



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:24 PM UTC

PDB ID : 7FTZ / pdb_00007ftz
Title : Crystal Structure of human cyclic GMP-AMP synthase in complex with 2-[2-[3-[[[(5-bromo-2-hydroxybenzoyl)amino]methyl]anilino]-1,3-thiazol-4-yl]acetic acid
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Deposited on : 2023-02-08
Resolution : 2.50 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

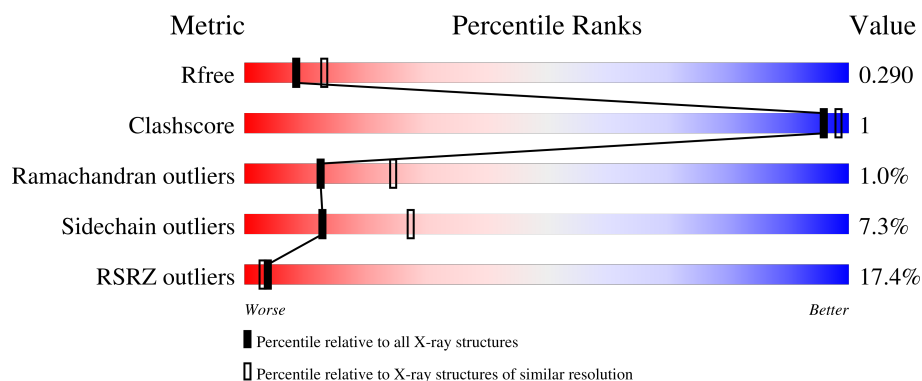
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	362	17% 83% 13% ..
1	B	362	17% 83% 13% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5912 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.

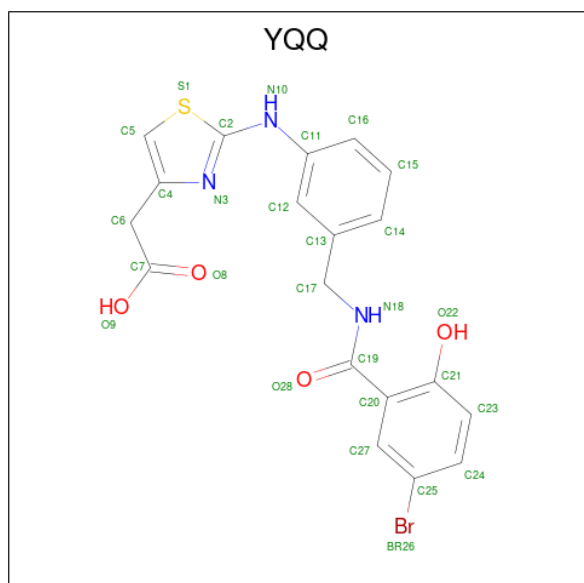
- Molecule 1 is a protein called Cyclic GMP-AMP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	0	0
			2885	1846	497	527	15			
1	B	351	Total	C	N	O	S	0	0	0
			2885	1846	497	527	15			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (2-{3-[(5-bromo-2-hydroxybenzamido)methyl]anilino}-1,3-thiazol-4-yl)acetic acid (CCD ID: YQQ) (formula: C₁₉H₁₆BrN₃O₄S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	Br	C	N	O	S	0	0
			28	1	19	3	4	1		
3	B	1	Total	Br	C	N	O	S	0	0
			28	1	19	3	4	1		

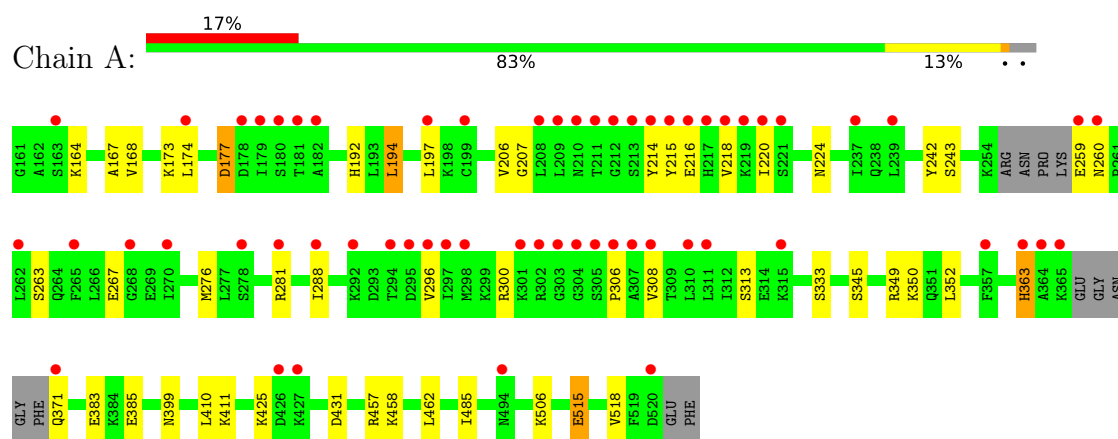
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	49	Total	O	0	0
			49	49		
4	B	35	Total	O	0	0
			35	35		

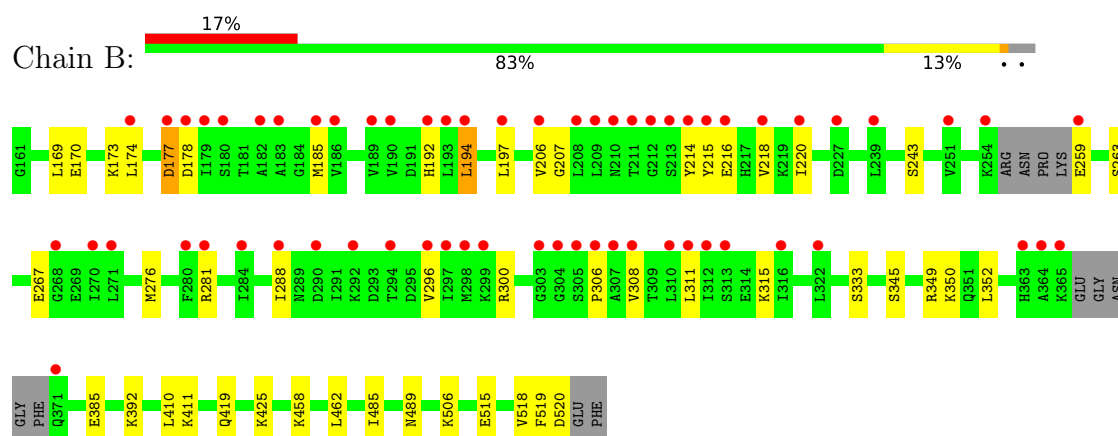
3 Residue-property plots [i](#)

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: Cyclic GMP-AMP synthase



• Molecule 1: Cyclic GMP-AMP synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	219.01Å 46.95Å 88.97Å 90.00° 110.62° 90.00°	Depositor
Resolution (Å)	83.27 – 2.50 83.27 – 2.50	Depositor EDS
% Data completeness (in resolution range)	82.1 (83.27-2.50) 82.2 (83.27-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.223 , 0.272 0.237 , 0.290	Depositor DCC
R_{free} test set	1271 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtrriage
Anisotropy	(Not available)	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 73.4	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5912	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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: ZN, YQQ

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	0.87	0/2940	1.43	27/3940 (0.7%)
1	B	0.88	0/2940	1.44	25/3940 (0.6%)
All	All	0.87	0/5880	1.43	52/7880 (0.7%)

There are no bond length outliers.

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	178	ASP	CA-CB-CG	8.50	121.10	112.60
1	B	177	ASP	CA-CB-CG	6.89	119.49	112.60
1	A	177	ASP	CA-CB-CG	6.67	119.27	112.60
1	B	169	LEU	CA-C-N	6.64	129.07	120.44
1	B	169	LEU	C-N-CA	6.64	129.07	120.44
1	A	214	TYR	N-CA-C	-6.38	103.58	112.45
1	B	214	TYR	N-CA-C	-6.30	103.69	112.45
1	A	399	ASN	CA-C-N	6.19	128.57	120.28
1	A	399	ASN	C-N-CA	6.19	128.57	120.28
1	B	462	LEU	CA-C-N	6.12	128.40	120.44
1	B	462	LEU	C-N-CA	6.12	128.40	120.44
1	A	462	LEU	CA-C-N	6.07	128.33	120.44
1	A	462	LEU	C-N-CA	6.07	128.33	120.44
1	A	425	LYS	CA-C-N	6.03	128.64	120.44
1	A	425	LYS	C-N-CA	6.03	128.64	120.44
1	B	519	PHE	CA-C-N	6.01	132.52	121.70
1	B	519	PHE	C-N-CA	6.01	132.52	121.70
1	B	306	PRO	CA-C-N	5.97	128.60	120.54
1	B	306	PRO	C-N-CA	5.97	128.60	120.54
1	B	425	LYS	CA-C-N	5.97	128.77	120.29
1	B	425	LYS	C-N-CA	5.97	128.77	120.29
1	A	306	PRO	CA-C-N	5.79	128.36	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	PRO	C-N-CA	5.79	128.36	120.54
1	A	218	VAL	N-CA-C	-5.58	104.92	111.00
1	A	410	LEU	CA-C-N	5.53	128.00	120.54
1	A	410	LEU	C-N-CA	5.53	128.00	120.54
1	B	392	LYS	N-CA-C	-5.50	105.29	112.23
1	B	170	GLU	CA-C-N	5.41	128.31	120.79
1	B	170	GLU	C-N-CA	5.41	128.31	120.79
1	B	410	LEU	CA-C-N	5.39	127.81	120.54
1	B	410	LEU	C-N-CA	5.39	127.81	120.54
1	A	281	ARG	CA-C-N	5.38	127.80	120.54
1	A	281	ARG	C-N-CA	5.38	127.80	120.54
1	B	281	ARG	CA-C-N	5.32	127.72	120.54
1	B	281	ARG	C-N-CA	5.32	127.72	120.54
1	A	224	ASN	CA-CB-CG	5.30	117.90	112.60
1	B	349	ARG	CA-C-N	5.29	127.32	120.44
1	B	349	ARG	C-N-CA	5.29	127.32	120.44
1	A	515	GLU	CB-CG-CD	5.28	121.58	112.60
1	B	489	ASN	CA-C-N	5.23	127.81	120.28
1	B	489	ASN	C-N-CA	5.23	127.81	120.28
1	B	218	VAL	N-CA-C	-5.22	105.31	111.00
1	A	167	ALA	CA-C-N	5.21	127.60	120.46
1	A	167	ALA	C-N-CA	5.21	127.60	120.46
1	A	349	ARG	CA-C-N	5.20	127.20	120.44
1	A	349	ARG	C-N-CA	5.20	127.20	120.44
1	A	431	ASP	CA-CB-CG	5.15	117.75	112.60
1	A	242	TYR	CA-C-N	5.07	131.22	121.54
1	A	242	TYR	C-N-CA	5.07	131.22	121.54
1	A	383	GLU	CA-C-N	5.05	127.04	120.28
1	A	383	GLU	C-N-CA	5.05	127.04	120.28
1	A	260	ASN	CA-CB-CG	5.04	117.64	112.60

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	2885	0	2931	7	0
1	B	2885	0	2931	6	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	28	0	0	0	0
3	B	28	0	0	0	0
4	A	49	0	0	0	0
4	B	35	0	0	0	0
All	All	5912	0	5862	13	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 (13) 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:352:LEU:HD21	1:A:385:GLU:HG2	1.90	0.53
1:B:352:LEU:HD21	1:B:385:GLU:HG2	1.90	0.53
1:B:419:GLN:HB3	1:B:518:VAL:HG22	1.95	0.48
1:A:164:LYS:O	1:A:168:VAL:HG23	2.14	0.47
1:B:194:LEU:HD21	1:B:207:GLY:HA2	2.00	0.42
1:B:288:ILE:HG23	1:B:296:VAL:HB	2.02	0.42
1:A:288:ILE:HG23	1:A:296:VAL:HB	2.02	0.41
1:A:333:SER:HB2	1:A:485:ILE:HG23	2.02	0.41
1:A:363:HIS:NE2	1:A:371:GLN:HB3	2.33	0.41
1:A:194:LEU:HD21	1:A:207:GLY:HA2	2.01	0.41
1:B:333:SER:HB2	1:B:485:ILE:HG23	2.03	0.41
1:B:352:LEU:CD2	1:B:385:GLU:HG2	2.49	0.41
1:A:352:LEU:CD2	1:A:385:GLU:HG2	2.49	0.41

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	345/362 (95%)	326 (94%)	15 (4%)	4 (1%)	10	20
1	B	345/362 (95%)	324 (94%)	18 (5%)	3 (1%)	14	27
All	All	690/724 (95%)	650 (94%)	33 (5%)	7 (1%)	12	24

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	SER
1	A	215	TYR
1	A	345	SER
1	B	215	TYR
1	B	345	SER
1	A	313	SER
1	B	243	SER

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	324/334 (97%)	301 (93%)	23 (7%)	13	29
1	B	324/334 (97%)	300 (93%)	24 (7%)	13	27
All	All	648/668 (97%)	601 (93%)	47 (7%)	13	27

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	173	LYS
1	A	174	LEU
1	A	177	ASP
1	A	192	HIS
1	A	194	LEU
1	A	197	LEU
1	A	206	VAL
1	A	216	GLU
1	A	220	ILE

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Mol	Chain	Res	Type
1	A	259	GLU
1	A	263	SER
1	A	267	GLU
1	A	276	MET
1	A	300	ARG
1	A	308	VAL
1	A	350	LYS
1	A	363	HIS
1	A	411	LYS
1	A	457	ARG
1	A	458	LYS
1	A	506	LYS
1	A	515	GLU
1	A	518	VAL
1	B	173	LYS
1	B	174	LEU
1	B	177	ASP
1	B	185	MET
1	B	192	HIS
1	B	194	LEU
1	B	197	LEU
1	B	206	VAL
1	B	216	GLU
1	B	220	ILE
1	B	259	GLU
1	B	263	SER
1	B	267	GLU
1	B	276	MET
1	B	300	ARG
1	B	308	VAL
1	B	311	LEU
1	B	315	LYS
1	B	350	LYS
1	B	411	LYS
1	B	458	LYS
1	B	506	LYS
1	B	515	GLU
1	B	520	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	224	ASN
1	A	238	GLN
1	A	389	ASN
1	A	451	GLN
1	A	514	ASN
1	B	224	ASN
1	B	238	GLN
1	B	451	GLN
1	B	514	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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
3	YQQ	A	602	-	30,30,30	1.90	7 (23%)	41,41,41	1.25	6 (14%)
3	YQQ	B	602	-	30,30,30	1.91	9 (30%)	41,41,41	1.20	4 (9%)

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	YQQ	A	602	-	-	1/17/17/17	0/3/3/3
3	YQQ	B	602	-	-	2/17/17/17	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	YQQ	C2-S1	4.84	1.80	1.74
3	A	602	YQQ	C2-S1	4.55	1.80	1.74
3	A	602	YQQ	C12-C11	3.26	1.44	1.39
3	A	602	YQQ	C19-N18	3.23	1.40	1.33
3	B	602	YQQ	BR26-C25	3.03	1.96	1.90
3	A	602	YQQ	C12-C13	2.81	1.44	1.39
3	B	602	YQQ	C12-C13	2.79	1.44	1.39
3	B	602	YQQ	C19-N18	2.77	1.39	1.33
3	B	602	YQQ	C12-C11	2.73	1.43	1.39
3	B	602	YQQ	O9-C7	2.69	1.39	1.30
3	A	602	YQQ	C23-C21	2.64	1.44	1.39
3	B	602	YQQ	C23-C21	2.42	1.43	1.39
3	A	602	YQQ	BR26-C25	2.40	1.95	1.90
3	A	602	YQQ	O9-C7	2.29	1.38	1.30
3	B	602	YQQ	O28-C19	2.10	1.28	1.23
3	B	602	YQQ	C2-N3	2.00	1.34	1.31

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	YQQ	C7-C6-C4	-3.04	107.25	113.25
3	B	602	YQQ	C7-C6-C4	-2.98	107.37	113.25
3	A	602	YQQ	S1-C2-N10	2.37	128.28	121.37
3	A	602	YQQ	O28-C19-C20	-2.26	116.89	121.03
3	B	602	YQQ	S1-C2-N10	2.19	127.75	121.37
3	B	602	YQQ	O22-C21-C23	-2.18	113.51	119.36
3	B	602	YQQ	O22-C21-C20	2.11	125.61	121.71
3	A	602	YQQ	BR26-C25-C27	-2.05	116.33	119.23
3	A	602	YQQ	O22-C21-C23	-2.04	113.88	119.36
3	A	602	YQQ	O8-C7-C6	2.01	127.85	122.11

There are no chirality outliers.

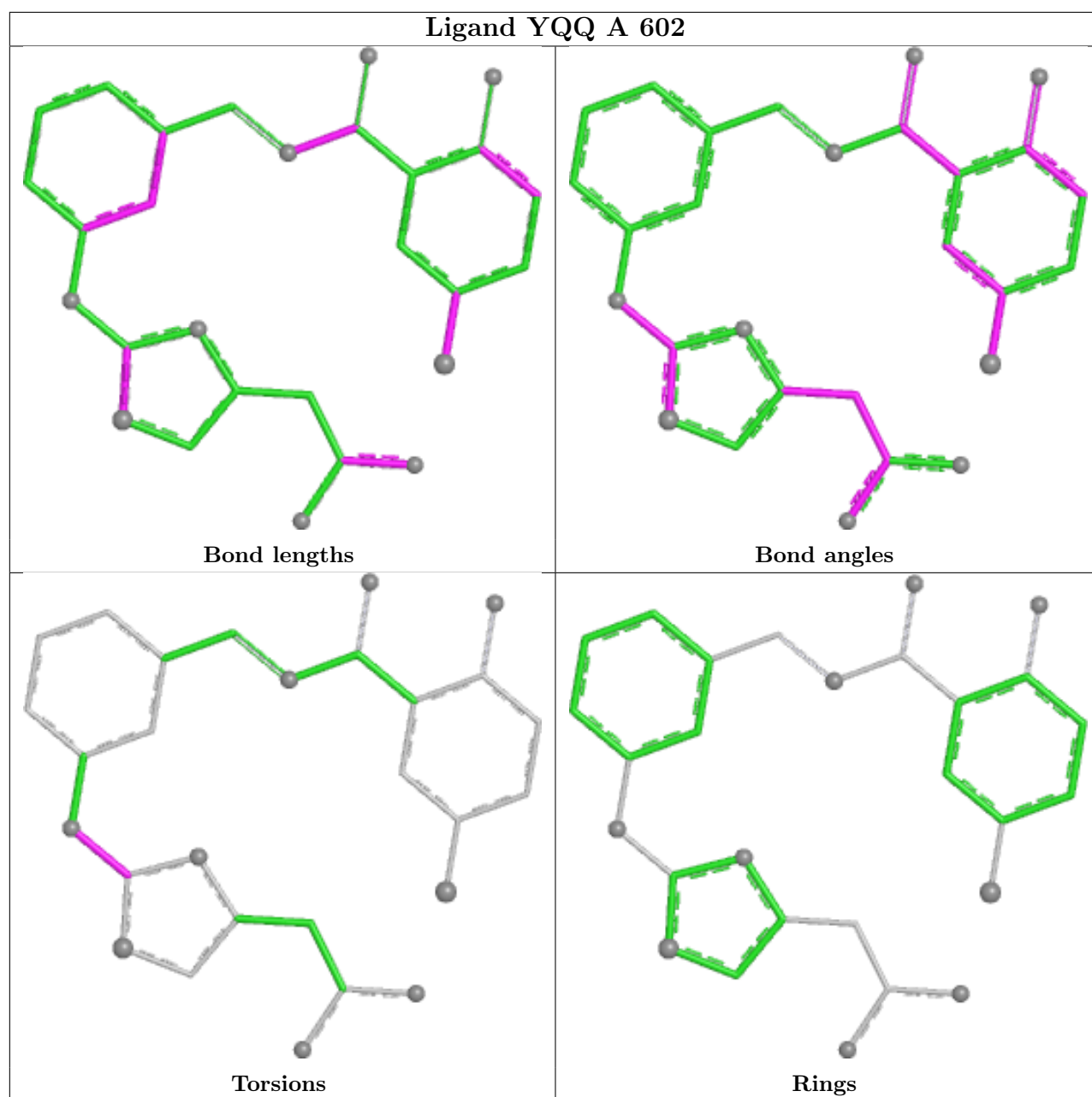
All (3) torsion outliers are listed below:

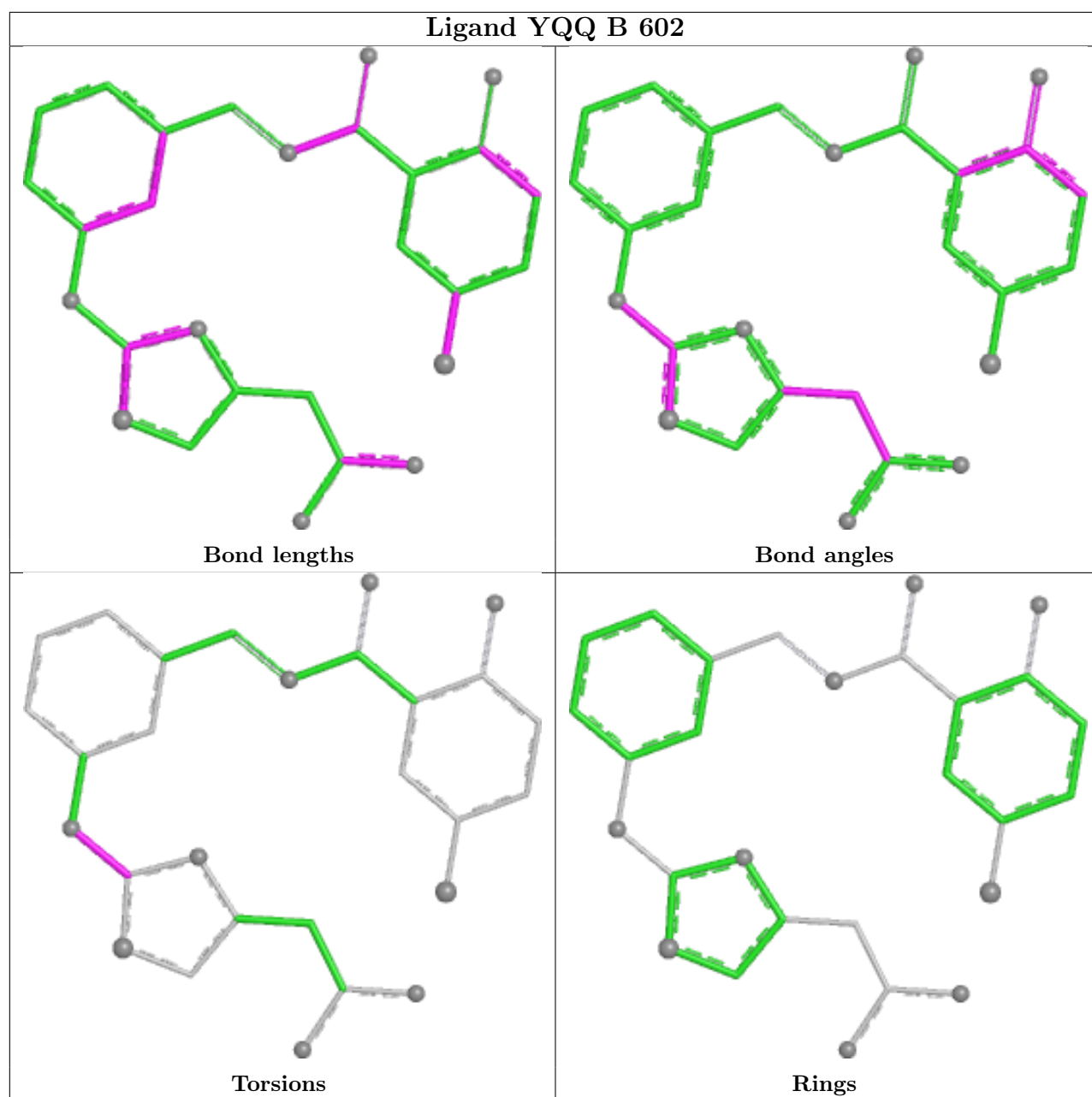
Mol	Chain	Res	Type	Atoms
3	B	602	YQQ	S1-C2-N10-C11
3	A	602	YQQ	N3-C2-N10-C11
3	B	602	YQQ	N3-C2-N10-C11

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	351/362 (96%)	0.89	60 (17%) 4 3	34, 65, 114, 140	0
1	B	351/362 (96%)	0.96	62 (17%) 4 3	31, 70, 124, 143	0
All	All	702/724 (96%)	0.92	122 (17%) 4 3	31, 68, 120, 143	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	179	ILE	5.6
1	A	179	ILE	5.4
1	A	305	SER	5.1
1	A	214	TYR	4.7
1	A	211	THR	4.7
1	B	183	ALA	4.5
1	A	302	ARG	4.3
1	A	364	ALA	4.2
1	B	214	TYR	4.2
1	B	211	THR	4.1
1	A	371	GLN	4.0
1	B	306	PRO	3.9
1	B	215	TYR	3.9
1	B	218	VAL	3.9
1	A	218	VAL	3.8
1	A	304	GLY	3.8
1	B	303	GLY	3.8
1	B	270	ILE	3.8
1	B	307	ALA	3.6
1	B	180	SER	3.6
1	B	298	MET	3.6
1	B	305	SER	3.6
1	A	307	ALA	3.6
1	A	215	TYR	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	294	THR	3.5
1	A	182	ALA	3.5
1	B	297	ILE	3.5
1	A	180	SER	3.5
1	A	178	ASP	3.4
1	B	178	ASP	3.3
1	B	371	GLN	3.3
1	B	194	LEU	3.3
1	B	208	LEU	3.3
1	A	213	SER	3.3
1	A	294	THR	3.2
1	B	186	VAL	3.2
1	A	288	ILE	3.1
1	B	311	LEU	3.1
1	A	306	PRO	3.1
1	A	260	ASN	3.1
1	B	292	LYS	3.0
1	A	181	THR	3.0
1	A	311	LEU	3.0
1	A	210	ASN	3.0
1	B	322	LEU	2.9
1	B	365	LYS	2.9
1	A	212	GLY	2.9
1	A	292	LYS	2.8
1	A	315	LYS	2.8
1	B	316	ILE	2.8
1	B	304	GLY	2.8
1	A	296	VAL	2.8
1	B	299	LYS	2.8
1	B	213	SER	2.7
1	B	193	LEU	2.7
1	A	270	ILE	2.7
1	B	280	PHE	2.7
1	B	210	ASN	2.7
1	A	197	LEU	2.7
1	B	185	MET	2.6
1	B	239	LEU	2.6
1	A	259	GLU	2.6
1	B	284	ILE	2.6
1	B	189	VAL	2.6
1	B	364	ALA	2.6
1	B	209	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	303	GLY	2.5
1	B	212	GLY	2.5
1	A	216	GLU	2.5
1	A	174	LEU	2.5
1	B	296	VAL	2.5
1	B	308	VAL	2.5
1	A	278	SER	2.5
1	A	237	ILE	2.5
1	B	216	GLU	2.5
1	B	288	ILE	2.5
1	B	190	VAL	2.5
1	B	251	VAL	2.5
1	B	254	LYS	2.5
1	B	310	LEU	2.5
1	B	206	VAL	2.5
1	B	182	ALA	2.4
1	A	365	LYS	2.4
1	B	174	LEU	2.4
1	B	290	ASP	2.4
1	B	281	ARG	2.3
1	A	308	VAL	2.3
1	A	520	ASP	2.3
1	B	177	ASP	2.3
1	A	219	LYS	2.3
1	A	268	GLY	2.3
1	B	268	GLY	2.3
1	A	220	ILE	2.3
1	B	259	GLU	2.3
1	A	297	ILE	2.2
1	B	197	LEU	2.2
1	B	312	ILE	2.2
1	A	265	PHE	2.2
1	B	192	HIS	2.2
1	A	209	LEU	2.1
1	A	310	LEU	2.1
1	A	295	ASP	2.1
1	A	301	LYS	2.1
1	A	262	LEU	2.1
1	A	217	HIS	2.1
1	A	357	PHE	2.1
1	B	227	ASP	2.1
1	A	208	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	239	LEU	2.1
1	A	281	ARG	2.1
1	B	220	ILE	2.1
1	A	426	ASP	2.1
1	A	199	CYS	2.1
1	A	494	ASN	2.0
1	B	271	LEU	2.0
1	A	221	SER	2.0
1	A	427	LYS	2.0
1	A	298	MET	2.0
1	A	163	SER	2.0
1	A	363	HIS	2.0
1	B	313	SER	2.0
1	B	363	HIS	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](#)

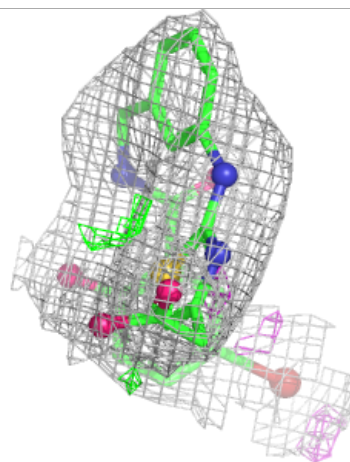
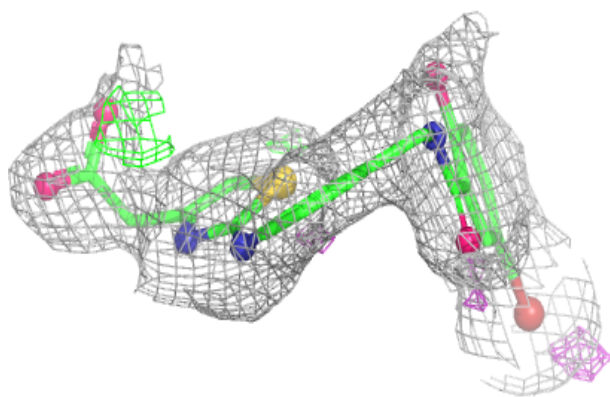
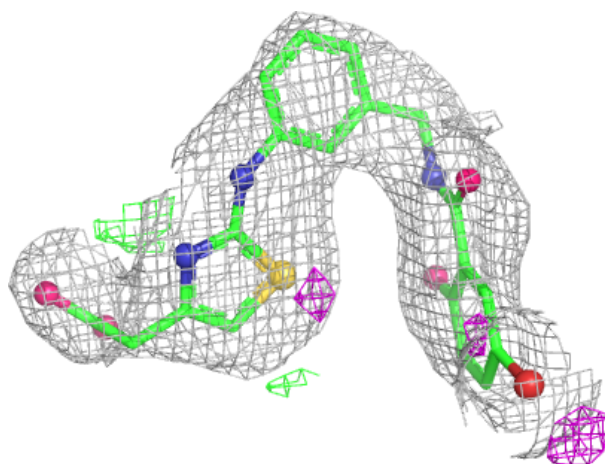
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	YQQ	A	602	28/28	0.95	0.10	48,54,72,77	0
3	YQQ	B	602	28/28	0.95	0.09	50,54,72,74	0
2	ZN	B	601	1/1	0.98	0.09	53,53,53,53	0
2	ZN	A	601	1/1	0.99	0.06	48,48,48,48	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.

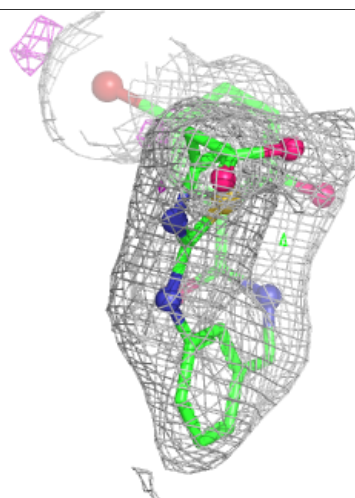
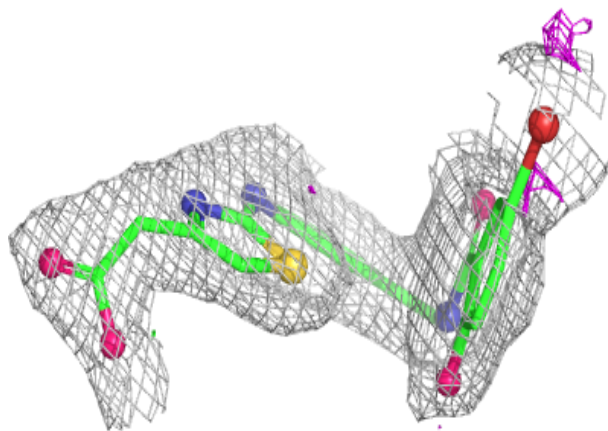
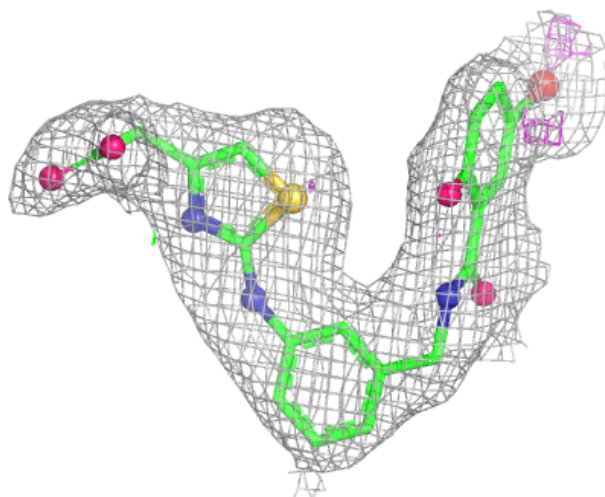
Electron density around YQQ A 602:

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



Electron density around YQQ B 602:

$2mF_o-DF_c$ (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.