



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:23 AM UTC

PDB ID : 8DZ0 / pdb_00008dz0
Title : Crystal Structure of SARS-CoV-2 Main protease in complex with Ensitrelvir
Authors : Noske, G.D.; Oliva, G.; Godoy, A.S.
Deposited on : 2022-08-06
Resolution : 2.29 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

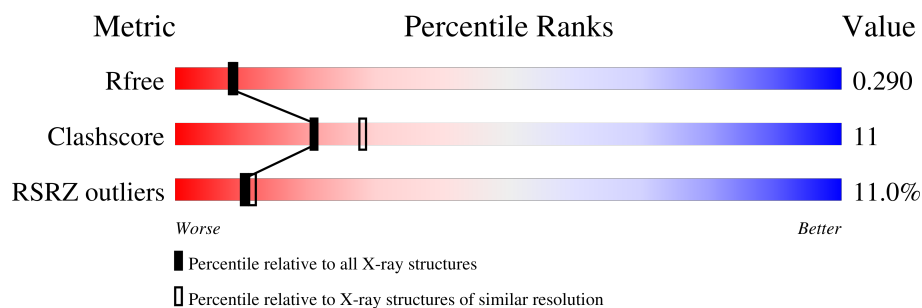
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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)
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 $\leq 5\%$. 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	306	9% 75% 24%
1	B	306	13% 75% 25%

2 Entry composition [i](#)

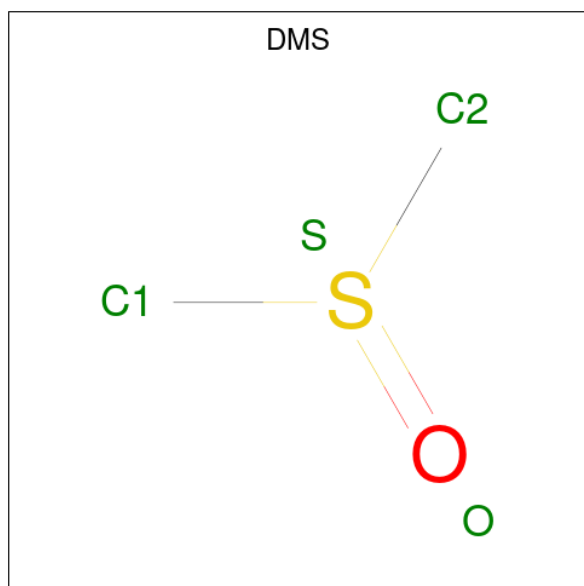
There are 4 unique types of molecules in this entry. The entry contains 4848 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	302	Total	C	N	O	S	0	1	0
			2338	1480	397	438	23			
1	B	306	Total	C	N	O	S	4	3	0
			2380	1511	402	444	23			

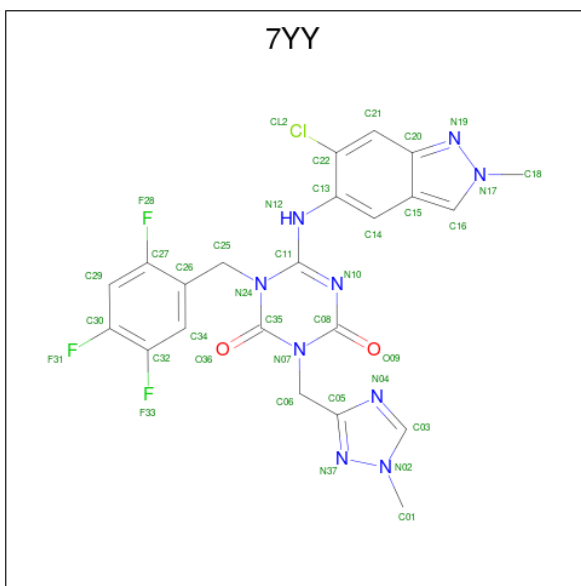
- Molecule 2 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



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

- Molecule 3 is 6-[(6-chloranyl-2-methyl-indazol-5-yl)amino]-3-[(1-methyl-1,2,4-triazol-3-yl)methyl]-1-[[2,4,5-tris(fluoranyl)phenyl)methyl]-1,3,5-triazine-2,4-dione (CCD ID: 7YY)

(formula: $C_{22}H_{17}ClF_3N_9O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	0	0
			37	22	1	3	9	2		
3	B	1	Total	C	Cl	F	N	O	0	0
			37	22	1	3	9	2		

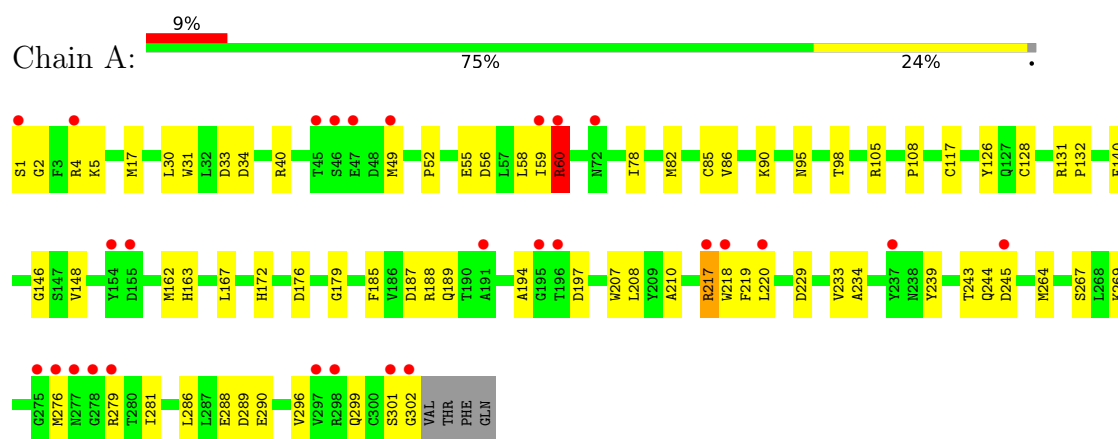
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	12	Total	O	0	0
			12	12		

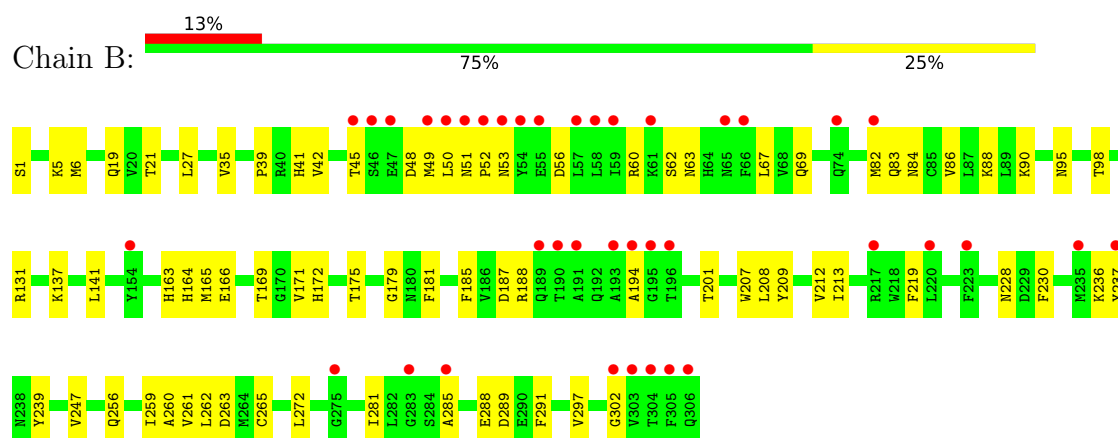
3 Residue-property plots [i](#)

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: 3C-like proteinase nsp5



- Molecule 1: 3C-like proteinase nsp5



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.91Å 99.02Å 103.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.42 – 2.29 71.42 – 2.29	Depositor EDS
% Data completeness (in resolution range)	47.8 (71.42-2.29) 48.0 (71.42-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.219 , 0.275 (Not available) , 0.290	Depositor DCC
R_{free} test set	810 reflections (2.52%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.039 for -h,l,k	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4848	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 4.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 7YY, DMS

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	0.31	1/2393 (0.0%)	0.45	2/3251 (0.1%)
1	B	0.15	0/2442	0.38	0/3319
All	All	0.24	1/4835 (0.0%)	0.42	2/6570 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	217	ARG	CB-CG	-6.93	1.31	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ARG	CB-CG-CD	-5.56	98.52	111.30
1	A	217	ARG	CB-CG-CD	-5.29	99.13	111.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	60	ARG	Sidechain

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	2338	0	2290	52	0
1	B	2380	0	2341	59	0
2	A	8	0	12	0	0
2	B	4	0	6	1	0
3	A	37	0	0	0	0
3	B	37	0	0	1	0
4	A	32	0	0	2	0
4	B	12	0	0	2	0
All	All	4848	0	4649	105	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 11.

All (105) 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:4:ARG:NH1	1:B:137:LYS:O	2.14	0.79
1:A:40:ARG:HD3	1:A:85:CYS:HA	1.62	0.79
1:A:58:LEU:C	1:A:60:ARG:H	1.93	0.75
1:B:49:MET:O	1:B:52:PRO:HD3	1.88	0.74
1:A:301:SER:OG	1:A:302:GLY:N	2.18	0.72
1:B:247:VAL:HG13	1:B:261:VAL:HG11	1.74	0.70
1:A:217:ARG:HD3	1:A:217:ARG:O	1.92	0.69
1:A:52:PRO:HD2	1:A:188:ARG:HG2	1.74	0.68
1:B:164:HIS:CD2	1:B:175:THR:HG23	2.29	0.67
1:A:49:MET:HB3	1:A:189:GLN:HG3	1.76	0.67
1:A:58:LEU:C	1:A:60:ARG:N	2.54	0.65
1:B:50:LEU:O	1:B:188:ARG:NE	2.29	0.65
1:B:56:ASP:OD1	1:B:60:ARG:NH2	2.31	0.64
1:B:137:LYS:NZ	4:B:502:HOH:O	2.30	0.64
1:B:288:GLU:HG2	1:B:291:PHE:HE1	1.65	0.62
1:B:259:ILE:HG22	1:B:260:ALA:H	1.66	0.60
1:B:163:HIS:HE1	1:B:172:HIS:HB3	1.64	0.60
1:A:140:PHE:O	1:B:1:SER:N	2.32	0.57
1:B:256:GLN:OE1	4:B:501:HOH:O	2.17	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:LEU:HG	1:B:285:ALA:HB2	1.85	0.57
1:B:131:ARG:NH2	1:B:289:ASP:OD2	2.29	0.57
1:A:17:MET:HG3	1:A:117:CYS:SG	2.45	0.56
1:A:30:LEU:HD22	1:A:148:VAL:HG11	1.88	0.55
1:A:1:SER:HB3	1:B:166:GLU:OE2	2.05	0.55
1:B:41:HIS:O	1:B:41:HIS:ND1	2.40	0.55
1:B:86[A]:VAL:HG13	1:B:179:GLY:HA2	1.88	0.55
1:A:167:LEU:HD21	1:A:185:PHE:CE2	2.41	0.55
1:A:56:ASP:O	1:A:60:ARG:HB2	2.07	0.55
1:B:175:THR:HG22	1:B:181:PHE:HA	1.88	0.54
1:B:288:GLU:HG2	1:B:291:PHE:CE1	2.43	0.54
1:B:19:GLN:HB3	1:B:69:GLN:HB2	1.89	0.53
1:A:55:GLU:O	1:A:59:ILE:HG22	2.08	0.53
1:B:21:THR:HB	1:B:67:LEU:HD12	1.89	0.53
1:A:1:SER:OG	1:A:2:GLY:N	2.42	0.53
1:A:210:ALA:HB2	1:A:296:VAL:HG13	1.90	0.52
1:B:185:PHE:CZ	1:B:194:ALA:HB2	2.45	0.52
1:A:131:ARG:HH22	1:A:289:ASP:CG	2.18	0.52
1:B:163:HIS:CE1	1:B:172:HIS:HB3	2.45	0.52
1:B:39:PRO:HB3	1:B:164:HIS:CE1	2.45	0.51
1:B:45:THR:HG23	1:B:48:ASP:H	1.76	0.51
1:A:58:LEU:HD22	1:A:82:MET:HB2	1.93	0.51
1:A:301:SER:O	4:A:501:HOH:O	2.19	0.50
1:A:131:ARG:HD3	1:A:197:ASP:OD2	2.11	0.50
1:A:299:GLN:O	1:B:141:LEU:HD21	2.11	0.50
1:A:163:HIS:CE1	1:A:172:HIS:HB3	2.47	0.50
1:A:217:ARG:HG2	1:A:220:LEU:HD13	1.93	0.49
1:A:31:TRP:CD2	1:A:95:ASN:HB2	2.48	0.49
1:A:187:ASP:N	1:A:187:ASP:OD1	2.40	0.49
1:B:236:LYS:C	1:B:236:LYS:HD3	2.37	0.49
1:A:243:THR:HG22	1:A:245:ASP:H	1.77	0.49
1:B:27:LEU:HD21	1:B:42:VAL:HB	1.94	0.49
1:B:239:TYR:CZ	1:B:272:LEU:HD21	2.48	0.48
1:A:59:ILE:HG13	1:A:59:ILE:O	2.14	0.48
1:A:207:TRP:CD2	1:A:288:GLU:HB2	2.48	0.48
1:B:169:THR:HG23	1:B:171:VAL:HG22	1.96	0.48
1:A:218:TRP:HH2	1:A:276:MET:HA	1.79	0.47
1:A:229:ASP:O	1:A:233:VAL:HG23	2.14	0.47
1:B:95:ASN:HB3	1:B:98:THR:OG1	2.14	0.47
1:B:228:ASN:OD1	1:B:228:ASN:N	2.46	0.47
1:A:78:ILE:HG13	1:A:90:LYS:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:LEU:O	1:B:212:VAL:HG23	2.14	0.47
1:B:230:PHE:CD2	1:B:265:CYS:HB3	2.50	0.46
1:B:83:GLN:HB2	1:B:88:LYS:HE2	1.96	0.46
1:A:146:GLY:HA2	1:A:162:MET:HE2	1.98	0.46
1:A:194:ALA:O	4:A:502:HOH:O	2.21	0.46
1:B:49:MET:O	1:B:51:ASN:N	2.49	0.46
1:A:219:PHE:O	1:A:267:SER:OG	2.33	0.45
1:A:234:ALA:O	1:A:239:TYR:HB2	2.17	0.45
1:A:86:VAL:HG13	1:A:179:GLY:HA2	1.99	0.45
1:A:208:LEU:HB3	1:A:264:MET:HE2	1.99	0.44
1:B:207:TRP:CH2	1:B:281:ILE:HB	2.53	0.44
1:B:236:LYS:HD3	1:B:236:LYS:O	2.17	0.44
1:B:261:VAL:O	1:B:262:LEU:HB2	2.18	0.44
1:A:233:VAL:HG21	1:A:269:LYS:HD2	1.99	0.44
1:B:56:ASP:O	1:B:60:ARG:HG2	2.17	0.44
1:B:187:ASP:OD1	1:B:187:ASP:N	2.46	0.44
1:B:82:MET:HE3	1:B:84:ASN:O	2.18	0.43
1:B:60:ARG:HG2	1:B:60:ARG:H	1.69	0.43
1:B:51:ASN:OD1	1:B:188:ARG:NH1	2.51	0.43
1:A:279:ARG:HG2	1:A:279:ARG:HH11	1.84	0.43
1:A:5:LYS:NZ	1:A:290:GLU:HB2	2.34	0.43
1:A:126:TYR:HD2	1:B:6:MET:HG3	1.82	0.43
1:B:260:ALA:O	1:B:263:ASP:HB3	2.19	0.43
1:A:217:ARG:HG2	1:A:220:LEU:CD1	2.49	0.43
1:B:62:SER:OG	1:B:63:ASN:N	2.51	0.43
1:A:108:PRO:HB3	1:A:132:PRO:HA	2.02	0.42
1:B:5:LYS:NZ	1:B:288:GLU:OE2	2.51	0.42
1:B:209:TYR:O	1:B:213:ILE:HG13	2.20	0.42
1:A:105:ARG:HD2	1:A:176:ASP:OD2	2.19	0.42
1:A:55:GLU:H	1:A:55:GLU:CD	2.28	0.42
1:B:83:GLN:O	1:B:84:ASN:HB2	2.20	0.42
1:B:185:PHE:HZ	1:B:194:ALA:HB2	1.84	0.42
1:B:219:PHE:CD2	1:B:281:ILE:HD13	2.54	0.41
1:B:297:VAL:O	1:B:302:GLY:N	2.54	0.41
1:A:95:ASN:HB3	1:A:98:THR:OG1	2.20	0.41
1:B:272:LEU:HD23	1:B:272:LEU:HA	1.86	0.41
1:A:207:TRP:CH2	1:A:281:ILE:HB	2.55	0.41
1:B:35:VAL:HG22	1:B:90:LYS:HG3	2.01	0.41
1:A:33:ASP:HB3	1:A:34:ASP:H	1.77	0.41
1:B:165:MET:HA	3:B:402:7YY:O36	2.21	0.41
1:B:53:ASN:HA	2:B:401:DMS:H13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:THR:HG22	1:A:244:GLN:N	2.37	0.40
1:B:201:THR:HG23	1:B:239:TYR:HD2	1.87	0.40
1:A:126:TYR:HE1	1:A:128:CYS:SG	2.44	0.40
1:B:236:LYS:HB3	1:B:237:TYR:CD1	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

5 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	DMS	B	401	-	3,3,3	0.68	0	3,3,3	0.52	0
2	DMS	A	402	-	3,3,3	0.68	0	3,3,3	0.56	0
3	7YY	B	402	-	39,41,41	2.71	15 (38%)	48,61,61	3.70	18 (37%)
3	7YY	A	403	-	39,41,41	2.71	15 (38%)	48,61,61	3.80	18 (37%)
2	DMS	A	401	-	3,3,3	0.68	0	3,3,3	0.56	0

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	7YY	B	402	-	-	3/12/12/12	0/5/5/5
3	7YY	A	403	-	-	1/12/12/12	0/5/5/5

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	403	7YY	C11-N12	7.39	1.46	1.36
3	B	402	7YY	C11-N12	7.23	1.46	1.36
3	B	402	7YY	C06-C05	6.54	1.53	1.49
3	B	402	7YY	C11-N24	6.37	1.45	1.39
3	A	403	7YY	C11-N24	6.34	1.45	1.39
3	A	403	7YY	C06-C05	6.30	1.53	1.49
3	A	403	7YY	C11-N10	5.25	1.44	1.32
3	B	402	7YY	C11-N10	5.20	1.44	1.32
3	A	403	7YY	C08-N10	4.70	1.45	1.36
3	B	402	7YY	C35-N07	4.49	1.46	1.38
3	B	402	7YY	C08-N10	4.44	1.45	1.36
3	A	403	7YY	C35-N07	4.34	1.46	1.38
3	B	402	7YY	C35-N24	4.08	1.45	1.38
3	A	403	7YY	C35-N24	3.81	1.45	1.38
3	A	403	7YY	C20-C15	-3.53	1.36	1.43
3	B	402	7YY	C20-C15	-3.46	1.36	1.43
3	B	402	7YY	O09-C08	-2.66	1.18	1.23
3	A	403	7YY	O09-C08	-2.64	1.18	1.23
3	A	403	7YY	C08-N07	2.45	1.46	1.41
3	A	403	7YY	C13-N12	2.39	1.46	1.41
3	B	402	7YY	C08-N07	2.39	1.46	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	7YY	O36-C35	-2.38	1.18	1.22
3	A	403	7YY	O36-C35	-2.38	1.18	1.22
3	B	402	7YY	C22-CL2	2.30	1.79	1.73
3	A	403	7YY	C22-CL2	2.29	1.79	1.73
3	A	403	7YY	N17-N19	2.29	1.39	1.36
3	B	402	7YY	N17-N19	2.25	1.39	1.36
3	B	402	7YY	C13-N12	2.24	1.46	1.41
3	A	403	7YY	C05-N04	-2.16	1.33	1.37
3	B	402	7YY	C05-N04	-2.14	1.33	1.37

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	403	7YY	C13-C22-CL2	17.33	138.31	119.52
3	B	402	7YY	C13-C22-CL2	16.74	137.68	119.52
3	A	403	7YY	C21-C22-CL2	-10.49	101.28	118.45
3	B	402	7YY	C21-C22-CL2	-10.07	101.95	118.45
3	A	403	7YY	N24-C11-N10	-6.51	120.29	124.07
3	A	403	7YY	C01-N02-N37	6.37	126.05	119.57
3	B	402	7YY	C01-N02-N37	6.23	125.91	119.57
3	B	402	7YY	N24-C11-N10	-6.19	120.47	124.07
3	A	403	7YY	C03-N04-C05	5.23	107.83	102.88
3	A	403	7YY	N37-C05-N04	-5.17	108.39	115.26
3	B	402	7YY	C03-N04-C05	5.16	107.76	102.88
3	B	402	7YY	N37-C05-N04	-5.10	108.48	115.26
3	A	403	7YY	C16-N17-N19	-4.59	108.16	112.71
3	B	402	7YY	C16-N17-N19	-4.47	108.28	112.71
3	A	403	7YY	N24-C35-N07	3.57	120.10	115.75
3	A	403	7YY	C20-N19-N17	3.53	107.92	104.26
3	B	402	7YY	N24-C35-N07	3.49	120.01	115.75
3	B	402	7YY	C20-N19-N17	3.47	107.85	104.26
3	B	402	7YY	C06-N07-C35	3.45	120.87	116.71
3	B	402	7YY	C34-C26-C27	3.37	120.25	116.54
3	A	403	7YY	C18-N17-C16	3.36	131.31	127.94
3	A	403	7YY	C34-C26-C27	3.35	120.23	116.54
3	B	402	7YY	C26-C25-N24	-3.19	109.28	113.14
3	A	403	7YY	C26-C25-N24	-3.16	109.32	113.14
3	B	402	7YY	C18-N17-C16	2.95	130.90	127.94
3	B	402	7YY	C25-N24-C35	2.81	120.85	116.92
3	A	403	7YY	N02-C03-N04	-2.69	108.09	110.71
3	B	402	7YY	N02-C03-N04	-2.67	108.11	110.71
3	A	403	7YY	C06-N07-C35	2.53	119.77	116.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	403	7YY	C29-C27-C26	-2.50	119.91	123.76
3	B	402	7YY	C29-C27-C26	-2.42	120.03	123.76
3	A	403	7YY	C25-N24-C35	2.36	120.22	116.92
3	A	403	7YY	C01-N02-C03	-2.35	125.90	129.74
3	A	403	7YY	C06-C05-N04	2.32	125.19	122.77
3	B	402	7YY	C01-N02-C03	-2.28	126.01	129.74
3	B	402	7YY	C14-C13-C22	2.14	120.39	118.21

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	7YY	C22-C13-N12-C11
3	B	402	7YY	C22-C13-N12-C11
3	B	402	7YY	N04-C05-C06-N07
3	B	402	7YY	N37-C05-C06-N07

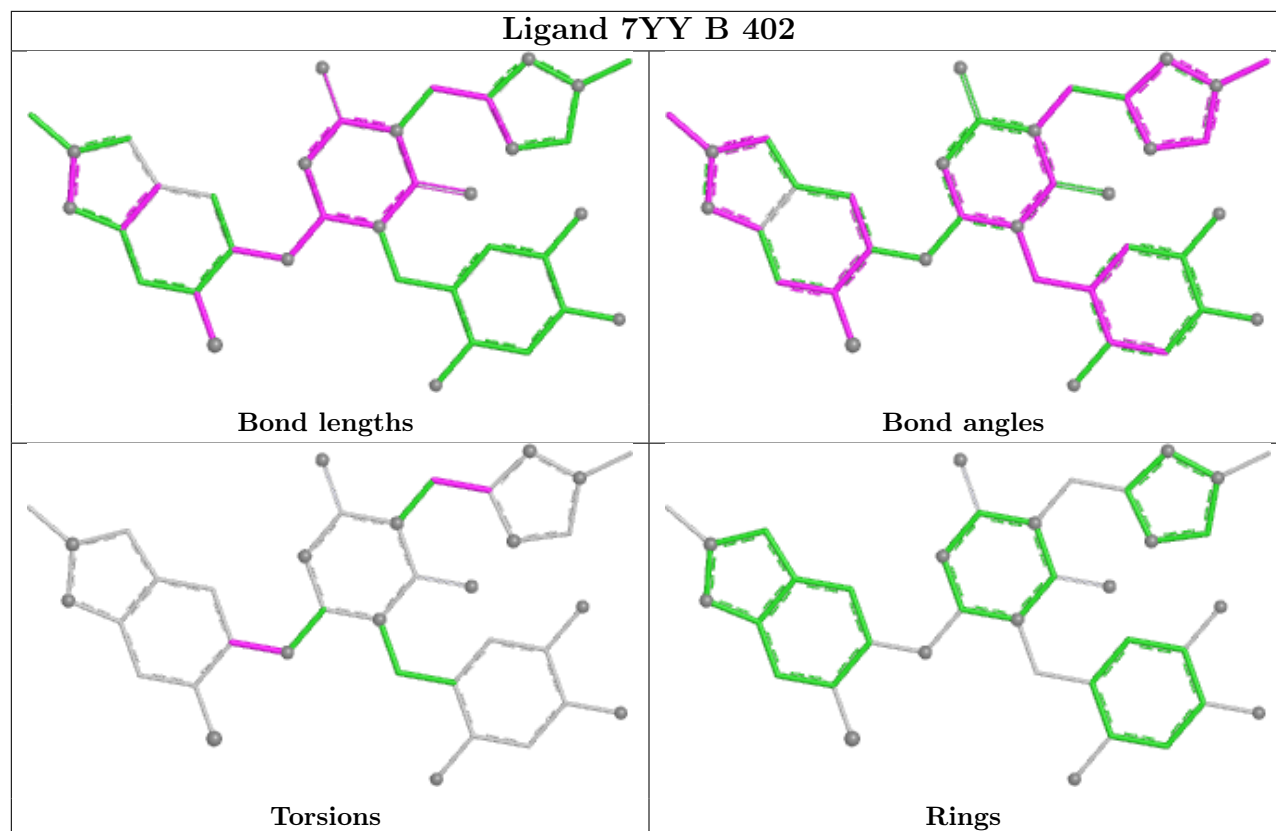
There are no ring outliers.

2 monomers are involved in 2 short contacts:

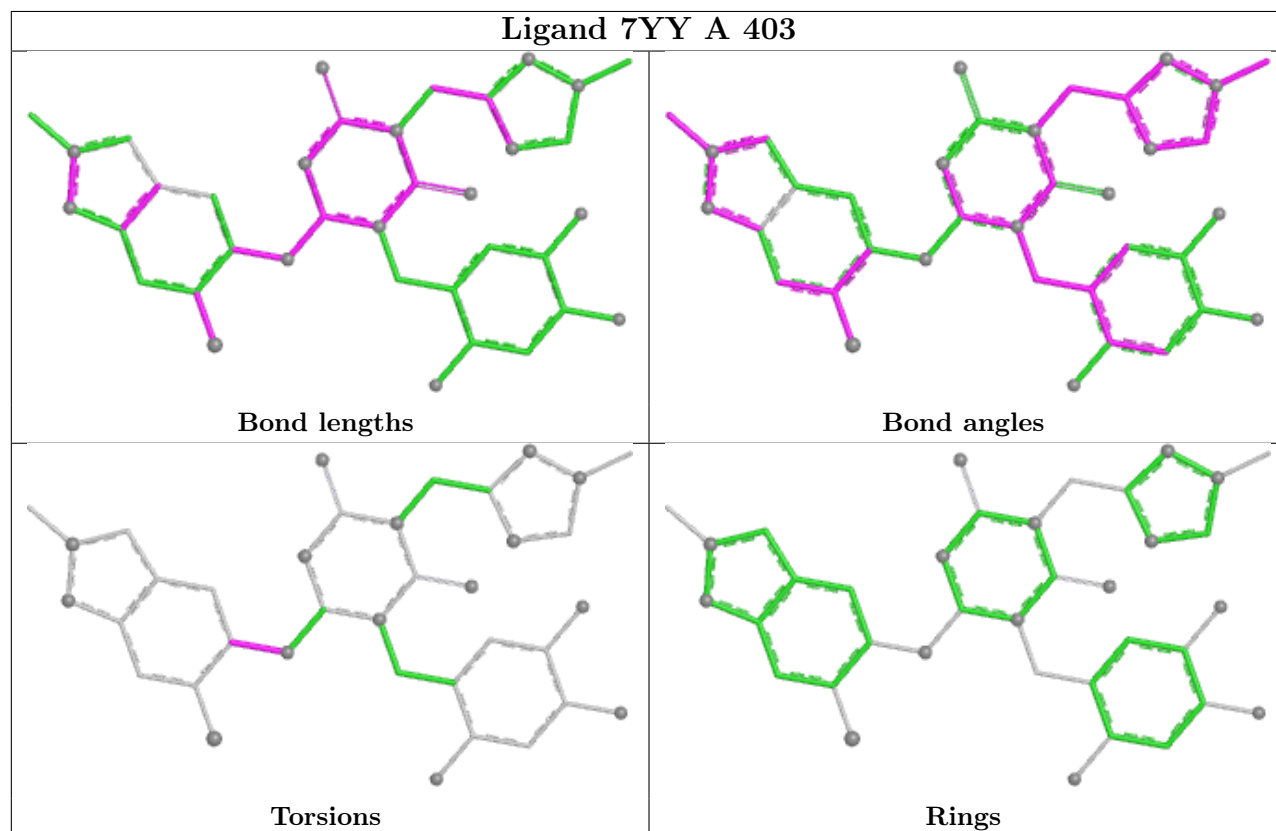
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	DMS	1	0
3	B	402	7YY	1	0

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.

Ligand 7YY B 402



Ligand 7YY A 403



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	302/306 (98%)	0.51	28 (9%) 14 16	18, 35, 69, 115	1 (0%)
1	B	306/306 (100%)	0.84	39 (12%) 8 8	21, 44, 90, 149	3 (0%)
All	All	608/612 (99%)	0.68	67 (11%) 10 11	18, 38, 80, 149	4 (0%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	50	LEU	8.1
1	A	60	ARG	5.9
1	A	218	TRP	5.8
1	B	305	PHE	5.5
1	A	217	ARG	5.3
1	A	49	MET	5.2
1	B	154	TYR	5.1
1	B	303	VAL	5.0
1	B	302	GLY	4.9
1	B	59	ILE	4.9
1	B	306	GLN	4.8
1	B	304	THR	4.7
1	A	46	SER	4.7
1	B	58	LEU	4.4
1	A	191	ALA	3.9
1	A	301	SER	3.9
1	B	235[A]	MET	3.9
1	B	190	THR	3.7
1	B	46	SER	3.7
1	B	49	MET	3.5
1	A	277	ASN	3.3
1	A	302	GLY	3.2
1	B	47	GLU	3.1
1	B	191	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	196	THR	2.9
1	B	51	ASN	2.9
1	B	283	GLY	2.9
1	A	298	ARG	2.9
1	A	278	GLY	2.8
1	B	61	LYS	2.8
1	A	47	GLU	2.8
1	A	275	GLY	2.7
1	B	74	GLN	2.7
1	A	196	THR	2.7
1	A	59	ILE	2.7
1	B	217	ARG	2.7
1	B	195	GLY	2.6
1	B	275	GLY	2.6
1	A	220	LEU	2.6
1	B	66	PHE	2.6
1	B	285	ALA	2.5
1	B	57	LEU	2.5
1	B	189	GLN	2.5
1	B	55	GLU	2.5
1	B	193	ALA	2.5
1	B	82	MET	2.4
1	A	245	ASP	2.4
1	B	220	LEU	2.4
1	A	155	ASP	2.3
1	A	154	TYR	2.3
1	A	279	ARG	2.3
1	A	276	MET	2.3
1	B	194	ALA	2.3
1	B	53	ASN	2.2
1	B	65	ASN	2.2
1	A	1	SER	2.2
1	A	297	VAL	2.2
1	A	72	ASN	2.2
1	B	237	TYR	2.2
1	A	45	THR	2.1
1	B	223	PHE	2.1
1	A	195	GLY	2.1
1	B	54	TYR	2.1
1	A	4	ARG	2.1
1	A	237	TYR	2.0
1	B	45	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	52	PRO	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](#)

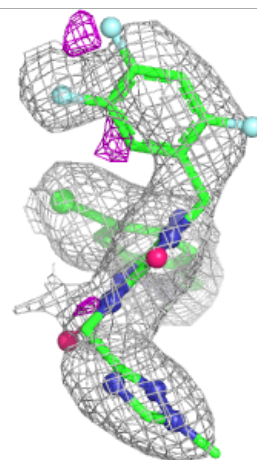
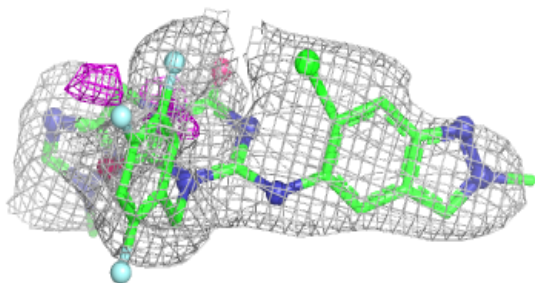
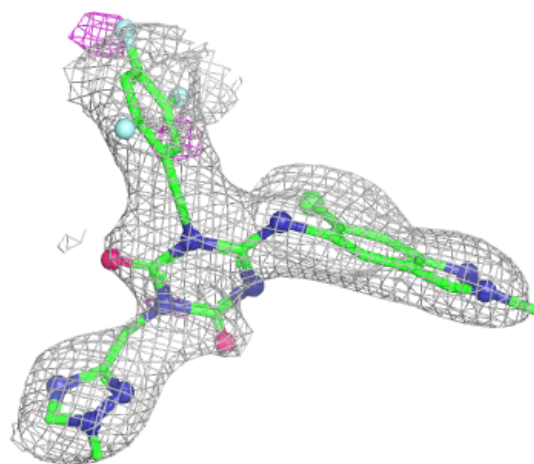
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	DMS	B	401	4/4	0.70	0.26	76,93,96,102	0
2	DMS	A	401	4/4	0.76	0.22	47,49,49,70	0
2	DMS	A	402	4/4	0.83	0.22	47,50,55,60	0
3	7YY	B	402	37/37	0.83	0.15	54,72,83,88	0
3	7YY	A	403	37/37	0.89	0.11	31,48,66,70	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.

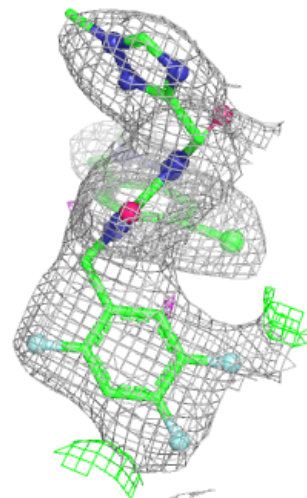
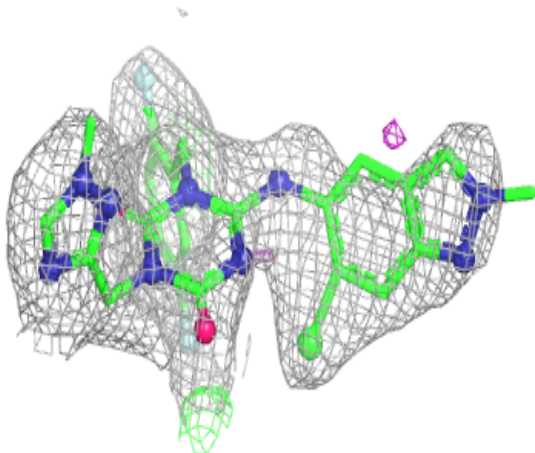
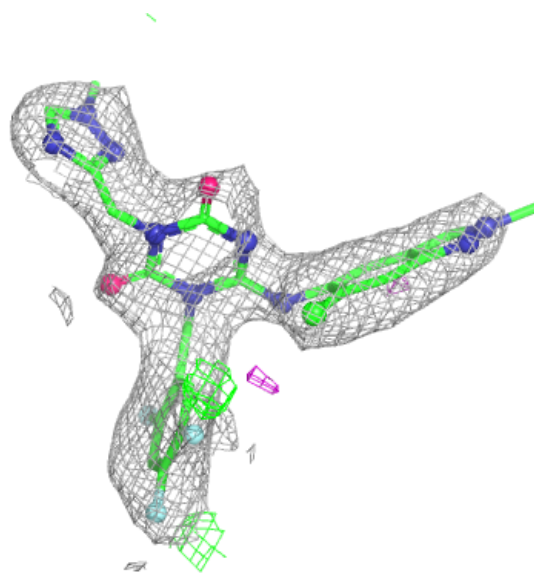
Electron density around 7YY B 402:

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



Electron density around 7YY A 403:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.