



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 12:18 PM EDT

PDB ID : 2Y50 / pdb_00002y50
Title : Crystal Structure of Collagenase G from Clostridium histolyticum at 2. 80 Angstrom Resolution
Authors : Eckhard, U.; Brandstetter, H.
Deposited on : 2011-01-11
Resolution : 2.80 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

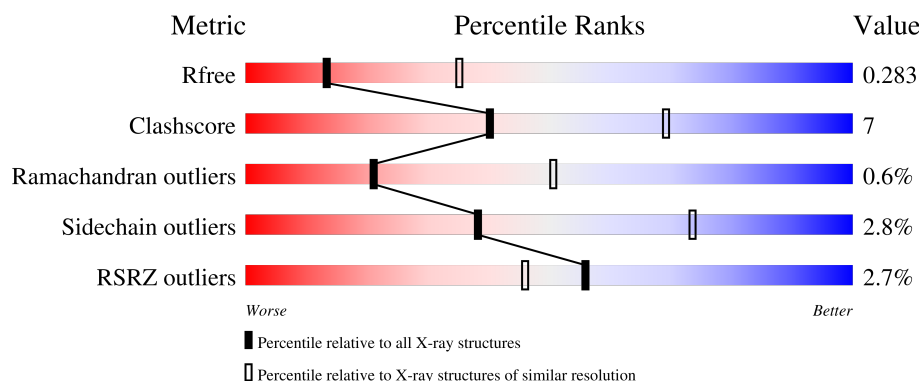
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 of 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 $\leq 5\%$. 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	785	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5458 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 COLLAGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	672	5300	3398	850	1039	13	0	0	0

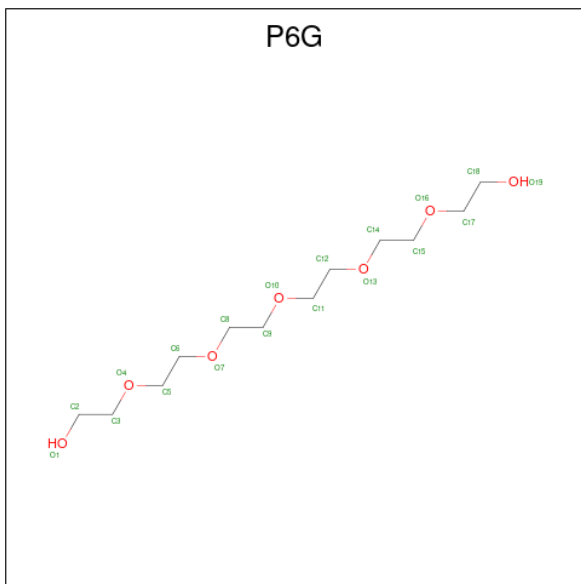
There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	96	MET	-	expression tag	UNP Q9X721
A	97	GLY	-	expression tag	UNP Q9X721
A	98	SER	-	expression tag	UNP Q9X721
A	99	SER	-	expression tag	UNP Q9X721
A	100	HIS	-	expression tag	UNP Q9X721
A	101	HIS	-	expression tag	UNP Q9X721
A	102	HIS	-	expression tag	UNP Q9X721
A	103	HIS	-	expression tag	UNP Q9X721
A	104	HIS	-	expression tag	UNP Q9X721
A	105	HIS	-	expression tag	UNP Q9X721
A	106	SER	-	expression tag	UNP Q9X721
A	107	SER	-	expression tag	UNP Q9X721
A	108	GLY	-	expression tag	UNP Q9X721
A	109	GLU	-	expression tag	UNP Q9X721
A	110	ASN	-	expression tag	UNP Q9X721
A	111	LEU	-	expression tag	UNP Q9X721
A	112	TYR	-	expression tag	UNP Q9X721
A	113	PHE	-	expression tag	UNP Q9X721
A	114	GLN	-	expression tag	UNP Q9X721
A	115	GLY	-	expression tag	UNP Q9X721
A	116	GLY	-	expression tag	UNP Q9X721
A	117	THR	-	expression tag	UNP Q9X721
A	118	MET	-	expression tag	UNP Q9X721
A	575	LEU	SER	SEE REMARK 999	UNP Q9X721
A	665	VAL	ALA	SEE REMARK 999	UNP Q9X721
A	790	GLY	ALA	SEE REMARK 999	UNP Q9X721

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

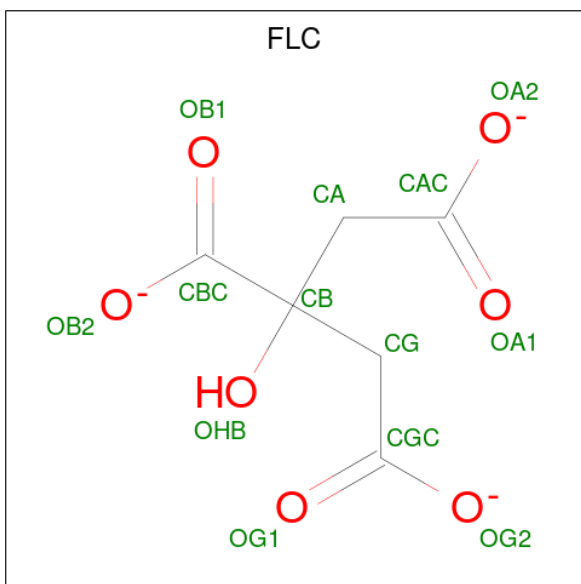
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: C₁₂H₂₆O₇).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			19	12 7		

- Molecule 4 is CITRATE ANION (CCD ID: FLC) (formula: C₆H₅O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	6	7		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	125	Total	O	0	0
			125	125		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.26Å 108.70Å 181.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.83 – 2.80 34.83 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.0 (34.83-2.80) 98.9 (34.83-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.232 , 0.282 0.237 , 0.283	Depositor DCC
R_{free} test set	1433 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.414	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 69.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5458	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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, 0.333 respectively for untwinned datasets, and 0.375, 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, FLC, P6G

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	0.44	1/5425 (0.0%)	0.82	1/7367 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	263	VAL	CA-CB	5.16	1.56	1.54

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	502	THR	N-CA-C	6.25	119.13	109.07

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	5300	0	4905	68	0
2	A	1	0	0	0	0
3	A	19	0	26	0	0
4	A	13	0	5	1	0
5	A	125	0	0	5	0
All	All	5458	0	4936	68	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 7.

All (68) 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:508:ARG:HG2	1:A:508:ARG:HH11	1.11	1.09
1:A:508:ARG:HD3	1:A:513:SER:HB3	1.58	0.84
1:A:508:ARG:HH11	1:A:508:ARG:CG	1.93	0.77
1:A:615:HIS:HD2	1:A:660:HIS:CE1	2.05	0.74
1:A:426:ALA:HB1	1:A:430:VAL:HG11	1.70	0.71
1:A:475:SER:HB2	1:A:476:PRO:HD2	1.74	0.70
1:A:508:ARG:HG2	1:A:508:ARG:NH1	1.92	0.70
1:A:495:LEU:HD21	1:A:497:ILE:HD11	1.75	0.68
1:A:615:HIS:HD2	1:A:660:HIS:HE1	1.42	0.67
1:A:370:TYR:O	1:A:376:ALA:HB2	1.94	0.66
1:A:189:PHE:O	5:A:2005:HOH:O	2.13	0.65
1:A:450:HIS:CE1	1:A:456:ASP:HA	2.32	0.64
1:A:255:ASN:OD1	1:A:258:VAL:HG12	2.00	0.60
1:A:416:THR:HG22	1:A:420:GLY:HA2	1.83	0.59
1:A:449:PHE:HA	1:A:561:PHE:CE1	2.38	0.59
1:A:506:TYR:O	1:A:508:ARG:NH1	2.36	0.58
1:A:770:VAL:HG13	1:A:774:ASN:HA	1.85	0.58
1:A:539:TRP:HZ3	5:A:2023:HOH:O	1.86	0.58
1:A:480:LYS:O	1:A:483:THR:HG22	2.04	0.57
1:A:282:LYS:N	5:A:2017:HOH:O	2.37	0.56
1:A:586:ASP:O	1:A:588:ARG:N	2.39	0.56
1:A:676:VAL:HG13	1:A:680:TYR:CD2	2.42	0.55
1:A:255:ASN:O	1:A:256:ALA:CB	2.56	0.54
1:A:433:GLU:O	1:A:437:ARG:HG3	2.09	0.53
1:A:255:ASN:O	1:A:256:ALA:HB2	2.09	0.53
1:A:222:MET:HE2	1:A:258:VAL:HG21	1.91	0.53
1:A:585:VAL:O	1:A:586:ASP:C	2.53	0.52
1:A:622:MET:HE3	1:A:626:ILE:HD11	1.92	0.52
1:A:712:SER:HB2	1:A:715:PHE:O	2.09	0.51
1:A:582:LYS:HE2	1:A:605:MET:HE2	1.92	0.51
1:A:426:ALA:HB1	1:A:430:VAL:CG1	2.37	0.50
1:A:309:MET:HE2	1:A:311:TRP:CH2	2.46	0.50
1:A:522:ARG:NH2	1:A:564:SER:O	2.45	0.50
1:A:508:ARG:CG	1:A:508:ARG:NH1	2.62	0.50
1:A:508:ARG:HD2	1:A:515:PHE:O	2.12	0.49
1:A:437:ARG:NH2	1:A:678:ASP:OD1	2.46	0.48
1:A:553:PHE:CE2	1:A:628:MET:HE3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:THR:CG2	1:A:420:GLY:HA2	2.42	0.48
1:A:279:ASP:C	5:A:2017:HOH:O	2.56	0.48
1:A:505:THR:HG23	1:A:506:TYR:N	2.28	0.47
1:A:615:HIS:CD2	1:A:660:HIS:HE1	2.27	0.47
1:A:164:VAL:HA	1:A:167:ILE:HD12	1.98	0.46
1:A:315:ILE:HD11	1:A:352:TYR:HB2	1.98	0.46
1:A:763:ALA:HA	1:A:781:VAL:O	2.16	0.46
1:A:657:TYR:CZ	1:A:661:ILE:HD11	2.51	0.45
1:A:427:GLY:O	1:A:430:VAL:HG12	2.16	0.45
1:A:189:PHE:C	5:A:2005:HOH:O	2.56	0.45
1:A:511:GLN:CD	1:A:511:GLN:N	2.75	0.44
1:A:553:PHE:HE2	1:A:628:MET:HE3	1.82	0.44
1:A:620:LYS:HD3	1:A:660:HIS:CD2	2.52	0.44
1:A:582:LYS:HE3	1:A:609:TYR:OH	2.18	0.44
1:A:380:MET:HE3	1:A:399:HIS:ND1	2.33	0.43
1:A:256:ALA:HB2	1:A:311:TRP:HB3	2.01	0.43
1:A:282:LYS:H	1:A:282:LYS:HG2	1.64	0.42
1:A:736:TRP:HZ3	4:A:1793:FLC:HA1	1.84	0.42
1:A:508:ARG:HD3	1:A:513:SER:CB	2.39	0.42
1:A:573:ARG:HH21	1:A:576:ILE:HD11	1.84	0.42
1:A:222:MET:HE3	1:A:249:ILE:HG12	2.02	0.42
1:A:555:GLU:O	1:A:559:GLU:HG2	2.19	0.42
1:A:277:ALA:HB3	1:A:335:THR:HG22	2.01	0.41
1:A:470:MET:CE	1:A:521:PHE:CE1	3.03	0.41
1:A:718:PHE:HB3	1:A:786:LEU:HD13	2.02	0.41
1:A:372:TYR:O	1:A:410:TYR:OH	2.24	0.41
1:A:617:LEU:O	1:A:621:ASP:C	2.64	0.41
1:A:443:ARG:HD3	1:A:443:ARG:HA	1.82	0.41
1:A:523:HIS:C	1:A:523:HIS:CD2	2.98	0.40
1:A:470:MET:HE1	1:A:521:PHE:CE1	2.56	0.40
1:A:481:PHE:HA	1:A:484:ASN:HD22	1.87	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	668/785 (85%)	636 (95%)	28 (4%)	4 (1%)	21	51

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	587	HIS
1	A	256	ALA
1	A	586	ASP
1	A	234	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	532/672 (79%)	517 (97%)	15 (3%)	38	73

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	151	SER
1	A	416	THR
1	A	430	VAL
1	A	446	LYS
1	A	461	VAL
1	A	490	THR
1	A	498	GLU
1	A	508	ARG
1	A	591	LEU
1	A	655	THR
1	A	676	VAL
1	A	740	SER
1	A	749	SER
1	A	770	VAL

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

Mol	Chain	Res	Type
1	A	230	ASN
1	A	238	GLN
1	A	320	ASN
1	A	448	GLN
1	A	484	ASN
1	A	596	ASN
1	A	608	ASN
1	A	615	HIS
1	A	634	ASN
1	A	660	HIS
1	A	767	ASN
1	A	777	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
3	P6G	A	1792	-	18,18,18	0.48	0	17,17,17	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FLC	A	1793	-	12,12,12	1.03	0	17,17,17	1.39	1 (5%)

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	P6G	A	1792	-	-	10/16/16/16	-
4	FLC	A	1793	-	-	10/16/16/16	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1793	FLC	OB2-CBC-CB	3.34	119.54	113.14

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1793	FLC	CAC-CA-CB-CBC
4	A	1793	FLC	CAC-CA-CB-CG
4	A	1793	FLC	CAC-CA-CB-OHB
4	A	1793	FLC	CG-CB-CBC-OB1
4	A	1793	FLC	CG-CB-CBC-OB2
4	A	1793	FLC	OHB-CB-CBC-OB1
4	A	1793	FLC	OHB-CB-CBC-OB2
3	A	1792	P6G	O7-C8-C9-O10
3	A	1792	P6G	O1-C2-C3-O4
3	A	1792	P6G	C2-C3-O4-C5
3	A	1792	P6G	C14-C15-O16-C17
3	A	1792	P6G	O13-C14-C15-O16
4	A	1793	FLC	CA-CB-CBC-OB1
4	A	1793	FLC	CA-CB-CBC-OB2
3	A	1792	P6G	C18-C17-O16-C15
3	A	1792	P6G	C15-C14-O13-C12
3	A	1792	P6G	C5-C6-O7-C8
3	A	1792	P6G	C11-C12-O13-C14
3	A	1792	P6G	C8-C9-O10-C11

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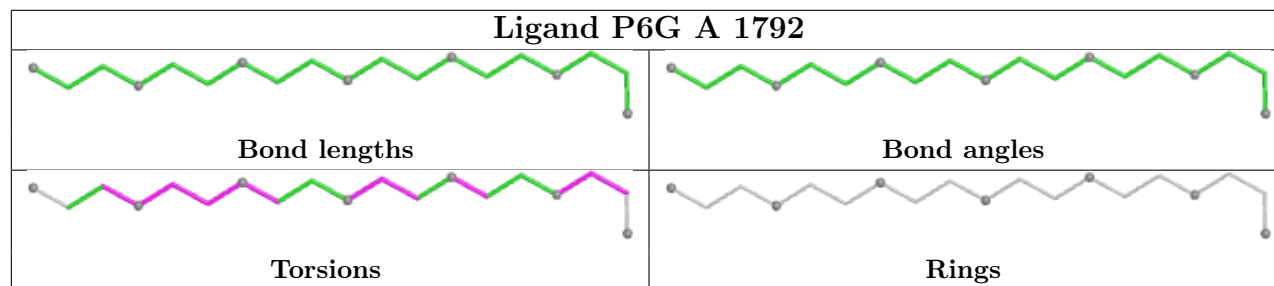
Mol	Chain	Res	Type	Atoms
4	A	1793	FLC	CB-CG-CGC-OG1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1793	FLC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 ⓘ

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	672/785 (85%)	0.37	18 (2%) 56 46	39, 71, 105, 138	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	586	ASP	6.3
1	A	202	TYR	3.9
1	A	585	VAL	3.9
1	A	601	ASP	3.3
1	A	182	ASP	3.2
1	A	587	HIS	2.9
1	A	597	SER	2.8
1	A	600	ASP	2.6
1	A	126	GLY	2.6
1	A	181	ASN	2.5
1	A	683	ASP	2.4
1	A	117	THR	2.4
1	A	122	GLU	2.4
1	A	790	GLY	2.3
1	A	575	LEU	2.3
1	A	582	LYS	2.2
1	A	754	SER	2.2
1	A	679	ASP	2.0

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 [i](#)

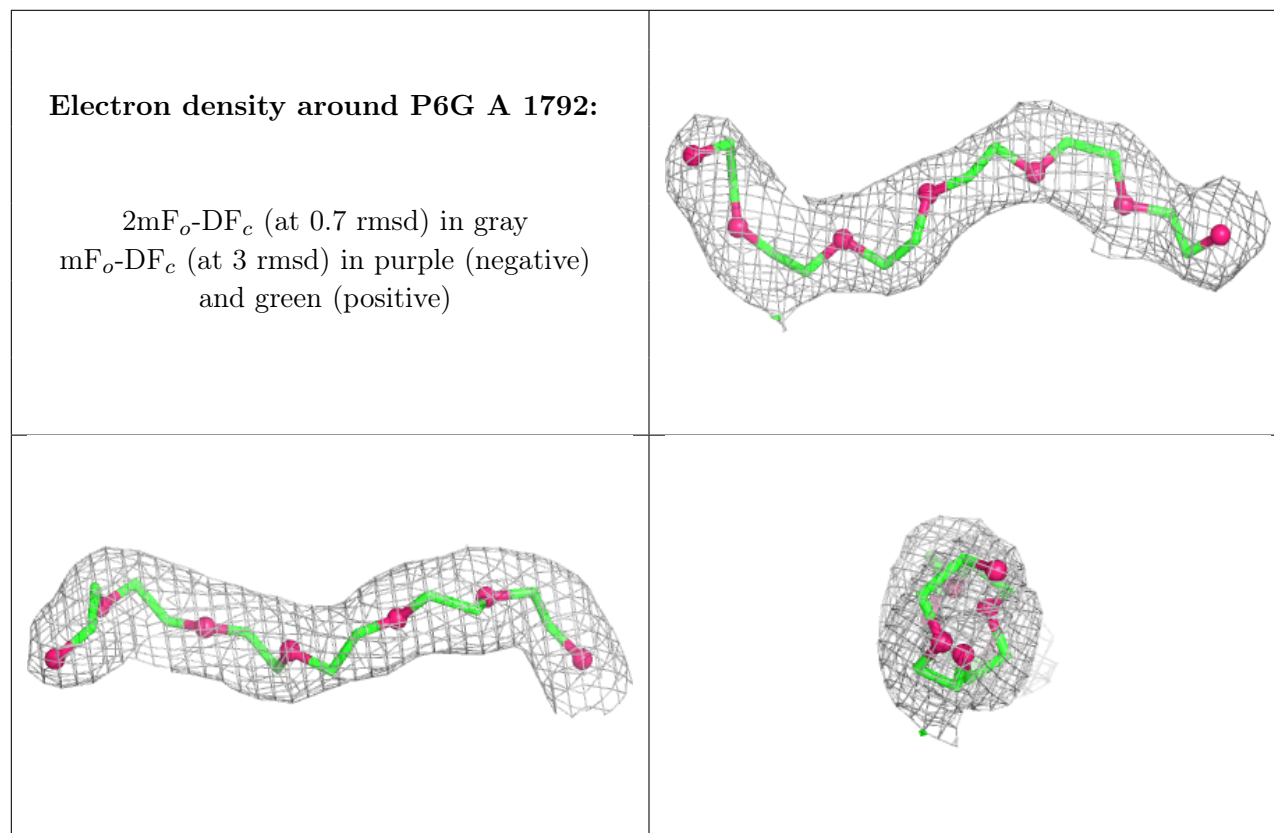
There are no oligosaccharides in this entry.

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. 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	B-factors($\text{\AA}^2$)	Q<0.9
4	FLC	A	1793	13/13	0.85	0.18	84,85,86,86	0
3	P6G	A	1792	19/19	0.92	0.11	72,73,75,76	0
2	ZN	A	1791	1/1	0.96	0.22	39,39,39,39	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.