Full wwPDB X-ray Structure Validation Report

May 22, 2020 – 01:12 am BST

PDB ID	:	2Y3U
Title	:	Crystal structure of apo collagenase G from Clostridium histolyticum at 2.55 Angstrom resolution
Authors	:	Eckhard, U.; Brandstetter, H.
Deposited on	:	2010-12-26
Resolution	:	2.55 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:

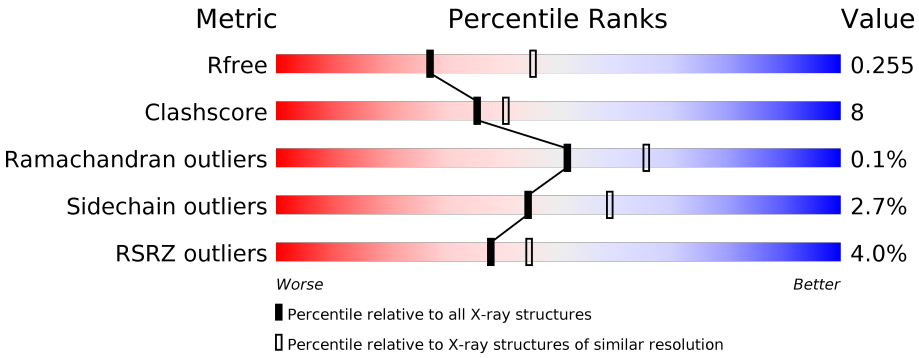
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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 2.55 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}	130704	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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 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% 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	3% 68% 17% • 14%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5615 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.

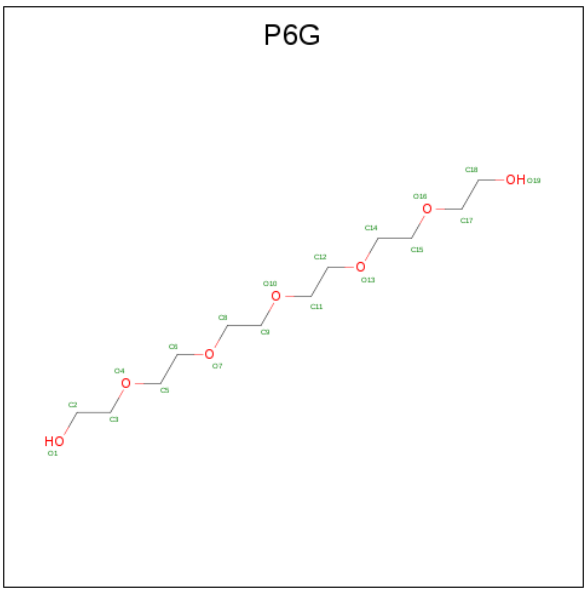
- Molecule 1 is a protein called COLLAGENASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	679	5490	3511	900	1066	2	11	128	0	0

There are 26 discrepancies between the modelled and reference sequences:

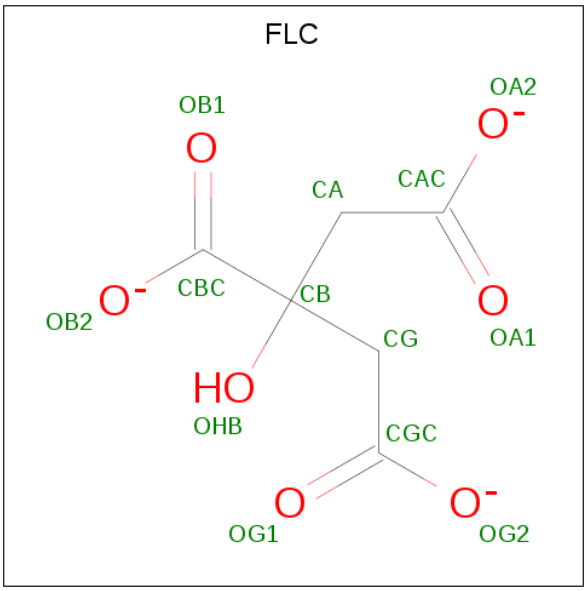
Chain	Residue	Modelled	Actual	Comment	Reference
A	96	MSE	-	expression tag	UNP Q9X721
A	97	GLY	-	expression tag	UNP Q9X721
A	98	SER	-	expression tag	UNP Q9X721
A	99	SER	-	expression tag	UNP Q9X721
A	100	HIS	-	expression tag	UNP Q9X721
A	101	HIS	-	expression tag	UNP Q9X721
A	102	HIS	-	expression tag	UNP 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	MSE	-	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 HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $C_{12}H_{26}O_7$).



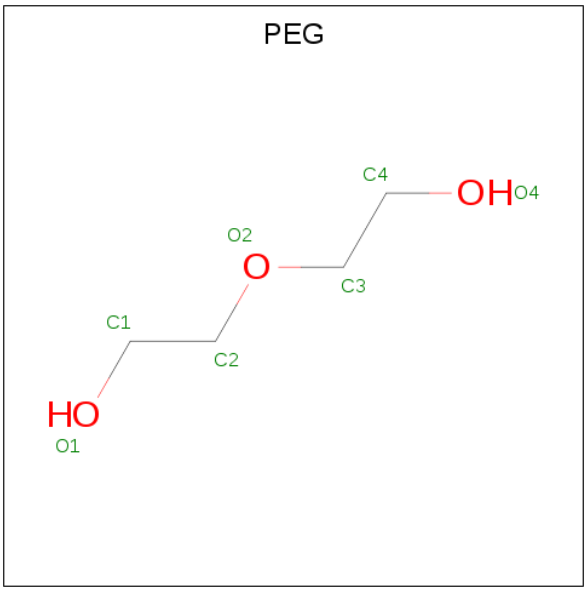
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 19 12 7	0	0

- Molecule 3 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7$).



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

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is water.

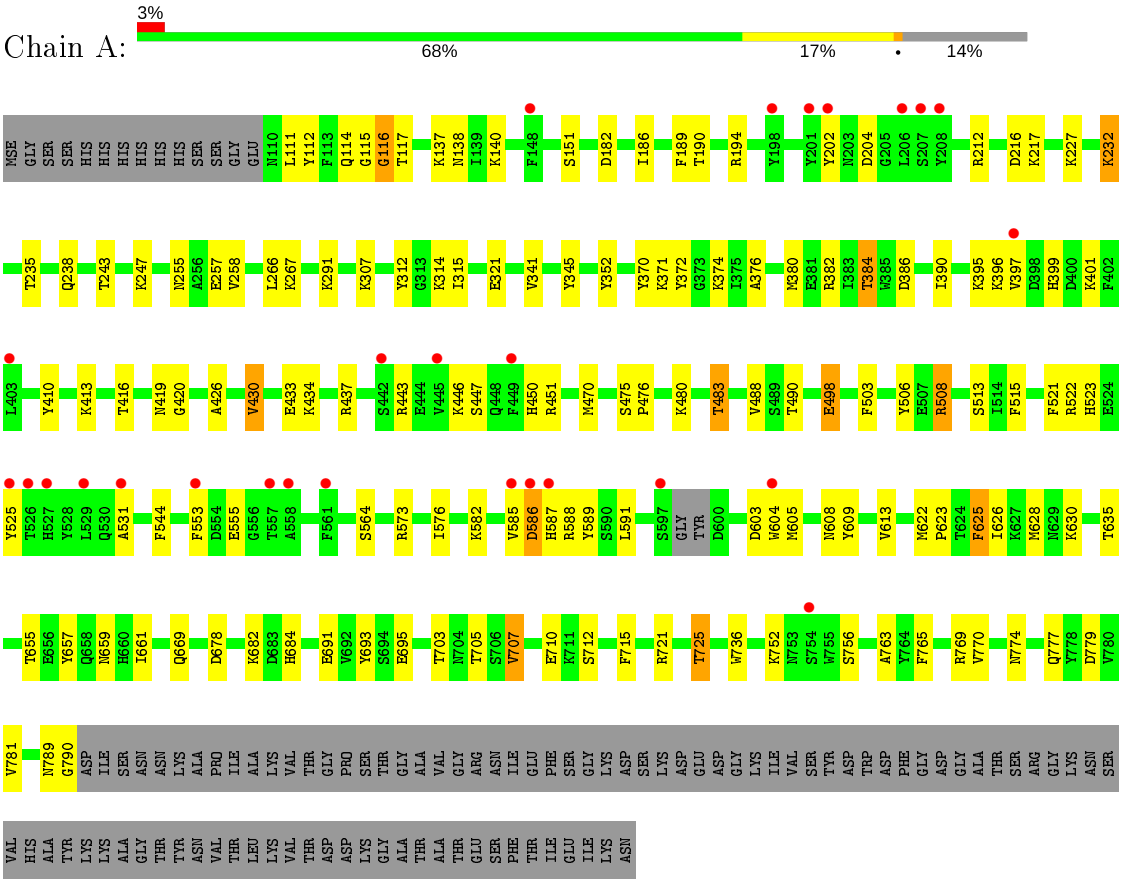
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	86	Total	O	0	0
			86	86		

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 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: COLLAGENASE

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	56.96Å 109.05Å 182.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.55 38.85 – 2.55	Depositor EDS
% Data completeness (in resolution range)	87.1 (40.00-2.55) 87.1 (38.85-2.55)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.72 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.208 , 0.256 0.209 , 0.255	Depositor DCC
R_{free} test set	1626 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å^2)	57.9	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 36.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5615	wwPDB-VP
Average B, all atoms (Å^2)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: FLC, PEG, 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	1.08	29/5607 (0.5%)	1.16	35/7562 (0.5%)

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	0	1

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	202	TYR	CG-CD2	40.15	1.91	1.39
1	A	202	TYR	CG-CD1	-26.38	1.04	1.39
1	A	371	LYS	CE-NZ	-20.17	0.98	1.49
1	A	112	TYR	CE2-CZ	18.24	1.62	1.38
1	A	112	TYR	CG-CD2	14.40	1.57	1.39
1	A	112	TYR	CG-CD1	14.11	1.57	1.39
1	A	401	LYS	CD-CE	12.82	1.83	1.51
1	A	116	GLY	N-CA	12.36	1.64	1.46
1	A	374	LYS	CD-CE	-11.40	1.22	1.51
1	A	390	ILE	CG1-CD1	-9.87	0.82	1.50
1	A	585	VAL	CB-CG2	-9.67	1.32	1.52
1	A	112	TYR	CE1-CZ	9.46	1.50	1.38
1	A	227	LYS	CD-CE	-8.87	1.29	1.51
1	A	111	LEU	CB-CG	8.81	1.78	1.52
1	A	116	GLY	C-N	-8.20	1.15	1.34
1	A	232	LYS	CG-CD	-7.53	1.26	1.52
1	A	267	LYS	CD-CE	7.11	1.69	1.51
1	A	396	LYS	CG-CD	-6.90	1.28	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	314	LYS	CB-CG	-6.75	1.34	1.52
1	A	669	GLN	CB-CG	6.68	1.70	1.52
1	A	682	LYS	CE-NZ	-6.38	1.33	1.49
1	A	413	LYS	CG-CD	-6.31	1.31	1.52
1	A	307	LYS	CG-CD	-6.22	1.31	1.52
1	A	752	LYS	CD-CE	-6.20	1.35	1.51
1	A	194	ARG	CB-CG	-5.75	1.37	1.52
1	A	291	LYS	CE-NZ	-5.43	1.35	1.49
1	A	182	ASP	CB-CG	-5.36	1.40	1.51
1	A	630	LYS	CG-CD	-5.23	1.34	1.52
1	A	204	ASP	CG-OD1	-5.14	1.13	1.25

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	202	TYR	CB-CG-CD1	52.67	152.60	121.00
1	A	202	TYR	CB-CG-CD2	-30.73	102.56	121.00
1	A	202	TYR	CG-CD1-CE1	24.43	140.84	121.30
1	A	204	ASP	CB-CG-OD1	-18.25	101.87	118.30
1	A	585	VAL	CA-CB-CG1	-17.11	85.24	110.90
1	A	371	LYS	CD-CE-NZ	13.67	143.14	111.70
1	A	204	ASP	OD1-CG-OD2	13.55	149.04	123.30
1	A	116	GLY	C-N-CA	13.05	154.34	121.70
1	A	202	TYR	CD1-CG-CD2	-12.90	103.71	117.90
1	A	182	ASP	CB-CG-OD1	-11.45	108.00	118.30
1	A	182	ASP	CB-CG-OD2	11.40	128.56	118.30
1	A	204	ASP	CB-CG-OD2	-10.72	108.65	118.30
1	A	395	LYS	CA-CB-CG	10.59	136.69	113.40
1	A	232	LYS	CB-CG-CD	9.93	137.41	111.60
1	A	116	GLY	O-C-N	-9.54	107.44	122.70
1	A	659	ASN	CB-CG-OD1	-9.28	103.04	121.60
1	A	140	LYS	CA-CB-CG	9.16	133.56	113.40
1	A	669	GLN	CA-CB-CG	-9.04	93.51	113.40
1	A	585	VAL	CA-CB-CG2	-8.96	97.46	110.90
1	A	401	LYS	CG-CD-CE	-8.49	86.42	111.90
1	A	112	TYR	CB-CG-CD2	-7.49	116.50	121.00
1	A	202	TYR	CG-CD2-CE2	-7.40	115.38	121.30
1	A	659	ASN	CB-CG-ND2	7.09	133.72	116.70
1	A	112	TYR	CD1-CG-CD2	6.75	125.32	117.90
1	A	390	ILE	CB-CG1-CD1	6.43	131.91	113.90
1	A	116	GLY	CA-C-N	6.41	131.30	117.20
1	A	232	LYS	CG-CD-CE	6.26	130.69	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	659	ASN	CA-CB-CG	5.56	125.63	113.40
1	A	395	LYS	CB-CG-CD	5.47	125.84	111.60
1	A	112	TYR	CG-CD1-CE1	-5.42	116.96	121.30
1	A	112	TYR	CG-CD2-CE2	-5.40	116.98	121.30
1	A	585	VAL	CG1-CB-CG2	5.40	119.54	110.90
1	A	217	LYS	CG-CD-CE	-5.32	95.94	111.90
1	A	267	LYS	CG-CD-CE	-5.19	96.32	111.90
1	A	140	LYS	CB-CG-CD	5.08	124.80	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	116	GLY	Peptide

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	5490	0	5295	83	0
2	A	19	0	26	0	0
3	A	13	0	5	3	0
4	A	7	0	10	0	0
5	A	86	0	0	3	0
All	All	5615	0	5336	85	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 8.

All (85) 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:HH11	1:A:508:ARG:HG2	0.97	1.11
1:A:508:ARG:HH11	1:A:508:ARG:CG	1.84	0.90
1:A:506:TYR:O	1:A:508:ARG:NH1	2.07	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:508:ARG:HG2	1:A:508:ARG:NH1	1.79	0.84
1:A:437:ARG:NH2	1:A:678:ASP:OD1	2.12	0.81
1:A:434:LYS:HE2	5:A:2043:HOH:O	1.80	0.80
1:A:370:TYR:O	1:A:376:ALA:HB2	1.87	0.74
1:A:426:ALA:HB1	1:A:430:VAL:HG11	1.74	0.70
1:A:114:GLN:HG3	1:A:115:GLY:H	1.56	0.69
1:A:684:HIS:HD2	1:A:756:SER:OG	1.74	0.69
1:A:483:THR:HB	1:A:488:VAL:O	1.93	0.69
1:A:255:ASN:HB2	1:A:258:VAL:HG23	1.74	0.68
1:A:604:TRP:CZ2	1:A:605:MSE:HE2	2.30	0.67
3:A:1793:FLC:OA1	3:A:1793:FLC:OHB	2.07	0.66
1:A:586:ASP:C	1:A:588:ARG:H	2.01	0.64
1:A:586:ASP:O	1:A:588:ARG:N	2.29	0.61
1:A:586:ASP:C	1:A:588:ARG:N	2.54	0.60
1:A:522:ARG:NH2	1:A:564:SER:O	2.35	0.59
1:A:508:ARG:HD3	1:A:513:SER:HB3	1.84	0.59
1:A:426:ALA:HB1	1:A:430:VAL:CG1	2.32	0.59
1:A:582:LYS:HD3	1:A:605:MSE:HE3	1.87	0.57
1:A:372:TYR:O	1:A:410:TYR:OH	2.13	0.56
1:A:380:MSE:HE1	1:A:397:VAL:CG1	2.36	0.56
1:A:380:MSE:HE2	1:A:399:HIS:HA	1.88	0.55
1:A:589:TYR:OH	1:A:603:ASP:OD1	2.25	0.55
1:A:712:SER:HB2	1:A:715:PHE:O	2.08	0.54
1:A:508:ARG:HD2	1:A:515:PHE:O	2.08	0.54
1:A:555:GLU:HG3	5:A:2051:HOH:O	2.08	0.54
1:A:266:LEU:HB3	1:A:321:GLU:HG3	1.90	0.54
1:A:380:MSE:HE1	1:A:397:VAL:HG12	1.90	0.53
1:A:622:MSE:HE3	1:A:626:ILE:HG13	1.89	0.52
1:A:770:VAL:HG13	1:A:774:ASN:HA	1.91	0.52
1:A:691:GLU:O	1:A:695:GLU:HG3	2.10	0.52
1:A:341:VAL:HG13	1:A:345:TYR:HD2	1.74	0.51
1:A:736:TRP:HZ3	3:A:1793:FLC:HA1	1.74	0.51
1:A:380:MSE:HG3	1:A:399:HIS:ND1	2.26	0.51
1:A:447:SER:O	1:A:451:ARG:HG3	2.11	0.49
1:A:470:MSE:HE3	1:A:503:PHE:CE2	2.47	0.49
1:A:315:ILE:HD11	1:A:352:TYR:HB2	1.95	0.49
1:A:470:MSE:HE3	1:A:503:PHE:HE2	1.77	0.49
1:A:341:VAL:HG13	1:A:345:TYR:CD2	2.47	0.49
1:A:380:MSE:O	1:A:384:THR:HG23	2.13	0.49
1:A:137:LYS:HG3	1:A:189:PHE:CE2	2.49	0.48
1:A:605:MSE:SE	1:A:608:ASN:HD22	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:573:ARG:HH21	1:A:576:ILE:HD11	1.79	0.47
1:A:609:TYR:O	1:A:613:VAL:HG23	2.15	0.47
1:A:480:LYS:O	1:A:483:THR:HG22	2.15	0.47
1:A:625:PHE:HD1	1:A:628:MSE:HE2	1.79	0.47
1:A:604:TRP:CH2	1:A:605:MSE:HE2	2.50	0.46
1:A:763:ALA:HA	1:A:781:VAL:O	2.16	0.46
1:A:498:GLU:HG3	1:A:531:ALA:HB1	1.97	0.46
1:A:573:ARG:NH2	1:A:576:ILE:HD11	2.31	0.46
1:A:705:THR:HA	1:A:721:ARG:O	2.17	0.45
3:A:1793:FLC:OG2	3:A:1793:FLC:CAC	2.64	0.45
1:A:446:LYS:HG2	1:A:450:HIS:CD2	2.52	0.45
1:A:523:HIS:HE1	5:A:2051:HOH:O	1.98	0.45
1:A:235:THR:HG22	1:A:238:GLN:OE1	2.17	0.45
1:A:725:THR:HB	1:A:777:GLN:HE21	1.81	0.45
1:A:765:PHE:HA	1:A:779:ASP:O	2.17	0.45
1:A:212:ARG:O	1:A:216:ASP:OD1	2.35	0.44
1:A:433:GLU:OE1	1:A:437:ARG:NH1	2.51	0.44
1:A:553:PHE:CE2	1:A:628:MSE:HE3	2.52	0.44
1:A:508:ARG:CG	1:A:508:ARG:NH1	2.54	0.43
1:A:443:ARG:HA	1:A:443:ARG:HD3	1.60	0.43
1:A:789:ASN:O	1:A:790:GLY:C	2.57	0.43
1:A:312:TYR:HD1	1:A:352:TYR:HA	1.84	0.43
1:A:544:PHE:HE2	1:A:635:THR:HA	1.83	0.43
1:A:186:ILE:O	1:A:190:THR:HG23	2.19	0.42
1:A:625:PHE:CD1	1:A:628:MSE:HE2	2.55	0.42
1:A:657:TYR:CZ	1:A:661:ILE:HD11	2.54	0.42
1:A:693:TYR:CE1	1:A:707:VAL:HG21	2.54	0.42
1:A:433:GLU:O	1:A:437:ARG:HG3	2.20	0.42
1:A:382:ARG:O	1:A:386:ASP:HB2	2.20	0.42
1:A:380:MSE:HG3	1:A:399:HIS:CE1	2.55	0.42
1:A:416:THR:HG23	1:A:420:GLY:HA2	2.01	0.42
1:A:475:SER:HB2	1:A:476:PRO:HD2	2.02	0.42
1:A:582:LYS:NZ	1:A:587:HIS:HB3	2.35	0.41
1:A:312:TYR:HA	1:A:352:TYR:HB3	2.02	0.41
1:A:622:MSE:N	1:A:623:PRO:CD	2.84	0.41
1:A:232:LYS:O	1:A:238:GLN:HB3	2.21	0.40
1:A:521:PHE:O	1:A:525:TYR:HB2	2.21	0.40
1:A:243:THR:O	1:A:247:LYS:HG3	2.22	0.40
1:A:255:ASN:HB3	1:A:257:GLU:H	1.86	0.40
1:A:419:ASN:ND2	1:A:710:GLU:HG2	2.36	0.40
1:A:635:THR:HG22	1:A:635:THR:O	2.21	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	675/785 (86%)	654 (97%)	20 (3%)	1 (0%)	51 65

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	117	THR

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	Rotameric	Outliers	Percentiles
1	A	587/660 (89%)	571 (97%)	16 (3%)	44 59

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	151	SER
1	A	384	THR
1	A	430	VAL
1	A	483	THR
1	A	490	THR
1	A	498	GLU

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Mol	Chain	Res	Type
1	A	508	ARG
1	A	586	ASP
1	A	591	LEU
1	A	625	PHE
1	A	655	THR
1	A	703	THR
1	A	707	VAL
1	A	725	THR
1	A	769	ARG

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	274	ASN
1	A	523	HIS
1	A	530	GLN
1	A	596	ASN
1	A	608	ASN
1	A	684	HIS
1	A	767	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)

3 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	$\#\|Z\| > 2$
3	FLC	A	1793	-	3,12,12	1.40	1 (33%)	3,17,17	3.41	2 (66%)
2	P6G	A	1792	-	18,18,18	0.56	0	17,17,17	0.51	0
4	PEG	A	1794	-	6,6,6	0.50	0	5,5,5	0.19	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
3	FLC	A	1793	-	-	4/6/16/16	-
2	P6G	A	1792	-	-	7/16/16/16	-
4	PEG	A	1794	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1793	FLC	CA-CB	-2.15	1.51	1.54

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1793	FLC	CB-CA-CAC	-5.49	106.19	114.98
3	A	1793	FLC	CG-CB-CA	2.06	114.85	109.33

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1793	FLC	CAC-CA-CB-CBC
3	A	1793	FLC	CAC-CA-CB-CG
3	A	1793	FLC	CAC-CA-CB-OHB
3	A	1793	FLC	CA-CB-CG-CGC
2	A	1792	P6G	O1-C2-C3-O4
2	A	1792	P6G	O16-C17-C18-O19

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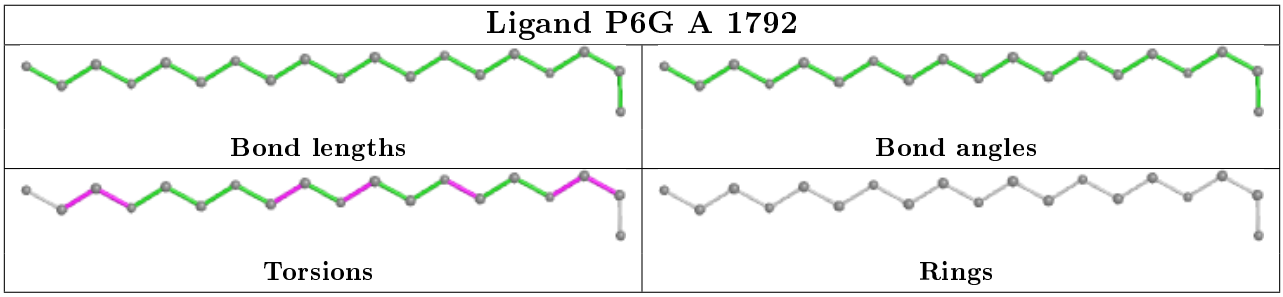
Mol	Chain	Res	Type	Atoms
4	A	1794	PEG	O1-C1-C2-O2
2	A	1792	P6G	O10-C11-C12-O13
2	A	1792	P6G	C5-C6-O7-C8
2	A	1792	P6G	C18-C17-O16-C15
2	A	1792	P6G	C8-C9-O10-C11
2	A	1792	P6G	C2-C3-O4-C5
4	A	1794	PEG	C1-C2-O2-C3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1793	FLC	3	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 then 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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	116:GLY	C	117:THR	N	1.15

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, 95^{th} 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	668/785 (85%)	0.13	27 (4%) 38 45	33, 57, 89, 172	49 (7%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	587	HIS	4.1
1	A	201	TYR	3.7
1	A	525	TYR	3.5
1	A	604	TRP	3.5
1	A	557	THR	3.2
1	A	529	LEU	3.1
1	A	558	ALA	3.1
1	A	526	THR	2.9
1	A	586	ASP	2.7
1	A	445	VAL	2.6
1	A	561	PHE	2.6
1	A	754	SER	2.5
1	A	585	VAL	2.4
1	A	553	PHE	2.4
1	A	527	HIS	2.3
1	A	208	TYR	2.3
1	A	442	SER	2.3
1	A	198	TYR	2.3
1	A	403	LEU	2.3
1	A	207	SER	2.2
1	A	397	VAL	2.2
1	A	449	PHE	2.2
1	A	531	ALA	2.1
1	A	202	TYR	2.1
1	A	148	PHE	2.1
1	A	597	SER	2.1
1	A	206	LEU	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)

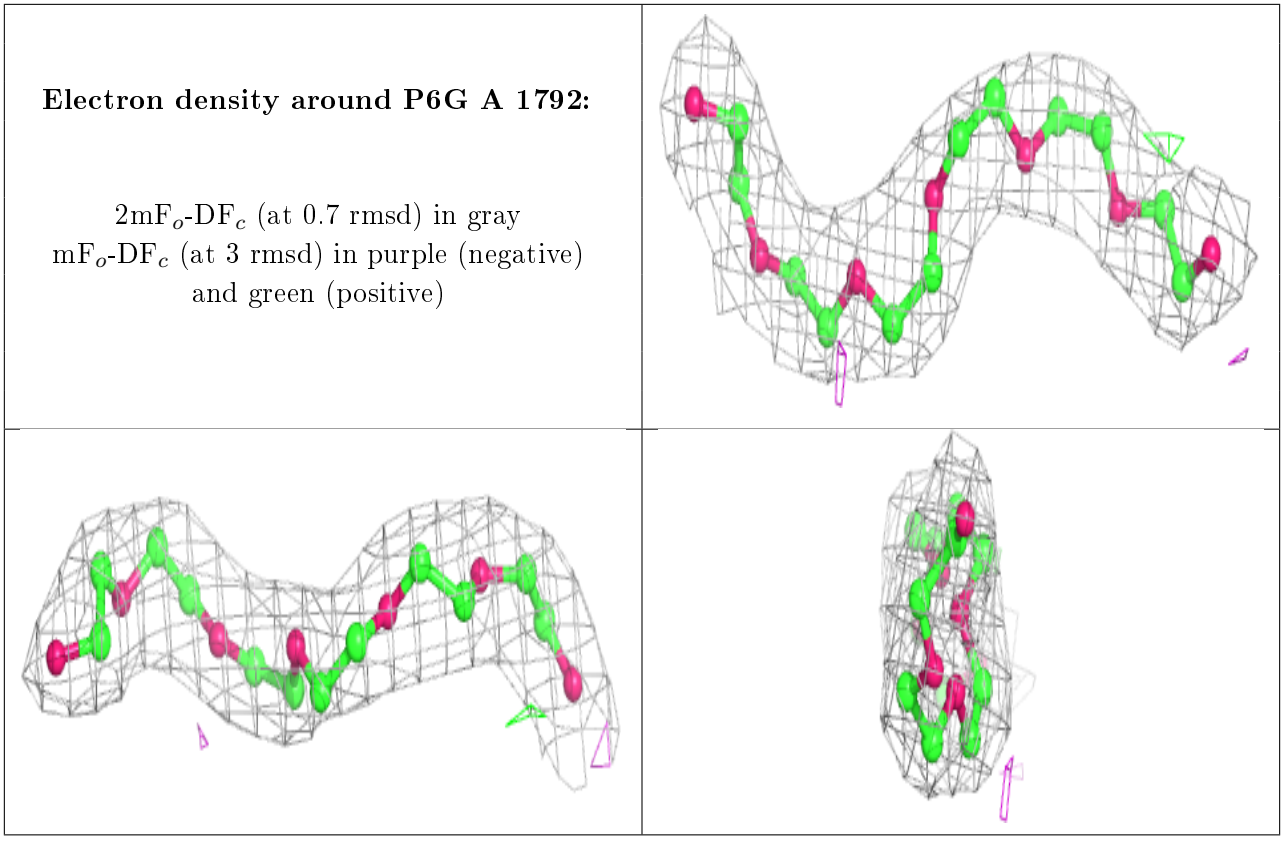
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	PEG	A	1794	7/7	0.65	0.25	90,92,93,93	0
3	FLC	A	1793	13/13	0.88	0.25	31,34,37,39	13
2	P6G	A	1792	19/19	0.93	0.13	55,58,62,62	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.

Electron density around P6G A 1792:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers

There are no such residues in this entry.