

Full wwPDB X-ray Structure Validation Report (i)

May 15, 2020 – 11:21 pm BST

PDB ID : 1GCD

Title : REFINED CRYSTAL STRUCTURE OF "AGED" AND "NON-AGED"

ORGANOPHOSPHORYL CONJUGATES OF GAMMA-CHYMOTRYPSIN

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Deposited on : 1994-05-05

Resolution : 1.90 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 (i) symbol.

The following versions of software and data (see references (1)) 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)

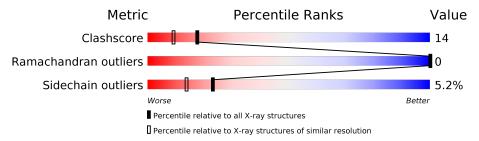
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain		
1	Λ	245	73%	20%	



2 Entry composition (i)

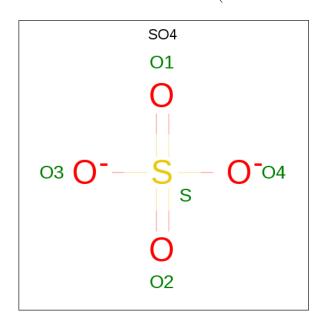
There are 4 unique types of molecules in this entry. The entry contains 2045 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GAMMA-CHYMOTRYPSIN.

Mol	Chain	Residues		Atoms		ZeroOcc	AltConf	Trace		
1	Λ	237	Total	С	N	О	S	0	0	0
1	A	237	1744	1096	295	341	12	0	U	0

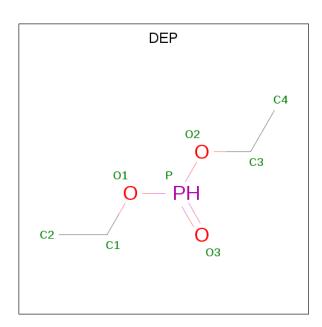
• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0

• Molecule 3 is DIETHYL PHOSPHONATE (three-letter code: DEP) (formula: C₄H₁₁O₃P).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C 1	3 ()	P 1	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	278	Total O 278 278	0	0

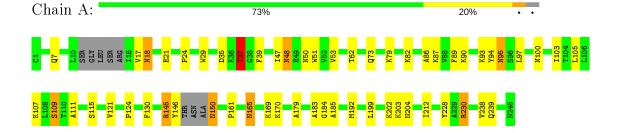


3 Residue-property plots (i)

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: GAMMA-CHYMOTRYPSIN





4 Data and refinement statistics (i)

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

Property	Value	Source	
Space group	P 42 21 2	Depositor	
Cell constants	69.00Å 69.00Å 95.30Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	6.00 - 1.90	Depositor	
% Data completeness	(Not available) (6.00-1.90)	Depositor	
(in resolution range)	(1101 available) (0.00 1.50)	Беровног	
R_{merge}	(Not available)	Depositor	
R_{sym}	(Not available)	Depositor	
Refinement program	PROLSQ	Depositor	
R, R_{free}	0.187 , (Not available)	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	2045	wwPDB-VP	
Average B, all atoms (Å ²)	13.0	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: DEP, SO4

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		
MIOI	Chain	RMSZ	# Z >5	RMSZ	# Z > 5	
1	A	0.99	0/1778	1.42	7/2422 (0.3%)	

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^o)$
1	Α	21	GLU	OE1-CD-OE2	9.14	134.27	123.30
1	A	145	ARG	RG NE-CZ-NH2		123.27	120.30
1	A	21	GLU	CG-CD-OE1	-5.81	106.69	118.30
1	A	37	THR	N-CA-CB	-5.27	100.29	110.30
1	A	230	ARG	NE-CZ-NH2	5.26	122.93	120.30
1	A	199	LEU	N-CA-CB	5.24	120.89	110.40
1	A	39	PHE	CB-CG-CD2	5.16	124.41	120.80

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	1744	0	1719	45	0
2	A	15	0	0	1	0
3	A	8	0	10	3	0
4	A	278	0	0	12	0

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\mathbf{Mol}	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
All	All	2045	0	1729	48	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 14.

All (48) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance} ({f A})$	overlap(A)
1:A:17:VAL:O	1:A:18:ASN:HB2	1.76	0.86
1:A:79:LYS:HG3	4:A:555:HOH:O	1.85	0.76
1:A:170:LYS:HE3	4:A:487:HOH:O	1.95	0.66
1:A:202:LYS:HE2	4:A:545:HOH:O	1.95	0.65
1:A:48:ASN:HD22	1:A:50:ASN:H	1.43	0.65
1:A:93:LYS:HD3	4:A:520:HOH:O	1.96	0.64
1:A:35:ASP:OD1	1:A:37:THR:HB	1.99	0.63
1:A:100:ASN:HD21	1:A:179:ALA:HB3	1.65	0.62
1:A:145:ARG:HB3	1:A:150:ASN:OD1	1.99	0.61
1:A:95:ASN:HD21	1:A:97:LEU:HB2	1.66	0.60
1:A:90:LYS:HD2	1:A:94:TYR:HD2	1.67	0.60
1:A:145:ARG:CZ	1:A:150:ASN:OD1	2.52	0.57
1:A:50:ASN:HD21	1:A:111:ALA:HB2	1.69	0.57
1:A:17:VAL:O	1:A:18:ASN:CB	2.51	0.57
1:A:203:LYS:O	1:A:204:ASN:C	2.45	0.55
1:A:87:LYS:HG3	1:A:89:PHE:CE2	2.42	0.55
1:A:146:TYR:HD1	1:A:192:MET:HE1	1.73	0.54
1:A:86:ALA:HB2	1:A:109:SER:HA	1.89	0.54
1:A:7:GLN:HG2	4:A:514:HOH:O	2.08	0.53
1:A:100:ASN:ND2	1:A:179:ALA:HB3	2.24	0.52
1:A:48:ASN:ND2	1:A:50:ASN:H	2.05	0.52
1:A:51:TRP:CZ2	1:A:107:LYS:HD2	2.45	0.52
3:A:290:DEP:H12	3:A:290:DEP:H32	1.92	0.51
1:A:105:LEU:HD11	1:A:238:VAL:HA	1.91	0.51
1:A:130:PHE:O	4:A:468:HOH:O	2.19	0.50
1:A:183:ALA:HB3	1:A:228:TYR:CE1	2.46	0.50
1:A:62:THR:HG23	4:A:568:HOH:O	2.12	0.50
1:A:165:ASN:O	1:A:169:LYS:HG2	2.12	0.49
1:A:90:LYS:HD2	1:A:94:TYR:CD2	2.47	0.49
1:A:115:SER:HB2	2:A:296:SO4:O3	2.13	0.48
1:A:73:GLN:NE2	4:A:333:HOH:O	2.44	0.47
1:A:202:LYS:CE	4:A:545:HOH:O	2.59	0.47
1:A:29:TRP:CG	1:A:121:VAL:HB	2.51	0.46

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Atom-1	Atom-2	Interatomic	Clash
1100111 1	1100111 =	$\operatorname{distance}\left(\mathrm{\AA} ight)$	overlap(A)
3:A:290:DEP:H21	4:A:564:HOH:O	2.14	0.46
1:A:48:ASN:C	1:A:48:ASN:HD22	2.17	0.46
1:A:48:ASN:HD21	1:A:51:TRP:HD1	1.62	0.46
1:A:95:ASN:C	1:A:95:ASN:HD22	2.20	0.45
1:A:24:PRO:HD3	4:A:310:HOH:O	2.16	0.45
1:A:165:ASN:ND2	1:A:230:ARG:HH21	2.15	0.45
3:A:290:DEP:H12	3:A:290:DEP:C3	2.46	0.44
1:A:146:TYR:CD1	1:A:192:MET:HE1	2.52	0.44
1:A:82:LYS:HE3	4:A:455:HOH:O	2.17	0.43
1:A:145:ARG:NE	1:A:150:ASN:OD1	2.52	0.43
1:A:103:ILE:HG13	1:A:212:ILE:CD1	2.48	0.43
1:A:184:GLY:O	1:A:185:ALA:HB3	2.19	0.42
1:A:47:ILE:CG2	1:A:53:VAL:HG23	2.50	0.41
1:A:47:ILE:HG23	1:A:53:VAL:HG23	2.03	0.41
1:A:145:ARG:HB3	1:A:150:ASN:CB	2.51	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	Percei	ntiles
1	A	231/245 (94%)	223 (96%)	8 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	Analysed Rotameric		Percentiles		
1	A	194/200 (97%)	184 (95%)	10 (5%)	23 14		

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	37	THR
1	A	48	ASN
1	A	95	ASN
1	A	109	SER
1	A	124	PRO
1	A	150	ASN
1	A	161	PRO
1	A	165	ASN
1	A	239	GLN

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	7	GLN
1	A	48	ASN
1	A	50	ASN
1	A	73	GLN
1	A	95	ASN
1	A	100	ASN
1	A	116	GLN
1	A	165	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

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	Tuno	Chain	Res	Res Link Bond lengths			Bond angles			
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
2	SO4	A	298	-	4,4,4	0.58	0	6,6,6	0.42	0
2	SO4	A	296	_	4,4,4	0.62	0	6,6,6	0.25	0
2	SO4	A	297	_	4,4,4	0.67	0	6,6,6	0.32	0
3	DEP	A	290	1	4,7,7	5.55	2 (50%)	2,7,7	1.67	1 (50%)

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
3	DEP	A	290	1	-	0/2/6/6	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
3	A	290	DEP	O2-C3	-9.98	1.20	1.44
3	A	290	DEP	O1-C1	-4.65	1.33	1.44

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
3	A	290	DEP	O2-C3-C4	2.36	128.69	110.88

There are no chirality outliers.

There are no torsion outliers.



There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	296	SO4	1	0
3	A	290	DEP	3	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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

