



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2024 – 03:08 AM EST

PDB ID : 3GR0
Title : Periplasmic domain of the T3SS inner membrane protein PrgH from *S.typhimurium* (fragment 170-362)
Authors : Yip, C.K.; Vockovic, M.; Yu, A.C.; Strynadka, N.C.J.
Deposited on : 2009-03-24
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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

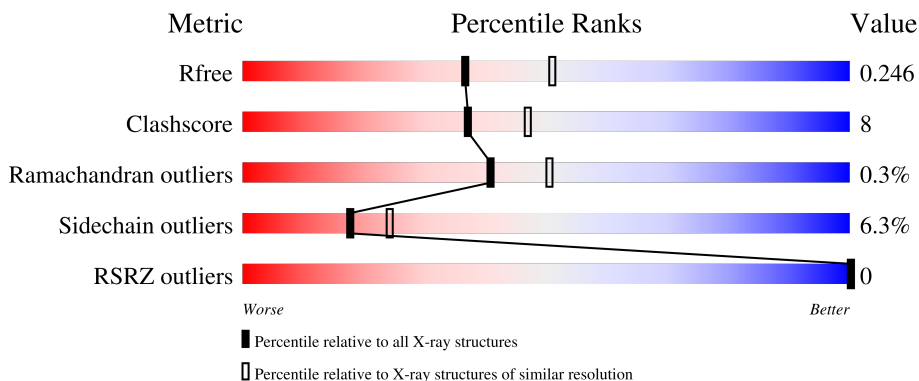
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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 of 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	197	 74% 19% • 6%
1	B	197	 82% 13% • •
1	C	197	 72% 16% • 11%
1	D	197	 71% 18% • 9%

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 6311 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 Protein prgH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	186	Total 1541	C 976	N 280	O 281	S 4	0	0	0
1	B	192	Total 1585	C 1003	N 286	O 292	S 4	0	0	0
1	C	176	Total 1461	C 928	N 268	O 261	S 4	0	0	0
1	D	180	Total 1493	C 947	N 272	O 270	S 4	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	166	GLY	-	expression tag	UNP P41783
A	167	SER	-	expression tag	UNP P41783
A	168	HIS	-	expression tag	UNP P41783
A	169	MET	-	expression tag	UNP P41783
B	166	GLY	-	expression tag	UNP P41783
B	167	SER	-	expression tag	UNP P41783
B	168	HIS	-	expression tag	UNP P41783
B	169	MET	-	expression tag	UNP P41783
C	166	GLY	-	expression tag	UNP P41783
C	167	SER	-	expression tag	UNP P41783
C	168	HIS	-	expression tag	UNP P41783
C	169	MET	-	expression tag	UNP P41783
D	166	GLY	-	expression tag	UNP P41783
D	167	SER	-	expression tag	UNP P41783
D	168	HIS	-	expression tag	UNP P41783
D	169	MET	-	expression tag	UNP P41783

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	66	Total 66	O 66	0	0
2	B	69	Total 69	O 69	0	0
2	C	41	Total 41	O 41	0	0
2	D	55	Total 55	O 55	0	0

4 Data and refinement statistics i

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	53.26Å 53.26Å 282.53Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.00 – 2.30 94.18 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (35.00-2.30) 94.1 (94.18-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.92 (at 2.10Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.206 , 0.255 0.200 , 0.246	Depositor DCC
R_{free} test set	2521 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtrriage
Anisotropy	0.343	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.024 for -h,-k,l 0.487 for h,-h-k,-l 0.027 for -k,-h,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6311	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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](#)

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	0.66	0/1573	0.70	0/2122
1	B	0.69	1/1617 (0.1%)	0.70	1/2182 (0.0%)
1	C	0.57	0/1491	0.65	0/2011
1	D	0.58	0/1525	0.69	1/2059 (0.0%)
All	All	0.63	1/6206 (0.0%)	0.69	2/8374 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	218	LYS	CE-NZ	10.62	1.75	1.49

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	353	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	B	307	LEU	CA-CB-CG	5.16	127.17	115.30

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	1541	0	1530	31	0
1	B	1585	0	1572	19	0
1	C	1461	0	1462	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1493	0	1483	27	0
2	A	66	0	0	0	0
2	B	69	0	0	1	0
2	C	41	0	0	0	0
2	D	55	0	0	3	0
All	All	6311	0	6047	95	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 8.

All (95) 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:218:LYS:CE	1:B:218:LYS:NZ	1.75	1.48
1:D:263:GLN:HE22	1:D:296:ASP:H	1.25	0.81
1:C:263:GLN:HE22	1:C:296:ASP:H	1.27	0.79
1:B:263:GLN:HE21	1:B:263:GLN:H	1.30	0.78
1:A:263:GLN:HE21	1:A:263:GLN:H	1.30	0.77
1:D:263:GLN:NE2	1:D:296:ASP:H	1.83	0.76
1:A:267:MET:HE2	1:A:271:GLU:HB3	1.70	0.74
1:C:209:GLN:O	1:C:213:ARG:HG2	1.88	0.73
1:C:263:GLN:NE2	1:C:296:ASP:H	1.85	0.73
1:A:263:GLN:HE22	1:A:296:ASP:H	1.38	0.70
1:C:205:LEU:HD11	1:D:266:THR:HA	1.74	0.69
1:A:304:GLU:HG2	1:A:325:PHE:CE2	2.28	0.69
1:B:263:GLN:HE22	1:B:296:ASP:H	1.40	0.68
1:A:304:GLU:HG2	1:A:325:PHE:HE2	1.58	0.67
1:B:304:GLU:HG2	1:B:325:PHE:CE2	2.29	0.66
1:A:196:VAL:HG13	1:A:222:VAL:HG22	1.76	0.66
1:A:361:LEU:HD13	1:C:343:VAL:HG21	1.79	0.64
1:D:201:GLU:H	1:D:201:GLU:CD	2.02	0.62
1:C:263:GLN:H	1:C:263:GLN:HE21	1.47	0.62
1:A:317:ARG:HH21	1:C:356:GLN:NE2	1.98	0.62
1:C:201:GLU:OE1	1:C:202:ARG:HD3	2.00	0.61
1:B:304:GLU:HG2	1:B:325:PHE:HE2	1.65	0.61
1:D:248:ILE:HG12	1:D:258:PHE:HD1	1.66	0.61
1:B:263:GLN:NE2	1:B:296:ASP:H	1.98	0.60
1:A:180:GLU:HG3	1:A:184:PHE:CD1	2.37	0.59
1:D:353:ARG:NH2	2:D:109:HOH:O	2.35	0.59
1:B:218:LYS:NZ	1:B:218:LYS:CD	2.61	0.58
1:A:187:LEU:HD11	1:A:253:PRO:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:ARG:HD3	1:B:318:ASN:OD1	2.04	0.57
1:A:196:VAL:CG1	1:A:222:VAL:HG22	2.34	0.57
1:D:304:GLU:HG2	1:D:325:PHE:CE2	2.40	0.56
1:D:304:GLU:HG2	1:D:325:PHE:HE2	1.71	0.56
1:D:263:GLN:HE21	1:D:263:GLN:H	1.53	0.55
1:A:263:GLN:NE2	1:A:296:ASP:H	2.03	0.55
1:B:225:GLU:HG3	1:B:250:PHE:CD1	2.42	0.54
1:C:248:ILE:HG12	1:C:258:PHE:HD1	1.72	0.54
1:D:263:GLN:HE22	1:D:296:ASP:N	2.03	0.53
1:B:225:GLU:HG3	1:B:250:PHE:HD1	1.73	0.53
1:D:232:ILE:HD13	1:D:248:ILE:CD1	2.39	0.53
1:D:243:LEU:HD11	1:D:260:LEU:HD11	1.91	0.53
1:D:248:ILE:HG12	1:D:258:PHE:CD1	2.44	0.52
1:A:201:GLU:O	1:A:205:LEU:HD13	2.09	0.52
1:C:304:GLU:HG3	1:C:325:PHE:CE2	2.45	0.52
1:C:306:GLY:HA3	1:C:342:PHE:CE2	2.46	0.50
1:A:317:ARG:HH21	1:C:356:GLN:HE21	1.57	0.50
1:A:201:GLU:OE2	1:A:224:ASN:ND2	2.45	0.49
1:C:234:ILE:O	1:C:237:ASP:HB2	2.12	0.49
1:D:312:LEU:HD21	1:D:339:ALA:HB2	1.94	0.49
1:B:177:LEU:HD21	1:B:186:VAL:HG23	1.95	0.48
1:B:187:LEU:HD11	1:B:253:PRO:HB2	1.94	0.48
1:B:235:TRP:CD2	1:B:282:LEU:HD11	2.48	0.48
1:B:319:HIS:ND1	2:B:366:HOH:O	2.35	0.48
1:A:306:GLY:HA3	1:A:342:PHE:CE2	2.49	0.48
1:A:361:LEU:HD21	1:C:340:ARG:HG2	1.96	0.48
1:C:202:ARG:NH2	1:C:205:LEU:HD23	2.28	0.48
1:D:240:TYR:HB3	1:D:243:LEU:HB2	1.95	0.48
1:C:339:ALA:O	1:C:343:VAL:HG23	2.13	0.47
1:C:191:ASP:C	1:C:192:LYS:HD2	2.35	0.47
1:A:313:PRO:HB3	1:D:287:ASP:HB3	1.97	0.47
1:C:228:GLU:O	1:C:232:ILE:HG12	2.16	0.47
1:A:180:GLU:HG3	1:A:184:PHE:HD1	1.80	0.46
1:D:280:ARG:HD2	1:D:288:SER:HA	1.97	0.46
1:A:263:GLN:HE21	1:A:263:GLN:N	2.05	0.46
1:C:254:ARG:HA	1:C:254:ARG:HD2	1.84	0.46
1:A:317:ARG:NH2	1:C:356:GLN:HE21	2.13	0.46
1:A:319:HIS:CE1	1:C:317:ARG:HB3	2.50	0.46
1:B:249:HIS:HE1	1:B:349:THR:HA	1.80	0.46
1:C:190:ARG:NH2	1:C:286:ALA:O	2.50	0.45
1:A:180:GLU:HG3	1:A:184:PHE:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:209:GLN:O	1:D:213:ARG:HG2	2.16	0.45
1:D:306:GLY:HA3	1:D:342:PHE:CE2	2.51	0.45
1:B:316:ARG:HG3	1:B:325:PHE:CE2	2.52	0.45
1:A:317:ARG:NH2	1:C:356:GLN:NE2	2.63	0.45
1:D:327:ILE:CG2	1:D:331:LEU:HD11	2.48	0.44
1:C:263:GLN:HE22	1:C:296:ASP:N	2.04	0.43
1:D:232:ILE:HD13	1:D:248:ILE:HD13	2.00	0.43
1:A:235:TRP:CD2	1:A:282:LEU:HD11	2.54	0.43
1:A:249:HIS:NE2	1:A:349:THR:O	2.50	0.43
1:B:263:GLN:H	1:B:263:GLN:NE2	2.08	0.43
1:D:262:ARG:NH2	2:D:363:HOH:O	2.50	0.43
1:D:207:ALA:HA	1:D:210:VAL:HG22	2.01	0.43
1:B:263:GLN:HE21	1:B:263:GLN:N	2.08	0.42
1:A:314:TYR:HA	1:A:326:VAL:O	2.19	0.42
1:A:263:GLN:H	1:A:263:GLN:NE2	2.09	0.42
1:C:248:ILE:HG12	1:C:258:PHE:CD1	2.53	0.42
1:A:248:ILE:HG12	1:A:258:PHE:CD1	2.55	0.41
1:A:320:LYS:HD3	1:A:320:LYS:HA	1.72	0.41
1:D:232:ILE:HG21	1:D:248:ILE:HD11	2.02	0.41
1:A:256:PRO:HD2	1:A:288:SER:O	2.21	0.41
1:D:185:GLN:HG2	1:D:187:LEU:HD11	2.02	0.41
1:B:306:GLY:HA3	1:B:342:PHE:CE2	2.56	0.40
1:D:356:GLN:HG3	2:D:368:HOH:O	2.20	0.40
1:C:240:TYR:HB3	1:C:243:LEU:HB2	2.03	0.40
1:C:263:GLN:HE22	1:C:295:ASP:HA	1.86	0.40
1:D:254:ARG:HA	1:D:254:ARG:HD2	1.99	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	184/197 (93%)	180 (98%)	4 (2%)	0	100	100
1	B	190/197 (96%)	185 (97%)	4 (2%)	1 (0%)	29	35
1	C	172/197 (87%)	170 (99%)	2 (1%)	0	100	100
1	D	178/197 (90%)	173 (97%)	4 (2%)	1 (1%)	25	31
All	All	724/788 (92%)	708 (98%)	14 (2%)	2 (0%)	41	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	215	ASP
1	B	320	LYS

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	160/168 (95%)	150 (94%)	10 (6%)	18	24
1	B	165/168 (98%)	153 (93%)	12 (7%)	14	18
1	C	152/168 (90%)	145 (95%)	7 (5%)	27	38
1	D	155/168 (92%)	144 (93%)	11 (7%)	14	19
All	All	632/672 (94%)	592 (94%)	40 (6%)	18	24

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	179	GLN
1	A	181	LYS
1	A	209	GLN
1	A	238	THR
1	A	263	GLN
1	A	277	GLN
1	A	307	LEU
1	A	345	SER
1	A	361	LEU

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Mol	Chain	Res	Type
1	A	362	LYS
1	B	177	LEU
1	B	181	LYS
1	B	201	GLU
1	B	226	ASN
1	B	252	GLU
1	B	263	GLN
1	B	307	LEU
1	B	309	GLN
1	B	316	ARG
1	B	336	ILE
1	B	338	ARG
1	B	361	LEU
1	C	192	LYS
1	C	194	LEU
1	C	202	ARG
1	C	263	GLN
1	C	307	LEU
1	C	338	ARG
1	C	360	GLU
1	D	192	LYS
1	D	194	LEU
1	D	211	LEU
1	D	213	ARG
1	D	255	LYS
1	D	263	GLN
1	D	264	ARG
1	D	307	LEU
1	D	309	GLN
1	D	312	LEU
1	D	320	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	226	ASN
1	A	263	GLN
1	A	328	GLN
1	A	341	GLN
1	B	185	GLN
1	B	199	GLN
1	B	242	GLN

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Mol	Chain	Res	Type
1	B	263	GLN
1	B	309	GLN
1	B	341	GLN
1	C	263	GLN
1	C	341	GLN
1	C	356	GLN
1	D	185	GLN
1	D	226	ASN
1	D	263	GLN

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](#)

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	186/197 (94%)	-0.34	0 100 100	10, 20, 35, 44	0
1	B	192/197 (97%)	-0.34	0 100 100	8, 22, 35, 51	0
1	C	176/197 (89%)	-0.33	0 100 100	14, 23, 34, 38	0
1	D	180/197 (91%)	-0.29	0 100 100	13, 24, 35, 40	0
All	All	734/788 (93%)	-0.32	0 100 100	8, 22, 35, 51	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.