



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 24, 2022 – 08:09 PM JST

PDB ID : 7E01
Title : Trans-3/4-proline-hydroxylase H11 in the sixth reaction state
Authors : Gong, W.G.; Yang, L.Y.
Deposited on : 2021-01-26
Resolution : 1.72 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 : 2.26
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) : 2.26

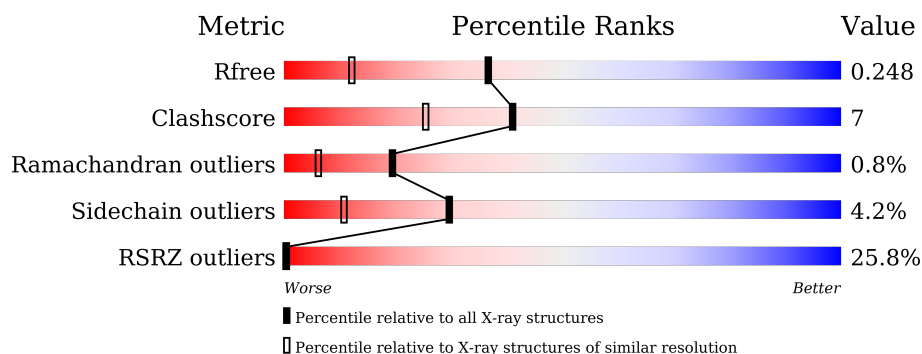
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 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	5722 (1.74-1.70)
Clashscore	141614	6152 (1.74-1.70)
Ramachandran outliers	138981	6051 (1.74-1.70)
Sidechain outliers	138945	6051 (1.74-1.70)
RSRZ outliers	127900	5629 (1.74-1.70)

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	266	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PRO	A	301	-	-	X	-
3	O	A	302	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	O	A	303	-	-	X	-

2 Entry composition [i](#)

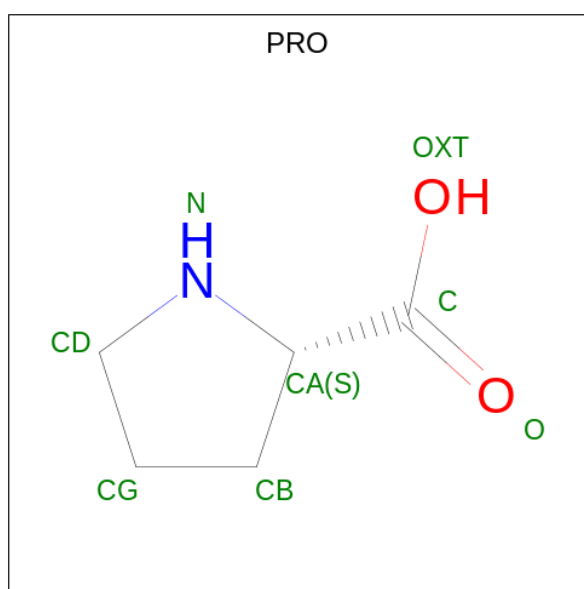
There are 4 unique types of molecules in this entry. The entry contains 2288 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Phytanoyl-CoA dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	2	0
			2088	1324	372	386	6			

- Molecule 2 is PROLINE (three-letter code: PRO) (formula: $C_5H_9NO_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			8	5	1	2		

- Molecule 3 is OXYGEN ATOM (three-letter code: O) (formula: O) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total	0	0
			2	2	

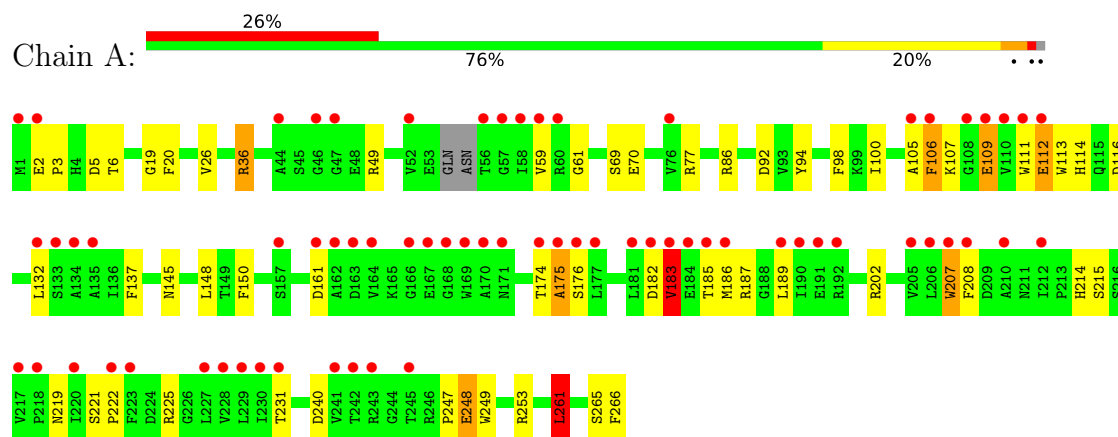
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	190	Total 190	O 190	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Phytanoyl-CoA dioxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	75.37Å 75.37Å 144.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.37 – 1.72 32.83 – 1.72	Depositor EDS
% Data completeness (in resolution range)	99.7 (75.37-1.72) 99.7 (32.83-1.72)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.20 (at 1.72Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.207 , 0.236 0.219 , 0.248	Depositor DCC
R_{free} test set	2204 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2288	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.53% 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

5.1 Standard geometry

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

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	1.42	11/2143 (0.5%)	1.43	28/2914 (1.0%)

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	3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	248	GLU	CD-OE2	11.14	1.37	1.25
1	A	207	TRP	CG-CD1	-7.27	1.26	1.36
1	A	94	TYR	CG-CD2	6.03	1.47	1.39
1	A	266	PHE	C-O	5.54	1.33	1.23
1	A	249	TRP	CB-CG	5.46	1.60	1.50
1	A	69	SER	CB-OG	-5.42	1.35	1.42
1	A	265	SER	CA-CB	5.36	1.60	1.52
1	A	94	TYR	CE1-CZ	5.30	1.45	1.38
1	A	247	PRO	CA-C	5.29	1.63	1.52
1	A	19	GLY	CA-C	5.22	1.60	1.51
1	A	208	PHE	CG-CD2	-5.09	1.31	1.38

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36[A]	ARG	NE-CZ-NH2	-14.28	113.16	120.30
1	A	36[B]	ARG	NE-CZ-NH2	-14.28	113.16	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	TYR	CB-CG-CD1	10.33	127.20	121.00
1	A	36[A]	ARG	NE-CZ-NH1	8.94	124.77	120.30
1	A	36[B]	ARG	NE-CZ-NH1	8.94	124.77	120.30
1	A	183	VAL	N-CA-C	8.94	135.14	111.00
1	A	225	ARG	NE-CZ-NH1	8.69	124.64	120.30
1	A	92	ASP	CB-CG-OD2	-8.36	110.78	118.30
1	A	86	ARG	NE-CZ-NH2	-8.02	116.29	120.30
1	A	94	TYR	CB-CG-CD2	-7.51	116.49	121.00
1	A	225	ARG	NE-CZ-NH2	-7.19	116.70	120.30
1	A	175	ALA	N-CA-C	-6.68	92.96	111.00
1	A	261	LEU	CB-CG-CD1	6.40	121.89	111.00
1	A	49	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	A	6	THR	N-CA-CB	5.84	121.39	110.30
1	A	137	PHE	CB-CG-CD2	-5.75	116.77	120.80
1	A	36[A]	ARG	CG-CD-NE	-5.70	99.84	111.80
1	A	36[B]	ARG	CG-CD-NE	-5.70	99.84	111.80
1	A	261	LEU	CB-CA-C	-5.59	99.59	110.20
1	A	92	ASP	CB-CG-OD1	5.52	123.27	118.30
1	A	148	LEU	CB-CG-CD2	5.40	120.18	111.00
1	A	231	THR	O-C-N	5.35	131.26	122.70
1	A	26	VAL	CG1-CB-CG2	-5.29	102.44	110.90
1	A	202	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	A	240	ASP	CB-CG-OD2	-5.19	113.63	118.30
1	A	70	GLU	OE1-CD-OE2	5.16	129.50	123.30
1	A	187	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	A	98	PHE	CB-CG-CD1	5.07	124.35	120.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	174	THR	Peptide
1	A	175	ALA	Peptide
1	A	36[A]	ARG	Sidechain

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	2088	0	2001	27	1
2	A	8	0	7	6	0
3	A	2	0	0	6	0
4	A	190	0	0	3	0
All	All	2288	0	2008	29	1

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 (29) 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:214:HIS:NE2	3:A:302:O:O	1.78	1.16
2:A:301:PRO:HG3	3:A:303:O:O	1.62	0.98
1:A:116[B]:ASP:OD1	2:A:301:PRO:HB3	1.78	0.82
1:A:116[B]:ASP:OD1	2:A:301:PRO:CB	2.27	0.82
1:A:214:HIS:CE1	3:A:302:O:O	2.32	0.81
1:A:116[B]:ASP:OD1	2:A:301:PRO:CG	2.36	0.72
1:A:185:THR:OG1	1:A:186:MET:N	2.23	0.70
2:A:301:PRO:CG	3:A:303:O:O	2.38	0.69
1:A:183:VAL:HG12	1:A:183:VAL:O	2.01	0.59
1:A:105:ALA:HB1	1:A:222:PRO:HA	1.87	0.57
1:A:114:HIS:NE2	3:A:302:O:O	2.39	0.56
1:A:5:ASP:HA	4:A:550:HOH:O	2.06	0.56
1:A:77:ARG:HB2	1:A:261:LEU:HD13	1.88	0.54
1:A:253:ARG:NH1	4:A:402:HOH:O	2.33	0.53
1:A:113:TRP:CD2	1:A:215:SER:HB2	2.45	0.51
1:A:145:ASN:HA	1:A:219:ASN:HA	1.93	0.50
1:A:116[B]:ASP:OD1	2:A:301:PRO:HG3	2.11	0.50
1:A:109:GLU:HG2	1:A:111:TRP:CD1	2.49	0.47
1:A:132:LEU:C	1:A:132:LEU:HD12	2.36	0.46
1:A:61:GLY:HA2	1:A:100:ILE:O	2.17	0.45
1:A:20:PHE:HA	1:A:207:TRP:O	2.17	0.45
1:A:116[A]:ASP:OD1	3:A:302:O:O	2.35	0.44
1:A:77:ARG:NH2	4:A:408:HOH:O	2.47	0.43
1:A:2:GLU:O	1:A:3:PRO:C	2.56	0.43
1:A:150:PHE:CD1	1:A:214:HIS:HB3	2.54	0.42
1:A:112:GLU:H	1:A:112:GLU:HG2	1.47	0.42
1:A:105:ALA:HB1	1:A:222:PRO:C	2.40	0.42
1:A:106:PHE:N	1:A:221:SER:O	2.46	0.42
1:A:59:VAL:O	1:A:59:VAL:HG23	2.21	0.40

All (1) 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)
1:A:106:PHE:CE2	1:A:183:VAL:CG1[5_455]	1.94	0.26

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	262/266 (98%)	250 (95%)	10 (4%)	2 (1%)	19 6

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	106	PHE
1	A	182	ASP

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	217/223 (97%)	208 (96%)	9 (4%)	30 11

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	107	LYS

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Mol	Chain	Res	Type
1	A	109	GLU
1	A	112	GLU
1	A	161	ASP
1	A	176	SER
1	A	183	VAL
1	A	189	LEU
1	A	248	GLU
1	A	261	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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
2	PRO	A	301	-	5,8,8	2.00	2 (40%)	6,10,10	1.38	1 (16%)

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	PRO	A	301	-	-	0/0/11/11	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	PRO	CB-CA	-3.45	1.50	1.54
2	A	301	PRO	CD-N	-2.65	1.39	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	PRO	CD-N-CA	2.75	114.43	107.06

There are no chirality outliers.

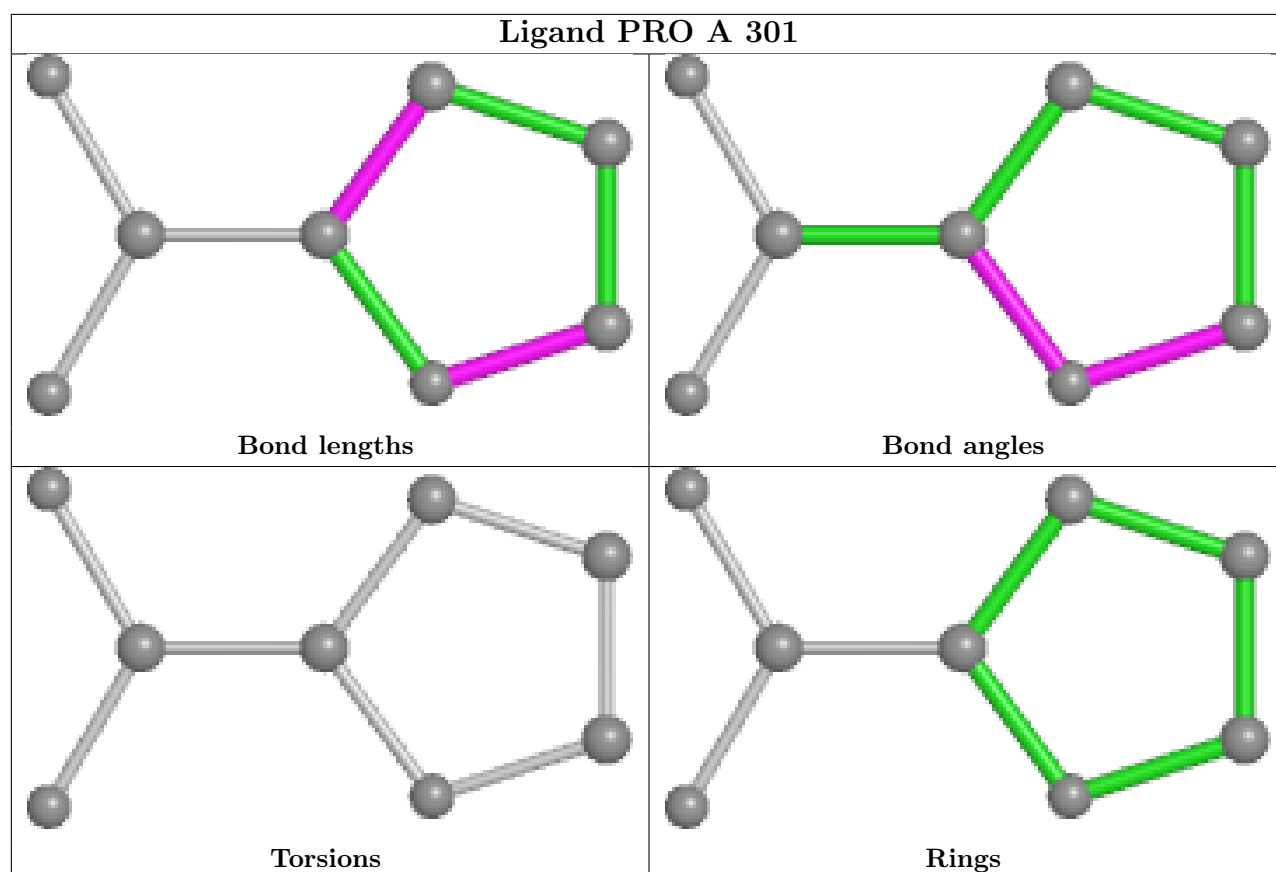
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	PRO	6	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	264/266 (99%)	1.11	68 (25%) 0 0	16, 31, 77, 94	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	56	THR	9.4
1	A	58	ILE	8.0
1	A	222	PRO	6.2
1	A	57	GLY	5.9
1	A	220	ILE	5.5
1	A	168	GLY	5.4
1	A	106	PHE	5.4
1	A	183	VAL	5.1
1	A	170	ALA	4.9
1	A	111	TRP	4.9
1	A	157	SER	4.7
1	A	2	GLU	4.5
1	A	245	THR	4.4
1	A	161	ASP	4.4
1	A	223	PHE	4.4
1	A	59	VAL	4.3
1	A	184	GLU	4.1
1	A	230	ILE	4.0
1	A	110	VAL	4.0
1	A	189	LEU	3.7
1	A	181	LEU	3.7
1	A	182	ASP	3.7
1	A	208	PHE	3.7
1	A	1	MET	3.7
1	A	167	GLU	3.6
1	A	169	TRP	3.6
1	A	134	ALA	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	229	LEU	3.4
1	A	177	LEU	3.3
1	A	163	ASP	3.2
1	A	206	LEU	3.2
1	A	132	LEU	3.1
1	A	46	GLY	3.1
1	A	212	ILE	2.9
1	A	176	SER	2.9
1	A	44	ALA	2.8
1	A	162	ALA	2.7
1	A	60	ARG	2.6
1	A	112	GLU	2.6
1	A	108	GLY	2.6
1	A	133	SER	2.6
1	A	243	ARG	2.6
1	A	228	VAL	2.5
1	A	166	GLY	2.5
1	A	171	ASN	2.5
1	A	109	GLU	2.5
1	A	47	GLY	2.5
1	A	192	ARG	2.5
1	A	207	TRP	2.4
1	A	242	THR	2.4
1	A	164	VAL	2.3
1	A	175	ALA	2.3
1	A	135	ALA	2.3
1	A	218	PRO	2.2
1	A	190	ILE	2.2
1	A	191	GLU	2.2
1	A	231	THR	2.2
1	A	217	VAL	2.2
1	A	210	ALA	2.2
1	A	174	THR	2.1
1	A	185	THR	2.1
1	A	227	LEU	2.1
1	A	52	VAL	2.1
1	A	105	ALA	2.1
1	A	186	MET	2.1
1	A	241	VAL	2.1
1	A	76	VAL	2.0
1	A	205	VAL	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 monosaccharides in this entry.

6.4 Ligands [i](#)

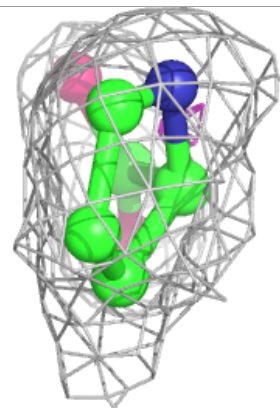
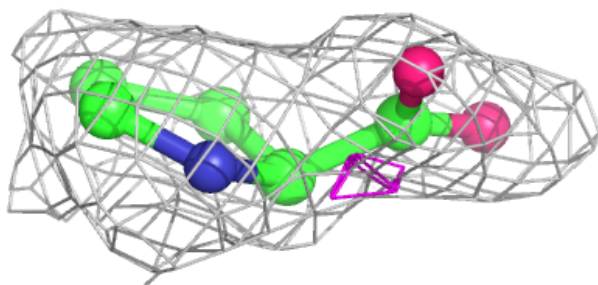
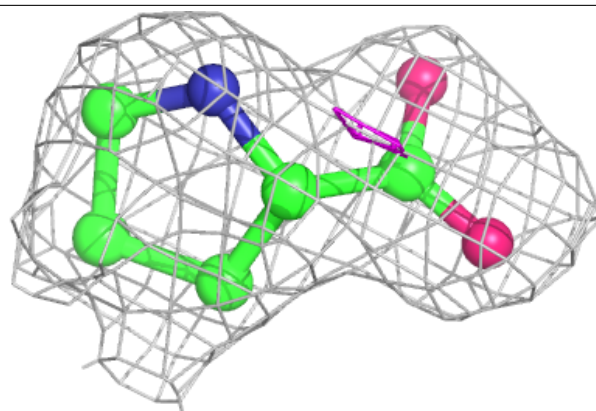
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PRO	A	301	8/8	0.87	0.16	35,40,45,46	0
3	O	A	302	1/1	0.87	0.17	39,39,39,39	0
3	O	A	303	1/1	0.95	0.63	54,54,54,54	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.

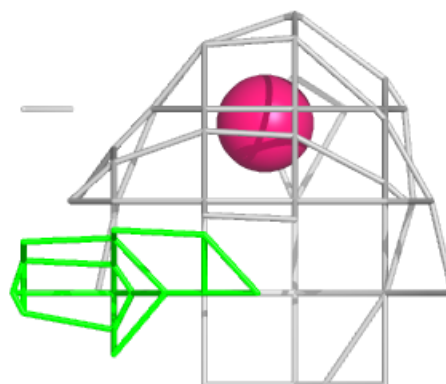
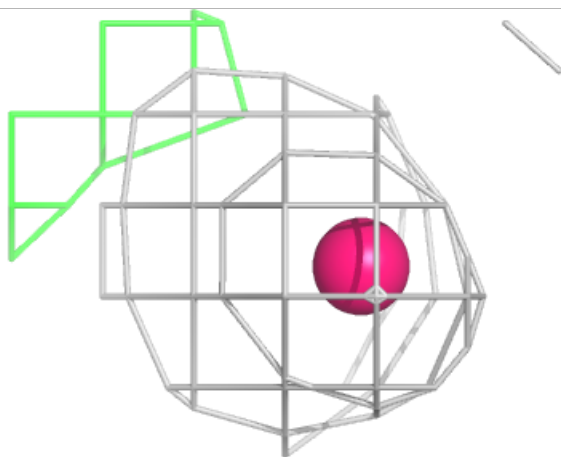
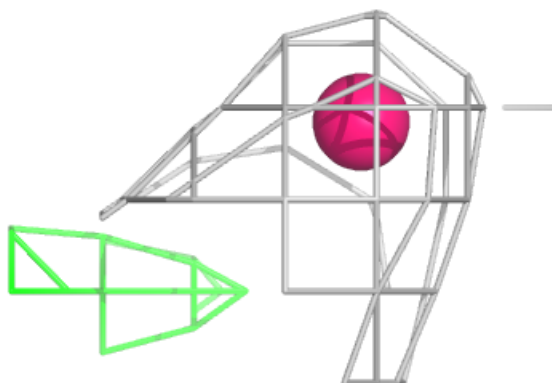
Electron density around PRO A 301:

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



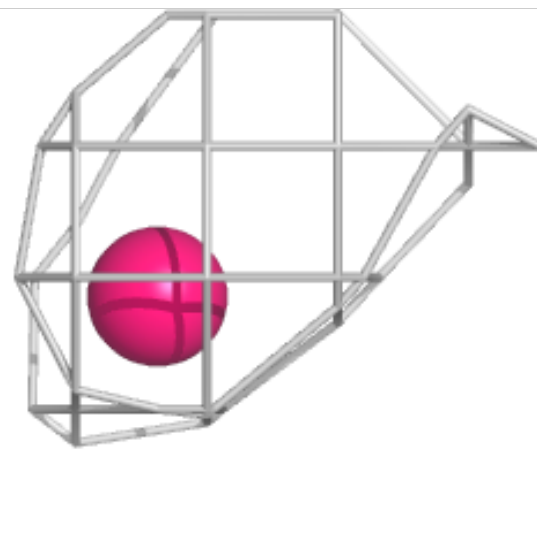
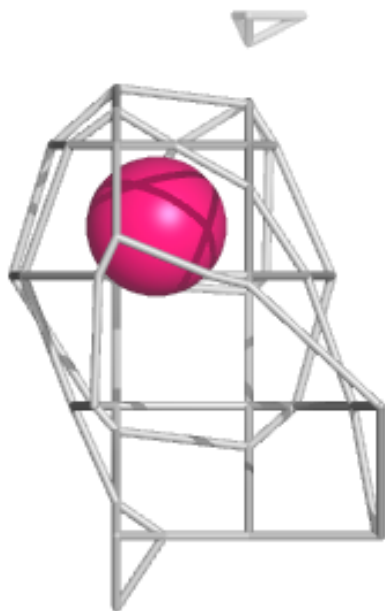
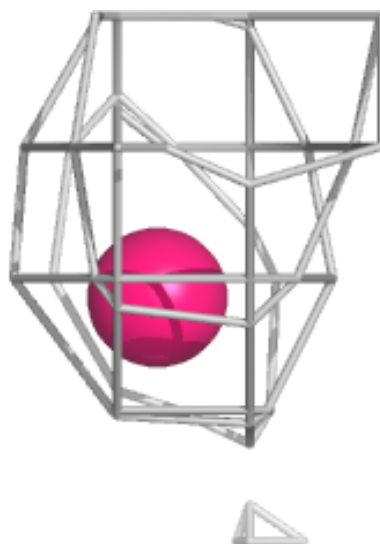
Electron density around O A 302:

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



Electron density around O A 303:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.