



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:58 AM GMT

PDB ID : 2F31
Title : Crystal structure of the autoinhibitory switch in Formin mDia1; the DID/DAD complex
Authors : Nezami, A.G.; Poy, F.; Eck, M.J.
Deposited on : 2005-11-18
Resolution : 2.10 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

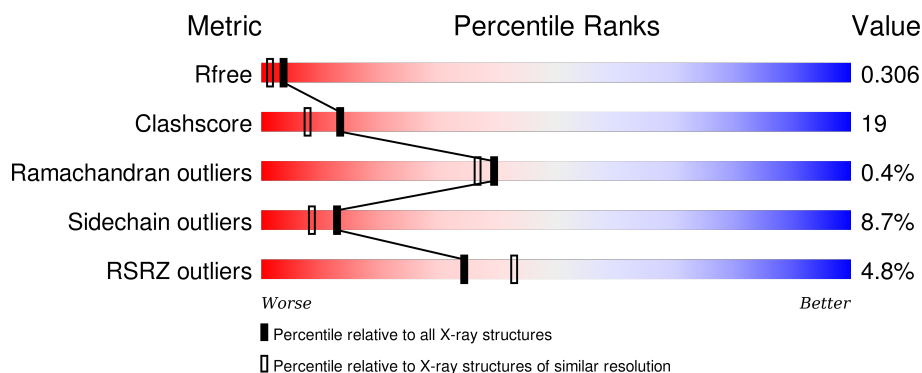
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


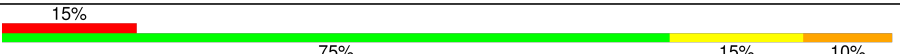
The reported resolution of this entry is 2.10 Å.

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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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 $\leq 5\%$. 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	233	
2	B	20	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2087 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 Diaphanous protein homolog 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1816	1140	311	347	18			

- Molecule 2 is a protein called Diaphanous protein homolog 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	20	Total	C	N	O	S	0	0	0
			146	89	24	32	1			

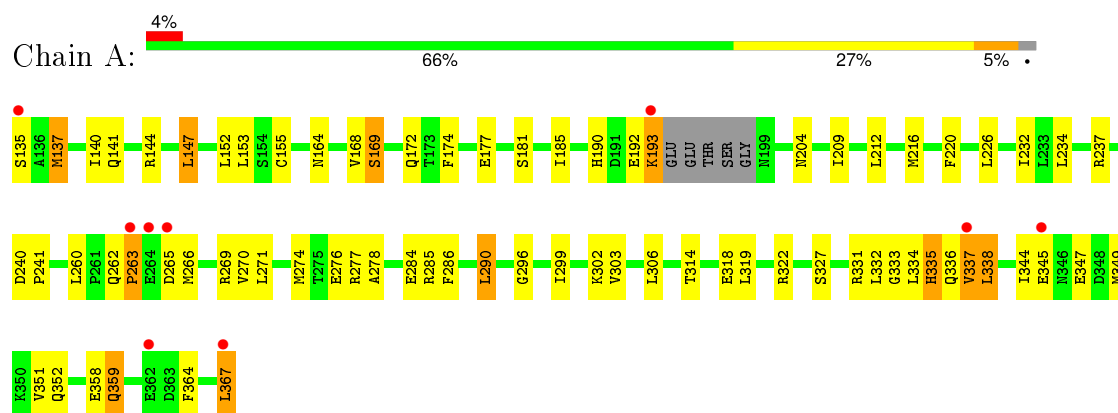
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	117	Total	O	0	0
			117	117		
3	B	8	Total	O	0	0
			8	8		

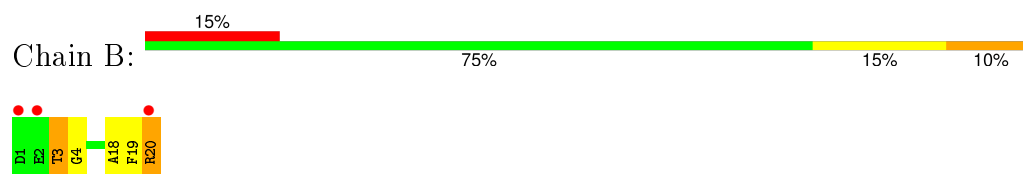
3 Residue-property plots

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: Diaphanous protein homolog 1



• Molecule 2: Diaphanous protein homolog 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.16 Å 65.79 Å 46.86 Å 90.00° 104.08° 90.00°	Depositor
Resolution (Å)	7.97 – 2.10 7.97 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (7.97-2.10) 96.9 (7.97-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.10 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.229 , 0.312 0.227 , 0.306	Depositor DCC
R_{free} test set	840 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.155	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.53 , 78.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 16811 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2087	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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	0.31	0/1838	0.77	0/2473
2	B	0.32	0/146	0.69	0/195
All	All	0.31	0/1984	0.77	0/2668

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
2	B	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	262	GLN	Peptide
1	A	263	PRO	Peptide
2	B	19	PHE	Peptide

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	1816	0	1833	69	0
2	B	146	0	141	5	0
3	A	117	0	0	22	0
3	B	8	0	0	0	0
All	All	2087	0	1974	74	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 19.

All (74) 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:A:192:GLU:O	1:A:193:LYS:HG3	1.33	1.23
1:A:344:ILE:HD12	3:A:442:HOH:O	1.42	1.16
1:A:140:ILE:HD11	3:A:474:HOH:O	1.50	1.09
1:A:240:ASP:HB3	3:A:476:HOH:O	1.59	1.01
1:A:164:ASN:HB2	3:A:470:HOH:O	1.63	0.97
1:A:192:GLU:O	1:A:193:LYS:CG	2.16	0.93
1:A:269:ARG:HD2	3:A:441:HOH:O	1.69	0.93
1:A:359:GLN:HG3	3:A:385:HOH:O	1.70	0.92
1:A:192:GLU:C	1:A:193:LYS:HG3	1.93	0.87
1:A:296:GLY:C	3:A:480:HOH:O	2.24	0.75
1:A:335:HIS:H	1:A:335:HIS:CD2	2.04	0.74
1:A:306:LEU:HD12	1:A:352:GLN:HB3	1.69	0.73
1:A:274:MET:HE1	1:A:286:PHE:HE1	1.53	0.72
1:A:226:LEU:HD22	1:A:266:MET:HE1	1.72	0.70
1:A:314:THR:OG1	3:A:385:HOH:O	2.11	0.69
1:A:240:ASP:CB	3:A:476:HOH:O	2.30	0.69
1:A:296:GLY:HA3	3:A:480:HOH:O	1.94	0.66
1:A:181:SER:O	1:A:185:ILE:HG13	1.95	0.66
1:A:212:LEU:O	1:A:216:MET:HG3	1.97	0.65
1:A:174:PHE:HA	3:A:474:HOH:O	1.97	0.64
1:A:296:GLY:CA	3:A:480:HOH:O	2.46	0.63
1:A:333:GLY:O	1:A:337:VAL:HG13	1.99	0.62
1:A:335:HIS:CD2	1:A:335:HIS:N	2.68	0.61
1:A:302:LYS:HB3	1:A:349:MET:HE1	1.83	0.61
1:A:349:MET:HG2	3:A:442:HOH:O	2.01	0.60
1:A:153:LEU:CA	1:A:204:ASN:HD21	2.14	0.60
1:A:226:LEU:HD22	1:A:266:MET:CE	2.31	0.60
1:A:306:LEU:HD13	1:A:306:LEU:O	2.02	0.60
1:A:302:LYS:HB3	1:A:349:MET:CE	2.32	0.59
2:B:3:THR:CG2	2:B:4:GLY:N	2.66	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:ILE:O	1:A:144:ARG:HG3	2.04	0.57
1:A:335:HIS:H	1:A:335:HIS:HD2	1.49	0.57
1:A:299:ILE:O	1:A:303:VAL:HG23	2.04	0.57
1:A:153:LEU:N	1:A:204:ASN:HD21	2.05	0.55
1:A:237:ARG:NH1	3:A:379:HOH:O	2.38	0.54
1:A:226:LEU:CD2	1:A:232:ILE:HD11	2.38	0.53
1:A:137:MET:O	1:A:141:GLN:HG3	2.07	0.53
2:B:18:ALA:C	2:B:20:ARG:HA	2.30	0.52
1:A:153:LEU:HA	1:A:204:ASN:HD21	1.74	0.52
1:A:226:LEU:HD23	1:A:232:ILE:HD11	1.91	0.52
1:A:290:LEU:HD13	1:A:332:LEU:HD23	1.91	0.52
1:A:168:VAL:O	1:A:172:GLN:HG3	2.09	0.52
1:A:153:LEU:HA	1:A:204:ASN:ND2	2.25	0.51
1:A:347:GLU:O	1:A:351:VAL:HG23	2.10	0.51
1:A:240:ASP:N	3:A:476:HOH:O	2.43	0.51
1:A:147:LEU:N	1:A:147:LEU:HD12	2.26	0.49
1:A:135:SER:N	3:A:437:HOH:O	2.45	0.49
1:A:334:LEU:O	1:A:338:LEU:HD22	2.13	0.49
1:A:177:GLU:HB3	3:A:446:HOH:O	2.12	0.48
1:A:271:LEU:HA	1:A:274:MET:HE2	1.96	0.48
1:A:172:GLN:HA	1:A:220:PHE:CE2	2.49	0.48
2:B:3:THR:HG23	2:B:4:GLY:H	1.79	0.47
1:A:299:ILE:HA	1:A:302:LYS:HD2	1.97	0.47
1:A:318:GLU:O	1:A:322:ARG:HG3	2.13	0.47
1:A:164:ASN:CB	3:A:470:HOH:O	2.37	0.46
1:A:190:HIS:HD2	1:A:234:LEU:HD23	1.81	0.46
2:B:3:THR:HG23	2:B:4:GLY:N	2.30	0.46
1:A:269:ARG:CD	3:A:441:HOH:O	2.41	0.45
1:A:278:ALA:HB2	1:A:285:ARG:HA	1.98	0.45
1:A:265:ASP:OD1	3:A:388:HOH:O	2.21	0.45
1:A:349:MET:HE3	3:A:442:HOH:O	2.18	0.44
1:A:277:ARG:HG2	1:A:277:ARG:O	2.17	0.44
1:A:266:MET:SD	1:A:269:ARG:HD3	2.58	0.44
1:A:190:HIS:HD2	1:A:234:LEU:CD2	2.30	0.43
2:B:18:ALA:O	2:B:20:ARG:HA	2.18	0.43
1:A:319:LEU:HD21	1:A:367:LEU:HB3	2.00	0.43
1:A:147:LEU:HD22	1:A:152:LEU:HA	2.00	0.43
1:A:240:ASP:HA	1:A:241:PRO:HD2	1.82	0.42
1:A:284:GLU:OE2	3:A:479:HOH:O	2.22	0.42
1:A:169:SER:HA	1:A:172:GLN:HE21	1.85	0.41
1:A:345:GLU:O	1:A:345:GLU:CG	2.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:SER:O	1:A:331:ARG:HG3	2.20	0.41
1:A:274:MET:CE	1:A:286:PHE:HE1	2.28	0.41
1:A:266:MET:O	1:A:270:VAL:HG23	2.20	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	224/233 (96%)	216 (96%)	7 (3%)	1 (0%)	39	37
2	B	18/20 (90%)	17 (94%)	1 (6%)	0	100	100
All	All	242/253 (96%)	233 (96%)	8 (3%)	1 (0%)	39	37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	263	PRO

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	204/208 (98%)	187 (92%)	17 (8%)	14	9
2	B	15/15 (100%)	13 (87%)	2 (13%)	5	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	219/223 (98%)	200 (91%)	19 (9%)	13 8

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	MET
1	A	147	LEU
1	A	155	CYS
1	A	169	SER
1	A	193	LYS
1	A	209	ILE
1	A	260	LEU
1	A	276	GLU
1	A	290	LEU
1	A	335	HIS
1	A	336	GLN
1	A	337	VAL
1	A	338	LEU
1	A	358	GLU
1	A	359	GLN
1	A	364	PHE
1	A	367	LEU
2	B	3	THR
2	B	20	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	172	GLN
1	A	204	ASN
1	A	205	GLN
1	A	335	HIS
2	B	14	GLN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	228/233 (97%)	0.15	9 (3%) 43 52	17, 33, 49, 62	0
2	B	20/20 (100%)	0.67	3 (15%) 3 5	28, 40, 63, 65	0
All	All	248/253 (98%)	0.19	12 (4%) 34 43	17, 33, 51, 65	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	263	PRO	3.5
2	B	20	ARG	3.1
2	B	2	GLU	2.7
1	A	264	GLU	2.7
1	A	345	GLU	2.5
2	B	1	ASP	2.4
1	A	362	GLU	2.3
1	A	193	LYS	2.3
1	A	135	SER	2.2
1	A	265	ASP	2.2
1	A	337	VAL	2.1
1	A	367	LEU	2.1

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 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.