



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:40 AM GMT

PDB ID : 2BEC
Title : Crystal structure of CHP2 in complex with its binding region in NHE1 and insights into the mechanism of pH regulation
Authors : Ben Ammar, Y.; Takeda, S.; Hisamitsu, T.; Mori, H.; Wakabayashi, S.
Deposited on : 2005-10-24
Resolution : 2.70 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

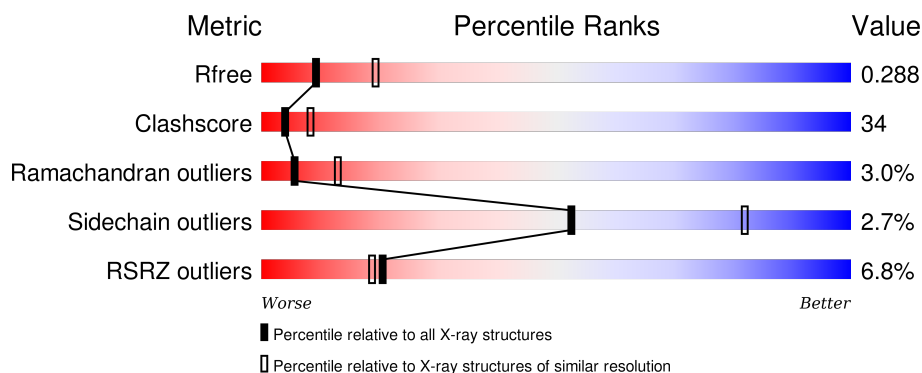
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	202	
2	B	43	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1674 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 Calcineurin B homologous protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1470	917	269	279	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	197	HIS	-	EXPRESSION TAG	UNP O43745
A	198	HIS	-	EXPRESSION TAG	UNP O43745
A	199	HIS	-	EXPRESSION TAG	UNP O43745
A	200	HIS	-	EXPRESSION TAG	UNP O43745
A	201	HIS	-	EXPRESSION TAG	UNP O43745
A	202	HIS	-	EXPRESSION TAG	UNP O43745

- Molecule 2 is a protein called Sodium/hydrogen exchanger 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	25	Total	C	N	O	S	0	0	0
			202	125	36	40	1			

- Molecule 3 is YTTRIUM (III) ION (three-letter code: YT3) (formula: Y).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Y	0	0
			2	2		

4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	49.26 Å 49.26 Å 102.59 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 28.82 – 2.67	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.70) 97.3 (28.82-2.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.34 (at 2.68 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.218 , 0.287 0.222 , 0.288	Depositor DCC
R_{free} test set	350 reflections (5.55%)	DCC
Wilson B-factor (Å ²)	60.3	Xtriage
Anisotropy	0.408	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 58.7	EDS
Estimated twinning fraction	0.074 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 6827 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	1674	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.79% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: YT3

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	0.30	0/1493	0.52	0/2004
2	B	0.34	0/205	0.49	0/276
All	All	0.31	0/1698	0.51	0/2280

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	1470	0	1456	110	0
2	B	202	0	189	16	0
3	A	2	0	0	0	0
All	All	1674	0	1645	113	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 34.

All (113) 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:21:THR:HG21	1:A:84:VAL:HG12	1.45	0.99
1:A:82:GLY:HA2	1:A:85:ARG:NH1	1.87	0.89
1:A:18:ARG:HG3	1:A:23:PHE:O	1.78	0.83
1:A:79:ASP:OD2	1:A:81:PRO:HG2	1.80	0.81
1:A:140:LEU:HD21	2:B:522:ILE:HD12	1.63	0.80
1:A:149:THR:HG23	1:A:152:GLN:HE21	1.47	0.79
1:A:57:LEU:HD22	2:B:537:ILE:HD11	1.65	0.79
1:A:136:MET:CE	2:B:522:ILE:HG21	2.15	0.76
1:A:43:LYS:HB3	1:A:45:TYR:CE1	2.20	0.76
1:A:60:ASN:HD22	1:A:62:LEU:H	1.38	0.72
1:A:32:HIS:CD2	1:A:36:ARG:HH21	2.08	0.71
1:A:136:MET:HE1	2:B:522:ILE:HG21	1.72	0.70
1:A:48:ARG:HG2	1:A:48:ARG:HH11	1.56	0.69
1:A:134:HIS:O	1:A:138:GLN:HG2	1.92	0.69
1:A:95:ASP:O	1:A:99:GLU:HG3	1.92	0.69
1:A:149:THR:OG1	1:A:152:GLN:HG3	1.92	0.69
1:A:193:ARG:HD2	2:B:535:GLU:OE1	1.92	0.69
1:A:173:SER:OG	1:A:176:GLU:HG3	1.96	0.65
1:A:185:ASP:O	1:A:189:LYS:HG3	1.98	0.64
1:A:60:ASN:ND2	1:A:62:LEU:H	1.96	0.64
1:A:146:VAL:O	1:A:148:VAL:HG23	1.98	0.64
1:A:152:GLN:HA	1:A:155:ASN:HD22	1.64	0.63
1:A:92:PRO:HB3	1:A:196:LYS:HE3	1.79	0.62
1:A:194:ILE:HG13	1:A:195:LEU:N	2.15	0.60
1:A:48:ARG:O	1:A:52:GLN:HG3	2.00	0.60
1:A:63:GLY:O	1:A:67:ILE:HG12	2.02	0.59
1:A:48:ARG:H	1:A:48:ARG:HD2	1.67	0.59
1:A:80:PHE:N	1:A:81:PRO:HD2	2.18	0.59
1:A:159:ARG:HA	1:A:162:GLN:HE21	1.68	0.58
1:A:112:ARG:HB2	1:A:187:GLU:OE1	2.03	0.58
1:A:82:GLY:HA2	1:A:85:ARG:HH12	1.66	0.57
1:A:24:SER:O	1:A:27:SER:HB2	2.03	0.57
1:A:38:LEU:HD22	1:A:54:ILE:HD11	1.86	0.57
1:A:57:LEU:HD23	1:A:67:ILE:HD11	1.86	0.57
1:A:174:PHE:O	1:A:177:PHE:HB3	2.05	0.57
1:A:60:ASN:HD22	1:A:60:ASN:C	2.08	0.57
2:B:516:ARG:HA	2:B:520:GLU:OE2	2.05	0.57
1:A:110:ASN:O	1:A:115:LYS:HE2	2.05	0.57
1:A:46:LEU:O	1:A:77:ARG:HA	2.05	0.56
1:A:118:TYR:HE1	2:B:530:LEU:HD11	1.70	0.56
1:A:164:ALA:HB2	2:B:519:ASN:ND2	2.20	0.56
1:A:90:PHE:HB3	1:A:192:ILE:HG13	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLU:HG3	1:A:21:THR:HG23	1.88	0.55
1:A:117:HIS:O	1:A:121:GLN:HG3	2.06	0.55
1:A:43:LYS:HB3	1:A:45:TYR:CD1	2.42	0.55
1:A:23:PHE:CE2	1:A:84:VAL:HG13	2.42	0.55
1:A:92:PRO:CB	1:A:196:LYS:HE3	2.36	0.55
1:A:35:PHE:CE2	1:A:46:LEU:HD13	2.42	0.55
1:A:72:PRO:HD2	1:A:75:SER:HB3	1.90	0.54
1:A:179:LYS:HD2	1:A:182:GLU:OE1	2.08	0.54
1:A:149:THR:HG23	1:A:152:GLN:NE2	2.21	0.54
1:A:79:ASP:N	1:A:79:ASP:OD1	2.41	0.53
1:A:32:HIS:HA	1:A:80:PHE:CE2	2.43	0.53
1:A:164:ALA:HB2	2:B:519:ASN:HD21	1.74	0.53
1:A:189:LYS:NZ	1:A:189:LYS:HB3	2.22	0.53
1:A:11:ILE:HG22	1:A:11:ILE:O	2.08	0.52
1:A:21:THR:HG21	1:A:84:VAL:CG1	2.28	0.52
1:A:34:ARG:O	1:A:38:LEU:HG	2.10	0.52
1:A:148:VAL:CG1	1:A:152:GLN:HB2	2.40	0.52
1:A:48:ARG:HG2	1:A:48:ARG:NH1	2.24	0.52
1:A:159:ARG:HD2	2:B:518:ILE:HG13	1.91	0.51
1:A:26:ALA:HA	1:A:29:LEU:HD12	1.91	0.51
1:A:62:LEU:O	1:A:66:ILE:HG13	2.11	0.50
1:A:157:ALA:O	1:A:161:VAL:HG23	2.11	0.50
1:A:90:PHE:CB	1:A:192:ILE:HG13	2.42	0.50
1:A:57:LEU:HD22	2:B:537:ILE:CD1	2.38	0.49
1:A:144:VAL:O	1:A:144:VAL:HG23	2.13	0.48
1:A:40:ARG:HH11	1:A:40:ARG:HG2	1.77	0.48
1:A:130:LYS:HB2	1:A:171:ALA:HB1	1.94	0.48
1:A:38:LEU:CD2	1:A:54:ILE:HD11	2.43	0.48
1:A:17:ILE:CD1	1:A:81:PRO:HB3	2.43	0.48
1:A:21:THR:HG1	1:A:23:PHE:HD2	1.61	0.47
1:A:190:MET:SD	2:B:527:LEU:HD22	2.54	0.47
1:A:70:PHE:O	1:A:78:VAL:HG22	2.14	0.47
1:A:14:GLY:HA2	1:A:17:ILE:HB	1.96	0.47
1:A:57:LEU:CD2	1:A:67:ILE:HD11	2.44	0.47
1:A:138:GLN:O	1:A:142:LEU:HG	2.14	0.47
1:A:19:ARG:HG3	1:A:19:ARG:HH11	1.79	0.47
1:A:12:PRO:C	1:A:14:GLY:H	2.18	0.46
2:B:533:GLY:O	2:B:537:ILE:HG12	2.16	0.46
1:A:13:ASP:O	1:A:13:ASP:OD1	2.34	0.46
1:A:148:VAL:HG12	1:A:152:GLN:HB2	1.98	0.46
1:A:27:SER:O	1:A:31:LEU:HG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:SER:O	1:A:48:ARG:C	2.55	0.45
1:A:131:ILE:HB	1:A:172:VAL:HB	1.98	0.45
1:A:64:ASP:O	1:A:68:GLU:HG2	2.17	0.45
1:A:14:GLY:O	1:A:18:ARG:NH1	2.49	0.45
1:A:40:ARG:O	1:A:42:LYS:HG2	2.17	0.45
1:A:18:ARG:HE	1:A:25:GLN:HA	1.81	0.44
1:A:34:ARG:NH2	1:A:54:ILE:HD12	2.31	0.44
1:A:17:ILE:HD12	1:A:81:PRO:HB3	2.00	0.43
1:A:82:GLY:CA	1:A:85:ARG:HH12	2.29	0.43
1:A:57:LEU:CD2	2:B:537:ILE:HD11	2.40	0.43
1:A:60:ASN:HD22	1:A:61:PRO:N	2.15	0.43
1:A:96:GLU:O	1:A:105:LYS:NZ	2.51	0.43
1:A:81:PRO:HB2	1:A:85:ARG:NH2	2.33	0.43
1:A:70:PHE:HB3	1:A:78:VAL:HG21	1.99	0.43
1:A:72:PRO:O	1:A:73:ASP:C	2.57	0.43
1:A:165:ASP:OD2	1:A:168:GLY:HA2	2.19	0.42
1:A:122:LEU:HB2	2:B:526:PHE:HZ	1.83	0.42
1:A:28:LEU:HD12	1:A:28:LEU:H	1.83	0.42
1:A:140:LEU:O	1:A:144:VAL:HG13	2.20	0.42
1:A:179:LYS:HD2	1:A:182:GLU:CD	2.40	0.42
1:A:189:LYS:HZ3	1:A:189:LYS:HB3	1.83	0.42
1:A:56:ALA:O	1:A:59:VAL:HG12	2.19	0.42
1:A:45:TYR:O	1:A:46:LEU:HD12	2.20	0.42
1:A:131:ILE:HA	1:A:135:GLU:OE2	2.20	0.41
1:A:46:LEU:HD23	1:A:51:LEU:HD21	2.01	0.41
1:A:18:ARG:NH2	1:A:25:GLN:OE1	2.54	0.41
2:B:518:ILE:O	2:B:522:ILE:HG12	2.21	0.41
1:A:93:VAL:HA	1:A:110:ASN:HD21	1.86	0.40
1:A:90:PHE:HZ	1:A:118:TYR:CE1	2.39	0.40
1:A:149:THR:CG2	1:A:152:GLN:HE21	2.25	0.40

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	177/202 (88%)	156 (88%)	15 (8%)	6 (3%)	5	10
2	B	23/43 (54%)	21 (91%)	2 (9%)	0	100	100
All	All	200/245 (82%)	177 (88%)	17 (8%)	6 (3%)	5	13

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	ASP
1	A	72	PRO
1	A	14	GLY
1	A	194	ILE
1	A	168	GLY
1	A	145	GLY

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	161/179 (90%)	156 (97%)	5 (3%)	47	78
2	B	23/39 (59%)	23 (100%)	0	100	100
All	All	184/218 (84%)	179 (97%)	5 (3%)	52	82

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

Mol	Chain	Res	Type
1	A	40	ARG
1	A	45	TYR
1	A	48	ARG
1	A	60	ASN
1	A	117	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	60	ASN
1	A	76	GLN
1	A	152	GLN
1	A	155	ASN
1	A	162	GLN
2	B	519	ASN

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 ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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 [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	181/202 (89%)	0.40	14 (7%) 16 14	33, 70, 111, 128	0
2	B	25/43 (58%)	0.03	0 100 100	38, 58, 85, 96	0
All	All	206/245 (84%)	0.36	14 (6%) 20 19	33, 69, 111, 128	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	99	GLU	3.6
1	A	49	MET	3.3
1	A	196	LYS	3.2
1	A	149	THR	3.2
1	A	146	VAL	3.2
1	A	14	GLY	2.5
1	A	105	LYS	2.2
1	A	187	GLU	2.2
1	A	147	GLN	2.2
1	A	40	ARG	2.2
1	A	106	PRO	2.1
1	A	168	GLY	2.1
1	A	39	ASP	2.1
1	A	186	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 [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	YT3	A	1202	1/1	1.00	0.13	-1.17	47,47,47,47	0
3	YT3	A	1201	1/1	0.93	0.06	-2.47	93,93,93,93	0

6.5 Other polymers [i](#)

There are no such residues in this entry.