



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:47 PM GMT

PDB ID : 1YME
Title : STRUCTURE OF CARBOXYPEPTIDASE
Authors : Greenblatt, H.M.; Tucker, P.A.; Shoham, G.
Deposited on : 1996-07-15
Resolution : 1.53 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

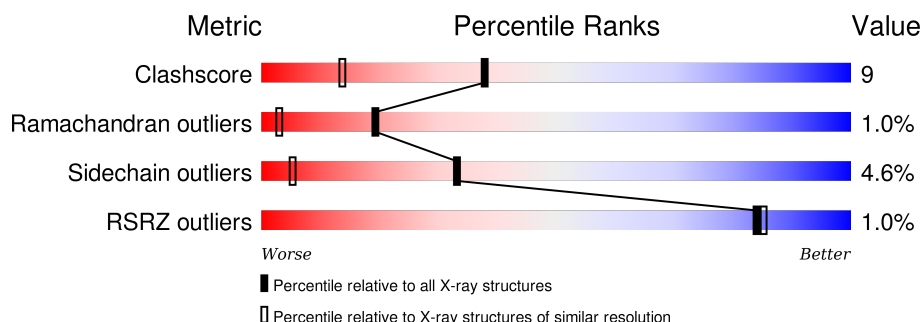
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.53 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	102246	1627 (1.56-1.52)
Ramachandran outliers	100387	1594 (1.56-1.52)
Sidechain outliers	100360	1592 (1.56-1.52)
RSRZ outliers	91569	1555 (1.56-1.52)

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	309	<div> <div></div> <div>74%</div> <div>21%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2650 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 CARBOXYPEPTIDASE A ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2436	1561	403	467	5			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

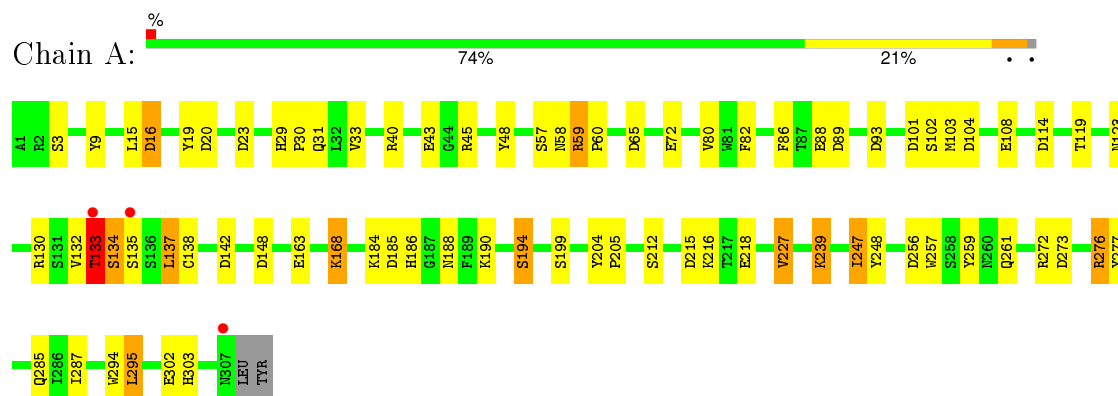
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	213	Total	O	0	0
			213	213		

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: CARBOXYPEPTIDASE A ALPHA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.62Å 60.17Å 47.17Å 90.00° 97.40° 90.00°	Depositor
Resolution (Å)	5.00 – 1.53 60.17 – 1.53	Depositor EDS
% Data completeness (in resolution range)	(Not available) (5.00-1.53) 79.9 (60.17-1.53)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 1.53Å)	Xtriage
Refinement program	PROLSQ, TNT	Depositor
R, R_{free}	0.148 , (Not available) 0.140 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	13.7	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 62.2	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	1 of 35158 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	2650	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.34% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	1.18	7/2502 (0.3%)	1.62	47/3402 (1.4%)

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	1	0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	88	GLU	CD-OE1	6.96	1.33	1.25
1	A	108	GLU	CD-OE1	-6.72	1.18	1.25
1	A	302	GLU	CD-OE2	6.44	1.32	1.25
1	A	218	GLU	CD-OE2	5.41	1.31	1.25
1	A	88	GLU	CD-OE2	-5.40	1.19	1.25
1	A	72	GLU	CD-OE2	-5.36	1.19	1.25
1	A	43	GLU	CD-OE1	5.20	1.31	1.25

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	ARG	NE-CZ-NH1	-8.88	115.86	120.30
1	A	142	ASP	CB-CG-OD2	-8.72	110.45	118.30
1	A	142	ASP	CB-CG-OD1	8.67	126.10	118.30
1	A	130	ARG	NE-CZ-NH2	-8.58	116.01	120.30
1	A	89	ASP	CB-CG-OD1	8.41	125.87	118.30
1	A	185	ASP	CB-CG-OD1	8.14	125.63	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	ASP	CB-CG-OD2	7.89	125.40	118.30
1	A	130	ARG	NE-CZ-NH1	7.81	124.20	120.30
1	A	185	ASP	CB-CG-OD2	-7.74	111.33	118.30
1	A	93	ASP	CB-CG-OD1	7.68	125.21	118.30
1	A	215	ASP	CB-CG-OD1	-7.55	111.50	118.30
1	A	227	VAL	CA-CB-CG1	7.48	122.12	110.90
1	A	93	ASP	CB-CG-OD2	-7.46	111.59	118.30
1	A	135	SER	N-CA-CB	7.44	121.66	110.50
1	A	82	PHE	CB-CG-CD1	7.42	126.00	120.80
1	A	45	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	A	256	ASP	CB-CG-OD1	6.82	124.44	118.30
1	A	89	ASP	CB-CG-OD2	-6.73	112.24	118.30
1	A	16	ASP	CB-CG-OD2	6.68	124.31	118.30
1	A	20	ASP	CB-CG-OD1	6.63	124.27	118.30
1	A	16	ASP	CB-CG-OD1	-6.60	112.36	118.30
1	A	40	ARG	NE-CZ-NH2	-6.58	117.01	120.30
1	A	101	ASP	CB-CG-OD1	6.55	124.20	118.30
1	A	65	ASP	CB-CG-OD1	6.48	124.14	118.30
1	A	23	ASP	CB-CG-OD2	6.44	124.10	118.30
1	A	194	SER	N-CA-CB	6.39	120.08	110.50
1	A	80	VAL	CA-CB-CG2	-6.27	101.50	110.90
1	A	114	ASP	CB-CG-OD2	6.26	123.93	118.30
1	A	104	ASP	CB-CG-OD2	-6.10	112.81	118.30
1	A	148	ASP	CB-CG-OD1	6.05	123.74	118.30
1	A	101	ASP	CB-CG-OD2	-6.04	112.86	118.30
1	A	295	LEU	CB-CG-CD1	6.03	121.25	111.00
1	A	104	ASP	CB-CG-OD1	5.94	123.64	118.30
1	A	259	TYR	CB-CG-CD2	-5.89	117.47	121.00
1	A	59	ARG	NE-CZ-NH2	5.88	123.24	120.30
1	A	65	ASP	CB-CG-OD2	-5.79	113.09	118.30
1	A	138	CYS	CA-CB-SG	-5.57	103.97	114.00
1	A	9	TYR	CB-CG-CD2	-5.44	117.74	121.00
1	A	82	PHE	CB-CG-CD2	-5.36	117.05	120.80
1	A	103	MET	CG-SD-CE	-5.27	91.77	100.20
1	A	102	SER	CB-CA-C	-5.26	100.11	110.10
1	A	227	VAL	CG1-CB-CG2	5.22	119.25	110.90
1	A	48	TYR	CB-CG-CD2	-5.18	117.89	121.00
1	A	134	SER	N-CA-C	5.16	124.92	111.00
1	A	273	ASP	CB-CG-OD2	5.14	122.93	118.30
1	A	3	SER	N-CA-CB	-5.08	102.87	110.50
1	A	133	THR	CB-CA-C	-5.05	97.98	111.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	134	SER	CA

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	2436	0	2346	44	0
2	A	1	0	0	0	0
3	A	213	0	0	10	0
All	All	2650	0	2346	44	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 9.

All (44) 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:137:LEU:HD22	1:A:137:LEU:H	1.25	1.00
1:A:186:HIS:HD2	1:A:188:ASN:H	1.03	0.99
1:A:15:LEU:HD21	1:A:19:TYR:CE2	2.18	0.79
1:A:186:HIS:CD2	1:A:188:ASN:H	1.96	0.76
1:A:303:HIS:ND1	3:A:610:HOH:O	2.23	0.71
1:A:186:HIS:HD2	1:A:188:ASN:N	1.84	0.70
1:A:168:LYS:HG2	3:A:596:HOH:O	1.91	0.69
1:A:137:LEU:N	1:A:137:LEU:HD13	2.08	0.68
1:A:212:SER:HB3	1:A:216:LYS:NZ	2.09	0.68
1:A:272:ARG:HH11	1:A:285:GLN:HE21	1.44	0.66
1:A:132:VAL:HG23	1:A:133:THR:N	2.11	0.65
1:A:15:LEU:O	1:A:15:LEU:HD23	1.97	0.65
1:A:137:LEU:N	1:A:137:LEU:HD22	2.07	0.59
1:A:15:LEU:HD21	1:A:19:TYR:CD2	2.43	0.54
1:A:186:HIS:HE1	3:A:449:HOH:O	1.93	0.52
1:A:15:LEU:HD21	1:A:19:TYR:CZ	2.44	0.51
1:A:239:LYS:HD3	3:A:499:HOH:O	2.11	0.50
1:A:272:ARG:HH11	1:A:285:GLN:NE2	2.08	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:ASP:HA	3:A:592:HOH:O	2.13	0.48
1:A:137:LEU:CD2	1:A:137:LEU:H	2.03	0.48
1:A:303:HIS:HB2	3:A:610:HOH:O	2.12	0.48
1:A:31:GLN:HB2	3:A:552:HOH:O	2.13	0.47
1:A:119:THR:HA	1:A:123:ASN:O	2.15	0.47
1:A:239:LYS:CB	1:A:239:LYS:HZ2	2.28	0.46
1:A:204:TYR:HB2	1:A:205:PRO:CD	2.45	0.46
1:A:212:SER:HB3	1:A:216:LYS:HZ3	1.80	0.46
1:A:239:LYS:CB	1:A:239:LYS:NZ	2.78	0.46
1:A:239:LYS:HZ2	1:A:239:LYS:HB3	1.80	0.46
1:A:59:ARG:HB3	1:A:60:PRO:HD2	1.97	0.45
1:A:163:GLU:HG3	3:A:598:HOH:O	2.17	0.45
1:A:190:LYS:HA	1:A:190:LYS:HD3	1.67	0.45
1:A:86:PHE:HE1	1:A:294:TRP:CZ3	2.35	0.45
1:A:15:LEU:C	1:A:15:LEU:HD23	2.36	0.44
1:A:15:LEU:HD22	3:A:592:HOH:O	2.16	0.44
1:A:212:SER:CB	1:A:216:LYS:HZ1	2.30	0.44
1:A:257:TRP:O	1:A:261:GLN:HG2	2.17	0.44
1:A:247:ILE:HG22	1:A:248:TYR:N	2.34	0.43
1:A:29:HIS:N	1:A:30:PRO:HD3	2.34	0.43
1:A:133:THR:HB	1:A:134:SER:H	0.90	0.42
1:A:204:TYR:HB2	1:A:205:PRO:HD2	2.00	0.42
1:A:276:ARG:HG3	1:A:277:TYR:CE2	2.54	0.42
1:A:15:LEU:CD2	1:A:19:TYR:CD2	3.02	0.42
1:A:132:VAL:O	1:A:133:THR:HG23	2.21	0.41
1:A:216:LYS:HE3	3:A:522:HOH:O	2.22	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	305/309 (99%)	291 (95%)	11 (4%)	3 (1%)	19 3

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	SER
1	A	133	THR
1	A	247	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	263/265 (99%)	251 (95%)	12 (5%)	33 5

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	VAL
1	A	57	SER
1	A	58	ASN
1	A	137	LEU
1	A	168	LYS
1	A	184	LYS
1	A	194	SER
1	A	227	VAL
1	A	239	LYS
1	A	276	ARG
1	A	287	ILE
1	A	295	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN

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Mol	Chain	Res	Type
1	A	37	GLN
1	A	92	GLN
1	A	122	GLN
1	A	186	HIS
1	A	220	ASN
1	A	249	GLN
1	A	285	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 ⓘ

Of 1 ligands modelled in this entry, 1 is 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	307/309 (99%)	-0.53	3 (0%) 84 85	10, 17, 43, 94	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	THR	6.3
1	A	135	SER	2.6
1	A	307	ASN	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 [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(Å ²)	Q<0.9
2	ZN	A	310	1/1	1.00	0.05	-2.53	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.