



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 07:06 PM UTC

PDB ID : 8FY2 / pdb_00008fy2
Title : E3:PROTAC:target ternary complex structure (VCB/WH244/BCL-2)
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Deposited on : 2023-01-25
Resolution : 2.98 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 : 2022.3.0, CSD as543be (2022)
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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

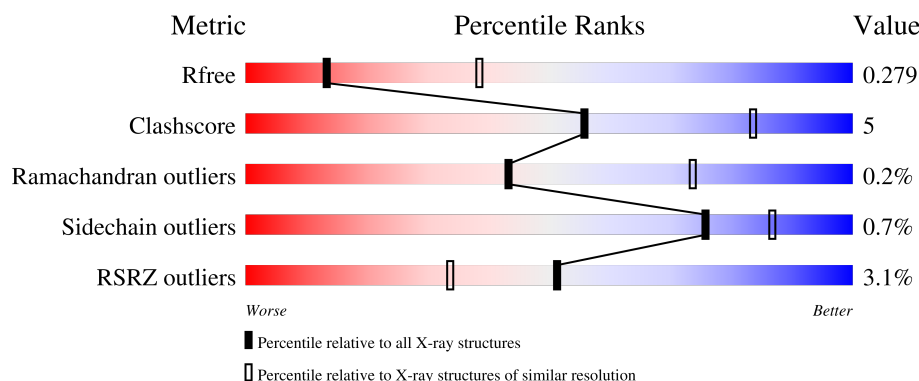
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	180	
2	B	118	
3	C	96	
4	D	209	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4022 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 von Hippel-Lindau disease tumor suppressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1223	776	227	218	2			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	MET	-	expression tag	UNP P40337
A	35	GLY	-	expression tag	UNP P40337
A	36	SER	-	expression tag	UNP P40337
A	37	SER	-	expression tag	UNP P40337
A	38	HIS	-	expression tag	UNP P40337
A	39	HIS	-	expression tag	UNP P40337
A	40	HIS	-	expression tag	UNP P40337
A	41	HIS	-	expression tag	UNP P40337
A	42	HIS	-	expression tag	UNP P40337
A	43	HIS	-	expression tag	UNP P40337
A	44	SER	-	expression tag	UNP P40337
A	45	SER	-	expression tag	UNP P40337
A	46	GLY	-	expression tag	UNP P40337
A	47	LEU	-	expression tag	UNP P40337
A	48	VAL	-	expression tag	UNP P40337
A	49	PRO	-	expression tag	UNP P40337
A	50	ARG	-	expression tag	UNP P40337
A	51	GLY	-	expression tag	UNP P40337
A	52	SER	-	expression tag	UNP P40337
A	53	HIS	-	expression tag	UNP P40337

- Molecule 2 is a protein called Elongin-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	105	Total	C	N	O	S	0	0	0
			829	525	139	160	5			

- Molecule 3 is a protein called Elongin-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	85	Total	C	N	O	S	0	0	0
			681	440	109	126	6			

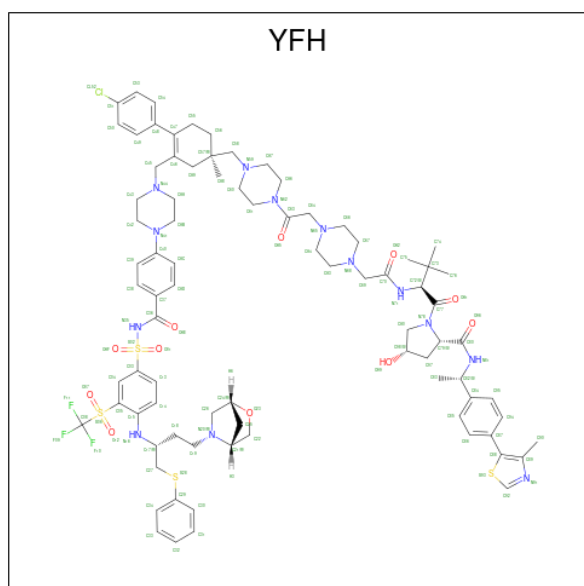
- Molecule 4 is a protein called Apoptosis regulator Bcl-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	141	Total	C	N	O	S	0	0	0
			1174	753	207	208	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	GLY	-	expression tag	UNP P10415
D	0	SER	-	expression tag	UNP P10415
D	8	VAL	GLY	conflict	UNP P10415
D	87	GLY	SER	conflict	UNP P10415

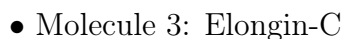
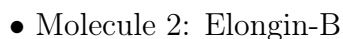
- Molecule 5 is N-({4-[2-(4-{[(4R)-4'-chloro-4-methyl-6-{[4-(4-{[4-{[(2R)-4-[(1R,4R)-2-oxa-5-azabicyclo[2.2.1]heptan-5-yl]-1-(phenylsulfanyl)butan-2-yl]amino}-3-(trifluoromethanesulfonyl)benzene-1-sulfonyl]carbamoyl}phenyl)piperazin-1-yl]methyl}-2,3,4,5-tetrahydro[1,1'-biphenyl]-4-yl]methyl}piperazin-1-yl)-2-oxoethyl]piperazin-1-yl]acetyl)-3-methyl-L-valyl-(4S)-4-hydroxy-N-{(1S)-1-[4-(4-methyl-1,3-thiazol-5-yl)phenyl]ethyl}-L-prolinamide (CCD ID: YFH) (formula: C₈₃H₁₀₅ClF₃N₁₃O₁₁S₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
5	D	1	Total	C	Cl	F	N	O	S	0	0
			115	83	1	3	13	11	4		

i

- Molecule 1: von Hippel-Lindau disease tumor suppressor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.28Å 102.60Å 167.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	83.68 – 2.98 83.68 – 2.98	Depositor EDS
% Data completeness (in resolution range)	99.4 (83.68-2.98) 89.7 (83.68-2.98)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.95 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.229 , 0.279 0.230 , 0.279	Depositor DCC
R_{free} test set	1733 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	80.9	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 54.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4022	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.64% 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: YFH

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	0.08	0/1255	0.23	0/1712
2	B	0.07	0/846	0.22	0/1144
3	C	0.07	0/695	0.21	0/937
4	D	0.07	0/1208	0.20	0/1637
All	All	0.08	0/4004	0.21	0/5430

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	1223	0	1224	14	0
2	B	829	0	831	6	0
3	C	681	0	683	6	0
4	D	1174	0	1113	11	0
5	D	115	0	0	3	0
All	All	4022	0	3851	37	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 5.

All (37) 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)
3:C:80:LYS:HE2	3:C:84:THR:HG21	1.78	0.65
1:A:84:VAL:HG11	1:A:151:ILE:HD13	1.79	0.64
2:B:78:ALA:HB1	2:B:86:GLU:HB3	1.79	0.64
2:B:9:ARG:HB2	2:B:77:LEU:HB3	1.82	0.61
3:C:20:LYS:HB3	3:C:59:GLU:HG2	1.83	0.60
4:D:119:LEU:HD21	4:D:157:MET:HG3	1.84	0.60
3:C:72:LYS:HA	3:C:75:MET:HE3	1.84	0.60
4:D:130:PHE:HB2	4:D:157:MET:HE1	1.84	0.59
2:B:79:PHE:O	2:B:86:GLU:HB2	2.04	0.57
1:A:182:ARG:NH1	1:A:186:GLU:OE2	2.41	0.54
2:B:6:MET:HE2	2:B:72:PRO:HG2	1.91	0.53
1:A:102:PRO:HD2	1:A:105:THR:HG21	1.92	0.51
1:A:87:VAL:HB	1:A:118:LEU:HB2	1.92	0.51
1:A:190:ASP:OD1	1:A:196:LYS:NZ	2.42	0.50
2:B:43:ARG:HB2	2:B:78:ALA:HB3	1.95	0.49
4:D:17:LYS:O	4:D:98:ARG:NH1	2.39	0.48
1:A:165:VAL:HG21	3:C:95:ILE:HB	1.96	0.47
4:D:188:TRP:O	4:D:192:ASN:ND2	2.35	0.47
4:D:12:ARG:O	4:D:16:MET:HG2	2.15	0.47
1:A:129:LEU:HG	1:A:154:PRO:HB3	1.95	0.47
2:B:23:THR:HA	2:B:56:THR:HA	1.97	0.47
4:D:21:TYR:CG	4:D:98:ARG:HD3	2.51	0.46
1:A:120:ARG:NH2	1:A:197:ASP:OD2	2.42	0.46
3:C:18:TYR:CE2	3:C:32:LYS:HG3	2.50	0.46
1:A:112:TYR:HB2	1:A:115:HIS:CE1	2.52	0.45
4:D:181:LEU:HA	4:D:185:LEU:HB2	1.99	0.44
4:D:145:GLY:HA3	5:D:301:YFH:N35	2.32	0.44
4:D:94:HIS:O	4:D:98:ARG:HG3	2.17	0.44
1:A:90:ASN:HD21	1:A:94:GLU:HB2	1.83	0.44
1:A:116:LEU:HD12	1:A:116:LEU:HA	1.88	0.43
1:A:60:ARG:HB2	1:A:61:PRO:HD3	2.02	0.42
1:A:109:ILE:HG23	5:D:301:YFH:C94	2.50	0.41
3:C:21:LEU:HD12	3:C:62:PHE:HE2	1.84	0.41
5:D:301:YFH:O07	5:D:301:YFH:N16	2.53	0.41
4:D:96:THR:HG21	4:D:199:VAL:HA	2.03	0.41
1:A:90:ASN:OD1	1:A:94:GLU:N	2.52	0.40
4:D:122:THR:HB	4:D:123:PRO:HD2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	147/180 (82%)	142 (97%)	5 (3%)	0	100	100
2	B	103/118 (87%)	94 (91%)	8 (8%)	1 (1%)	12	42
3	C	81/96 (84%)	80 (99%)	1 (1%)	0	100	100
4	D	137/209 (66%)	129 (94%)	8 (6%)	0	100	100
All	All	468/603 (78%)	445 (95%)	22 (5%)	1 (0%)	43	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	47	ASP

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	139/164 (85%)	138 (99%)	1 (1%)	76	87
2	B	93/103 (90%)	91 (98%)	2 (2%)	45	73
3	C	77/85 (91%)	77 (100%)	0	100	100
4	D	123/165 (74%)	123 (100%)	0	100	100
All	All	432/517 (84%)	429 (99%)	3 (1%)	76	87

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

Mol	Chain	Res	Type
1	A	130	VAL
2	B	80	ARG
2	B	85	PHE

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

Mol	Chain	Res	Type
1	A	150	ASN
4	D	120	HIS
4	D	190	GLN

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](#)

1 ligand is 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	$\# Z > 2$
5	YFH	D	301	-	125,127,127	1.17	12 (9%)	166,188,188	1.40	17 (10%)

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
5	YFH	D	301	-	-	27/107/181/181	2/14/13/13

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	301	YFH	C89-N91	5.05	1.45	1.38
5	D	301	YFH	C39-C38	2.53	1.42	1.38
5	D	301	YFH	C86-C87	2.35	1.43	1.39
5	D	301	YFH	C92-S93	2.24	1.76	1.72
5	D	301	YFH	C45-C46	2.23	1.57	1.49
5	D	301	YFH	C14-C15	2.20	1.43	1.39
5	D	301	YFH	CAD-C37	2.11	1.42	1.39
5	D	301	YFH	CAD-CAC	2.09	1.42	1.38
5	D	301	YFH	C21-N20	2.03	1.52	1.47
5	D	301	YFH	C19-N20	2.02	1.50	1.47
5	D	301	YFH	CA9-C57	2.02	1.56	1.54
5	D	301	YFH	C22-C21	2.01	1.55	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	301	YFH	C92-S93-C88	6.81	92.53	88.96
5	D	301	YFH	C89-C88-S93	-4.42	106.93	109.64
5	D	301	YFH	C03-S02-N35	4.18	112.72	106.06
5	D	301	YFH	S93-C92-N91	-4.08	112.58	116.08
5	D	301	YFH	C37-C36-N35	3.57	120.34	116.08
5	D	301	YFH	OAF-S02-C03	-3.44	103.65	107.98
5	D	301	YFH	C87-C88-C89	3.25	135.23	130.81
5	D	301	YFH	C05-C15-N16	-3.10	119.19	122.18
5	D	301	YFH	C86-C87-C88	2.73	125.19	120.84
5	D	301	YFH	C24-C25-C21	-2.59	96.13	102.36
5	D	301	YFH	C94-C87-C86	-2.47	115.43	118.57
5	D	301	YFH	CAB-N41-C40	2.46	124.83	118.11
5	D	301	YFH	C25-C21-N20	-2.27	98.68	103.74
5	D	301	YFH	C42-N41-C40	2.22	124.15	118.11
5	D	301	YFH	C56-C57-CA9	2.21	109.97	107.73
5	D	301	YFH	C55-C56-C57	2.13	115.86	112.60
5	D	301	YFH	CAB-CAA-N44	2.05	114.79	110.65

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	301	YFH	C27-C17-C18-C19
5	D	301	YFH	CA8-C57-C58-N59
5	D	301	YFH	C63-C64-N65-C66
5	D	301	YFH	C36-N35-S02-C03
5	D	301	YFH	C36-N35-S02-OAF
5	D	301	YFH	C77-C72-C73-C75
5	D	301	YFH	C77-C72-C73-C76
5	D	301	YFH	C36-N35-S02-O01
5	D	301	YFH	C77-C72-C73-C74
5	D	301	YFH	N71-C72-C73-C76
5	D	301	YFH	F09-C08-S06-O12
5	D	301	YFH	F10-C08-S06-O12
5	D	301	YFH	F11-C08-S06-O12
5	D	301	YFH	N71-C72-C73-C75
5	D	301	YFH	F09-C08-S06-O07
5	D	301	YFH	F10-C08-S06-O07
5	D	301	YFH	F11-C08-S06-O07
5	D	301	YFH	N16-C17-C18-C19
5	D	301	YFH	C56-C57-C58-N59
5	D	301	YFH	CA9-C57-C58-N59
5	D	301	YFH	F09-C08-S06-C05
5	D	301	YFH	F10-C08-S06-C05
5	D	301	YFH	F11-C08-S06-C05
5	D	301	YFH	C17-C18-C19-N20
5	D	301	YFH	N71-C72-C73-C74
5	D	301	YFH	C30-C29-S28-C27
5	D	301	YFH	C34-C29-S28-C27

All (2) ring outliers are listed below:

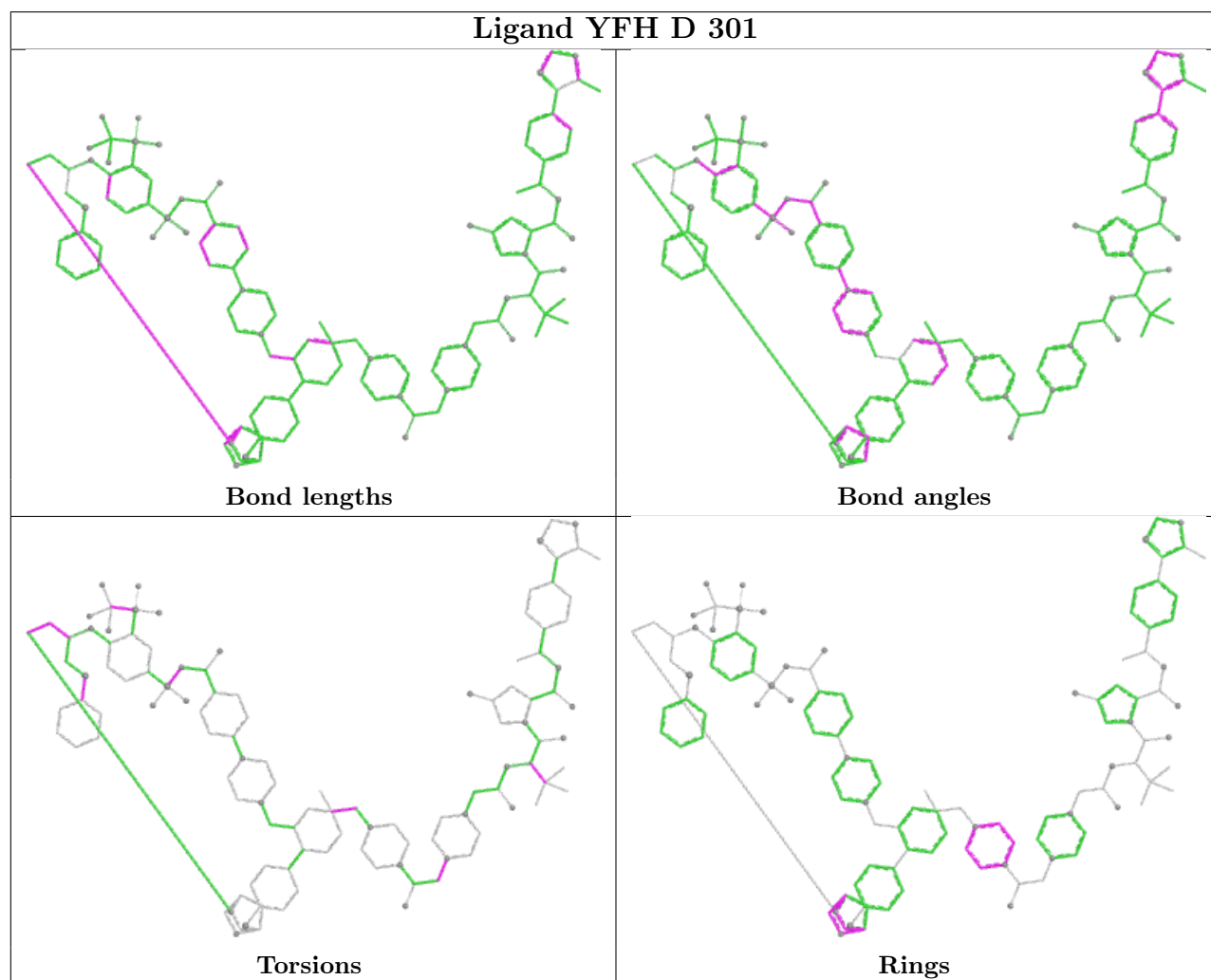
Mol	Chain	Res	Type	Atoms
5	D	301	YFH	C60-C61-CA6-CA7-N59-N62
5	D	301	YFH	C21-C22-C24-C26-N20-O23

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	301	YFH	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	149/180 (82%)	0.29	4 (2%)	56 37	63, 88, 146, 172	0
2	B	105/118 (88%)	0.16	2 (1%)	66 47	63, 84, 121, 150	0
3	C	85/96 (88%)	0.06	2 (2%)	59 40	68, 81, 109, 142	0
4	D	141/209 (67%)	0.59	7 (4%)	34 20	80, 108, 141, 176	0
All	All	480/603 (79%)	0.31	15 (3%)	51 34	63, 91, 137, 176	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	8	VAL	4.2
4	D	28	TYR	3.2
2	B	105	PRO	3.2
4	D	88	PRO	3.1
1	A	114	GLY	3.0
4	D	92	VAL	2.6
4	D	140	ASP	2.6
1	A	208	HIS	2.5
4	D	124	PHE	2.5
4	D	203	GLY	2.4
1	A	144	GLY	2.4
2	B	86	GLU	2.2
3	C	112	CYS	2.1
1	A	63	LEU	2.1
3	C	46	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

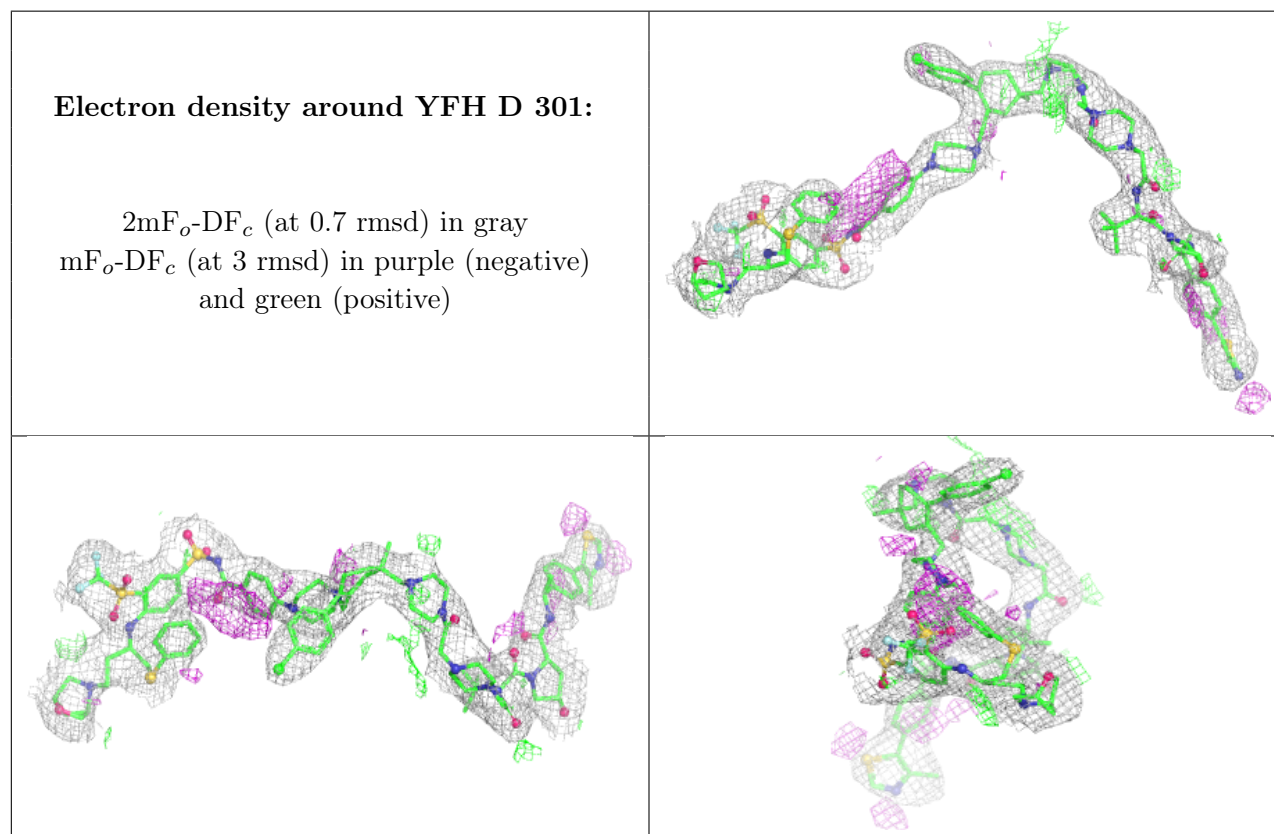
There are no oligosaccharides in this entry.

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
5	YFH	D	301	115/115	0.90	0.14	48,84,116,128	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.