



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 08:57 pm BST

PDB ID : 1AH8
Title : STRUCTURE OF THE ORTHORHOMBIC FORM OF THE N-TERMINAL
DOMAIN OF THE YEAST HSP90 CHAPERONE
Authors : Prodromou, C.; Roe, S.M.; Pearl, L.H.
Deposited on : 1997-04-14
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

<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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

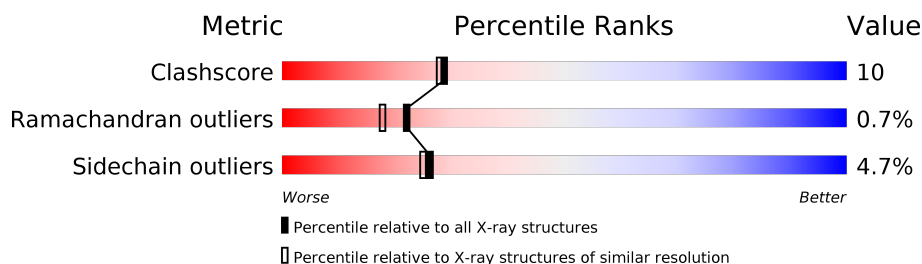
1 Overall quality at a glance

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))
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (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 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	220	
1	B	220	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3986 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 HEAT SHOCK PROTEIN 90.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1699	1080	278	337	4			
1	B	218	Total	C	N	O	S	0	0	0
			1733	1105	284	340	4			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	266	Total 266	O 266	0	0
3	B	264	Total 264	O 264	0	0

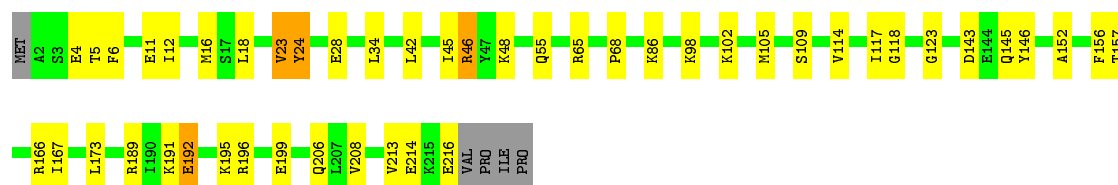
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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. 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. 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.

Note EDS was not executed.

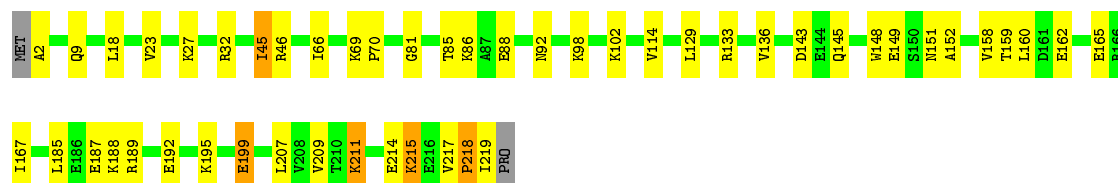
• Molecule 1: HEAT SHOCK PROTEIN 90

Chain A: 



• Molecule 1: HEAT SHOCK PROTEIN 90

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	115.54Å 112.90Å 44.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10	Depositor
% Data completeness (in resolution range)	99.0 (8.00-2.10)	Depositor
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.196 , 0.255	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3986	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

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: GOL

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	0.35	0/1723	0.61	0/2325
1	B	0.35	0/1758	0.63	0/2370
All	All	0.35	0/3481	0.62	0/4695

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 [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	1699	0	1698	38	0
1	B	1733	0	1758	33	0
2	A	6	0	8	1	0
2	B	18	0	24	0	0
3	A	266	0	0	11	6
3	B	264	0	0	6	6
All	All	3986	0	3488	67	10

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 10.

All (67) 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:145:GLN:HE22	1:A:167:ILE:H	1.20	0.89
1:A:11:GLU:HG2	1:A:102:LYS:HB2	1.57	0.86
1:A:199:GLU:O	1:B:215:LYS:HE2	1.83	0.77
1:B:145:GLN:HE22	1:B:167:ILE:H	1.36	0.72
1:B:2:ALA:HB2	3:B:399:HOH:O	1.92	0.69
1:A:16:MET:HE1	1:A:152:ALA:HA	1.80	0.64
1:A:48:LYS:HZ1	1:A:55:GLN:NE2	1.97	0.62
1:B:2:ALA:N	1:B:159:THR:HG1	1.97	0.62
1:B:218:PRO:O	1:B:219:ILE:HB	2.00	0.62
1:B:23:VAL:HG22	3:B:324:HOH:O	1.99	0.61
1:A:46:ARG:NH2	3:A:382:HOH:O	2.04	0.60
1:A:48:LYS:NZ	1:A:55:GLN:NE2	2.51	0.59
1:A:192:GLU:OE2	1:A:196:ARG:NH2	2.37	0.58
1:A:199:GLU:OE2	1:B:215:LYS:HG2	2.04	0.57
1:A:195:LYS:HE3	1:B:214:GLU:OE2	2.05	0.56
1:A:114:VAL:HG23	3:A:448:HOH:O	2.07	0.55
1:A:213:VAL:HG12	3:A:368:HOH:O	2.06	0.55
1:B:86:LYS:HG3	1:B:158:VAL:HG11	1.88	0.53
1:A:166:ARG:HD2	3:A:465:HOH:O	2.08	0.53
1:A:86:LYS:HD2	1:A:146:TYR:CE2	2.42	0.53
1:B:102:LYS:HD2	3:B:435:HOH:O	2.09	0.52
1:A:34:LEU:HD22	1:A:173:LEU:CD2	2.39	0.52
1:B:45:ILE:HD11	1:B:81:GLY:HA2	1.91	0.52
1:A:145:GLN:NE2	1:A:167:ILE:H	1.98	0.51
1:A:23:VAL:O	1:A:24:TYR:C	2.49	0.50
1:B:46:ARG:NE	3:B:393:HOH:O	2.43	0.50
1:A:18:LEU:HD13	1:A:105:MET:SD	2.53	0.49
1:B:114:VAL:HG23	3:B:409:HOH:O	2.11	0.49
1:B:218:PRO:O	1:B:219:ILE:CB	2.60	0.49
1:A:68:PRO:HD2	1:A:208:VAL:O	2.13	0.49
1:B:207:LEU:HG	1:B:209:VAL:CG2	2.43	0.48
1:A:23:VAL:O	1:A:24:TYR:O	2.30	0.48
1:A:189:ARG:HH12	1:A:196:ARG:HH12	1.61	0.47
1:A:98:LYS:HD3	3:A:336:HOH:O	2.14	0.47
1:A:145:GLN:HE22	1:A:167:ILE:N	1.99	0.47
3:A:371:HOH:O	1:B:195:LYS:HG3	2.15	0.47
1:B:189:ARG:O	1:B:192:GLU:HG2	2.14	0.46
1:B:70:PRO:HA	1:B:185:LEU:HD13	1.96	0.46
1:A:173:LEU:HD11	2:A:221:GOL:H11	1.98	0.46
1:A:46:ARG:NE	3:A:382:HOH:O	2.45	0.46
1:A:143:ASP:HB3	3:A:240:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:LYS:HG2	1:B:129:LEU:HB3	1.97	0.46
1:B:151:ASN:O	1:B:152:ALA:HB3	2.15	0.46
1:A:6:PHE:HB2	1:A:156:PHE:CZ	2.51	0.45
1:B:145:GLN:NE2	1:B:167:ILE:H	2.10	0.45
1:B:187:GLU:HG3	1:B:207:LEU:HD21	1.99	0.45
1:B:18:LEU:HD23	1:B:18:LEU:C	2.37	0.44
1:B:195:LYS:HA	1:B:199:GLU:HG3	1.99	0.44
1:B:9:GLN:HE21	1:B:98:LYS:NZ	2.15	0.43
1:A:46:ARG:NH1	3:A:484:HOH:O	2.43	0.43
1:A:28:GLU:HB3	3:A:406:HOH:O	2.19	0.42
1:A:4:GLU:O	1:A:157:THR:HA	2.19	0.42
1:A:117:ILE:HG23	1:A:118:GLY:N	2.35	0.42
1:B:143:ASP:HB3	3:B:234:HOH:O	2.20	0.42
1:B:85:THR:HG22	1:B:143:ASP:HB2	2.02	0.41
1:A:23:VAL:HG13	1:A:23:VAL:O	2.20	0.41
1:B:136:VAL:HB	1:B:148:TRP:HB3	2.02	0.41
1:A:195:LYS:HE3	1:B:214:GLU:CD	2.40	0.41
1:A:42:LEU:O	1:A:45:ILE:HG22	2.21	0.41
1:B:18:LEU:O	1:B:18:LEU:HD23	2.21	0.41
1:A:117:ILE:HG12	1:A:123:GLY:HA2	2.03	0.41
1:A:5:THR:OG1	1:A:157:THR:HG22	2.21	0.41
3:A:473:HOH:O	1:B:211:LYS:HB3	2.20	0.41
1:B:133:ARG:HD3	1:B:149:GLU:OE2	2.21	0.41
1:B:88:GLU:O	1:B:92:ASN:HB2	2.21	0.41
1:A:65:ARG:HD3	1:A:206:GLN:OE1	2.22	0.40
1:A:12:ILE:O	1:A:16:MET:HG3	2.22	0.40

All (10) 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	Interatomic distance (Å)	Clash overlap (Å)
3:B:485:HOH:O	3:B:486:HOH:O[1_556]	1.50	0.70
3:B:227:HOH:O	3:B:365:HOH:O[2_575]	1.80	0.40
3:A:301:HOH:O	3:A:326:HOH:O[2_564]	1.80	0.40
3:A:331:HOH:O	3:A:391:HOH:O[2_565]	1.85	0.35
3:B:295:HOH:O	3:B:378:HOH:O[1_554]	1.91	0.29
3:A:460:HOH:O	3:B:430:HOH:O[1_556]	1.95	0.25
3:A:288:HOH:O	3:A:438:HOH:O[1_556]	2.01	0.19
3:B:368:HOH:O	3:B:462:HOH:O[1_554]	2.02	0.18
3:A:403:HOH:O	3:B:414:HOH:O[3_544]	2.06	0.14
3:A:305:HOH:O	3:A:467:HOH:O[1_556]	2.17	0.03

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	213/220 (97%)	199 (93%)	12 (6%)	2 (1%)	17	12
1	B	216/220 (98%)	208 (96%)	7 (3%)	1 (0%)	29	26
All	All	429/440 (98%)	407 (95%)	19 (4%)	3 (1%)	22	18

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24	TYR
1	B	218	PRO
1	A	214	GLU

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	187/195 (96%)	181 (97%)	6 (3%)	39	41
1	B	193/195 (99%)	181 (94%)	12 (6%)	18	15
All	All	380/390 (97%)	362 (95%)	18 (5%)	26	25

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	VAL
1	A	46	ARG
1	A	109	SER

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Mol	Chain	Res	Type
1	A	191	LYS
1	A	192	GLU
1	A	216	GLU
1	B	32	ARG
1	B	45	ILE
1	B	66	ILE
1	B	69	LYS
1	B	160	LEU
1	B	162	GLU
1	B	165	GLU
1	B	188	LYS
1	B	199	GLU
1	B	211	LYS
1	B	215	LYS
1	B	217	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	55	GLN
1	A	72	GLN
1	A	91	ASN
1	A	92	ASN
1	A	141	ASN
1	A	145	GLN
1	A	181	GLN
1	B	9	GLN
1	B	92	ASN
1	B	145	GLN
1	B	181	GLN

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

4 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	GOL	B	221	-	5,5,5	0.94	0	5,5,5	0.25	0
2	GOL	A	221	-	5,5,5	0.95	0	5,5,5	0.23	0
2	GOL	B	223	-	5,5,5	1.08	0	5,5,5	0.29	0
2	GOL	B	222	-	5,5,5	0.83	0	5,5,5	0.21	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	B	221	-	-	4/4/4/4	-
2	GOL	A	221	-	-	4/4/4/4	-
2	GOL	B	223	-	-	2/4/4/4	-
2	GOL	B	222	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	221	GOL	O1-C1-C2-C3
2	A	221	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
2	A	221	GOL	C1-C2-C3-O3
2	B	223	GOL	O1-C1-C2-C3
2	B	222	GOL	O1-C1-C2-C3
2	B	221	GOL	O1-C1-C2-O2
2	B	222	GOL	O1-C1-C2-O2
2	A	221	GOL	O1-C1-C2-O2
2	B	223	GOL	O1-C1-C2-O2
2	A	221	GOL	O2-C2-C3-O3
2	B	221	GOL	C1-C2-C3-O3
2	B	221	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	221	GOL	1	0

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 ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.