



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 01:57 AM GMT

PDB ID : 3E29
Title : X-Ray structure of the protein Q7WE92_BORBR from thioesterase superfamily. Northeast Structural Genomics Consortium Target BoR214A.
Authors : Kuzin, A.P.; Chen, Y.; Setharaman, J.; Wang, D.; Mao, L.; Foote, E.L.; Xiao, R.; Nair, R.; Everett, J.K.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2008-08-05
Resolution : 2.40 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>

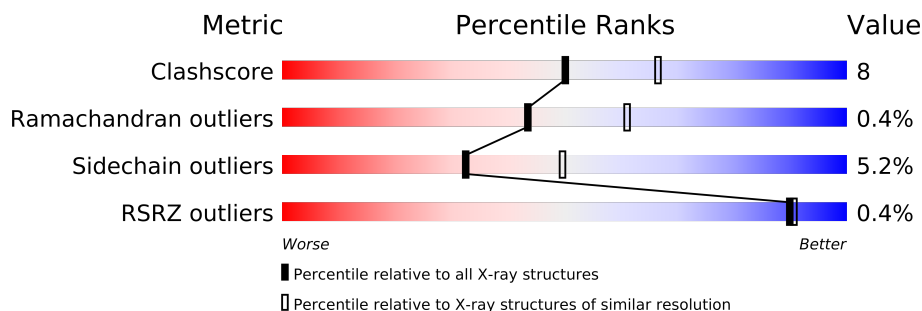
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
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 2.40 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	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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.

Mol	Chain	Length	Quality of chain
1	A	144	
1	B	144	
1	C	144	
1	D	144	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4321 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 uncharacterized protein Q7WE92_BORBR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	134	Total	C	N	O	Se	8	2	0
			1043	656	195	186	6			
1	B	135	Total	C	N	O	Se	8	1	0
			1040	654	193	187	6			
1	C	133	Total	C	N	O	Se	8	2	0
			1035	651	194	185	5			
1	D	132	Total	C	N	O	Se	0	1	0
			1021	643	190	183	5			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	137	LEU	-	expression tag	UNP Q7WE92
A	138	GLU	-	expression tag	UNP Q7WE92
A	139	HIS	-	expression tag	UNP Q7WE92
A	140	HIS	-	expression tag	UNP Q7WE92
A	141	HIS	-	expression tag	UNP Q7WE92
A	142	HIS	-	expression tag	UNP Q7WE92
A	143	HIS	-	expression tag	UNP Q7WE92
A	144	HIS	-	expression tag	UNP Q7WE92
B	137	LEU	-	expression tag	UNP Q7WE92
B	138	GLU	-	expression tag	UNP Q7WE92
B	139	HIS	-	expression tag	UNP Q7WE92
B	140	HIS	-	expression tag	UNP Q7WE92
B	141	HIS	-	expression tag	UNP Q7WE92
B	142	HIS	-	expression tag	UNP Q7WE92
B	143	HIS	-	expression tag	UNP Q7WE92
B	144	HIS	-	expression tag	UNP Q7WE92
C	137	LEU	-	expression tag	UNP Q7WE92
C	138	GLU	-	expression tag	UNP Q7WE92
C	139	HIS	-	expression tag	UNP Q7WE92
C	140	HIS	-	expression tag	UNP Q7WE92
C	141	HIS	-	expression tag	UNP Q7WE92

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Chain	Residue	Modelled	Actual	Comment	Reference
C	142	HIS	-	expression tag	UNP Q7WE92
C	143	HIS	-	expression tag	UNP Q7WE92
C	144	HIS	-	expression tag	UNP Q7WE92
D	137	LEU	-	expression tag	UNP Q7WE92
D	138	GLU	-	expression tag	UNP Q7WE92
D	139	HIS	-	expression tag	UNP Q7WE92
D	140	HIS	-	expression tag	UNP Q7WE92
D	141	HIS	-	expression tag	UNP Q7WE92
D	142	HIS	-	expression tag	UNP Q7WE92
D	143	HIS	-	expression tag	UNP Q7WE92
D	144	HIS	-	expression tag	UNP Q7WE92

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	43	Total O 43 43	0	0
2	B	41	Total O 41 41	0	0
2	C	49	Total O 49 49	0	0
2	D	49	Total O 49 49	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	112.52Å 112.52Å 119.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.59 – 2.40 28.60 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (28.59-2.40) 99.3 (28.60-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.27 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.221 , 0.258 0.212 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.532	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 26.8	EDS
Estimated twinning fraction	0.480 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 33459 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4321	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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 57.13 % 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 2.4741e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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	0.53	0/1065	0.69	1/1431 (0.1%)
1	B	0.54	0/1059	0.67	1/1424 (0.1%)
1	C	0.51	0/1057	0.62	0/1421
1	D	0.48	0/1040	0.64	0/1399
All	All	0.52	0/4221	0.66	2/5675 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	128	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	118	ASP	CB-CG-OD1	5.02	122.82	118.30

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	1043	0	1054	27	1
1	B	1040	0	1047	19	4
1	C	1035	0	1042	11	0
1	D	1021	0	1025	9	0
2	A	43	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	41	0	0	2	0
2	C	49	0	0	1	0
2	D	49	0	0	1	3
All	All	4321	0	4168	62	4

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 8.

All (62) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1:MSE:HB3	1:A:30:GLY:HA3	1.32	1.07
1:C:103:ILE:HG13	1:C:113:HIS:CE1	2.23	0.73
1:B:108:ARG:HH21	1:D:108:ARG:NH1	1.88	0.71
1:A:103:ILE:HG13	1:A:113:HIS:CE1	2.26	0.69
1:C:97:ARG:HD3	2:C:159:HOH:O	1.92	0.68
1:C:88:HIS:HE1	1:C:126:SER:OG	1.77	0.66
1:D:88:HIS:HE1	1:D:126:SER:OG	1.79	0.64
1:A:60:THR:OG1	1:B:56:GLY:HA3	1.97	0.64
1:A:108[A]:ARG:HE	1:A:108[A]:ARG:H	1.49	0.59
1:B:103:ILE:HG13	1:B:113:HIS:CE1	2.37	0.59
1:B:115:ARG:NH1	2:B:174:HOH:O	2.30	0.59
1:A:106:GLY:HA3	1:A:108[A]:ARG:NH2	2.19	0.58
1:A:108[A]:ARG:NE	1:A:108[A]:ARG:H	2.02	0.58
1:A:108[A]:ARG:HE	1:A:109:PHE:H	1.53	0.57
1:C:103:ILE:HG13	1:C:113:HIS:ND1	2.19	0.57
1:D:41:ARG:H	1:D:44:LEU:HD22	1.69	0.57
1:B:115:ARG:HD3	2:B:174:HOH:O	2.05	0.56
1:A:113:HIS:CD2	1:A:128:ARG:HB3	2.41	0.56
1:B:134:ARG:HG2	1:B:135:ALA:H	1.70	0.55
1:A:108[A]:ARG:HG2	1:A:109:PHE:CD2	2.42	0.55
1:B:108:ARG:HH21	1:D:108:ARG:HH11	1.56	0.53
1:B:92:THR:HB	1:B:93:PRO:HD2	1.90	0.53
1:C:41:ARG:H	1:C:44:LEU:HD22	1.74	0.53
1:B:88:HIS:HE1	1:B:126:SER:OG	1.91	0.53
1:D:46:SER:HB2	1:D:54:HIS:HA	1.92	0.52
1:A:1:MSE:HB3	1:A:30:GLY:CA	2.23	0.51
1:A:56:GLY:HA3	1:B:60:THR:OG1	2.11	0.50
1:A:83:MSE:HE1	1:A:85:VAL:HB	1.94	0.50
1:B:133:ILE:O	1:B:134:ARG:HB2	2.12	0.50
1:A:108[A]:ARG:CD	1:A:108[A]:ARG:H	2.27	0.48
1:B:74:THR:HB	1:B:76:HIS:CE1	2.50	0.47
1:C:46:SER:HB2	1:C:54:HIS:HA	1.97	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:81:MSE:HG3	1:A:130:LEU:HG	1.97	0.46
1:D:99:GLU:O	1:D:114:ALA:HA	2.15	0.46
1:B:113:HIS:CD2	1:B:128:ARG:HB3	2.50	0.45
1:A:83:MSE:CE	1:A:85:VAL:HB	2.46	0.45
1:B:92:THR:HB	1:B:93:PRO:CD	2.47	0.44
1:B:41:ARG:O	1:B:44:LEU:HB2	2.17	0.44
1:C:3:SER:O	1:C:7:GLU:HB2	2.18	0.44
1:B:46:SER:HB2	1:B:54:HIS:HA	2.00	0.44
1:C:81:MSE:HE3	1:C:132:LEU:HB2	1.99	0.43
1:A:88:HIS:HE1	1:A:126:SER:OG	2.01	0.43
1:A:106:GLY:HA3	1:A:108[A]:ARG:HH21	1.84	0.42
1:D:103:ILE:HG12	1:D:113:HIS:CE1	2.55	0.42
1:C:88:HIS:CE1	1:C:126:SER:OG	2.65	0.42
1:A:103:ILE:HA	1:A:103:ILE:HD13	1.86	0.42
1:A:7:GLU:HG3	1:A:8:MSE:N	2.34	0.42
1:B:24:MSE:HA	1:B:37:GLY:O	2.20	0.42
1:C:118:ASP:OD1	1:C:122:ASN:HB2	2.20	0.41
1:A:113:HIS:HD2	1:A:128:ARG:HB3	1.85	0.41
1:A:88:HIS:HD2	2:D:181:HOH:O	2.02	0.41
1:A:108[A]:ARG:HG2	1:A:109:PHE:CE2	2.55	0.41
1:A:92:THR:HB	1:A:93:PRO:CD	2.50	0.41
1:D:92:THR:HB	1:D:93:PRO:CD	2.51	0.41
1:B:134:ARG:HG2	1:B:135:ALA:N	2.35	0.40
1:D:63:ASP:HB2	1:D:83:MSE:SE	2.71	0.40
1:C:92:THR:HB	1:C:93:PRO:CD	2.52	0.40
1:A:24:MSE:HA	1:A:37:GLY:O	2.21	0.40
1:A:6:LEU:O	1:A:10:SER:HB2	2.22	0.40

All (4) 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:B:51[A]:ARG:CZ	2:D:183:HOH:O[6_555]	1.60	0.60
1:B:51[A]:ARG:NH1	2:D:183:HOH:O[6_555]	1.81	0.39
1:B:51[A]:ARG:NH2	2:D:183:HOH:O[6_555]	1.98	0.22
1:A:7:GLU:OE2	1:B:11:ARG:NH1[3_444]	2.15	0.05

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	134/144 (93%)	133 (99%)	1 (1%)	0	100	100
1	B	134/144 (93%)	131 (98%)	2 (2%)	1 (1%)	30	43
1	C	133/144 (92%)	129 (97%)	3 (2%)	1 (1%)	27	39
1	D	131/144 (91%)	129 (98%)	2 (2%)	0	100	100
All	All	532/576 (92%)	522 (98%)	8 (2%)	2 (0%)	43	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	134	ARG
1	C	133	ILE

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	108/109 (99%)	100 (93%)	8 (7%)	20	30
1	B	107/109 (98%)	102 (95%)	5 (5%)	36	54
1	C	107/109 (98%)	102 (95%)	5 (5%)	36	54
1	D	105/109 (96%)	100 (95%)	5 (5%)	35	53
All	All	427/436 (98%)	404 (95%)	23 (5%)	32	47

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU

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Mol	Chain	Res	Type
1	A	32	GLN
1	A	44	LEU
1	A	49	GLU
1	A	107	LYS
1	A	108[A]	ARG
1	A	108[B]	ARG
1	A	128	ARG
1	B	1	MSE
1	B	44	LEU
1	B	107	LYS
1	B	124	VAL
1	B	128	ARG
1	C	6	LEU
1	C	44	LEU
1	C	49	GLU
1	C	117	LEU
1	C	128	ARG
1	D	6	LEU
1	D	41	ARG
1	D	44	LEU
1	D	47	SER
1	D	128	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	84	HIS
1	A	88	HIS
1	B	19	ASN
1	B	76	HIS
1	B	84	HIS
1	B	88	HIS
1	C	88	HIS
1	D	88	HIS

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 ⓘ

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, 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	134/144 (93%)	-0.19	0 100 100	22, 28, 39, 55	1 (0%)
1	B	135/144 (93%)	-0.16	1 (0%) 84 84	22, 29, 40, 56	1 (0%)
1	C	133/144 (92%)	-0.14	0 100 100	23, 30, 43, 49	1 (0%)
1	D	132/144 (91%)	-0.11	1 (0%) 83 82	22, 30, 41, 47	1 (0%)
All	All	534/576 (92%)	-0.15	2 (0%) 90 90	22, 29, 41, 56	4 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	134	ARG	3.5
1	B	2	SER	2.3

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

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.