



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 9WS1 / pdb_00009ws1
Title : ToFrO 15E_cryo
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Deposited on : 2025-09-12
Resolution : 1.80 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 : 1.8.5 (274361), CSD as541be (2020)
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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

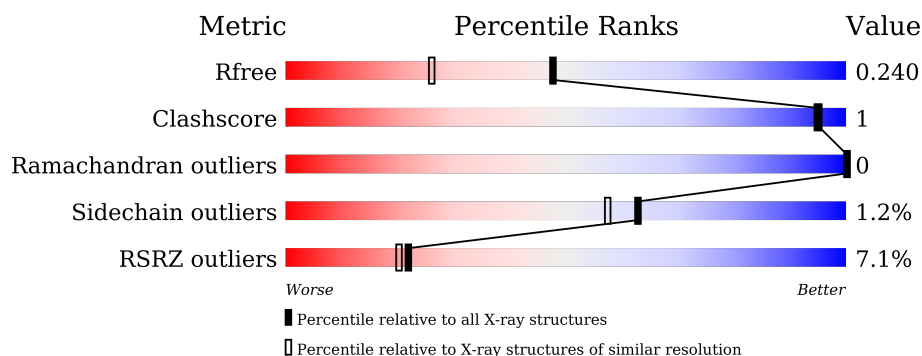
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	178	5% 84% 13%
1	B	178	7% 84% 13%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2740 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.

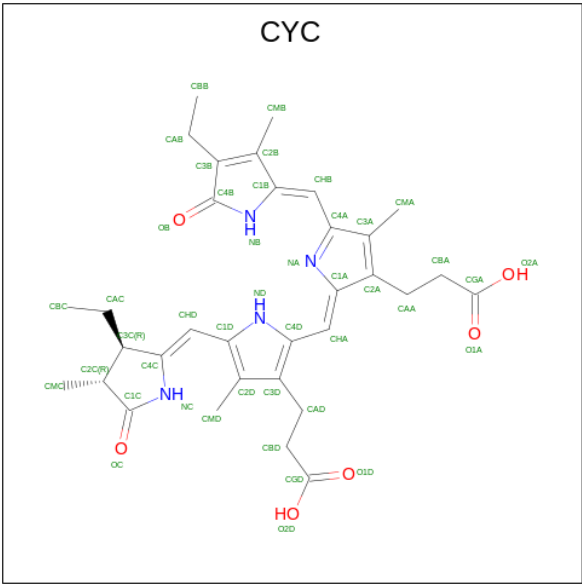
- Molecule 1 is a protein called histidine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	155	Total	C	N	O	S	0	0	0
			1209	770	212	221	6			
1	A	155	Total	C	N	O	S	0	0	0
			1191	760	207	218	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	554	MET	-	initiating methionine	UNP A0A6G9SQI3
B	555	ASP	-	expression tag	UNP A0A6G9SQI3
B	556	PRO	-	expression tag	UNP A0A6G9SQI3
B	557	SER	-	expression tag	UNP A0A6G9SQI3
B	558	SER	-	expression tag	UNP A0A6G9SQI3
B	559	SER	-	expression tag	UNP A0A6G9SQI3
B	726	HIS	-	expression tag	UNP A0A6G9SQI3
B	727	HIS	-	expression tag	UNP A0A6G9SQI3
B	728	HIS	-	expression tag	UNP A0A6G9SQI3
B	729	HIS	-	expression tag	UNP A0A6G9SQI3
B	730	HIS	-	expression tag	UNP A0A6G9SQI3
B	731	HIS	-	expression tag	UNP A0A6G9SQI3
A	554	MET	-	initiating methionine	UNP A0A6G9SQI3
A	555	ASP	-	expression tag	UNP A0A6G9SQI3
A	556	PRO	-	expression tag	UNP A0A6G9SQI3
A	557	SER	-	expression tag	UNP A0A6G9SQI3
A	558	SER	-	expression tag	UNP A0A6G9SQI3
A	559	SER	-	expression tag	UNP A0A6G9SQI3
A	726	HIS	-	expression tag	UNP A0A6G9SQI3
A	727	HIS	-	expression tag	UNP A0A6G9SQI3
A	728	HIS	-	expression tag	UNP A0A6G9SQI3
A	729	HIS	-	expression tag	UNP A0A6G9SQI3
A	730	HIS	-	expression tag	UNP A0A6G9SQI3
A	731	HIS	-	expression tag	UNP A0A6G9SQI3

- Molecule 2 is PHYCOCYANOBILIN (CCD ID: CYC) (formula: C₃₃H₄₀N₄O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			43	33	4	6		
2	A	1	Total	C	N	O	0	0
			43	33	4	6		

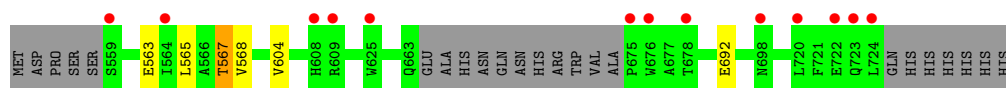
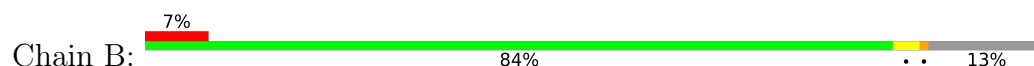
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	146	Total	O	0	0
			146	146		
3	A	108	Total	O	0	0
			108	108		

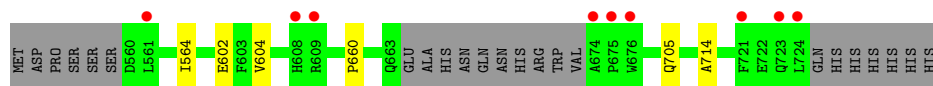
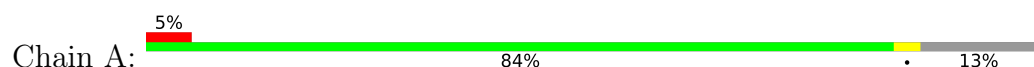
3 Residue-property plots [i](#)

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: histidine kinase



- Molecule 1: histidine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.04Å 53.87Å 79.95Å 90.00° 101.15° 90.00°	Depositor
Resolution (Å)	27.87 – 1.80 27.87 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.2 (27.87-1.80) 96.2 (27.87-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.220 , 0.242 0.218 , 0.240	Depositor DCC
R_{free} test set	1834 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	12.1	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2740	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.86% 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: CYC

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	0.27	0/1212	0.42	0/1653
1	B	0.30	0/1231	0.48	0/1677
All	All	0.29	0/2443	0.45	0/3330

There are no bond length outliers.

There are no bond angle outliers.

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	1191	0	1146	3	0
1	B	1209	0	1169	3	0
2	A	43	0	37	0	0
2	B	43	0	37	0	0
3	A	108	0	0	1	0
3	B	146	0	0	1	0
All	All	2740	0	2389	6	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 1.

All (6) 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:B:563:GLU:O	1:B:567:THR:HG23	2.04	0.57
1:B:604:VAL:HG12	3:B:940:HOH:O	2.05	0.57
1:A:564:ILE:HD13	1:A:714:ALA:HB1	1.95	0.48
1:B:565:LEU:HA	1:B:568:VAL:HG12	1.97	0.45
1:A:604:VAL:HG12	3:A:927:HOH:O	2.17	0.44
1:A:660:PRO:HD2	1:A:705:GLN:OE1	2.20	0.42

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	151/178 (85%)	150 (99%)	1 (1%)	0	100	100
1	B	151/178 (85%)	149 (99%)	2 (1%)	0	100	100
All	All	302/356 (85%)	299 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	121/160 (76%)	120 (99%)	1 (1%)	73	70

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	125/160 (78%)	123 (98%)	2 (2%)	55	47
All	All	246/320 (77%)	243 (99%)	3 (1%)	63	57

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	567	THR
1	B	692	GLU
1	A	602	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	618	GLN
1	B	679	ASN
1	A	654	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	CYC	B	801	1	46,46,46	6.15	28 (60%)	61,67,67	3.28	21 (34%)
2	CYC	A	801	1	46,46,46	6.31	28 (60%)	61,67,67	3.26	18 (29%)

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	CYC	B	801	1	-	8/26/74/74	0/4/4/4
2	CYC	A	801	1	-	5/26/74/74	0/4/4/4

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	CYC	C2C-C1C	-22.42	1.32	1.52
2	B	801	CYC	C2C-C1C	-21.83	1.32	1.52
2	A	801	CYC	C1C-NC	21.44	1.65	1.37
2	B	801	CYC	C1C-NC	20.58	1.64	1.37
2	B	801	CYC	OB-C4B	11.39	1.45	1.23
2	A	801	CYC	OB-C4B	11.36	1.45	1.23
2	A	801	CYC	CHD-C4C	10.64	1.57	1.36
2	B	801	CYC	CHD-C4C	10.01	1.56	1.36
2	A	801	CYC	CHA-C4D	8.52	1.60	1.40
2	B	801	CYC	CHA-C4D	8.24	1.59	1.40
2	B	801	CYC	CHD-C1D	8.19	1.59	1.40
2	A	801	CYC	OC-C1C	8.11	1.39	1.23
2	A	801	CYC	CHD-C1D	8.04	1.59	1.40
2	A	801	CYC	CHA-C1A	7.78	1.53	1.38
2	A	801	CYC	CHB-C1B	7.76	1.56	1.38
2	B	801	CYC	CHA-C1A	7.70	1.53	1.38
2	B	801	CYC	CHB-C1B	7.54	1.56	1.38
2	B	801	CYC	OC-C1C	7.26	1.37	1.23
2	A	801	CYC	CHB-C4A	6.83	1.56	1.40
2	B	801	CYC	CHB-C4A	6.34	1.55	1.40
2	A	801	CYC	CMC-C2C	6.08	1.66	1.53
2	B	801	CYC	CMC-C2C	5.94	1.66	1.53
2	B	801	CYC	C3C-C4C	5.38	1.58	1.50
2	A	801	CYC	CAC-C3C	-5.02	1.44	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	CYC	C3C-C4C	4.98	1.58	1.50
2	B	801	CYC	C4B-C3B	-4.78	1.39	1.48
2	B	801	CYC	CAC-C3C	-4.73	1.44	1.54
2	B	801	CYC	C4C-NC	-4.58	1.27	1.37
2	A	801	CYC	C4B-C3B	-4.28	1.40	1.48
2	A	801	CYC	CAB-C3B	4.26	1.62	1.51
2	B	801	CYC	CAB-C3B	4.06	1.61	1.51
2	A	801	CYC	C4C-NC	-3.98	1.29	1.37
2	B	801	CYC	CMA-C3A	3.60	1.58	1.50
2	B	801	CYC	C1B-NB	-3.49	1.32	1.37
2	A	801	CYC	CMA-C3A	3.38	1.57	1.50
2	A	801	CYC	C1B-NB	-3.36	1.32	1.37
2	B	801	CYC	C4D-ND	-3.29	1.32	1.37
2	B	801	CYC	C1A-C2A	-3.20	1.40	1.45
2	A	801	CYC	C1A-C2A	-3.19	1.40	1.45
2	B	801	CYC	C4B-NB	-3.02	1.31	1.38
2	A	801	CYC	C4D-ND	-2.97	1.32	1.37
2	A	801	CYC	C4B-NB	-2.94	1.31	1.38
2	A	801	CYC	C1B-C2B	-2.77	1.40	1.45
2	B	801	CYC	C1B-C2B	-2.73	1.40	1.45
2	A	801	CYC	C4A-C3A	-2.60	1.40	1.45
2	B	801	CYC	C4A-C3A	-2.47	1.40	1.45
2	B	801	CYC	CBA-CGA	2.45	1.56	1.50
2	A	801	CYC	CBA-CGA	2.35	1.56	1.50
2	A	801	CYC	CAD-C3D	2.30	1.57	1.51
2	B	801	CYC	CAD-C3D	2.29	1.57	1.51
2	A	801	CYC	CMD-C2D	2.27	1.55	1.50
2	B	801	CYC	C1A-NA	-2.19	1.33	1.38
2	A	801	CYC	C1A-NA	-2.17	1.33	1.38
2	B	801	CYC	CMD-C2D	2.14	1.55	1.50
2	B	801	CYC	CAA-C2A	2.11	1.56	1.51
2	A	801	CYC	CAA-C2A	2.07	1.56	1.51

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	CYC	OC-C1C-NC	-16.07	105.48	124.94
2	A	801	CYC	OC-C1C-C2C	-14.91	114.31	126.17
2	A	801	CYC	OC-C1C-NC	-13.66	108.40	124.94
2	B	801	CYC	OC-C1C-C2C	-11.68	116.89	126.17
2	B	801	CYC	CHD-C4C-NC	-8.38	115.25	125.20
2	A	801	CYC	CHD-C4C-NC	-7.75	115.99	125.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	CYC	CHA-C4D-C3D	-3.96	118.21	127.15
2	B	801	CYC	C1B-CHB-C4A	-3.95	118.42	128.08
2	A	801	CYC	CHA-C4D-C3D	-3.94	118.26	127.15
2	B	801	CYC	CMB-C2B-C1B	3.66	128.74	124.17
2	B	801	CYC	CHD-C1D-C2D	-3.59	118.69	127.49
2	A	801	CYC	CMB-C2B-C1B	3.56	128.61	124.17
2	A	801	CYC	CAB-C3B-C4B	3.49	126.90	121.38
2	A	801	CYC	C1B-CHB-C4A	-3.29	120.03	128.08
2	A	801	CYC	CHB-C1B-NB	-3.00	119.63	126.06
2	B	801	CYC	C3B-C4B-NB	2.98	109.18	106.78
2	B	801	CYC	C3D-C4D-ND	2.97	111.58	107.57
2	A	801	CYC	C3B-C4B-NB	2.94	109.15	106.78
2	A	801	CYC	C2C-C1C-NC	-2.93	105.75	108.27
2	A	801	CYC	CHD-C1D-C2D	-2.93	120.31	127.49
2	B	801	CYC	CHB-C1B-NB	-2.91	119.80	126.06
2	A	801	CYC	CHA-C1A-NA	2.87	130.55	124.58
2	B	801	CYC	C2D-C1D-ND	2.79	111.32	107.40
2	A	801	CYC	C3D-C4D-ND	2.67	111.17	107.57
2	A	801	CYC	CHA-C1A-C2A	-2.63	119.24	125.32
2	B	801	CYC	CHA-C1A-C2A	-2.60	119.31	125.32
2	A	801	CYC	C2D-C1D-ND	2.58	111.03	107.40
2	B	801	CYC	CHA-C1A-NA	2.58	129.95	124.58
2	B	801	CYC	CAC-C3C-C4C	-2.54	106.14	112.67
2	B	801	CYC	CAB-C3B-C4B	2.47	125.27	121.38
2	A	801	CYC	CMA-C3A-C4A	2.44	128.82	125.06
2	B	801	CYC	C1B-C2B-C3B	-2.42	105.34	107.87
2	A	801	CYC	CHA-C4D-ND	2.29	130.56	125.32
2	A	801	CYC	CHB-C1B-C2B	2.29	131.48	126.95
2	B	801	CYC	CHA-C4D-ND	2.13	130.21	125.32
2	B	801	CYC	C4D-ND-C1D	-2.12	106.00	109.79
2	B	801	CYC	C4D-C3D-C2D	-2.08	105.16	107.62
2	B	801	CYC	CHD-C1D-ND	2.04	129.99	125.32
2	B	801	CYC	CHB-C1B-C2B	2.01	130.93	126.95

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	801	CYC	NA-C4A-CHB-C1B
2	B	801	CYC	C3A-C4A-CHB-C1B
2	B	801	CYC	NC-C4C-CHD-C1D
2	A	801	CYC	NA-C4A-CHB-C1B

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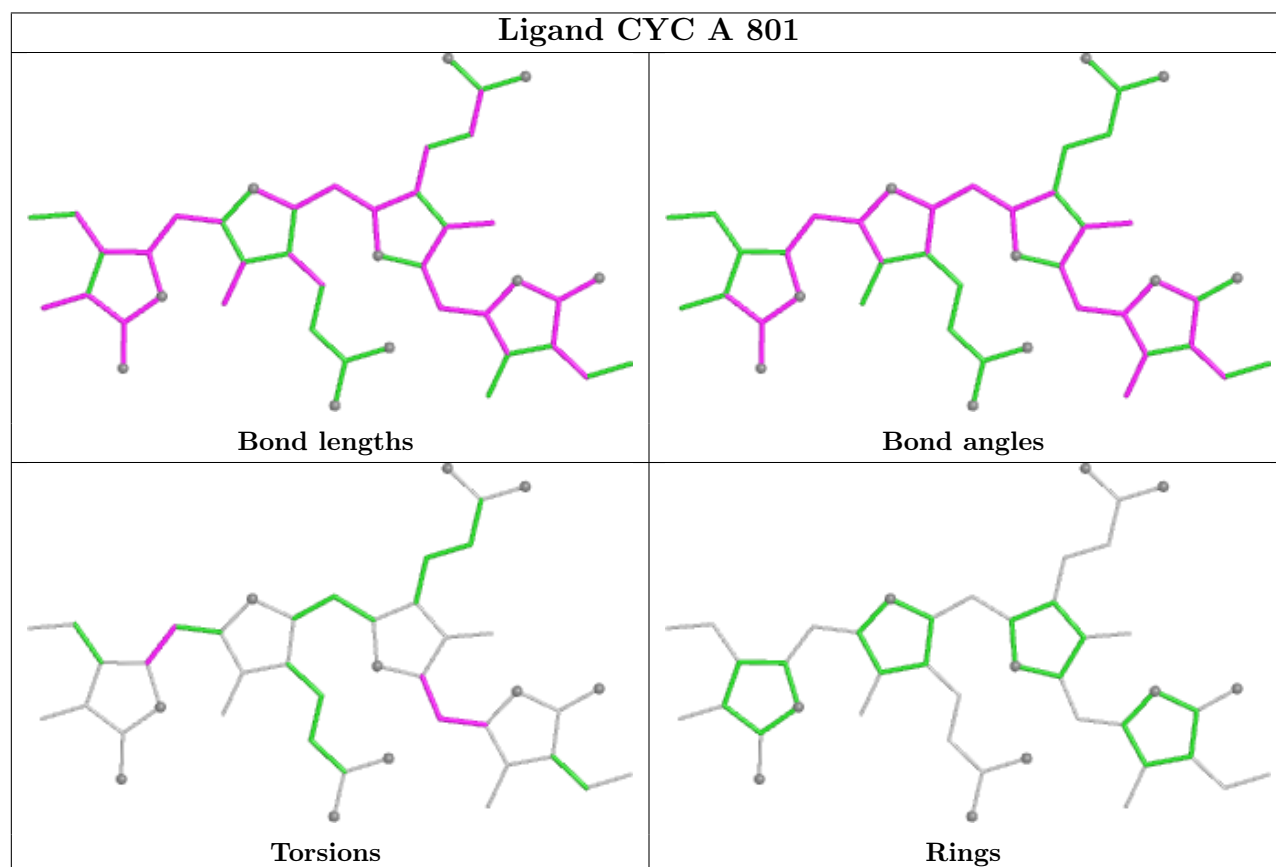
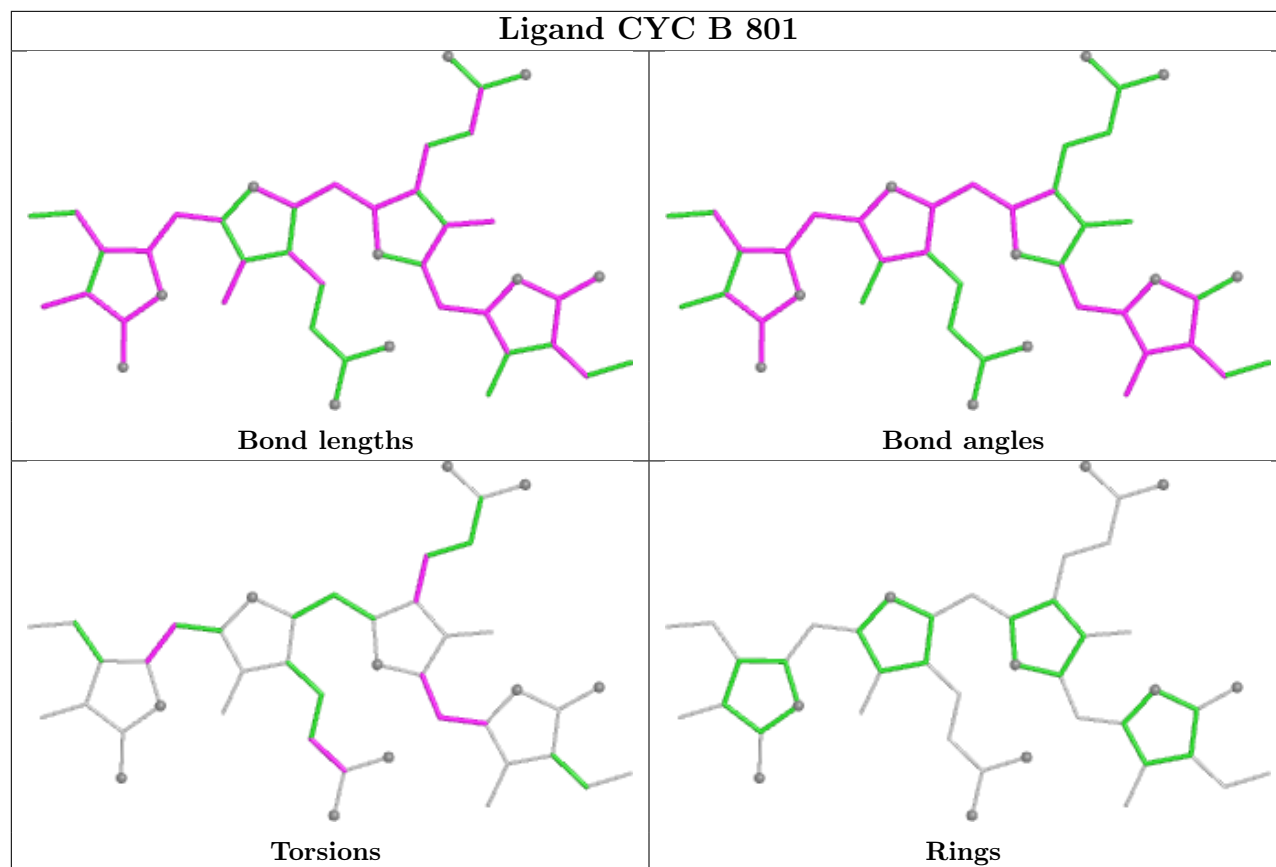
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Mol	Chain	Res	Type	Atoms
2	A	801	CYC	C3A-C4A-CHB-C1B
2	A	801	CYC	NC-C4C-CHD-C1D
2	B	801	CYC	C2B-C1B-CHB-C4A
2	A	801	CYC	C2B-C1B-CHB-C4A
2	B	801	CYC	NB-C1B-CHB-C4A
2	A	801	CYC	NB-C1B-CHB-C4A
2	B	801	CYC	CAD-CBD-CGD-O2D
2	B	801	CYC	CAD-CBD-CGD-O1D
2	B	801	CYC	C1A-C2A-CAA-CBA

There are no ring outliers.

No monomer is involved in short contacts.

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	155/178 (87%)	0.36	9 (5%) 29 27	7, 16, 33, 40	0
1	B	155/178 (87%)	0.24	13 (8%) 17 15	6, 11, 30, 38	0
All	All	310/356 (87%)	0.30	22 (7%) 22 20	6, 13, 32, 40	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	676	TRP	4.0
1	A	721	PHE	3.7
1	A	724	LEU	3.6
1	B	608	HIS	3.4
1	B	675	PRO	3.2
1	B	676	TRP	3.1
1	A	674	ALA	2.9
1	B	724	LEU	2.8
1	A	723	GLN	2.7
1	A	609	ARG	2.7
1	B	698	ASN	2.6
1	A	608	HIS	2.5
1	B	678	THR	2.5
1	B	609	ARG	2.5
1	B	722	GLU	2.5
1	A	561	LEU	2.4
1	B	625	TRP	2.4
1	B	723	GLN	2.3
1	B	559	SER	2.3
1	B	564	ILE	2.2
1	B	720	LEU	2.2
1	A	675	PRO	2.2

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

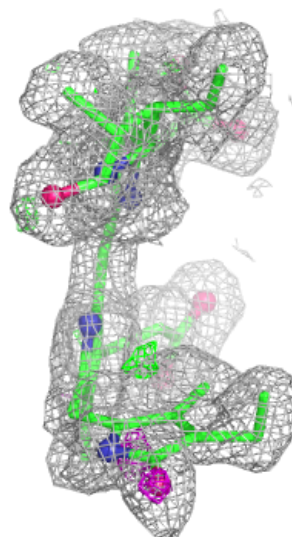
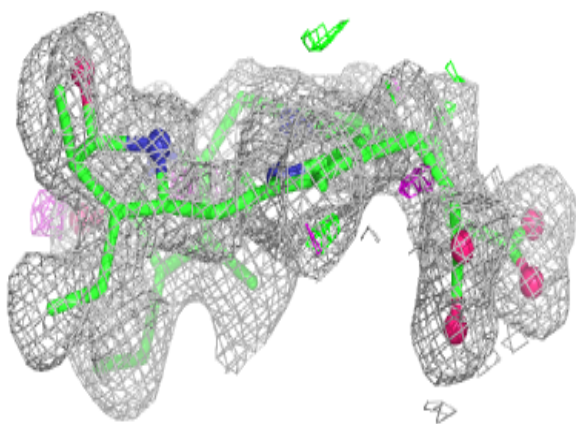
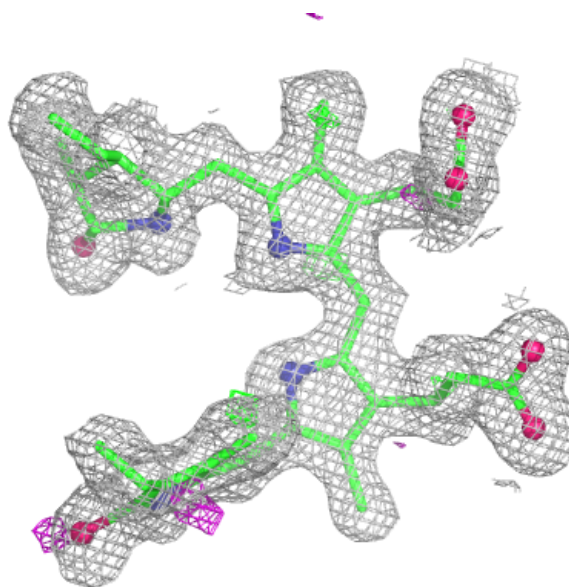
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	CYC	B	801	43/43	0.91	0.09	5,6,12,17	0
2	CYC	A	801	43/43	0.91	0.10	10,14,19,21	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.

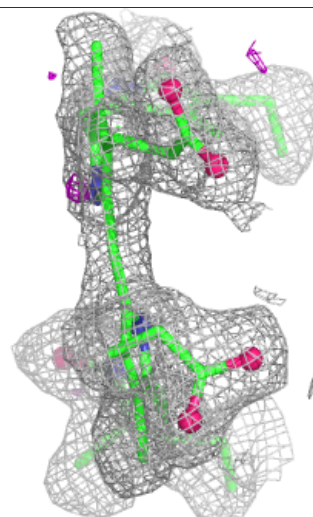
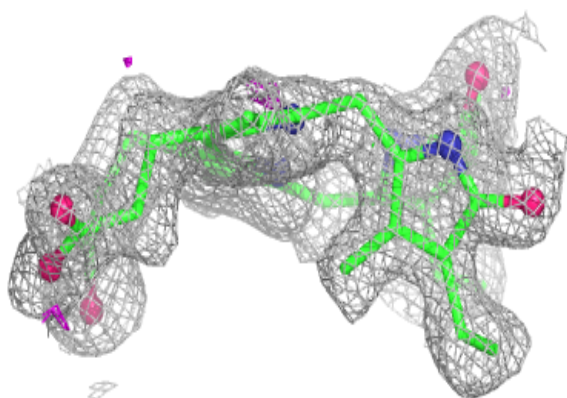
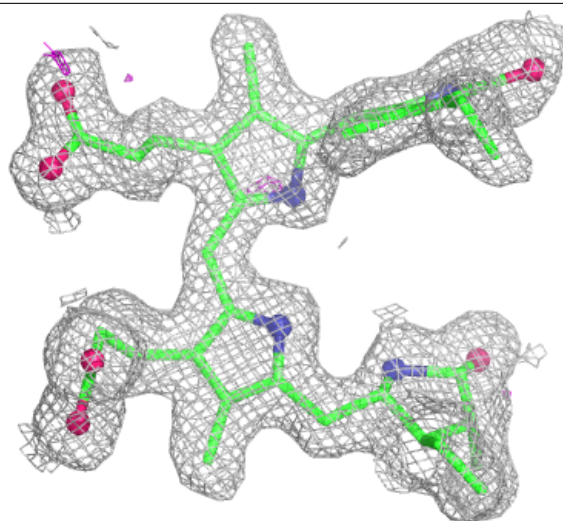
Electron density around CYC B 801:

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



Electron density around CYC A 801:

$2mF_o-DF_c$ (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.