Full wwPDB X-ray Structure Validation Report (i)

Sep 22, 2026 – 04:31 pm BST

PDB ID	:	5HDR / pdb_00005hdr
Title	:	Crystal structure of a bacterial fucosidase with iminocyclitol (2S,3S,4R,5S)-3, 4-dihydroxy-2-ethynyl-5-methylpyrrolidine
Authors	:	Wright, D.W.; Davies, G.J.
Deposited on	:	2016-01-05
Resolution	:	2.30 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 (i) 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 (i)) were used in the production of this report:

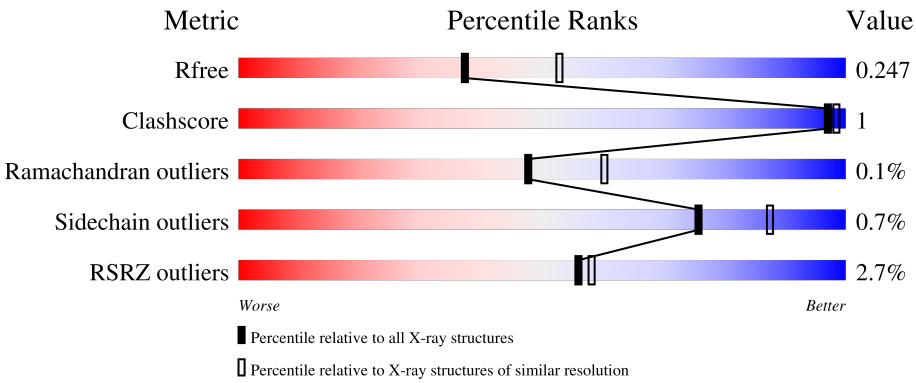
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 <=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	447	7% 95%
1	B	447	% 96%
1	C	447	% 95%
1	D	447	2% 95%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 14754 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 Alpha-L-fucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	440	3505	2250	587	652	16	0	0	0
1	B	440	3573	2298	599	659	17	0	2	0
1	C	441	3547	2283	593	655	16	0	1	0
1	D	440	3531	2271	590	654	16	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	ALA	-	expression tag	UNP Q8A3I4
A	473	LYS	-	expression tag	UNP Q8A3I4
A	474	GLY	-	expression tag	UNP Q8A3I4
A	475	THR	-	expression tag	UNP Q8A3I4
A	476	LYS	-	expression tag	UNP Q8A3I4
A	477	SER	-	expression tag	UNP Q8A3I4
A	478	ILE	-	expression tag	UNP Q8A3I4
A	479	TYR	-	expression tag	UNP Q8A3I4
A	480	ARG	-	expression tag	UNP Q8A3I4
B	34	ALA	-	expression tag	UNP Q8A3I4
B	473	LYS	-	expression tag	UNP Q8A3I4
B	474	GLY	-	expression tag	UNP Q8A3I4
B	475	THR	-	expression tag	UNP Q8A3I4
B	476	LYS	-	expression tag	UNP Q8A3I4
B	477	SER	-	expression tag	UNP Q8A3I4
B	478	ILE	-	expression tag	UNP Q8A3I4
B	479	TYR	-	expression tag	UNP Q8A3I4
B	480	ARG	-	expression tag	UNP Q8A3I4
C	34	ALA	-	expression tag	UNP Q8A3I4
C	473	LYS	-	expression tag	UNP Q8A3I4
C	474	GLY	-	expression tag	UNP Q8A3I4

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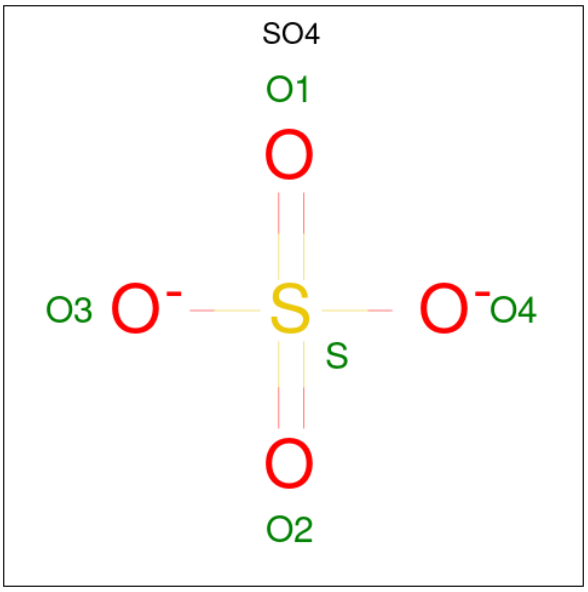
Chain	Residue	Modelled	Actual	Comment	Reference
C	475	THR	-	expression tag	UNP Q8A3I4
C	476	LYS	-	expression tag	UNP Q8A3I4
C	477	SER	-	expression tag	UNP Q8A3I4
C	478	ILE	-	expression tag	UNP Q8A3I4
C	479	TYR	-	expression tag	UNP Q8A3I4
C	480	ARG	-	expression tag	UNP Q8A3I4
D	34	ALA	-	expression tag	UNP Q8A3I4
D	473	LYS	-	expression tag	UNP Q8A3I4
D	474	GLY	-	expression tag	UNP Q8A3I4
D	475	THR	-	expression tag	UNP Q8A3I4
D	476	LYS	-	expression tag	UNP Q8A3I4
D	477	SER	-	expression tag	UNP Q8A3I4
D	478	ILE	-	expression tag	UNP Q8A3I4
D	479	TYR	-	expression tag	UNP Q8A3I4
D	480	ARG	-	expression tag	UNP Q8A3I4

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

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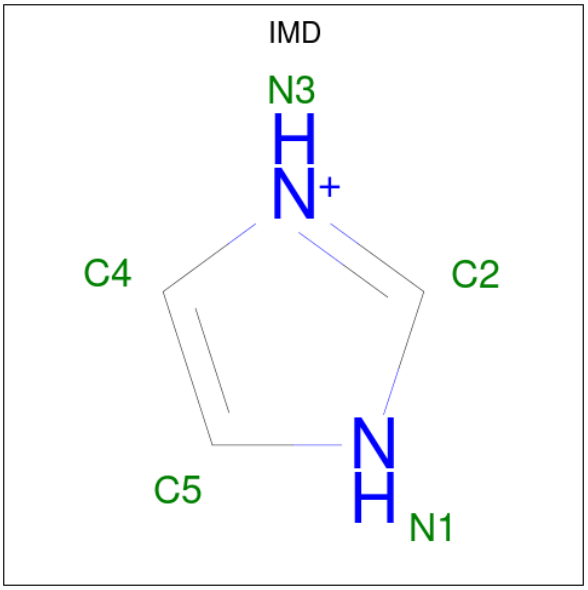
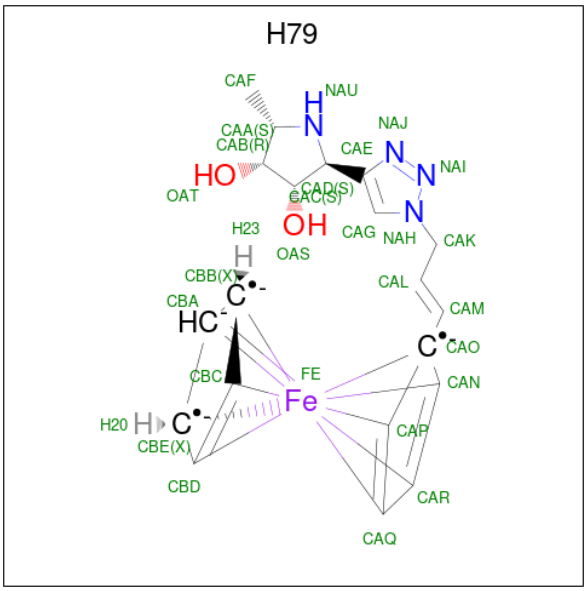
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	C	1	5	4	1	0	0

- Molecule 3 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	N		
3	A	1	5	3	2	0	0
3	B	1	5	3	2	0	0
3	C	1	5	3	2	0	0
3	D	1	5	3	2	0	0

- Molecule 4 is [(1,2,3,4,5-eta)-cyclopentadienyl][(1,2,3,4,5-eta)-1-{3-[4-(3,4-dihydroxy-5-methylpyrrolidin-2-yl)-1H-1,2,3-triazol-1-yl]prop-1-en-1-yl}cyclopentadienyl]iron (CCD ID: H79) (formula: $C_{20}H_{24}FeN_4O_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
4	A	1	27	20	1	4	2	0	0
4	B	1	27	20	1	4	2	0	0
4	C	1	27	20	1	4	2	0	0
4	D	1	27	20	1	4	2	0	0

- Molecule 5 is water.

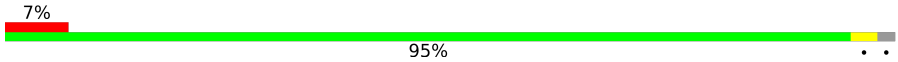
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	99	99	99	0	0
5	B	167	167	167	0	0
5	C	95	95	95	0	0
5	D	84	84	84	0	0

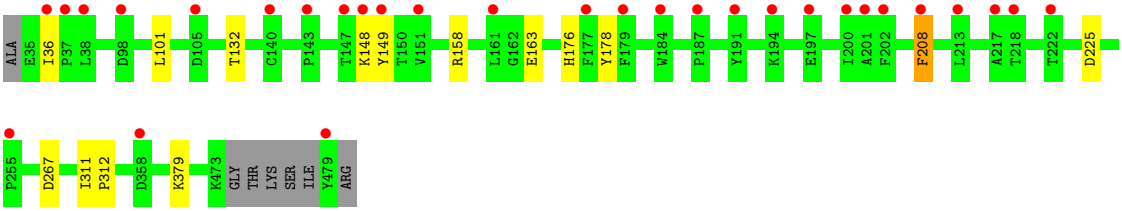
3 Residue-property plots (i)

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: Alpha-L-fucosidase

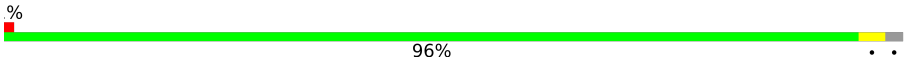
Chain A:

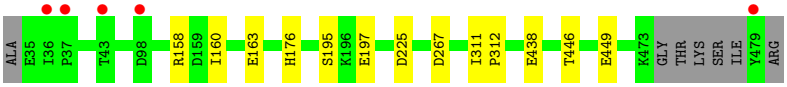




- Molecule 1: Alpha-L-fucosidase

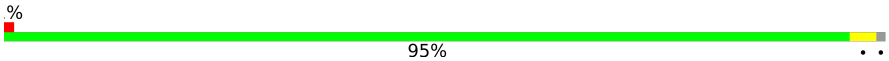
Chain B:





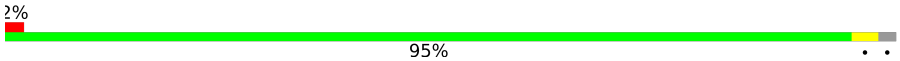
- Molecule 1: Alpha-L-fucosidase

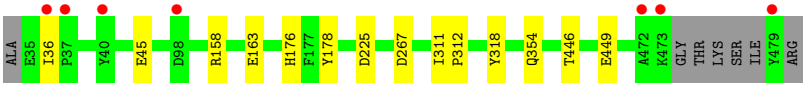
Chain C:



- Molecule 1: Alpha-L-fucosidase

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	56.50Å 188.76Å 98.14Å 90.00° 93.89° 90.00°	Depositor
Resolution (Å)	97.91 – 2.30 97.91 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.4 (97.91-2.30) 99.4 (97.91-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.51 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.201 , 0.243 0.206 , 0.247	Depositor DCC
R_{free} test set	4567 reflections (2.19%)	wwPDB-VP
Wilson B-factor (Å^2)	42.3	Xtriage
Anisotropy	0.851	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 27.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	14754	wwPDB-VP
Average B, all atoms (Å^2)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: IMD, SO4, H79

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	0.84	0/3612	0.90	1/4919 (0.0%)
1	B	0.85	0/3686	0.88	0/5004
1	C	0.83	0/3657	0.88	1/4971 (0.0%)
1	D	0.83	0/3638	0.89	2/4947 (0.0%)
All	All	0.84	0/14593	0.89	4/19841 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	36	ILE	N-CA-C	6.13	115.28	108.05
1	D	45	GLU	N-CA-C	-5.23	100.00	108.52
1	C	200	ILE	CB-CA-C	-5.13	105.40	111.97
1	A	36	ILE	N-CA-C	5.12	114.09	108.05

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	3505	0	3258	6	0
1	B	3573	0	3406	5	0
1	C	3547	0	3352	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3531	0	3322	7	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	5	0	0	0	0
3	A	5	0	5	0	0
3	B	5	0	5	0	0
3	C	5	0	5	0	0
3	D	5	0	5	0	0
4	A	27	0	15	0	0
4	B	27	0	15	0	0
4	C	27	0	15	0	0
4	D	27	0	15	0	0
5	A	99	0	0	0	0
5	B	167	0	0	0	0
5	C	95	0	0	0	0
5	D	84	0	0	0	0
All	All	14754	0	13418	24	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 (24) 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:C:446:THR:HG22	1:C:449:GLU:HB3	1.82	0.61
1:B:446:THR:HG22	1:B:449:GLU:HB3	1.82	0.60
1:C:176:HIS:CG	1:C:225:ASP:HB3	2.38	0.59
1:D:176:HIS:CG	1:D:225:ASP:HB3	2.39	0.57
1:D:446:THR:HG22	1:D:449:GLU:HB3	1.85	0.57
1:B:176:HIS:CG	1:B:225:ASP:HB3	2.41	0.56
1:A:176:HIS:CG	1:A:225:ASP:HB3	2.40	0.55
1:D:311:ILE:HB	1:D:312:PRO:HD3	1.95	0.49
1:A:208:PHE:CD1	1:A:208:PHE:C	2.91	0.49
1:D:318:TYR:CE2	1:D:354:GLN:HG2	2.49	0.48
1:A:311:ILE:HB	1:A:312:PRO:HD3	1.96	0.47
1:C:158:ARG:HD2	1:C:163:GLU:OE2	2.15	0.47
1:B:158:ARG:HD2	1:B:163:GLU:OE2	2.15	0.46
1:A:158:ARG:HD2	1:A:163:GLU:OE2	2.16	0.46
1:D:158:ARG:HD2	1:D:163:GLU:OE2	2.16	0.46
1:B:311:ILE:HB	1:B:312:PRO:HD3	1.98	0.45
1:C:311:ILE:HB	1:C:312:PRO:HD3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:178:TYR:CD1	1:D:178:TYR:C	2.97	0.43
1:A:132:THR:HA	1:A:178:TYR:HB3	2.01	0.43
1:B:195:SER:OG	1:B:197[A]:GLU:HG2	2.19	0.43
1:C:141:LEU:N	1:C:141:LEU:HD12	2.36	0.41
1:C:132:THR:HA	1:C:178:TYR:HB3	2.03	0.41
1:D:318:TYR:CD1	1:D:318:TYR:C	2.99	0.41
1:A:101:LEU:HD13	1:A:149:TYR:CE1	2.57	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	437/447 (98%)	424 (97%)	13 (3%)	0	100	100
1	B	439/447 (98%)	427 (97%)	12 (3%)	0	100	100
1	C	439/447 (98%)	426 (97%)	12 (3%)	1 (0%)	43	55
1	D	437/447 (98%)	424 (97%)	13 (3%)	0	100	100
All	All	1752/1788 (98%)	1701 (97%)	50 (3%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	60	ARG

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	351/380 (92%)	347 (99%)	4 (1%)	65	81
1	B	367/380 (97%)	364 (99%)	3 (1%)	73	86
1	C	359/380 (94%)	357 (99%)	2 (1%)	78	89
1	D	357/380 (94%)	356 (100%)	1 (0%)	86	93
All	All	1434/1520 (94%)	1424 (99%)	10 (1%)	76	87

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	148	LYS
1	A	208	PHE
1	A	267	ASP
1	A	379	LYS
1	B	160	ILE
1	B	267	ASP
1	B	438	GLU
1	C	54	GLN
1	C	267	ASP
1	D	267	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	245	HIS
1	C	375	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 oligosaccharides in this entry.

5.6 Ligand geometry (i)

13 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	SO4	A	502	-	4,4,4	0.35	0	6,6,6	0.55	0
2	SO4	B	501	-	4,4,4	0.49	0	6,6,6	0.21	0
3	IMD	D	501	-	5,5,5	0.76	0	5,5,5	0.44	0
2	SO4	B	502	-	4,4,4	0.45	0	6,6,6	0.14	0
4	H79	B	504	-	17,38,38	1.66	2 (11%)	14,103,103	3.83	7 (50%)
2	SO4	A	501	-	4,4,4	0.45	0	6,6,6	0.35	0
3	IMD	A	503	-	5,5,5	0.66	0	5,5,5	0.29	0
3	IMD	B	503	-	5,5,5	0.81	0	5,5,5	0.66	0
4	H79	A	504	-	17,38,38	1.76	4 (23%)	14,103,103	4.12	8 (57%)
4	H79	D	502	-	17,38,38	1.96	6 (35%)	14,103,103	3.63	7 (50%)
2	SO4	C	501	-	4,4,4	0.30	0	6,6,6	0.38	0
3	IMD	C	502	-	5,5,5	0.76	0	5,5,5	0.64	0
4	H79	C	503	-	17,38,38	1.73	4 (23%)	14,103,103	3.94	7 (50%)

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	IMD	D	501	-	-	-	0/1/1/1
4	H79	B	504	-	-	3/6/230/230	0/14/12/12
3	IMD	A	503	-	-	-	0/1/1/1
3	IMD	B	503	-	-	-	0/1/1/1
4	H79	A	504	-	-	1/6/230/230	0/14/12/12

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	H79	D	502	-	-	5/6/230/230	0/14/12/12
3	IMD	C	502	-	-	-	0/1/1/1
4	H79	C	503	-	-	4/6/230/230	0/14/12/12

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	504	H79	NAJ-NAI	4.08	1.38	1.32
4	D	502	H79	CAD-CAE	-3.91	1.47	1.51
4	B	504	H79	NAJ-NAI	3.77	1.38	1.32
4	C	503	H79	CAD-CAE	-3.76	1.47	1.51
4	C	503	H79	CAD-NAU	3.76	1.51	1.46
4	A	504	H79	CAG-CAE	3.55	1.41	1.36
4	C	503	H79	NAJ-NAI	3.49	1.37	1.32
4	A	504	H79	NAH-NAI	3.11	1.39	1.34
4	D	502	H79	NAJ-NAI	3.04	1.37	1.32
4	B	504	H79	CAG-CAE	2.96	1.40	1.36
4	A	504	H79	CAD-CAE	-2.75	1.48	1.51
4	D	502	H79	CAK-NAH	-2.71	1.45	1.47
4	D	502	H79	CAG-CAE	2.65	1.40	1.36
4	D	502	H79	CAB-CAA	-2.62	1.50	1.53
4	C	503	H79	NAH-NAI	2.28	1.38	1.34
4	D	502	H79	CAD-NAU	2.14	1.49	1.46

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	504	H79	CAE-NAJ-NAI	-10.10	100.86	109.24
4	C	503	H79	CAE-NAJ-NAI	-9.64	101.25	109.24
4	D	502	H79	CAE-NAJ-NAI	-9.24	101.58	109.24
4	B	504	H79	CAE-NAJ-NAI	-7.52	103.00	109.24
4	B	504	H79	OAS-CAC-CAB	7.05	134.63	111.82
4	A	504	H79	NAH-NAI-NAJ	6.67	113.07	107.01
4	B	504	H79	NAH-NAI-NAJ	5.86	112.34	107.01
4	C	503	H79	OAS-CAC-CAB	5.71	130.28	111.82
4	C	503	H79	NAH-NAI-NAJ	5.60	112.10	107.01
4	D	502	H79	NAH-NAI-NAJ	5.42	111.94	107.01
4	D	502	H79	CAG-CAE-NAJ	5.24	114.04	107.97
4	A	504	H79	CAG-CAE-NAJ	5.15	113.94	107.97
4	C	503	H79	CAG-CAE-NAJ	5.09	113.87	107.97
4	B	504	H79	CAK-NAH-CAG	4.59	136.99	128.26
4	B	504	H79	CAG-NAH-NAI	-4.53	106.09	110.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	D	502	H79	CAK-NAH-CAG	4.47	136.76	128.26
4	A	504	H79	CAK-NAH-CAG	4.33	136.48	128.26
4	A	504	H79	CAG-NAH-NAI	-4.20	106.43	110.73
4	A	504	H79	OAS-CAC-CAD	4.06	118.59	110.49
4	C	503	H79	CAG-NAH-NAI	-3.78	106.86	110.73
4	B	504	H79	CAG-CAE-NAJ	3.39	111.90	107.97
4	C	503	H79	OAS-CAC-CAD	2.94	116.36	110.49
4	D	502	H79	OAS-CAC-CAD	2.67	115.81	110.49
4	D	502	H79	CAG-NAH-NAI	-2.57	108.11	110.73
4	A	504	H79	CAF-CAA-NAU	2.39	114.50	111.70
4	C	503	H79	CAK-NAH-CAG	2.35	132.73	128.26
4	A	504	H79	CAL-CAK-NAH	-2.24	105.85	112.27
4	B	504	H79	CAF-CAA-NAU	2.15	114.22	111.70
4	D	502	H79	CAF-CAA-NAU	-2.04	109.32	111.70

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	504	H79	NAU-CAD-CAE-CAG
4	B	504	H79	CAC-CAD-CAE-CAG
4	B	504	H79	NAU-CAD-CAE-CAG
4	C	503	H79	CAC-CAD-CAE-CAG
4	C	503	H79	CAL-CAK-NAH-CAG
4	B	504	H79	NAH-CAK-CAL-CAM
4	D	502	H79	CAL-CAK-NAH-NAI
4	C	503	H79	NAH-CAK-CAL-CAM
4	D	502	H79	CAK-CAL-CAM-CAO
4	C	503	H79	NAU-CAD-CAE-CAG
4	D	502	H79	NAU-CAD-CAE-CAG
4	D	502	H79	CAL-CAK-NAH-CAG
4	D	502	H79	NAH-CAK-CAL-CAM

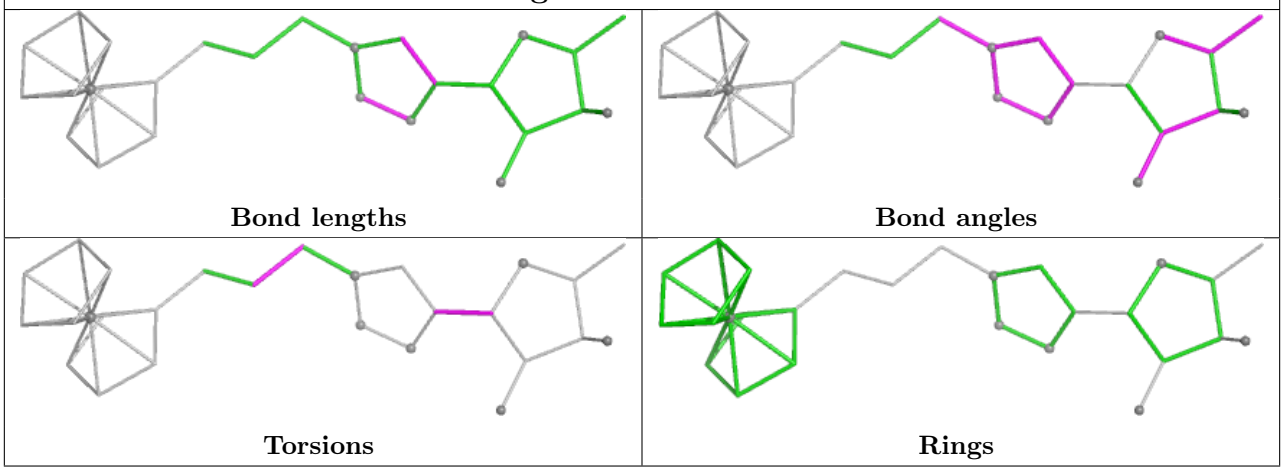
There are no ring outliers.

No monomer is involved in short contacts.

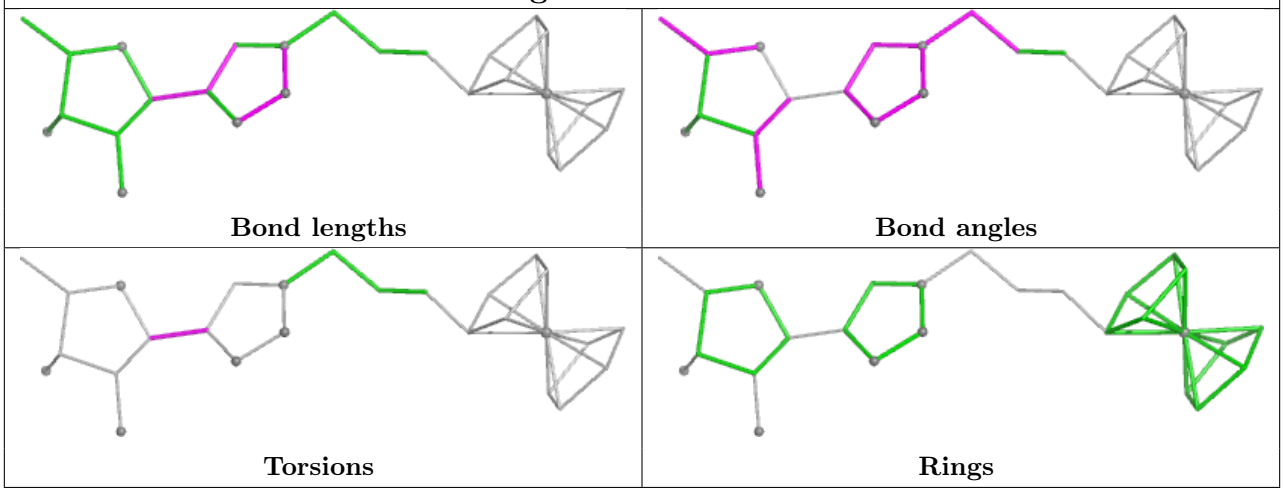
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

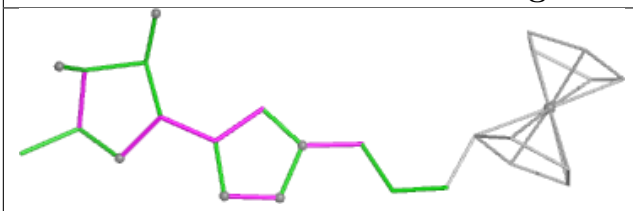
Ligand H79 B 504



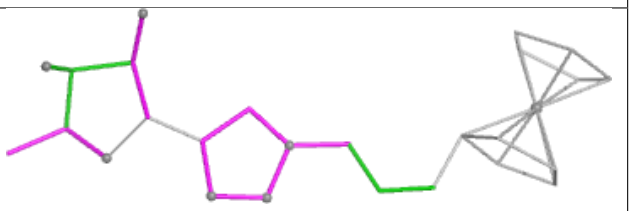
Ligand H79 A 504



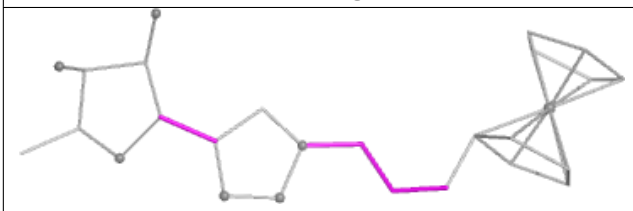
Ligand H79 D 502



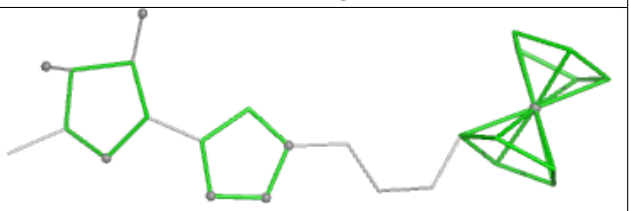
Bond lengths



Bond angles

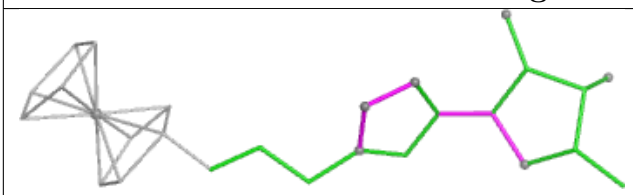


Torsions

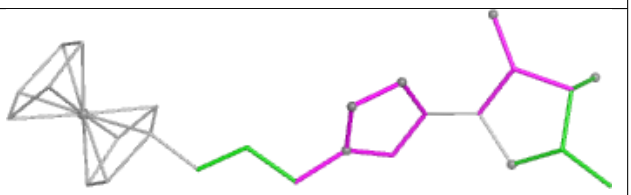


Rings

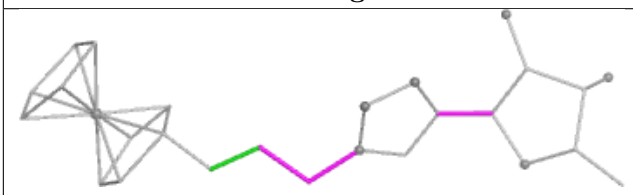
Ligand H79 C 503



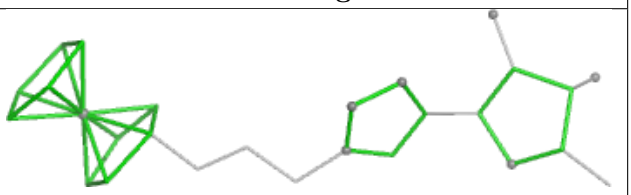
Bond lengths



Bond angles



Torsions



Rings

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	440/447 (98%)	0.49	30 (6%)	23	25	31, 58, 89, 126	1 (0%)
1	B	440/447 (98%)	-0.05	5 (1%)	78	79	28, 45, 72, 101	3 (0%)
1	C	441/447 (98%)	0.00	6 (1%)	73	75	32, 48, 70, 100	2 (0%)
1	D	440/447 (98%)	0.23	7 (1%)	70	72	33, 52, 79, 108	1 (0%)
All	All	1761/1788 (98%)	0.17	48 (2%)	56	58	28, 50, 80, 126	7 (0%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	ASP	4.7
1	B	37	PRO	3.7
1	A	179	PHE	3.6
1	A	218	THR	3.4
1	D	37	PRO	3.3
1	C	98	ASP	3.2
1	A	479	TYR	3.2
1	C	479	TYR	2.9
1	B	98	ASP	2.8
1	A	191	TYR	2.8
1	A	200	ILE	2.8
1	A	194	LYS	2.7
1	D	473	LYS	2.7
1	A	149	TYR	2.7
1	A	37	PRO	2.7
1	A	201	ALA	2.6
1	D	98	ASP	2.6
1	C	36	ILE	2.5
1	C	79	GLY	2.5
1	B	479	TYR	2.5
1	C	473	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	197	GLU	2.4
1	A	213	LEU	2.4
1	A	140	CYS	2.4
1	A	255	PRO	2.4
1	C	37	PRO	2.4
1	A	208	PHE	2.3
1	B	36	ILE	2.3
1	A	143	PRO	2.3
1	A	161	LEU	2.2
1	D	472	ALA	2.2
1	D	36	ILE	2.2
1	A	151	VAL	2.2
1	A	187	PRO	2.2
1	D	40	TYR	2.1
1	D	479	TYR	2.1
1	A	358	ASP	2.1
1	A	184	TRP	2.1
1	B	43	THR	2.1
1	A	217	ALA	2.1
1	A	36	ILE	2.1
1	A	147	THR	2.1
1	A	177	PHE	2.1
1	A	222	THR	2.0
1	A	38	LEU	2.0
1	A	202	PHE	2.0
1	A	105	ASP	2.0
1	A	148	LYS	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 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,

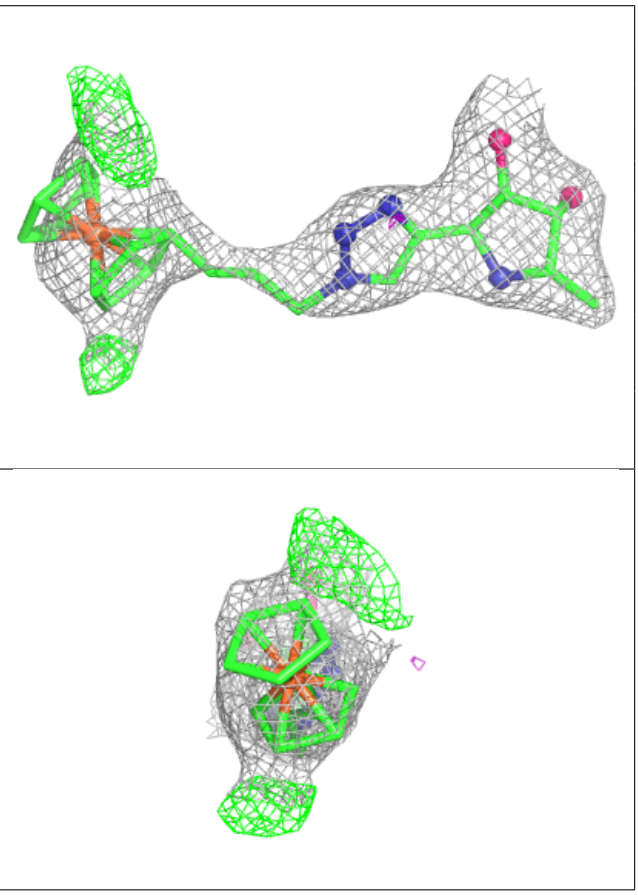
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	SO4	A	501	5/5	0.73	0.10	92,95,99,104	0
2	SO4	B	501	5/5	0.78	0.09	89,92,94,98	0
4	H79	A	504	27/27	0.81	0.16	57,92,208,212	0
2	SO4	C	501	5/5	0.85	0.15	48,50,58,59	5
2	SO4	A	502	5/5	0.85	0.14	45,45,46,52	5
2	SO4	B	502	5/5	0.86	0.09	88,90,95,99	0
4	H79	D	502	27/27	0.86	0.17	38,75,150,157	0
4	H79	C	503	27/27	0.87	0.17	39,75,199,200	0
4	H79	B	504	27/27	0.89	0.16	36,82,208,218	0
3	IMD	D	501	5/5	0.90	0.19	85,89,94,96	0
3	IMD	C	502	5/5	0.92	0.14	61,68,70,72	0
3	IMD	B	503	5/5	0.93	0.22	58,60,63,64	0
3	IMD	A	503	5/5	0.95	0.11	74,75,82,83	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.

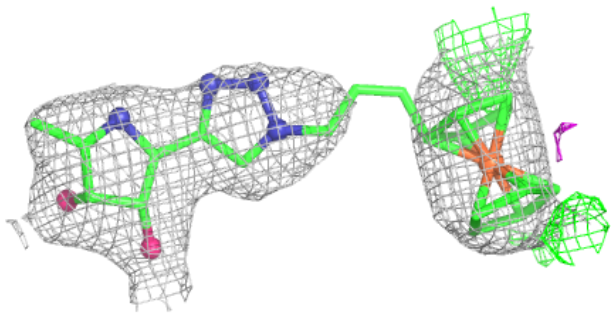
Electron density around H79 A 504:

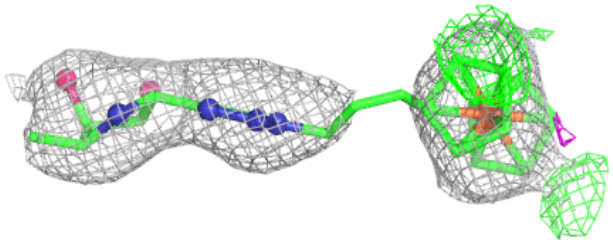
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

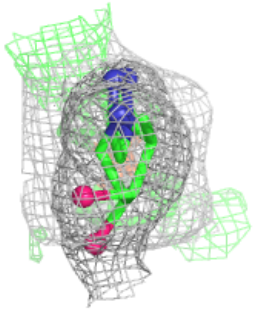


Electron density around H79 D 502:

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

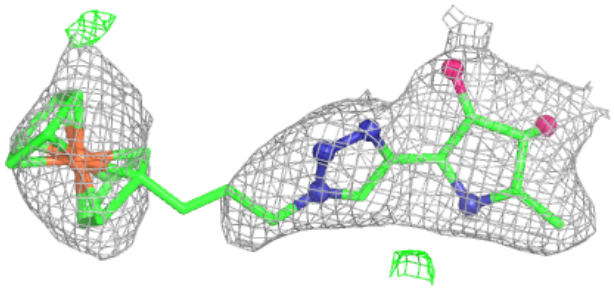


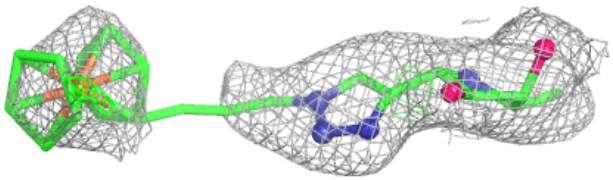


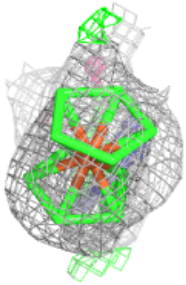


Electron density around H79 C 503:

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

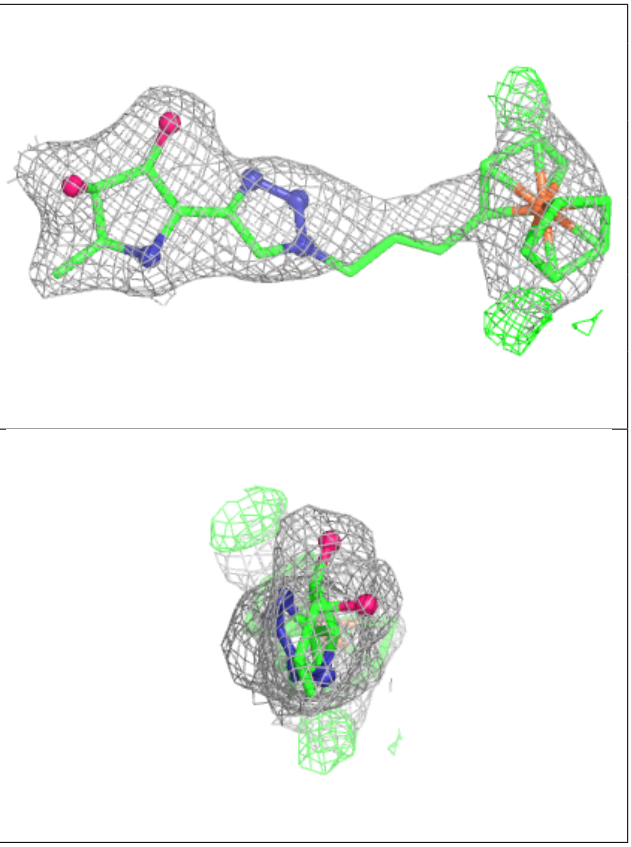






Electron density around H79 B 504:

$2mF_o$-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.