



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:09 PM UTC

PDB ID : 8WUR / pdb_00008wur
Title : Crystal structure of SARS-Cov-2 main protease D48N mutant in complex with shikonin
Authors : Zhao, Z.Y.; Li, W.W.; Zhang, J.; Li, J.
Deposited on : 2023-10-21
Resolution : 2.08 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

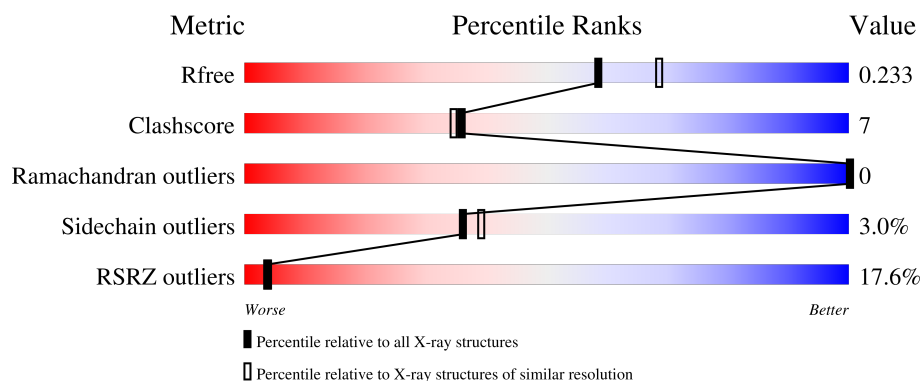
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	296	14% 65% 30% .
1	B	296	21% 70% 25% ..

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	FNO	A	301	X	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4510 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.

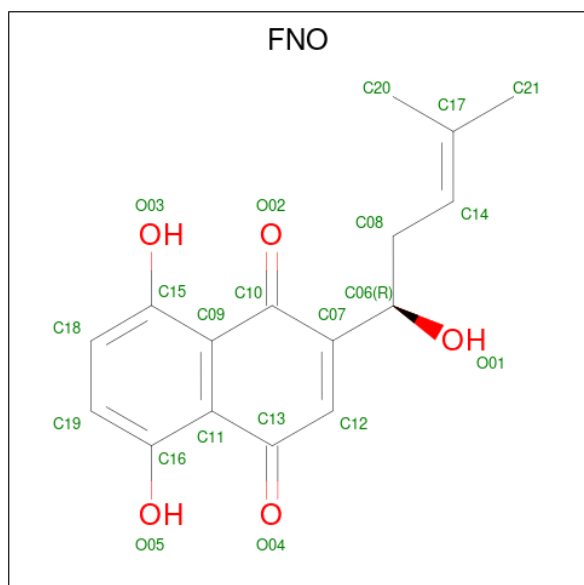
- Molecule 1 is a protein called 3C-like proteinase nsp5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	295	Total	C	N	O	S	0	0	0
			2229	1414	380	415	20			
1	B	291	Total	C	N	O	S	0	0	0
			2170	1378	366	407	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	48	ASN	ASP	engineered mutation	UNP P0DTC1
A	142	ALA	ASN	conflict	UNP P0DTC1
B	48	ASN	ASP	engineered mutation	UNP P0DTC1
B	142	ALA	ASN	conflict	UNP P0DTC1

- Molecule 2 is 2-[(1R)-4-methyl-1-oxidanyl-pent-3-enyl]-5,8-bis(oxidanyl)naphthalene-1,4-dione (CCD ID: FNO) (formula: C₁₆H₁₆O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			21	16	5		

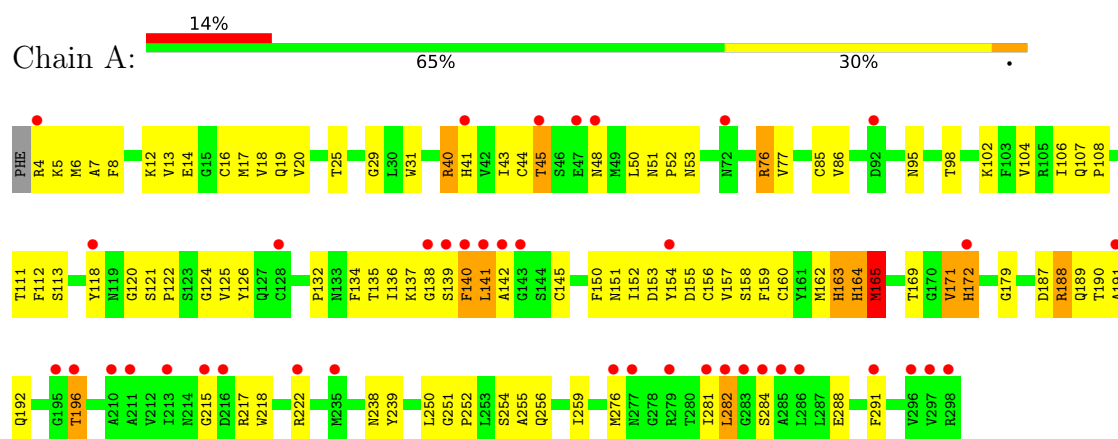
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	55	Total	O	0	0
			55	55		
3	B	35	Total	O	0	0
			35	35		

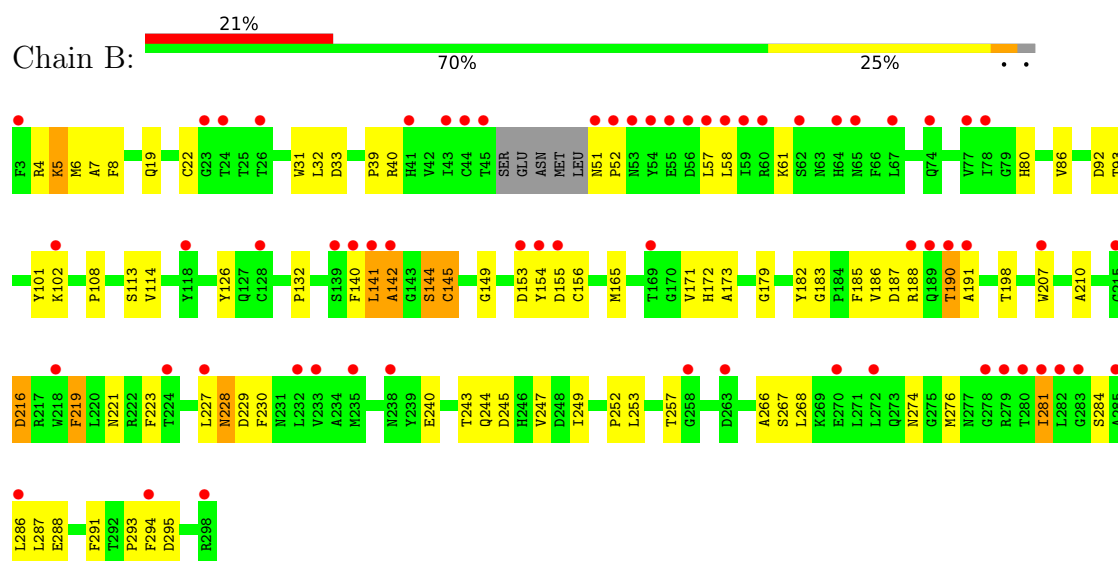
3 Residue-property plots

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: 3C-like proteinase nsp4



• Molecule 1: 3C-like proteinase nsp4



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.05Å 102.94Å 103.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.30 – 2.08 46.30 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.30-2.08) 93.7 (46.30-2.08)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.12 _2829	Depositor
R, R_{free}	0.211 , 0.238 0.216 , 0.233	Depositor DCC
R_{free} test set	2004 reflections (4.48%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.546	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4510	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% 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: FNO

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.76	79/2278 (3.5%)	1.04	11/3103 (0.4%)
1	B	1.53	38/2219 (1.7%)	0.93	7/3028 (0.2%)
All	All	1.65	117/4497 (2.6%)	0.98	18/6131 (0.3%)

All (117) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	12	LYS	C-O	-8.09	1.14	1.24
1	B	8	PHE	C-O	-7.77	1.14	1.23
1	B	267	SER	C-O	-7.69	1.15	1.24
1	A	136	ILE	C-O	-7.68	1.16	1.24
1	A	152	ILE	C-O	-7.66	1.15	1.24
1	A	14	GLU	C-O	-7.63	1.15	1.24
1	A	13	VAL	C-O	-7.47	1.15	1.24
1	B	39	PRO	C-O	-7.34	1.15	1.23
1	A	157	VAL	C-O	-7.33	1.15	1.24
1	A	158	SER	C-O	-7.32	1.15	1.24
1	B	210	ALA	C-O	-7.29	1.15	1.24
1	A	126	TYR	C-O	-7.11	1.15	1.23
1	A	29	GLY	C-O	-7.07	1.13	1.23
1	A	106	ILE	C-O	-7.03	1.16	1.23
1	B	294	PHE	C-O	-6.96	1.16	1.24
1	A	104	VAL	C-O	-6.96	1.16	1.23
1	A	77	VAL	C-O	-6.90	1.16	1.24
1	B	186	VAL	C-O	-6.83	1.17	1.24
1	A	25	THR	C-O	-6.82	1.15	1.24
1	A	112	PHE	C-O	-6.82	1.15	1.23
1	A	217	ARG	C-O	-6.76	1.14	1.24
1	A	43	ILE	C-O	-6.55	1.16	1.24
1	B	145	CYS	C-O	-6.53	1.16	1.23
1	A	163	HIS	C-O	-6.52	1.16	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	122	PRO	C-O	-6.51	1.16	1.23
1	A	153	ASP	C-O	-6.49	1.16	1.24
1	B	156	CYS	C-O	-6.48	1.16	1.24
1	B	288	GLU	C-O	-6.46	1.16	1.23
1	A	134	PHE	C-O	-6.45	1.15	1.23
1	A	188	ARG	C-O	-6.44	1.16	1.23
1	B	32	LEU	C-O	-6.44	1.16	1.24
1	A	16	CYS	C-O	-6.43	1.15	1.24
1	A	189	GLN	C-O	-6.40	1.15	1.23
1	A	156	CYS	C-O	-6.36	1.16	1.24
1	A	121	SER	C-O	-6.33	1.17	1.24
1	A	107	GLN	C-O	-6.33	1.17	1.24
1	A	18	VAL	C-O	-6.27	1.16	1.23
1	A	160	CYS	C-O	-6.26	1.15	1.24
1	A	172	HIS	C-O	-6.26	1.15	1.23
1	B	40	ARG	C-O	-6.25	1.16	1.24
1	A	118	TYR	C-O	-6.22	1.17	1.24
1	B	33	ASP	C-O	-6.19	1.18	1.24
1	B	219	PHE	C-O	-6.14	1.16	1.24
1	B	295	ASP	C-O	-6.14	1.17	1.24
1	A	6	MET	C-O	-6.11	1.16	1.23
1	A	218	TRP	C-O	-6.11	1.16	1.24
1	A	7	ALA	N-CA	-6.06	1.39	1.45
1	A	251	GLY	C-O	-6.06	1.18	1.24
1	B	284	SER	CA-C	-6.05	1.45	1.52
1	B	6	MET	C-O	-6.04	1.16	1.23
1	A	8	PHE	C-O	-6.03	1.16	1.23
1	A	162	MET	C-O	-6.01	1.17	1.24
1	B	230	PHE	C-O	-6.01	1.17	1.24
1	A	164	HIS	C-O	-5.93	1.16	1.24
1	A	44	CYS	C-O	-5.90	1.17	1.23
1	A	41	HIS	C-O	-5.89	1.16	1.24
1	A	187	ASP	C-O	-5.89	1.17	1.23
1	A	135	THR	C-O	-5.88	1.16	1.23
1	A	124	GLY	C-O	-5.87	1.17	1.24
1	B	171	VAL	C-O	-5.86	1.17	1.24
1	A	51	ASN	C-O	-5.85	1.17	1.24
1	A	196	THR	C-O	-5.83	1.16	1.23
1	B	153	ASP	N-CA	-5.82	1.39	1.46
1	A	19	GLN	C-O	-5.82	1.17	1.24
1	A	52	PRO	C-O	-5.81	1.17	1.23
1	B	268	LEU	C-O	-5.78	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	7	ALA	C-O	-5.75	1.16	1.23
1	A	5	LYS	C-O	-5.73	1.17	1.24
1	A	284	SER	CA-C	-5.73	1.46	1.52
1	A	151	ASN	C-O	-5.70	1.16	1.23
1	B	183	GLY	C-O	-5.64	1.16	1.24
1	A	255	ALA	C-O	-5.64	1.17	1.24
1	B	229	ASP	C-O	-5.63	1.17	1.24
1	A	165	MET	C-O	-5.63	1.17	1.23
1	A	76	ARG	C-O	-5.61	1.17	1.24
1	A	259	ILE	C-O	-5.58	1.18	1.24
1	A	238	ASN	C-O	-5.58	1.16	1.23
1	A	254	SER	C-O	-5.56	1.17	1.24
1	B	182	TYR	C-O	-5.55	1.17	1.24
1	A	138	GLY	C-O	-5.55	1.16	1.23
1	A	53	ASN	C-O	-5.44	1.17	1.24
1	A	102	LYS	C-O	-5.43	1.17	1.23
1	A	113	SER	C-O	-5.41	1.17	1.23
1	A	250	LEU	C-O	-5.39	1.16	1.24
1	A	150	PHE	C-O	-5.38	1.17	1.23
1	A	17	MET	C-O	-5.38	1.17	1.24
1	B	243	THR	C-O	-5.38	1.16	1.23
1	A	7	ALA	C-O	-5.36	1.17	1.23
1	B	31	TRP	C-O	-5.34	1.17	1.24
1	A	137	LYS	C-O	-5.34	1.17	1.23
1	A	120	GLY	C-O	-5.33	1.16	1.24
1	A	256	GLN	C-O	-5.33	1.17	1.24
1	A	282	LEU	CA-C	-5.33	1.46	1.53
1	A	40	ARG	C-O	-5.32	1.17	1.24
1	B	19	GLN	C-O	-5.32	1.17	1.24
1	A	31	TRP	C-O	-5.30	1.17	1.24
1	B	22	CYS	C-O	-5.30	1.18	1.24
1	B	252	PRO	C-O	-5.29	1.17	1.24
1	A	159	PHE	C-O	-5.28	1.17	1.24
1	A	125	VAL	C-O	-5.28	1.18	1.24
1	A	191	ALA	C-O	-5.25	1.17	1.24
1	B	221	ASN	C-O	-5.24	1.17	1.23
1	A	50	LEU	C-O	-5.19	1.18	1.24
1	A	20	VAL	C-O	-5.19	1.18	1.24
1	B	144	SER	C-O	-5.17	1.16	1.23
1	B	247	VAL	C-O	-5.15	1.18	1.24
1	B	185	PHE	C-O	-5.13	1.17	1.23
1	A	239	TYR	C-O	-5.12	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	253	LEU	C-O	-5.11	1.18	1.24
1	A	171	VAL	C-O	-5.11	1.18	1.24
1	B	257	THR	C-O	-5.10	1.17	1.24
1	A	111	THR	C-O	-5.08	1.17	1.23
1	B	266	ALA	C-O	-5.07	1.18	1.24
1	B	187	ASP	C-O	-5.06	1.16	1.23
1	A	282	LEU	CA-CB	-5.05	1.46	1.53
1	B	293	PRO	C-O	-5.02	1.17	1.24
1	A	136	ILE	N-CA	-5.01	1.40	1.46

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	GLY	N-CA-C	10.22	128.95	115.36
1	A	141	LEU	N-CA-C	-8.46	100.00	110.41
1	A	251	GLY	O-C-N	8.32	124.06	121.07
1	B	223	PHE	N-CA-C	8.27	121.72	110.55
1	B	274	ASN	N-CA-C	7.51	123.48	113.72
1	B	190	THR	N-CA-C	-7.50	99.71	110.59
1	A	165	MET	N-CA-C	7.11	119.67	108.79
1	A	150	PHE	N-CA-C	6.75	118.46	108.60
1	A	282	LEU	N-CA-C	-6.48	101.65	111.56
1	A	282	LEU	CA-CB-CG	-6.08	95.00	116.30
1	B	182	TYR	N-CA-C	-5.59	99.62	108.73
1	A	13	VAL	N-CA-C	5.55	116.33	110.72
1	A	142	ALA	N-CA-C	5.55	118.09	111.71
1	B	5	LYS	N-CA-C	-5.38	101.71	110.20
1	A	192	GLN	N-CA-C	-5.35	99.96	109.06
1	B	286	LEU	N-CA-C	-5.34	101.07	109.50
1	B	142	ALA	N-CA-C	-5.32	103.12	110.35
1	A	140	PHE	N-CA-C	5.11	117.57	109.24

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	2229	0	2156	33	0
1	B	2170	0	2054	32	0
2	A	21	0	0	10	0
3	A	55	0	0	0	0
3	B	35	0	0	0	0
All	All	4510	0	4210	63	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 7.

All (63) 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:165:MET:CE	2:A:301:FNO:O04	2.04	1.04
1:A:165:MET:HE2	2:A:301:FNO:C12	1.88	1.03
1:B:51:ASN:HA	1:B:188:ARG:HD2	1.43	0.99
1:A:165:MET:CE	2:A:301:FNO:C13	2.41	0.97
1:A:165:MET:HE3	2:A:301:FNO:O04	1.66	0.92
1:A:140:PHE:HD2	1:A:172:HIS:CD2	1.91	0.89
1:A:165:MET:CE	2:A:301:FNO:C12	2.51	0.88
1:B:216:ASP:OD2	1:B:281:ILE:O	1.90	0.88
1:A:76:ARG:HH11	1:A:76:ARG:HG3	1.40	0.87
1:B:141:LEU:HD12	1:B:142:ALA:O	1.74	0.87
1:A:45:THR:H	1:A:48:ASN:HD22	1.21	0.85
1:A:165:MET:HE2	2:A:301:FNO:C13	2.09	0.79
1:A:76:ARG:HG3	1:A:76:ARG:NH1	2.02	0.75
1:A:139:SER:HB2	1:B:4:ARG:HB2	1.72	0.72
1:B:52:PRO:HD2	1:B:188:ARG:HG2	1.75	0.69
1:A:169:THR:HG23	1:A:171:VAL:HG22	1.78	0.66
1:A:165:MET:HE1	2:A:301:FNO:O04	1.97	0.64
1:A:140:PHE:CD2	1:A:172:HIS:CD2	2.81	0.63
1:B:92:ASP:OD1	1:B:93:THR:OG1	2.08	0.61
1:A:4:ARG:HD2	1:B:126:TYR:CD2	2.38	0.59
1:B:198:THR:OG1	1:B:240:GLU:OE1	2.20	0.58
1:B:58:LEU:HD11	1:B:80:HIS:HD2	1.71	0.56
1:B:190:THR:HG22	1:B:191:ALA:N	2.21	0.55
1:B:140:PHE:HB2	1:B:172:HIS:CE1	2.42	0.55
1:A:45:THR:H	1:A:48:ASN:ND2	2.00	0.54
1:A:165:MET:HE3	2:A:301:FNO:C13	2.27	0.54
1:A:95:ASN:HB3	1:A:98:THR:OG1	2.08	0.54
1:B:52:PRO:HD2	1:B:188:ARG:HD2	1.89	0.53
1:A:154:TYR:O	1:A:155:ASP:HB2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:VAL:HG13	1:A:179:GLY:HA2	1.90	0.53
1:B:140:PHE:HB3	1:B:144:SER:OG	2.08	0.53
1:B:108:PRO:HB3	1:B:132:PRO:HA	1.92	0.52
1:B:245:ASP:O	1:B:249:ILE:HG12	2.09	0.52
1:B:228:ASN:OD1	1:B:228:ASN:N	2.42	0.51
1:A:40:ARG:HD3	1:A:85:CYS:HA	1.93	0.49
1:B:5:LYS:HA	1:B:291:PHE:CZ	2.48	0.49
1:A:108:PRO:HB3	1:A:132:PRO:HA	1.95	0.48
1:B:52:PRO:HD2	1:B:188:ARG:CG	2.43	0.48
1:A:163:HIS:CE1	1:A:172:HIS:HB3	2.49	0.48
1:A:140:PHE:CD2	1:A:172:HIS:CG	3.01	0.48
1:A:288:GLU:HG2	1:A:291:PHE:CE2	2.49	0.48
1:A:165:MET:HE2	1:A:165:MET:HB3	1.75	0.47
1:B:101:TYR:C	1:B:102:LYS:HD2	2.40	0.47
1:A:164:HIS:O	2:A:301:FNO:C16	2.62	0.46
1:B:52:PRO:HD2	1:B:188:ARG:CD	2.45	0.46
1:A:140:PHE:HD2	1:A:172:HIS:NE2	2.14	0.45
1:B:154:TYR:O	1:B:155:ASP:HB2	2.16	0.44
1:B:58:LEU:HD13	1:B:58:LEU:O	2.18	0.44
1:A:165:MET:HE1	2:A:301:FNO:C12	2.45	0.43
1:A:188:ARG:HG2	1:A:190:THR:HG23	2.00	0.43
1:B:190:THR:HG22	1:B:191:ALA:H	1.82	0.42
1:A:222:ARG:HG2	1:A:222:ARG:HH11	1.85	0.42
1:B:57:LEU:O	1:B:61:LYS:HG2	2.20	0.42
1:B:86:VAL:HG13	1:B:179:GLY:HA2	2.02	0.42
1:B:114:VAL:HG11	1:B:140:PHE:HZ	1.86	0.41
1:B:207:TRP:CZ3	1:B:287:LEU:HA	2.56	0.41
1:A:288:GLU:HG2	1:A:291:PHE:HE2	1.85	0.41
1:B:114:VAL:HG11	1:B:140:PHE:CZ	2.56	0.41
1:B:219:PHE:CD2	1:B:281:ILE:HD13	2.55	0.41
1:A:276:MET:HE3	1:A:281:ILE:HG13	2.02	0.41
1:B:113:SER:O	1:B:149:GLY:HA2	2.21	0.41
1:B:165:MET:HE2	1:B:173:ALA:HB3	2.03	0.40
1:B:227:LEU:HD12	1:B:227:LEU:HA	1.87	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	293/296 (99%)	286 (98%)	7 (2%)	0	100	100
1	B	287/296 (97%)	281 (98%)	6 (2%)	0	100	100
All	All	580/592 (98%)	567 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	240/254 (94%)	233 (97%)	7 (3%)	37	40
1	B	229/254 (90%)	222 (97%)	7 (3%)	35	38
All	All	469/508 (92%)	455 (97%)	14 (3%)	36	39

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	THR
1	A	141	LEU
1	A	145	CYS
1	A	165	MET
1	A	196	THR
1	A	252	PRO
1	A	282	LEU
1	B	141	LEU

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Mol	Chain	Res	Type
1	B	145	CYS
1	B	216	ASP
1	B	228	ASN
1	B	244	GLN
1	B	276	MET
1	B	281	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	B	244	GLN
1	B	256	GLN
1	B	277	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](#)

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
2	FNO	A	301	1	21,22,22	2.09	7 (33%)	29,32,32	2.16	8 (27%)

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	FNO	A	301	1	1/1/5/7	2/9/25/25	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	FNO	O02-C10	-4.07	1.14	1.23
2	A	301	FNO	C11-C09	-3.73	1.32	1.41
2	A	301	FNO	O04-C13	-3.63	1.16	1.24
2	A	301	FNO	C07-C10	-3.62	1.37	1.46
2	A	301	FNO	O01-C06	-3.03	1.36	1.42
2	A	301	FNO	C12-C13	-2.42	1.39	1.44
2	A	301	FNO	C09-C10	-2.04	1.41	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	FNO	O02-C10-C07	-5.66	116.94	122.52
2	A	301	FNO	C09-C10-C07	5.29	123.87	116.59
2	A	301	FNO	C11-C13-C12	3.32	122.55	116.31
2	A	301	FNO	C21-C17-C20	3.22	122.00	114.59
2	A	301	FNO	C18-C15-C09	3.17	124.19	120.15
2	A	301	FNO	C19-C18-C15	-2.58	117.92	120.50
2	A	301	FNO	O04-C13-C11	-2.14	118.61	122.10
2	A	301	FNO	C18-C19-C16	-2.08	118.42	120.50

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	301	FNO	C06

All (2) torsion outliers are listed below:

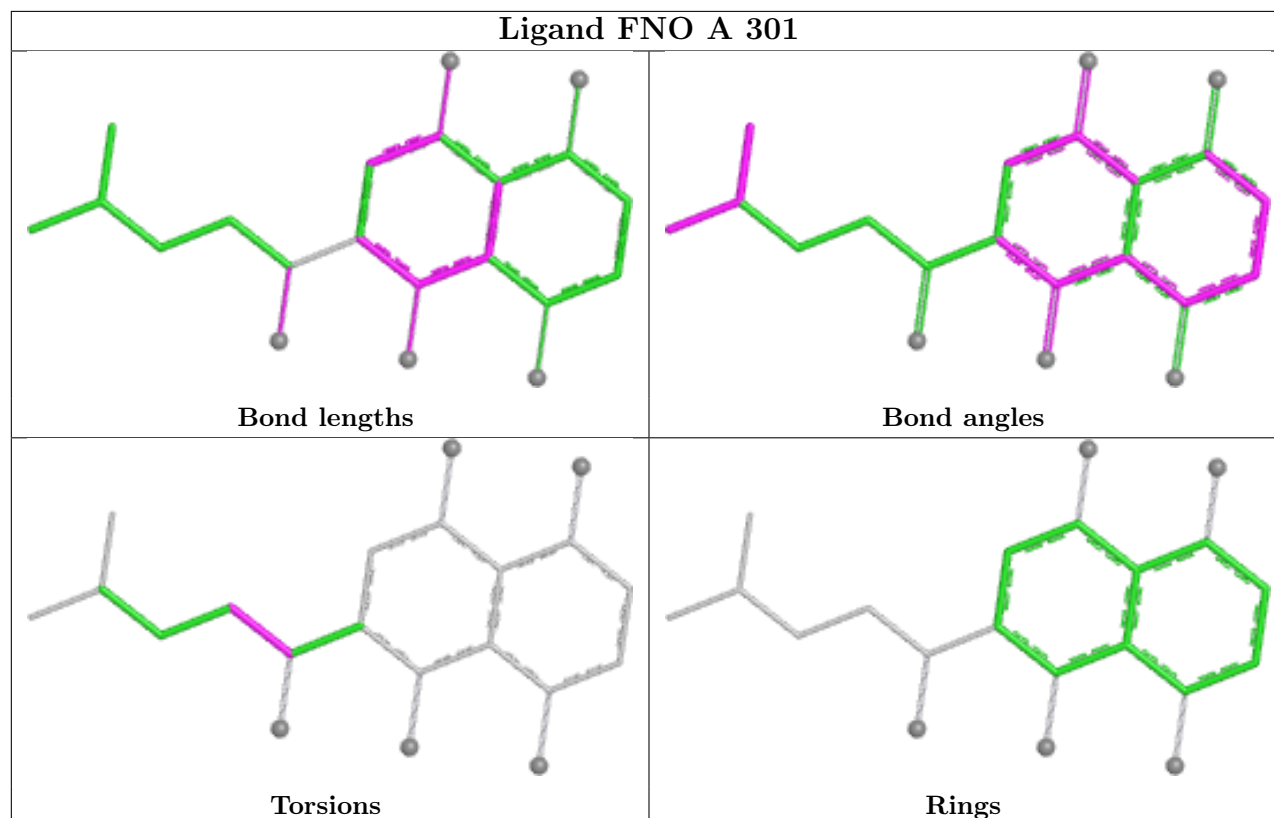
Mol	Chain	Res	Type	Atoms
2	A	301	FNO	O01-C06-C08-C14
2	A	301	FNO	C07-C06-C08-C14

There are no ring outliers.

1 monomer is involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	FNO	10	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	295/296 (99%)	0.71	40 (13%) 7 7	26, 38, 57, 68	0
1	B	291/296 (98%)	1.12	63 (21%) 2 2	27, 47, 73, 79	0
All	All	586/592 (98%)	0.92	103 (17%) 4 4	26, 42, 66, 79	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	140	PHE	6.7
1	A	142	ALA	6.1
1	A	282	LEU	4.9
1	B	140	PHE	4.2
1	A	277	ASN	4.1
1	B	154	TYR	4.1
1	B	45	THR	3.9
1	A	297	VAL	3.9
1	B	142	ALA	3.9
1	B	58	LEU	3.9
1	B	57	LEU	3.7
1	B	298	ARG	3.7
1	B	282	LEU	3.6
1	A	128	CYS	3.6
1	A	47	GLU	3.6
1	A	141	LEU	3.6
1	B	155	ASP	3.5
1	B	55	GLU	3.5
1	A	215	GLY	3.4
1	A	286	LEU	3.4
1	A	154	TYR	3.3
1	B	141	LEU	3.3
1	B	59	ILE	3.3
1	B	279	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	62	SER	3.3
1	B	128	CYS	3.2
1	B	263	ASP	3.2
1	A	195	GLY	3.1
1	B	232	LEU	3.1
1	B	286	LEU	3.1
1	B	169	THR	3.1
1	B	52	PRO	3.1
1	B	53	ASN	3.1
1	B	67	LEU	3.0
1	A	41	HIS	3.0
1	B	54	TYR	2.9
1	B	74	GLN	2.9
1	A	298	ARG	2.9
1	B	139	SER	2.9
1	B	191	ALA	2.9
1	A	172	HIS	2.9
1	A	222	ARG	2.8
1	A	291	PHE	2.8
1	B	26	THR	2.8
1	A	143	GLY	2.8
1	B	283	GLY	2.8
1	B	294	PHE	2.8
1	A	211	ALA	2.7
1	B	44	CYS	2.7
1	B	153	ASP	2.7
1	B	281	ILE	2.7
1	A	283	GLY	2.7
1	B	60	ARG	2.7
1	B	3	PHE	2.6
1	B	118	TYR	2.6
1	B	235	MET	2.6
1	A	216	ASP	2.6
1	A	276	MET	2.5
1	A	48	ASN	2.5
1	B	65	ASN	2.5
1	A	92	ASP	2.5
1	A	72	ASN	2.5
1	B	188	ARG	2.5
1	A	213	ILE	2.5
1	B	285	ALA	2.4
1	B	56	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	279	ARG	2.4
1	A	284	SER	2.4
1	B	224	THR	2.4
1	A	210	ALA	2.4
1	B	227	LEU	2.4
1	B	43	ILE	2.4
1	A	196	THR	2.4
1	B	24	THR	2.4
1	B	258	GLY	2.3
1	A	281	ILE	2.3
1	A	139	SER	2.3
1	B	78	ILE	2.3
1	B	23	GLY	2.3
1	B	189	GLN	2.3
1	A	285	ALA	2.3
1	A	296	VAL	2.2
1	A	138	GLY	2.2
1	B	207	TRP	2.2
1	B	280	THR	2.2
1	B	102	LYS	2.2
1	B	51	ASN	2.2
1	B	238	ASN	2.2
1	A	45	THR	2.1
1	B	190	THR	2.1
1	B	218	TRP	2.1
1	B	41	HIS	2.1
1	B	278	GLY	2.1
1	B	233	VAL	2.1
1	A	191	ALA	2.1
1	B	272	LEU	2.1
1	A	118	TYR	2.1
1	B	215	GLY	2.1
1	A	4	ARG	2.0
1	B	77	VAL	2.0
1	B	270	GLU	2.0
1	A	235	MET	2.0
1	B	64	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [i](#)

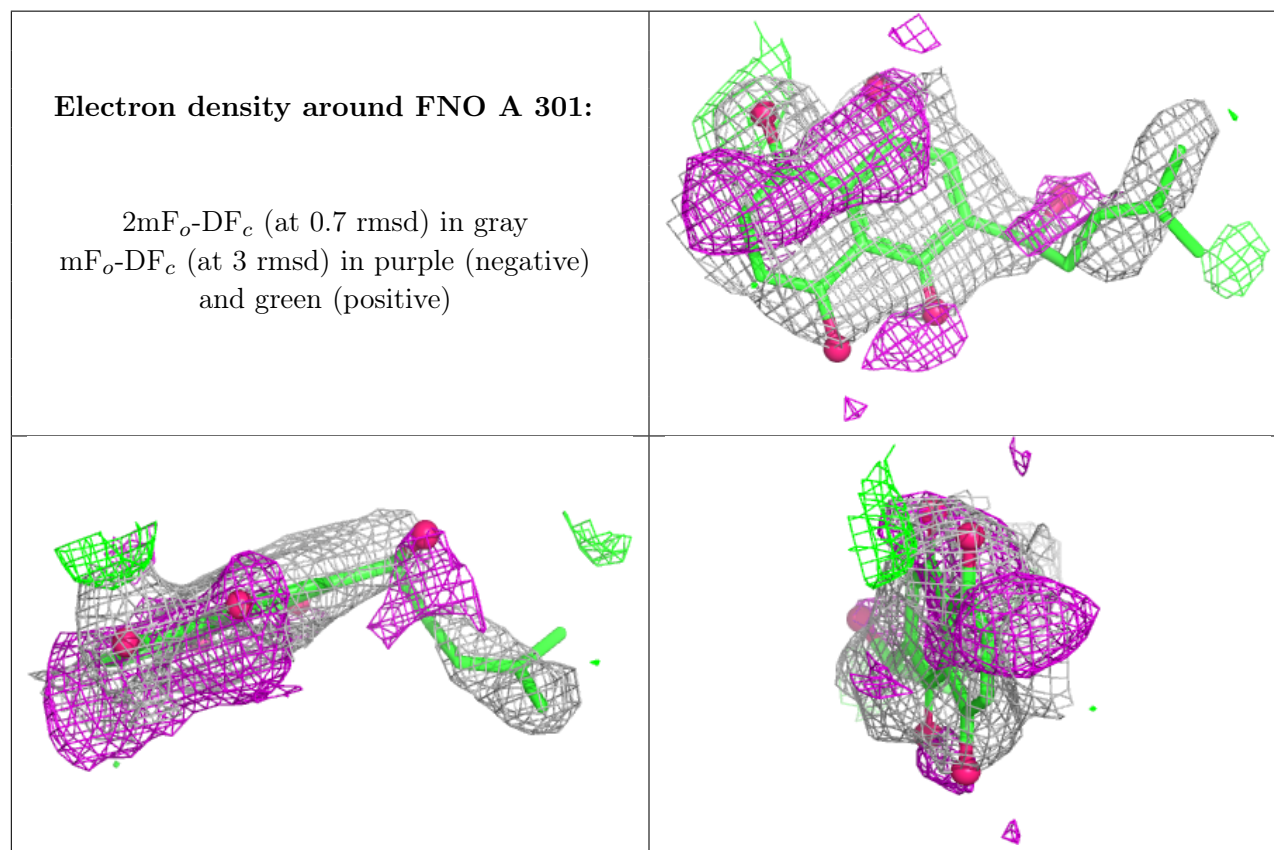
There are no oligosaccharides in this entry.

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	FNO	A	301	21/21	0.58	0.20	49,59,65,68	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.