



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 16, 2024 – 02:08 AM EST

PDB ID : 7RW5
Title : Crystal structure of human methionine adenosyltransferase 2A (MAT2A) in complex with SAM and allosteric inhibitor Compound 1
Authors : Jin, L.; Padyana, A.K.
Deposited on : 2021-08-19
Resolution : 2.48 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

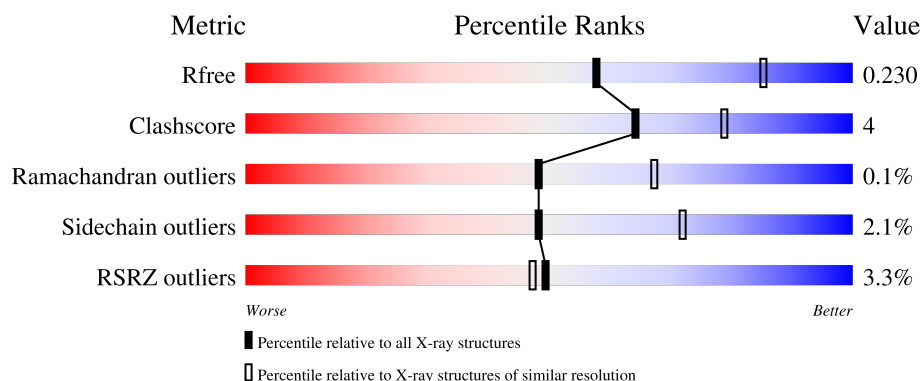
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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}	164625	7106 (2.50-2.46)
Clashscore	180529	7991 (2.50-2.46)
Ramachandran outliers	177936	7888 (2.50-2.46)
Sidechain outliers	177891	7890 (2.50-2.46)
RSRZ outliers	164620	7106 (2.50-2.46)

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	396	<div> <div>88%</div> <div>7%</div> <div>5%</div> </div>
1	B	396	<div> <div>87%</div> <div>7%</div> <div>6%</div> </div>
1	C	396	<div> <div>3%</div> <div>82%</div> <div>14%</div> <div>••</div> </div>
1	D	396	<div> <div>8%</div> <div>72%</div> <div>12%</div> <div>•</div> <div>14%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11985 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.

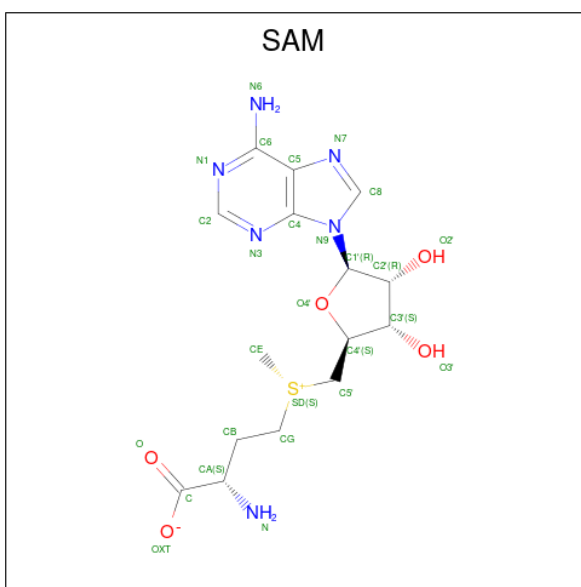
- Molecule 1 is a protein called S-adenosylmethionine synthase isoform type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	378	Total	C	N	O	S	0	2	0
			2944	1860	511	561	12			
1	B	373	Total	C	N	O	S	0	0	0
			2898	1838	501	548	11			
1	C	383	Total	C	N	O	S	0	0	0
			2977	1882	518	566	11			
1	D	339	Total	C	N	O	S	0	0	0
			2625	1661	456	497	11			

There are 4 discrepancies between the modelled and reference sequences:

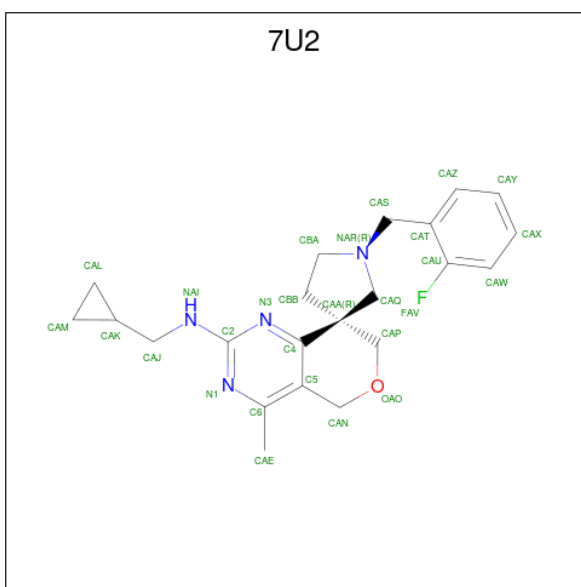
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP P31153
B	0	SER	-	expression tag	UNP P31153
C	0	SER	-	expression tag	UNP P31153
D	0	SER	-	expression tag	UNP P31153

- Molecule 2 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: C₁₅H₂₂N₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	D	1	Total	C	N	O	S	0	0
			27	15	6	5	1		

- Molecule 3 is (3'R)-N-(cyclopropylmethyl)-1'-[(2-fluorophenyl)methyl]-4-methyl-5H,7H-spiro[pyrano[4,3-d]pyrimidine-8,3'-pyrrolidin]-2-amine (three-letter code: 7U2) (formula: C₂₂H₂₇FN₄O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	F	N	O	0	0
			28	22	1	4	1		

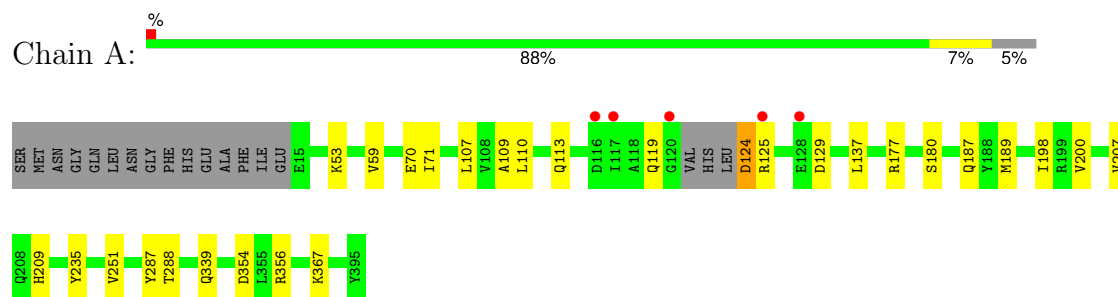
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	160	Total 160	O 160	0	0
4	B	142	Total 142	O 142	0	0
4	C	98	Total 98	O 98	0	0
4	D	59	Total 59	O 59	0	0

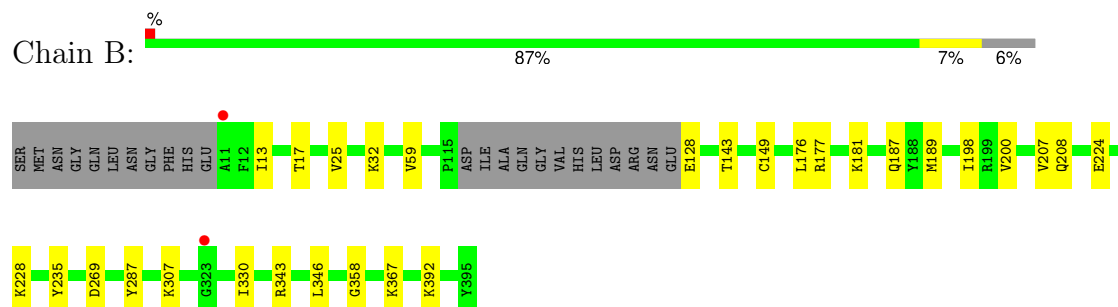
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

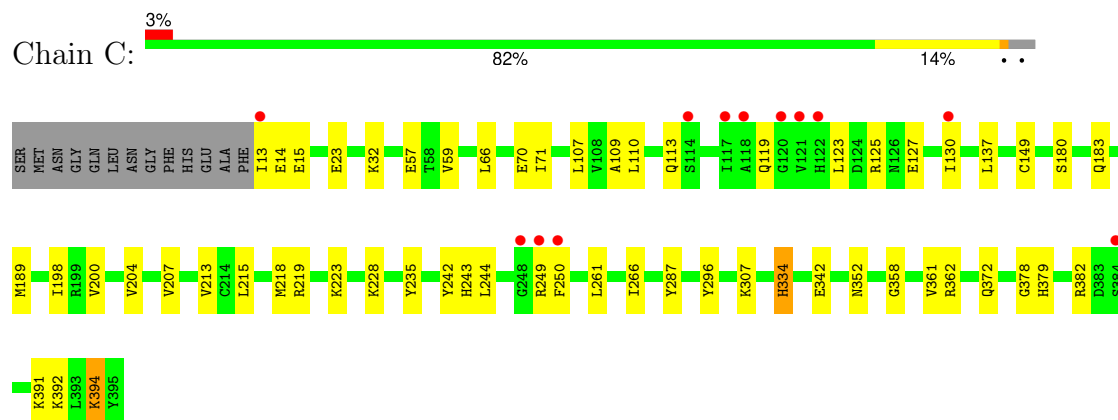
- Molecule 1: S-adenosylmethionine synthase isoform type-2



- Molecule 1: S-adenosylmethionine synthase isoform type-2

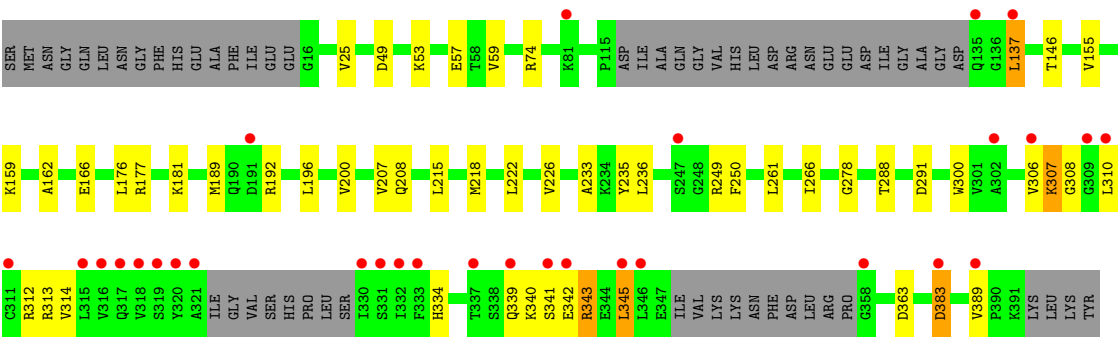


- Molecule 1: S-adenosylmethionine synthase isoform type-2



- Molecule 1: S-adenosylmethionine synthase isoform type-2





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.24Å 103.99Å 113.97Å 90.00° 93.47° 90.00°	Depositor
Resolution (Å)	37.92 – 2.48 37.92 – 2.48	Depositor EDS
% Data completeness (in resolution range)	96.3 (37.92-2.48) 96.5 (37.92-2.48)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.175 , 0.229 0.177 , 0.230	Depositor DCC
R_{free} test set	2849 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11985	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.30% 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: SAM, 7U2

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	0.24	0/3007	0.43	0/4065
1	B	0.24	0/2956	0.43	0/3997
1	C	0.24	0/3036	0.44	0/4107
1	D	0.24	0/2675	0.43	0/3618
All	All	0.24	0/11674	0.43	0/15787

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	2944	0	2934	16	0
1	B	2898	0	2894	20	0
1	C	2977	0	2968	33	0
1	D	2625	0	2608	27	0
2	B	27	0	21	1	0
2	D	27	0	22	0	0
3	B	28	0	0	0	0
4	A	160	0	0	1	0
4	B	142	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	98	0	0	1	0
4	D	59	0	0	0	0
All	All	11985	0	11447	88	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 4.

All (88) 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:B:149:CYS:SG	1:B:307:LYS:NZ	2.56	0.79
1:C:149:CYS:SG	1:C:307:LYS:NZ	2.61	0.73
1:C:352:ASN:O	1:C:391:LYS:NZ	2.25	0.69
1:B:189:MET:HB2	1:B:198:ILE:HD11	1.73	0.68
1:D:383:ASP:OD1	1:D:383:ASP:N	2.30	0.65
1:B:128:GLU:HB3	1:B:358:GLY:HA3	1.79	0.64
1:B:200:VAL:HG23	1:B:235:TYR:HB3	1.82	0.61
1:A:187:GLN:NE2	4:A:401:HOH:O	2.33	0.61
1:C:189:MET:HB2	1:C:198:ILE:HD11	1.81	0.60
1:C:13:ILE:HG13	1:C:15:GLU:H	1.65	0.60
1:C:200:VAL:HG23	1:C:235:TYR:HB3	1.84	0.59
1:D:177:ARG:HB2	1:D:208:GLN:HB3	1.83	0.59
1:D:314:VAL:HG23	1:D:334:HIS:HA	1.86	0.57
1:A:189:MET:HB2	1:A:198:ILE:HD11	1.86	0.56
1:D:200:VAL:HG23	1:D:235:TYR:HB3	1.88	0.56
1:C:59:VAL:HG21	1:D:59:VAL:HG21	1.87	0.56
1:A:177:ARG:NH1	1:A:251:VAL:O	2.38	0.55
1:D:310:LEU:O	1:D:339:GLN:NE2	2.40	0.53
1:C:372:GLN:NE2	4:C:402:HOH:O	2.39	0.52
1:C:57:GLU:HB3	1:D:261:LEU:HD11	1.92	0.52
1:C:70:GLU:HB3	1:C:113:GLN:HA	1.93	0.51
1:B:177:ARG:HB2	1:B:208:GLN:HB3	1.94	0.50
1:C:378:GLY:O	1:C:382:ARG:NH2	2.34	0.50
1:A:59:VAL:HG21	1:B:59:VAL:HG21	1.94	0.50
1:B:25:VAL:HG12	1:B:181:LYS:HG2	1.94	0.50
1:D:25:VAL:HG12	1:D:181:LYS:HG2	1.93	0.50
1:A:200:VAL:HG23	1:A:235:TYR:HB3	1.94	0.49
1:A:339:GLN:H	1:A:339:GLN:CD	2.16	0.49
1:C:14:GLU:HG3	1:D:343:ARG:HD2	1.95	0.48
1:B:32:LYS:NZ	1:B:269:ASP:OD1	2.37	0.48
1:C:71:ILE:HD12	1:C:110:LEU:HD13	1.93	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:LYS:HE3	1:A:288:THR:HG21	1.96	0.48
1:A:107:LEU:HD23	1:C:109:ALA:HB2	1.96	0.48
1:B:13:ILE:O	1:B:17:THR:HB	2.14	0.48
1:A:70:GLU:HB3	1:A:113:GLN:HA	1.95	0.47
1:C:23:GLU:HB3	1:C:183:GLN:HG3	1.97	0.47
1:D:155:VAL:HG13	1:D:159:LYS:HE2	1.97	0.47
1:D:222:LEU:HD23	1:D:226:VAL:HG21	1.97	0.47
1:C:394:LYS:O	1:C:394:LYS:HE3	2.14	0.47
1:C:127:GLU:O	1:C:358:GLY:HA3	2.14	0.47
1:B:17:THR:HG21	1:B:187:GLN:HE21	1.79	0.46
1:D:339:GLN:H	1:D:339:GLN:CD	2.18	0.46
1:A:109:ALA:HB2	1:C:107:LEU:HD23	1.97	0.46
1:B:287:TYR:CZ	1:B:367:LYS:HG2	2.51	0.46
1:B:330:ILE:HB	1:B:346:LEU:HD11	1.98	0.46
1:B:181:LYS:HD2	2:B:401:SAM:H4'	1.98	0.46
1:C:125:ARG:HD2	1:C:130:ILE:HA	1.98	0.46
1:C:180:SER:HB3	1:C:207:VAL:HG23	1.96	0.46
1:A:354:ASP:OD2	1:A:356:ARG:NH2	2.49	0.45
1:C:266:ILE:HD11	1:D:266:ILE:HD11	1.98	0.45
1:C:213:VAL:O	1:C:249:ARG:NH1	2.49	0.45
1:B:13:ILE:HD13	1:B:13:ILE:HA	1.78	0.45
1:D:189:MET:HB3	1:D:196:LEU:HB3	1.98	0.45
1:A:124:ASP:O	1:A:125:ARG:NE	2.44	0.45
1:D:233:ALA:HA	1:D:236:LEU:HD12	1.99	0.45
1:D:308:GLY:O	1:D:310:LEU:HD12	2.17	0.45
1:C:204:VAL:HB	1:C:243:HIS:HB2	1.99	0.44
1:A:180:SER:HB3	1:A:207:VAL:HG23	1.99	0.44
1:D:300:TRP:NE1	1:D:389:VAL:O	2.41	0.44
1:D:137:LEU:O	1:D:278:GLY:HA3	2.18	0.44
1:B:32:LYS:HA	1:B:32:LYS:HD3	1.69	0.44
1:C:261:LEU:HD11	1:D:57:GLU:HB3	1.99	0.44
1:C:334:HIS:CD2	1:C:342:GLU:HG3	2.53	0.43
1:D:176:LEU:HD22	1:D:207:VAL:HG21	2.00	0.43
1:C:287:TYR:CD1	1:C:361:VAL:HB	2.53	0.43
1:D:310:LEU:HD23	1:D:345:LEU:HD23	2.01	0.43
1:A:287:TYR:CE2	1:A:367:LYS:HG2	2.53	0.43
1:D:53:LYS:HE3	1:D:288:THR:HG21	1.99	0.43
1:D:306:VAL:HG23	1:D:314:VAL:HG12	2.00	0.43
1:B:343:ARG:HH22	1:B:346:LEU:HD23	1.84	0.43
1:C:218:MET:HE1	1:C:249:ARG:HG2	2.00	0.43
1:C:219:ARG:HG2	1:C:244:LEU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:296:TYR:OH	1:C:379:HIS:ND1	2.40	0.42
1:A:209:HIS:O	1:A:251:VAL:HA	2.19	0.42
1:B:392:LYS:HD3	1:B:392:LYS:HA	1.88	0.42
1:C:223:LYS:HE2	1:C:242:TYR:CG	2.55	0.42
1:D:49:ASP:OD1	1:D:74:ARG:NH2	2.53	0.42
1:D:162:ALA:O	1:D:166:GLU:HG2	2.19	0.42
1:B:176:LEU:HD22	1:B:207:VAL:HG21	2.01	0.42
1:C:32:LYS:HD2	1:C:32:LYS:HA	1.59	0.41
1:C:223:LYS:O	1:C:228:LYS:HD3	2.20	0.41
1:B:224:GLU:HA	1:B:228:LYS:HD2	2.01	0.41
1:D:215:LEU:HD23	1:D:218:MET:HE3	2.02	0.41
1:A:71:ILE:HD12	1:A:110:LEU:HD13	2.02	0.41
1:D:307:LYS:HA	1:D:307:LYS:HD2	1.85	0.41
1:B:143:THR:O	1:B:149:CYS:HA	2.22	0.40
1:C:215:LEU:HD13	1:C:249:ARG:HG3	2.03	0.40
1:C:119:GLN:HA	1:C:123:LEU:HB2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	376/396 (95%)	366 (97%)	10 (3%)	0	100	100
1	B	369/396 (93%)	361 (98%)	8 (2%)	0	100	100
1	C	381/396 (96%)	367 (96%)	14 (4%)	0	100	100
1	D	331/396 (84%)	318 (96%)	12 (4%)	1 (0%)	37	54
All	All	1457/1584 (92%)	1412 (97%)	44 (3%)	1 (0%)	48	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	313	ARG

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	314/328 (96%)	310 (99%)	4 (1%)	65	83
1	B	308/328 (94%)	308 (100%)	0	100	100
1	C	317/328 (97%)	310 (98%)	7 (2%)	47	70
1	D	278/328 (85%)	263 (95%)	15 (5%)	18	35
All	All	1217/1312 (93%)	1191 (98%)	26 (2%)	48	72

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	GLN
1	A	124	ASP
1	A	129	ASP
1	A	137	LEU
1	C	66	LEU
1	C	137	LEU
1	C	250	PHE
1	C	334	HIS
1	C	362	ARG
1	C	392	LYS
1	C	394	LYS
1	D	137	LEU
1	D	146	THR
1	D	192	ARG
1	D	249	ARG
1	D	250	PHE
1	D	291	ASP
1	D	307	LYS
1	D	312	ARG
1	D	340	LYS
1	D	341	SER

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Mol	Chain	Res	Type
1	D	342	GLU
1	D	343	ARG
1	D	345	LEU
1	D	363	ASP
1	D	383	ASP

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

Mol	Chain	Res	Type
1	B	161	ASN
1	D	161	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 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	SAM	B	401	-	23,29,29	1.27	3 (13%)	20,42,42	1.63	2 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	7U2	B	402	-	31,32,32	5.69	13 (41%)	36,47,47	5.92	12 (33%)
2	SAM	D	401	-	23,29,29	1.27	3 (13%)	20,42,42	1.65	2 (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
2	SAM	B	401	-	-	3/13/33/33	0/3/3/3
3	7U2	B	402	-	-	1/9/35/35	0/5/5/5
2	SAM	D	401	-	-	2/13/33/33	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	7U2	CAQ-NAR	-20.97	1.27	1.46
3	B	402	7U2	CAS-NAR	-15.51	1.17	1.47
3	B	402	7U2	CBB-CBA	-10.48	1.34	1.52
3	B	402	7U2	C2-NAI	8.00	1.47	1.34
3	B	402	7U2	CBA-NAR	6.50	1.60	1.47
3	B	402	7U2	CBB-CAA	-5.18	1.48	1.55
2	B	401	SAM	C2-N3	4.21	1.38	1.32
2	D	401	SAM	C2-N3	4.11	1.38	1.32
3	B	402	7U2	C5-C6	3.30	1.44	1.39
3	B	402	7U2	CAA-C4	3.30	1.58	1.52
3	B	402	7U2	CAE-C6	2.90	1.55	1.50
3	B	402	7U2	CAP-CAA	2.76	1.60	1.52
3	B	402	7U2	CAS-CAT	2.71	1.56	1.51
3	B	402	7U2	CAM-CAL	-2.64	1.39	1.48
2	B	401	SAM	C2-N1	2.58	1.38	1.33
2	D	401	SAM	C2-N1	2.53	1.38	1.33
3	B	402	7U2	C4-N3	2.49	1.37	1.33
2	D	401	SAM	OXT-C	-2.23	1.23	1.30
2	B	401	SAM	OXT-C	-2.14	1.23	1.30

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	7U2	CAM-CAK-CAJ	24.01	150.00	119.17
3	B	402	7U2	CAL-CAK-CAJ	23.94	149.91	119.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	SAM	N3-C2-N1	-6.32	120.10	128.67
2	B	401	SAM	N3-C2-N1	-6.15	120.32	128.67
3	B	402	7U2	CAA-C4-N3	4.10	121.05	117.21
3	B	402	7U2	CAJ-NAI-C2	-3.10	118.96	123.94
3	B	402	7U2	CAA-CAQ-NAR	3.08	106.45	103.64
3	B	402	7U2	CAS-NAR-CAQ	-2.94	109.00	112.89
3	B	402	7U2	N3-C2-N1	-2.75	121.76	126.26
2	B	401	SAM	OXT-C-O	-2.65	118.08	124.08
3	B	402	7U2	CAA-C4-C5	-2.64	119.94	123.26
3	B	402	7U2	C5-C4-N3	-2.60	119.01	123.66
2	D	401	SAM	OXT-C-O	-2.59	118.20	124.08
3	B	402	7U2	C5-C6-N1	-2.56	119.18	122.09
3	B	402	7U2	CAE-C6-N1	2.37	120.21	116.53
3	B	402	7U2	C2-N1-C6	2.33	120.52	116.00

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	402	7U2	NAI-CAJ-CAK-CAL
2	B	401	SAM	O-C-CA-CB
2	B	401	SAM	OXT-C-CA-CB
2	D	401	SAM	O-C-CA-CB
2	D	401	SAM	OXT-C-CA-CB
2	B	401	SAM	OXT-C-CA-N

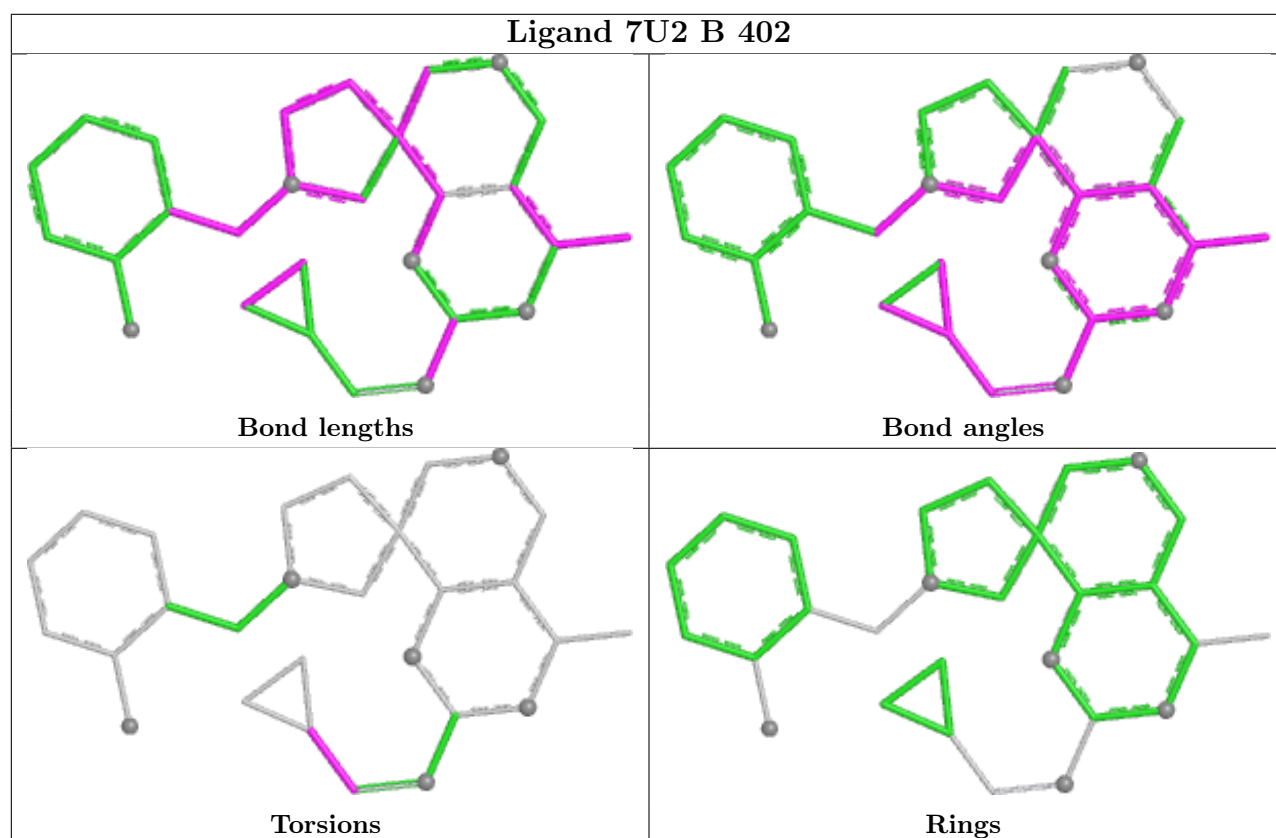
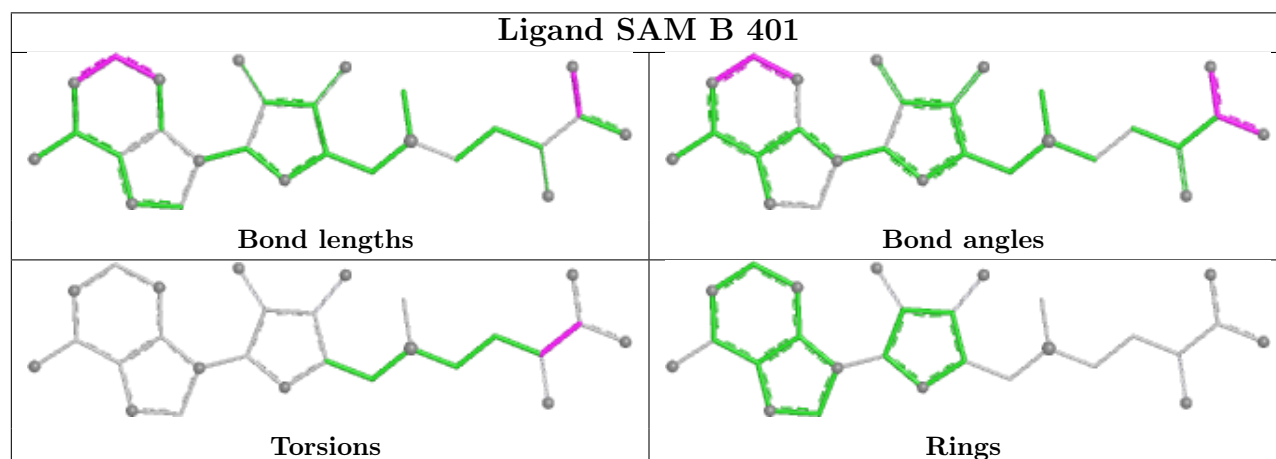
There are no ring outliers.

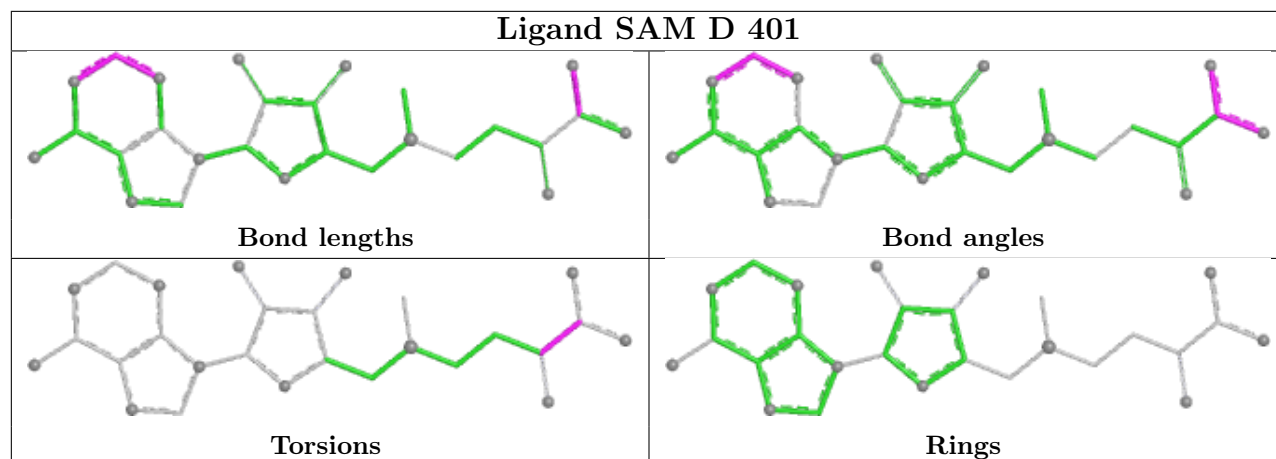
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	SAM	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.





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	378/396 (95%)	-0.41	5 (1%)	74 73	19, 33, 72, 128	2 (0%)
1	B	373/396 (94%)	-0.46	2 (0%)	87 85	20, 34, 69, 102	0
1	C	383/396 (96%)	-0.16	12 (3%)	51 49	21, 42, 78, 109	0
1	D	339/396 (85%)	0.25	30 (8%)	17 16	25, 55, 106, 159	0
All	All	1473/1584 (92%)	-0.20	49 (3%)	49 47	19, 39, 86, 159	2 (0%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	249	ARG	4.7
1	D	333	PHE	4.4
1	C	118	ALA	4.4
1	D	137	LEU	4.0
1	D	332	ILE	3.9
1	D	331	SER	3.7
1	C	13	ILE	3.7
1	A	117	ILE	3.6
1	D	330	ILE	3.5
1	C	114	SER	3.3
1	D	317	GLN	3.3
1	D	339	GLN	3.2
1	D	318	VAL	3.2
1	D	247	SER	3.2
1	C	122	HIS	3.1
1	D	358	GLY	3.1
1	C	248	GLY	3.0
1	D	311	CYS	2.9
1	C	121	VAL	2.9
1	D	321	ALA	2.9
1	B	323	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	316	VAL	2.9
1	D	319	SER	2.8
1	D	320	TYR	2.6
1	D	342	GLU	2.6
1	C	117	ILE	2.6
1	C	250	PHE	2.6
1	C	384	SER	2.6
1	D	389	VAL	2.6
1	D	309	GLY	2.6
1	D	306	VAL	2.5
1	D	81	LYS	2.5
1	A	120	GLY	2.5
1	D	135	GLN	2.4
1	D	341	SER	2.4
1	D	302	ALA	2.4
1	A	116	ASP	2.3
1	C	120	GLY	2.3
1	D	337	THR	2.3
1	B	11	ALA	2.3
1	D	315	LEU	2.3
1	C	130	ILE	2.3
1	D	345	LEU	2.3
1	D	346	LEU	2.2
1	D	383	ASP	2.2
1	D	310	LEU	2.1
1	A	125	ARG	2.1
1	D	191	ASP	2.1
1	A	128	GLU	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 monosaccharides in this entry.

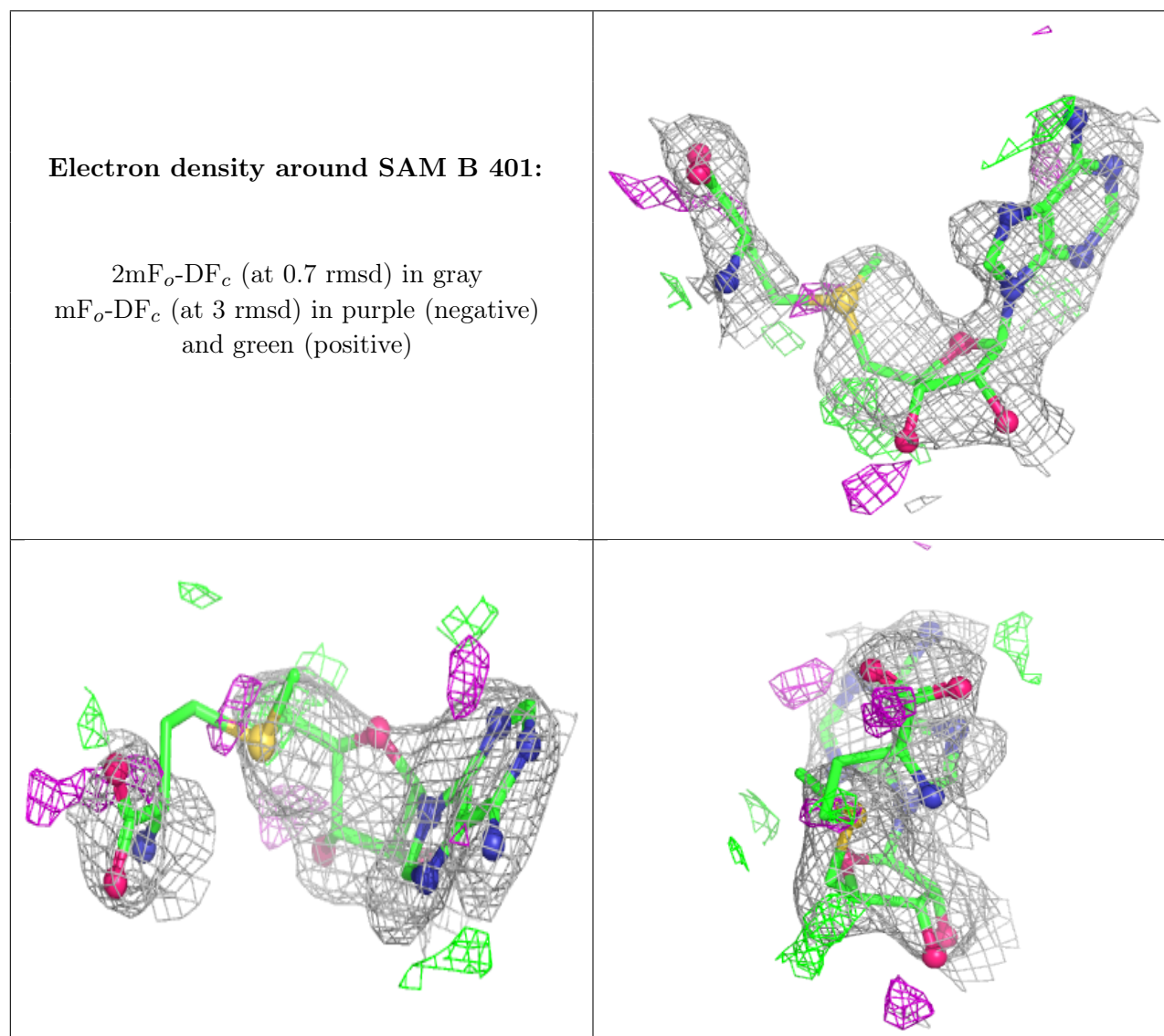
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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

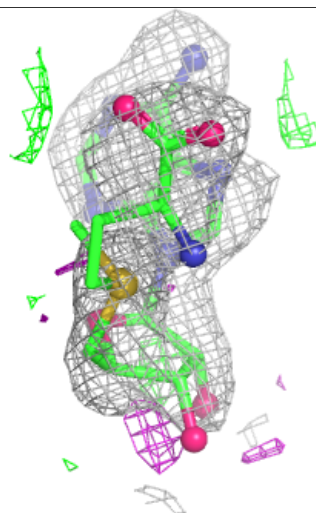
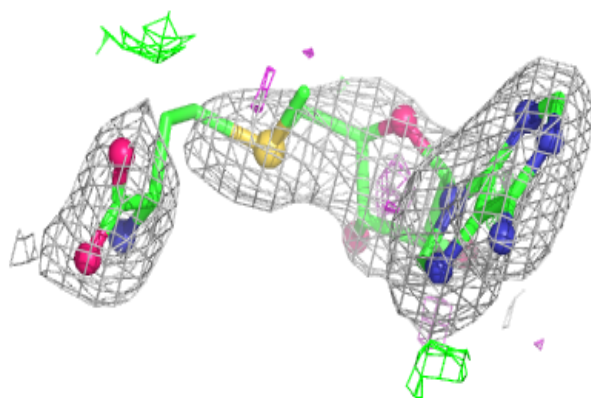
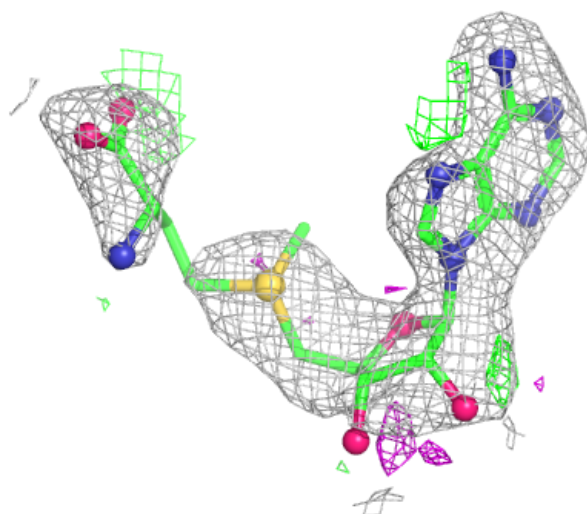
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SAM	B	401	27/27	0.78	0.19	58,74,87,90	0
2	SAM	D	401	27/27	0.85	0.14	50,71,79,80	0
3	7U2	B	402	28/28	0.93	0.09	37,46,53,54	0

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.



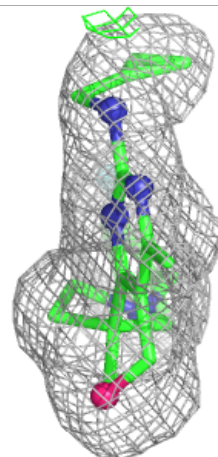
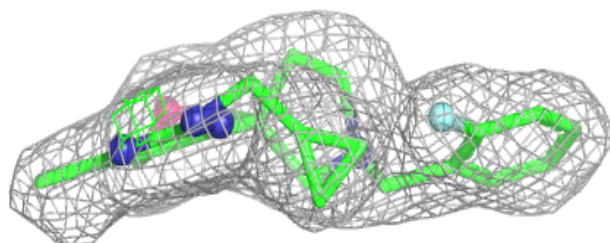
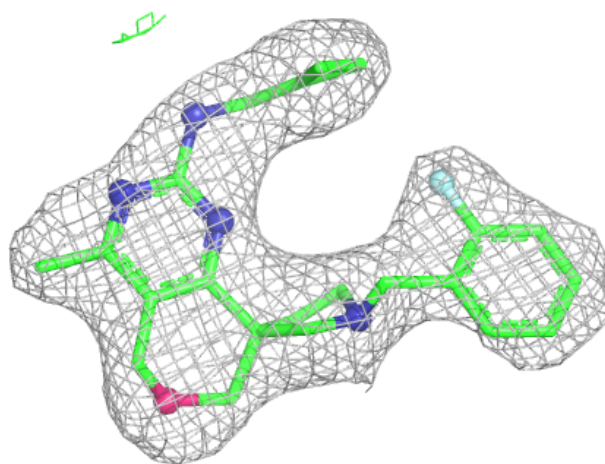
Electron density around SAM D 401:

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



Electron density around 7U2 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.