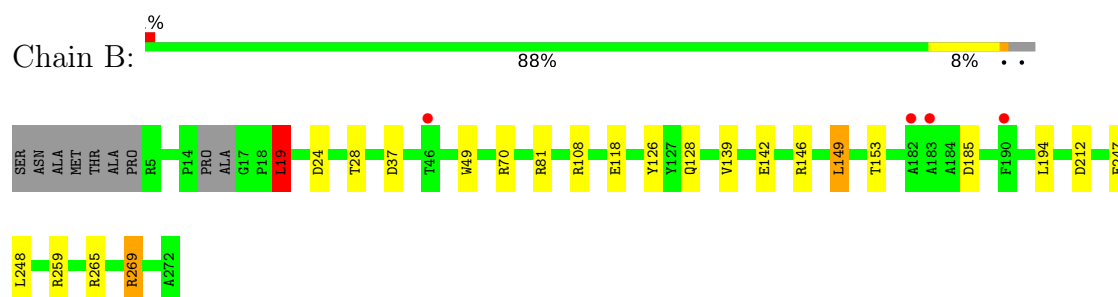
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Dieckmann cyclase, NcmC



- Molecule 1: Dieckmann cyclase, NcmC



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	80.36Å 80.36Å 78.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	80.36 – 1.55 56.82 – 1.55	Depositor EDS
% Data completeness (in resolution range)	99.4 (80.36-1.55) 99.4 (56.82-1.55)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.173 , 0.215 0.187 , 0.226	Depositor DCC
R_{free} test set	3570 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	18.0	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.004 for -h,-l,-k 0.000 for -h,l,k 0.000 for l,-k,h 0.007 for -l,-k,-h 0.038 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4465	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.00 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8573e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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: SO4

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	1.06	3/2011 (0.1%)	1.13	17/2743 (0.6%)
1	B	1.11	4/2003 (0.2%)	1.21	18/2732 (0.7%)
All	All	1.09	7/4014 (0.2%)	1.17	35/5475 (0.6%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	70	ARG	CZ-NH1	6.89	1.42	1.33
1	A	73	GLU	CD-OE1	-6.06	1.19	1.25
1	B	118	GLU	CD-OE1	5.90	1.32	1.25
1	B	247	GLU	CD-OE1	5.27	1.31	1.25
1	B	118	GLU	CD-OE2	5.14	1.31	1.25
1	A	199	ARG	CZ-NH2	5.13	1.39	1.33
1	A	218	SER	CB-OG	-5.12	1.35	1.42

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	259	ARG	NE-CZ-NH1	11.31	125.96	120.30
1	A	259	ARG	NE-CZ-NH1	10.38	125.49	120.30
1	B	108	ARG	NE-CZ-NH1	8.96	124.78	120.30
1	A	70	ARG	NE-CZ-NH2	-8.65	115.97	120.30
1	B	24	ASP	CB-CG-OD1	8.39	125.85	118.30
1	A	24	ASP	CB-CG-OD1	8.02	125.52	118.30
1	B	70	ARG	NE-CZ-NH1	7.96	124.28	120.30
1	A	259	ARG	NE-CZ-NH2	-7.82	116.39	120.30
1	B	108	ARG	NE-CZ-NH2	-7.75	116.42	120.30
1	B	70	ARG	NE-CZ-NH2	-7.72	116.44	120.30
1	A	19	LEU	CA-CB-CG	7.10	131.63	115.30
1	A	70	ARG	NE-CZ-NH1	6.98	123.79	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	81	ARG	NE-CZ-NH1	6.86	123.73	120.30
1	A	240	ARG	NE-CZ-NH2	-6.79	116.91	120.30
1	B	24	ASP	CB-CG-OD2	-6.79	112.19	118.30
1	B	259	ARG	NE-CZ-NH2	-6.64	116.98	120.30
1	A	101	ARG	NE-CZ-NH2	-6.43	117.08	120.30
1	B	269	ARG	NE-CZ-NH1	6.33	123.46	120.30
1	B	265	ARG	NE-CZ-NH2	-6.27	117.17	120.30
1	A	178	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	A	24	ASP	CB-CG-OD2	-6.19	112.73	118.30
1	B	19	LEU	CA-CB-CG	6.16	129.47	115.30
1	B	247	GLU	OE1-CD-OE2	-6.10	115.98	123.30
1	B	81	ARG	NE-CZ-NH2	-6.04	117.28	120.30
1	B	212	ASP	CB-CG-OD2	-5.97	112.93	118.30
1	A	104	ARG	NE-CZ-NH1	5.73	123.17	120.30
1	B	146	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	A	37	ASP	CB-CG-OD2	5.52	123.27	118.30
1	A	165	LEU	CB-CG-CD2	5.33	120.07	111.00
1	B	37	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	265	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	B	259	ARG	CD-NE-CZ	5.21	130.89	123.60
1	A	5	ARG	NE-CZ-NH1	5.19	122.90	120.30
1	A	66	ASP	CB-CG-OD2	-5.16	113.66	118.30
1	A	171	ARG	NE-CZ-NH2	-5.07	117.77	120.30

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	1972	0	1939	9	0
1	B	1964	0	1932	8	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	245	0	0	4	0
3	B	274	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4465	0	3871	17	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 2.

All (17) 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:126:TYR:OH	1:B:153:THR:HG22	1.58	1.04
1:B:149:LEU:O	1:B:153:THR:HG23	1.87	0.73
1:A:46:THR:HB	3:A:481:HOH:O	1.92	0.68
1:A:259:ARG:HG3	3:A:525:HOH:O	1.93	0.68
1:B:185:ASP:CB	3:B:634:HOH:O	2.42	0.66
1:A:154:ALA:O	1:A:158:ARG:HG2	1.99	0.62
1:B:139:VAL:HG11	1:B:194:LEU:HD11	1.83	0.60
1:A:46:THR:CB	3:A:481:HOH:O	2.49	0.57
1:B:126:TYR:HH	1:B:153:THR:HG22	1.70	0.55
1:B:19:LEU:HD11	1:B:49:TRP:CZ2	2.46	0.51
1:A:178:ARG:HH11	1:A:178:ARG:CG	2.25	0.50
1:A:128:GLN:O	1:A:132:VAL:HG23	2.14	0.46
1:B:19:LEU:HD11	1:B:49:TRP:CE2	2.55	0.42
1:A:147:GLU:HB3	3:A:586:HOH:O	2.20	0.41
1:A:181:CYS:SG	1:A:188:ASP:CB	3.08	0.41
1:B:28:THR:HG23	3:B:610:HOH:O	2.20	0.41
1:A:178:ARG:HH11	1:A:178:ARG:HG3	1.86	0.41

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	262/275 (95%)	259 (99%)	3 (1%)	0	100	100
1	B	262/275 (95%)	257 (98%)	5 (2%)	0	100	100
All	All	524/550 (95%)	516 (98%)	8 (2%)	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	193/199 (97%)	185 (96%)	8 (4%)	30	5
1	B	191/199 (96%)	185 (97%)	6 (3%)	40	11
All	All	384/398 (96%)	370 (96%)	14 (4%)	35	8

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ARG
1	A	19	LEU
1	A	108	ARG
1	A	131	LYS
1	A	139	VAL
1	A	158	ARG
1	A	178	ARG
1	A	188	ASP
1	B	19	LEU
1	B	128	GLN
1	B	142	GLU
1	B	149	LEU
1	B	248	LEU
1	B	269	ARG

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

Mol	Chain	Res	Type
1	B	253	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	SO4	A	301	-	4,4,4	0.47	0	6,6,6	0.69	0
2	SO4	B	301	-	4,4,4	0.29	0	6,6,6	1.25	0

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, 95th 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	266/275 (96%)	-0.04	16 (6%) 21 25	12, 21, 44, 66	0
1	B	266/275 (96%)	-0.24	4 (1%) 73 78	11, 19, 42, 54	0
All	All	532/550 (96%)	-0.14	20 (3%) 40 47	11, 20, 44, 66	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	183	ALA	6.9
1	A	187	ASP	5.5
1	A	127	TYR	4.1
1	A	182	ALA	4.1
1	B	182	ALA	4.0
1	A	137	ALA	3.4
1	A	188	ASP	3.4
1	A	190	PHE	2.8
1	A	149	LEU	2.8
1	B	46	THR	2.5
1	A	186	LEU	2.5
1	A	46	THR	2.5
1	A	140	LEU	2.4
1	B	190	PHE	2.3
1	A	189	GLU	2.1
1	A	18	PRO	2.1
1	A	185	ASP	2.1
1	A	181	CYS	2.1
1	A	152	GLY	2.0
1	A	44	PRO	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 monosaccharides 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, 95th 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(\AA^2)	Q<0.9
2	SO4	A	301	5/5	0.97	0.06	29,29,32,35	0
2	SO4	B	301	5/5	0.99	0.05	25,26,27,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.