



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 09:08 am BST

PDB ID : 6HJE
Title : Trypanosoma cruzi proline racemase in complex with inhibitor NG-P27
Authors : Saul, F.; Haouz, A.; Uriac, P.; Blondel, A.; Minoprio, P.
Deposited on : 2018-09-03
Resolution : 2.00 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

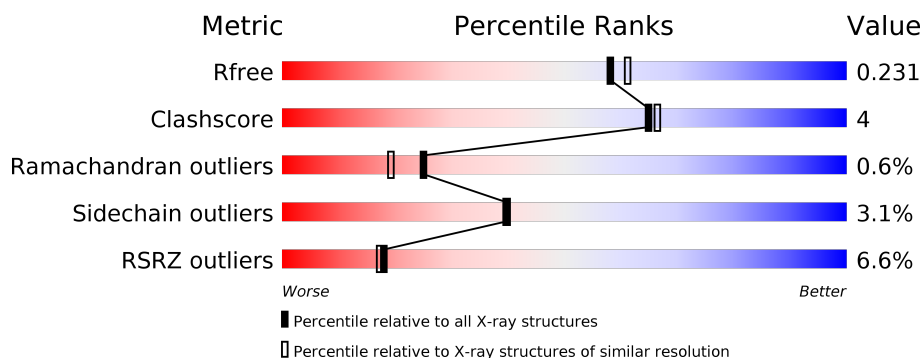
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 $\leq 5\%$. 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	414	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>6%</div> <div>14%</div> </div> </div>
1	B	414	<div> <div>7%</div> <div> <div></div> <div>77%</div> <div>9%</div> <div>14%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5872 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 Proline racemase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	0	3	0
			2734	1736	463	517	18			
1	B	356	Total	C	N	O	S	0	4	0
			2724	1730	460	517	17			

There are 46 discrepancies between the modelled and reference sequences:

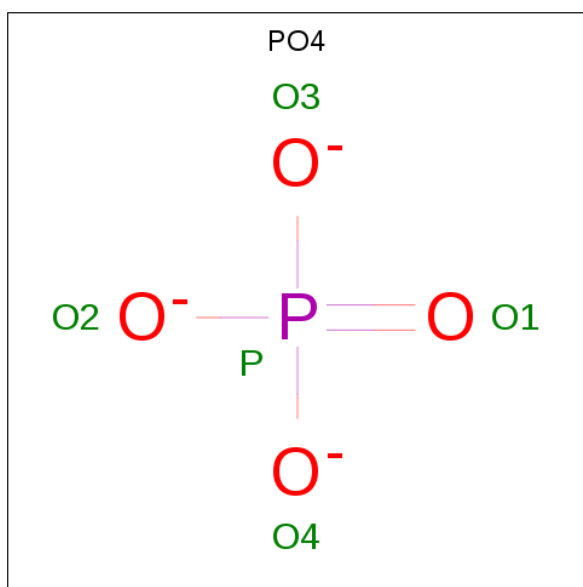
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q4DA80
A	118	ILE	MET	conflict	UNP Q4DA80
A	394	TYR	-	expression tag	UNP Q4DA80
A	395	ILE	-	expression tag	UNP Q4DA80
A	396	TRP	-	expression tag	UNP Q4DA80
A	397	SER	-	expression tag	UNP Q4DA80
A	398	SER	-	expression tag	UNP Q4DA80
A	399	SER	-	expression tag	UNP Q4DA80
A	400	VAL	-	expression tag	UNP Q4DA80
A	401	ASP	-	expression tag	UNP Q4DA80
A	402	LYS	-	expression tag	UNP Q4DA80
A	403	LEU	-	expression tag	UNP Q4DA80
A	404	ALA	-	expression tag	UNP Q4DA80
A	405	ALA	-	expression tag	UNP Q4DA80
A	406	ALA	-	expression tag	UNP Q4DA80
A	407	LEU	-	expression tag	UNP Q4DA80
A	408	GLU	-	expression tag	UNP Q4DA80
A	409	HIS	-	expression tag	UNP Q4DA80
A	410	HIS	-	expression tag	UNP Q4DA80
A	411	HIS	-	expression tag	UNP Q4DA80
A	412	HIS	-	expression tag	UNP Q4DA80
A	413	HIS	-	expression tag	UNP Q4DA80
A	414	HIS	-	expression tag	UNP Q4DA80
B	1	MET	-	initiating methionine	UNP Q4DA80
B	118	ILE	MET	conflict	UNP Q4DA80

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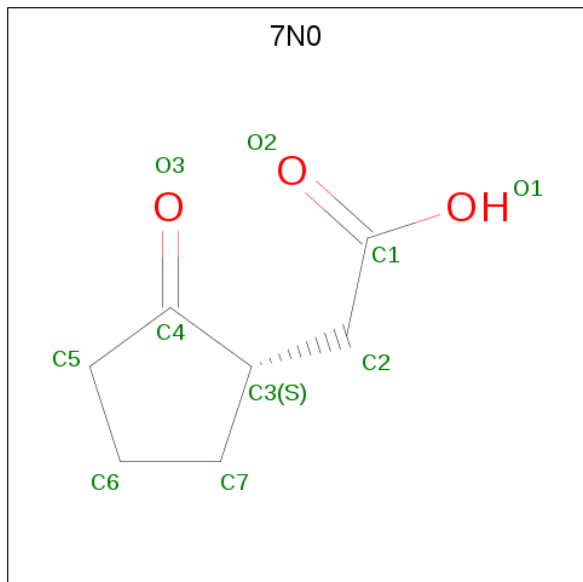
Chain	Residue	Modelled	Actual	Comment	Reference
B	394	TYR	-	expression tag	UNP Q4DA80
B	395	ILE	-	expression tag	UNP Q4DA80
B	396	TRP	-	expression tag	UNP Q4DA80
B	397	SER	-	expression tag	UNP Q4DA80
B	398	SER	-	expression tag	UNP Q4DA80
B	399	SER	-	expression tag	UNP Q4DA80
B	400	VAL	-	expression tag	UNP Q4DA80
B	401	ASP	-	expression tag	UNP Q4DA80
B	402	LYS	-	expression tag	UNP Q4DA80
B	403	LEU	-	expression tag	UNP Q4DA80
B	404	ALA	-	expression tag	UNP Q4DA80
B	405	ALA	-	expression tag	UNP Q4DA80
B	406	ALA	-	expression tag	UNP Q4DA80
B	407	LEU	-	expression tag	UNP Q4DA80
B	408	GLU	-	expression tag	UNP Q4DA80
B	409	HIS	-	expression tag	UNP Q4DA80
B	410	HIS	-	expression tag	UNP Q4DA80
B	411	HIS	-	expression tag	UNP Q4DA80
B	412	HIS	-	expression tag	UNP Q4DA80
B	413	HIS	-	expression tag	UNP Q4DA80
B	414	HIS	-	expression tag	UNP Q4DA80

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

- Molecule 3 is 2-[(1 {S})-2-oxidanylidencyclopentyl]ethanoic acid (three-letter code: 7N0) (formula: C₇H₁₀O₃) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	7	3		
3	B	1	Total	C	O	0	0
			10	7	3		

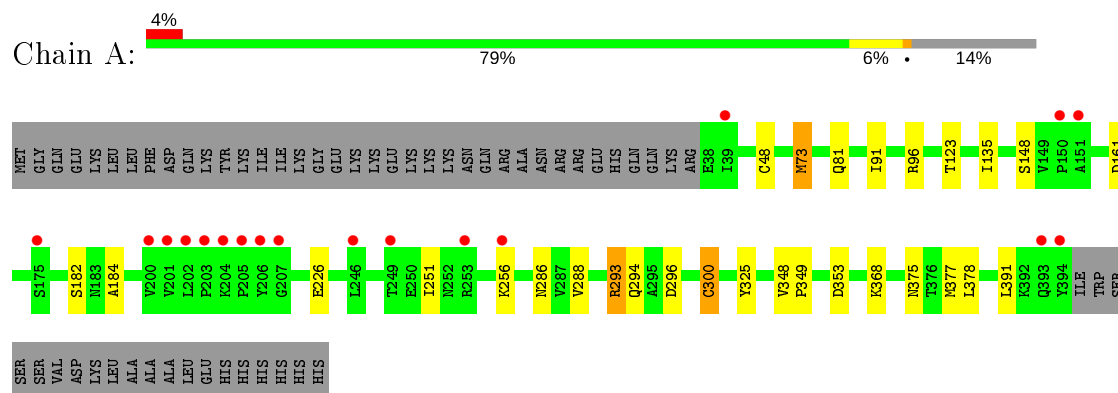
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	213	Total	O	0	0
			213	213		
4	B	176	Total	O	0	0
			176	176		

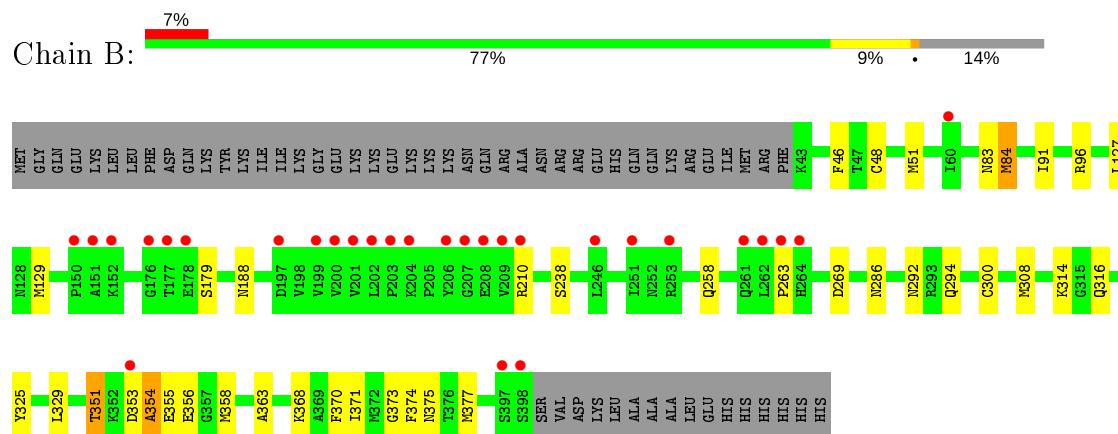
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Proline racemase A



• Molecule 1: Proline racemase A



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	133.23Å 90.64Å 85.39Å 90.00° 126.04° 90.00°	Depositor
Resolution (Å)	45.00 – 2.00 45.32 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.0 (45.00-2.00) 96.1 (45.32-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.177 , 0.230 0.186 , 0.231	Depositor DCC
R_{free} test set	1049 reflections (1.97%)	wwPDB-VP
Wilson B-factor (Å ²)	34.2	Xtriage
Anisotropy	0.432	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.3	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	0.97	EDS
Total number of atoms	5872	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% 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: PO4, 7N0

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	0.52	0/2798	0.72	1/3796 (0.0%)
1	B	0.52	0/2792	0.71	0/3791
All	All	0.52	0/5590	0.71	1/7587 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	296	ASP	CB-CG-OD1	5.14	122.93	118.30

There are no chirality outliers.

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	2734	0	2727	22	0
1	B	2724	0	2715	24	0
2	A	5	0	0	0	0
3	A	10	0	0	4	0
3	B	10	0	0	2	0
4	A	213	0	0	2	0
4	B	176	0	0	0	1
All	All	5872	0	5442	43	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 4.

All (43) 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:300:CYS:SG	3:A:502:7N0:C3	2.72	0.77
1:A:91:ILE:HA	1:A:377[A]:MET:HE1	1.70	0.72
1:A:391:LEU:HD21	1:B:329:LEU:HD21	1.79	0.65
1:B:46:PHE:CE2	1:B:91:ILE:HD11	2.32	0.64
1:B:368:LYS:HE3	1:B:370:PHE:CZ	2.38	0.59
1:A:48[B]:CYS:SG	1:A:377[B]:MET:CE	2.91	0.58
1:A:48[B]:CYS:SG	1:A:377[B]:MET:HE1	2.43	0.58
1:B:308:MET:HE3	1:B:363:ALA:HB1	1.86	0.58
1:A:123:THR:OG1	1:A:294:GLN:OE1	2.22	0.57
1:B:300:CYS:SG	3:B:501:7N0:C3	2.92	0.56
1:B:353:ASP:O	1:B:354:ALA:CB	2.55	0.55
1:B:258:GLN:NE2	1:B:263:PRO:HA	2.24	0.53
1:A:293[A]:ARG:HH11	1:A:293[A]:ARG:HG2	1.72	0.53
4:A:652:HOH:O	1:B:373:GLY:HA3	2.08	0.53
1:B:292:ASN:H	1:B:294:GLN:HE21	1.55	0.53
1:B:46:PHE:CD2	1:B:91:ILE:HD11	2.46	0.51
1:A:182:SER:O	1:A:368:LYS:HE3	2.11	0.50
1:A:81:GLN:NE2	4:A:602:HOH:O	2.35	0.49
1:B:314:LYS:HB2	1:B:316:GLN:HE21	1.78	0.49
1:B:48[B]:CYS:SG	1:B:375:ASN:HB3	2.53	0.49
1:A:288:VAL:HG11	3:A:502:7N0:C7	2.43	0.48
1:A:48[B]:CYS:SG	1:A:377[B]:MET:HE3	2.54	0.48
1:A:96:ARG:HD3	1:A:377[A]:MET:HE3	1.96	0.47
1:A:348:VAL:HB	1:A:349:PRO:HD2	1.97	0.46
1:B:91:ILE:HA	1:B:377[A]:MET:HE1	1.96	0.46
1:B:308:MET:CE	1:B:363:ALA:HB1	2.46	0.45
1:B:51:MET:SD	1:B:371:ILE:HG12	2.57	0.44
1:A:300:CYS:SG	3:A:502:7N0:C7	3.06	0.44
1:B:83:ASN:O	1:B:84:MET:HG2	2.18	0.43
1:B:351:THR:HG23	1:B:358:MET:HA	1.99	0.43
1:B:300:CYS:SG	3:B:501:7N0:C7	3.07	0.43
1:A:377[A]:MET:HE3	1:A:377[A]:MET:HB3	1.90	0.42
1:B:353:ASP:O	1:B:354:ALA:HB3	2.18	0.42
1:B:83:ASN:C	1:B:84:MET:HG2	2.39	0.42
1:A:378:LEU:HD12	1:B:371:ILE:HG22	2.01	0.42
1:A:293[A]:ARG:NH1	1:A:293[A]:ARG:HG2	2.35	0.41
1:A:300:CYS:SG	3:A:502:7N0:C1	3.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:MET:HG2	1:A:161:ASP:O	2.21	0.41
1:A:48[B]:CYS:SG	1:A:375:ASN:HB3	2.61	0.40
1:A:135:ILE:HG23	1:A:184:ALA:HB3	2.03	0.40
1:B:355:GLU:O	1:B:356:GLU:HG3	2.21	0.40
1:A:378:LEU:HD11	1:B:374:PHE:CE2	2.56	0.40
1:B:96:ARG:HD3	1:B:377[A]:MET:HE3	2.02	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)
4:B:775:HOH:O	4:B:775:HOH:O[2_555]	2.12	0.08

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	358/414 (86%)	349 (98%)	6 (2%)	3 (1%)	19	13
1	B	358/414 (86%)	340 (95%)	16 (4%)	2 (1%)	25	19
All	All	716/828 (86%)	689 (96%)	22 (3%)	5 (1%)	25	16

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	293[A]	ARG
1	A	293[B]	ARG
1	B	354	ALA
1	B	179	SER
1	A	300	CYS

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	297/347 (86%)	289 (97%)	8 (3%)	44	46
1	B	297/347 (86%)	287 (97%)	10 (3%)	37	36
All	All	594/694 (86%)	576 (97%)	18 (3%)	40	41

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	MET
1	A	148	SER
1	A	226	GLU
1	A	251	ILE
1	A	256	LYS
1	A	286	ASN
1	A	325	TYR
1	A	353	ASP
1	B	84	MET
1	B	127	LEU
1	B	129	MET
1	B	188	ASN
1	B	210	ARG
1	B	238	SER
1	B	269	ASP
1	B	286	ASN
1	B	325	TYR
1	B	351	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	ASN
1	A	155	ASN
1	A	174	GLN
1	A	196	GLN

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Mol	Chain	Res	Type
1	A	236	ASN
1	A	241	GLN
1	A	387	ASN
1	B	133	ASN
1	B	155	ASN
1	B	174	GLN
1	B	196	GLN
1	B	235	GLN
1	B	236	ASN
1	B	241	GLN
1	B	258	GLN
1	B	266	ASN
1	B	294	GLN
1	B	316	GLN
1	B	387	ASN

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 ⓘ

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
2	PO4	A	501	-	4,4,4	1.11	0	6,6,6	0.85	0
3	7N0	A	502	1	7,10,10	0.82	0	9,13,13	2.96	4 (44%)
3	7N0	B	501	1	7,10,10	0.59	0	9,13,13	2.85	2 (22%)

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	7N0	A	502	1	-	2/2/14/14	0/1/1/1
3	7N0	B	501	1	-	2/2/14/14	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	7N0	C5-C4-C3	6.77	113.74	108.89
3	A	502	7N0	C5-C4-C3	6.52	113.57	108.89
3	B	501	7N0	O3-C4-C3	-4.02	121.60	125.54
3	A	502	7N0	C7-C3-C2	-3.44	110.49	118.30
3	A	502	7N0	C6-C5-C4	-3.20	99.91	105.27
3	A	502	7N0	O3-C4-C3	-2.58	123.02	125.54

There are no chirality outliers.

All (4) torsion outliers are listed below:

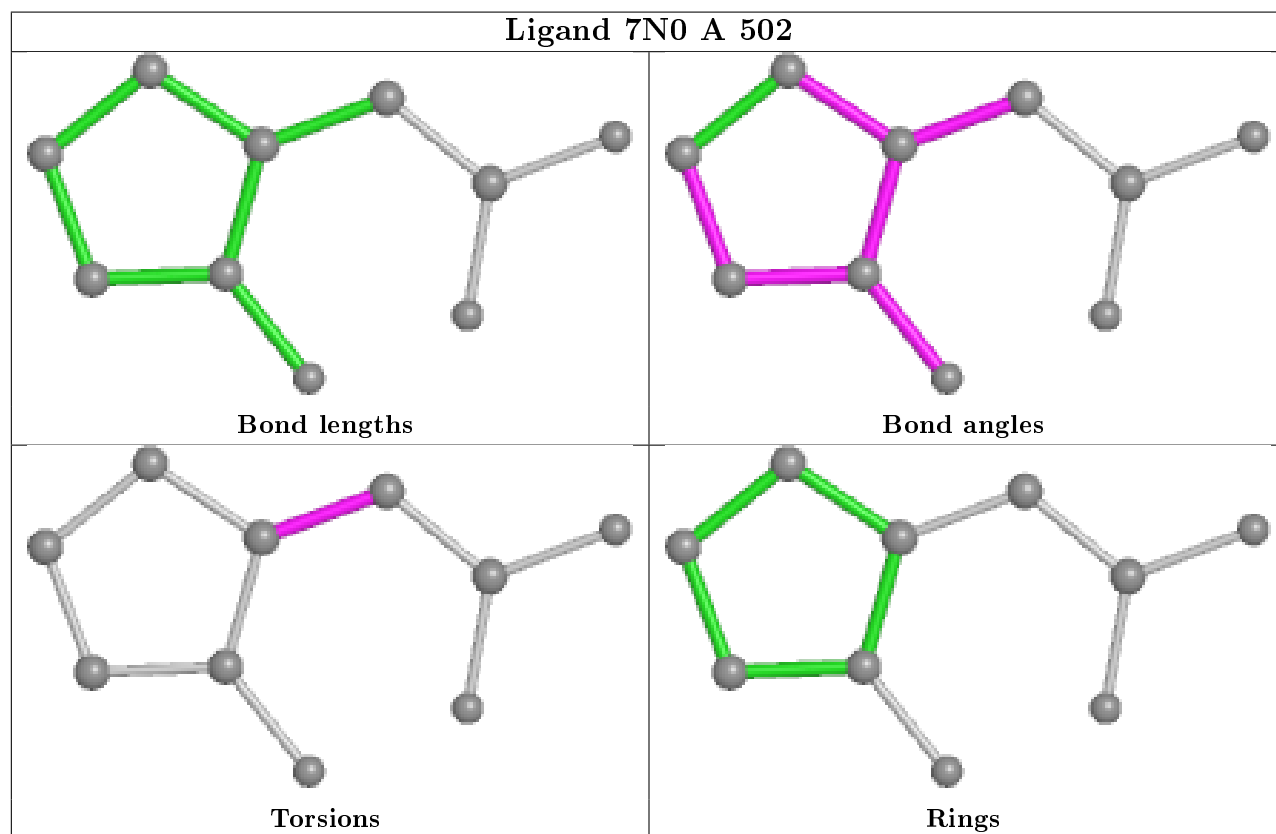
Mol	Chain	Res	Type	Atoms
3	A	502	7N0	C1-C2-C3-C7
3	B	501	7N0	C1-C2-C3-C7
3	A	502	7N0	C1-C2-C3-C4
3	B	501	7N0	C1-C2-C3-C4

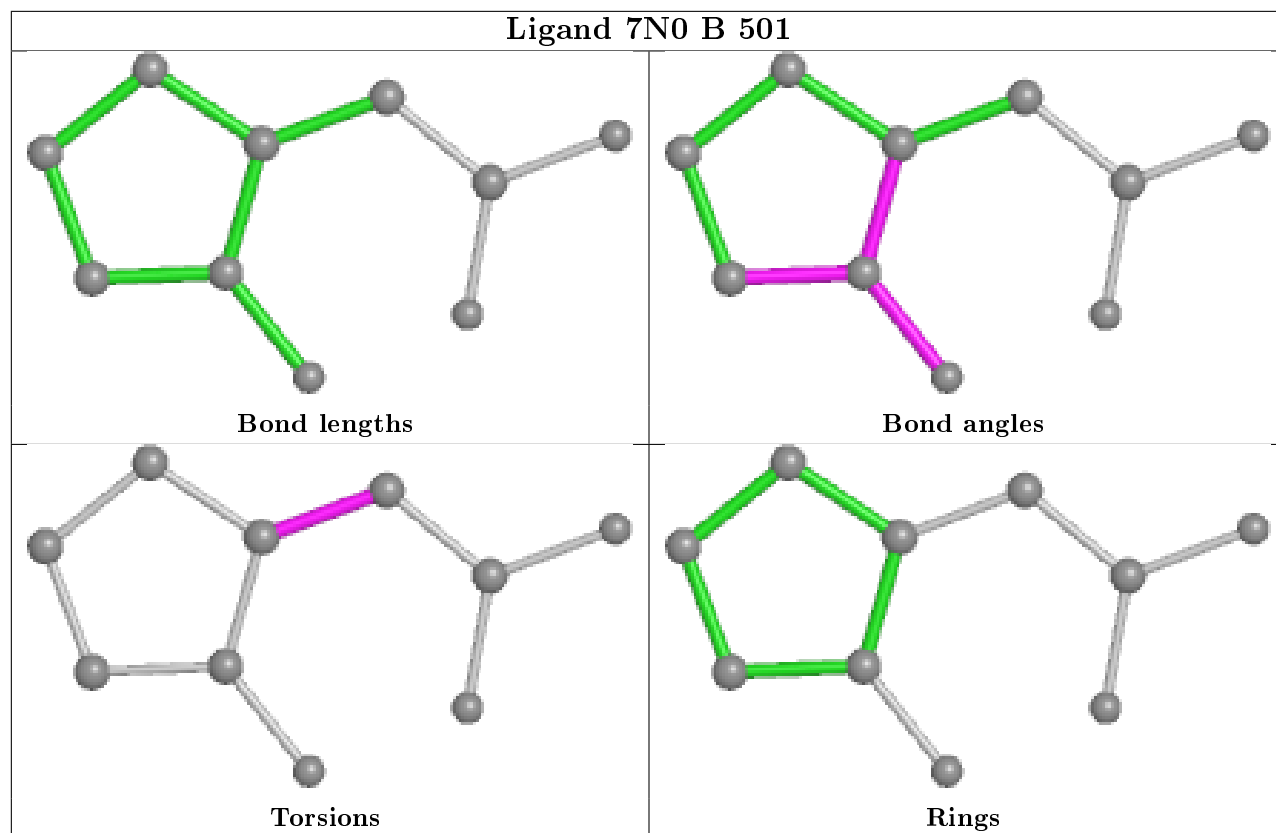
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	7N0	4	0
3	B	501	7N0	2	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	357/414 (86%)	-0.05	18 (5%) 28 28	26, 41, 82, 100	0
1	B	356/414 (85%)	0.18	29 (8%) 12 11	26, 45, 90, 100	0
All	All	713/828 (86%)	0.07	47 (6%) 18 17	26, 43, 87, 100	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	394	TYR	6.7
1	A	201	VAL	5.5
1	A	39	ILE	5.1
1	B	151	ALA	5.1
1	B	353	ASP	4.6
1	B	201	VAL	4.4
1	A	203	PRO	4.3
1	B	203	PRO	4.2
1	A	202	LEU	4.0
1	B	209	VAL	4.0
1	B	199	VAL	3.9
1	B	246	LEU	3.9
1	B	208	GLU	3.9
1	A	249	THR	3.7
1	B	210	ARG	3.2
1	B	202	LEU	3.2
1	A	393	GLN	3.2
1	A	175	SER	3.1
1	A	207	GLY	3.1
1	A	253	ARG	2.9
1	A	204	LYS	2.8
1	B	152	LYS	2.8
1	B	398	SER	2.8
1	B	264	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	263	PRO	2.8
1	B	177	THR	2.6
1	A	256	LYS	2.6
1	B	204	LYS	2.5
1	B	150	PRO	2.5
1	B	178	GLU	2.4
1	B	261	GLN	2.4
1	B	397	SER	2.4
1	B	262	LEU	2.3
1	A	200	VAL	2.3
1	A	205	PRO	2.3
1	B	200	VAL	2.3
1	B	207	GLY	2.3
1	B	251	ILE	2.3
1	B	176	GLY	2.2
1	A	151	ALA	2.2
1	A	246	LEU	2.2
1	B	253	ARG	2.2
1	B	197	ASP	2.2
1	A	206	TYR	2.1
1	A	150	PRO	2.1
1	B	206	TYR	2.1
1	B	60	ILE	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, 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.

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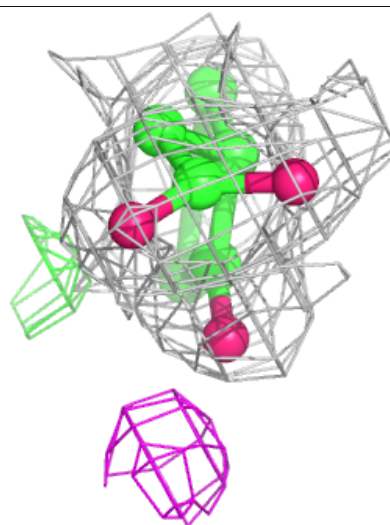
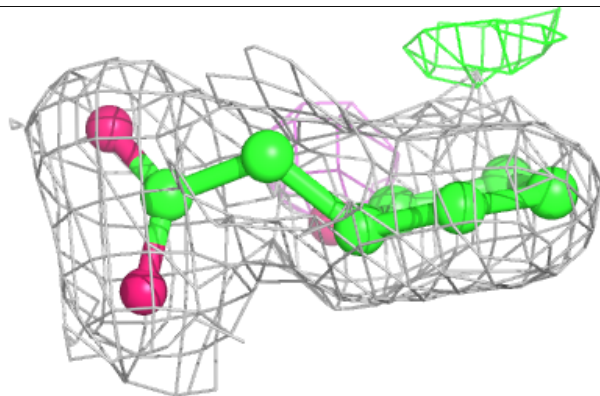
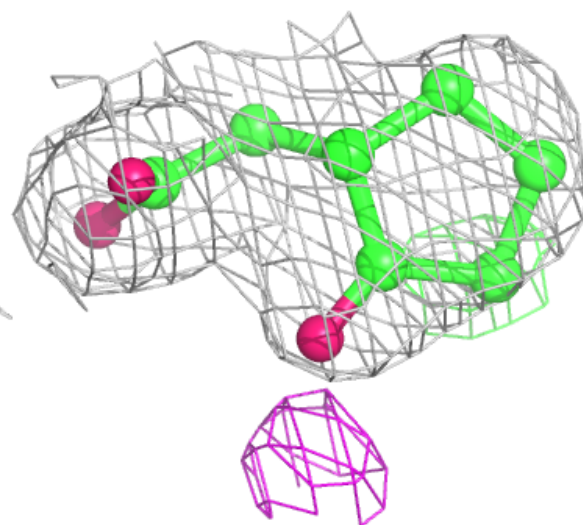
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	7N0	B	501	10/10	0.94	0.10	40,53,62,64	0
3	7N0	A	502	10/10	0.97	0.12	33,47,55,70	0
2	PO4	A	501	5/5	0.99	0.08	36,38,42,42	0

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.

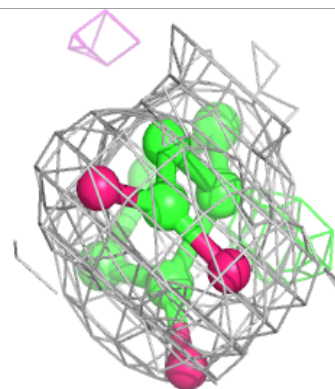
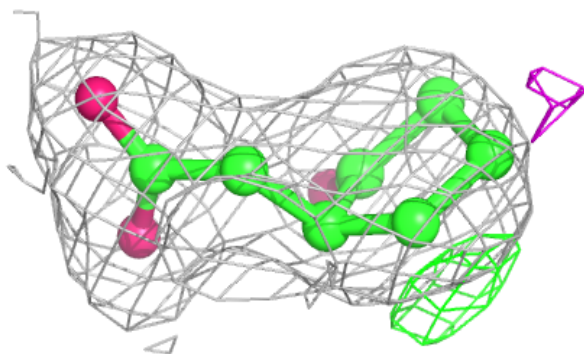
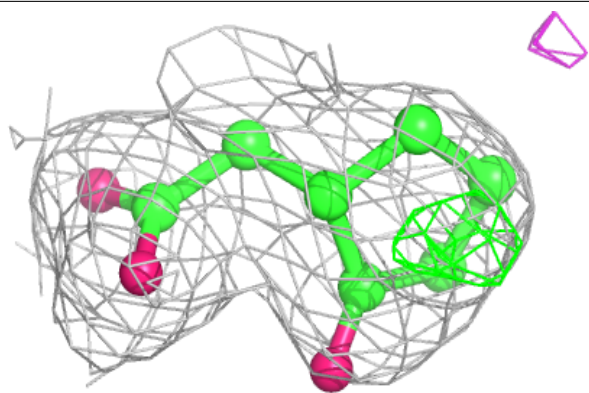
Electron density around 7N0 B 501:

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 7N0 A 502:

$2mF_o - DF_c$ (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.