



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:04 PM UTC

PDB ID : 7YRJ / pdb_00007yrj
Title : S-adenosylmethionine sensor upstream of mTORC1
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Deposited on : 2022-08-10
Resolution : 2.35 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

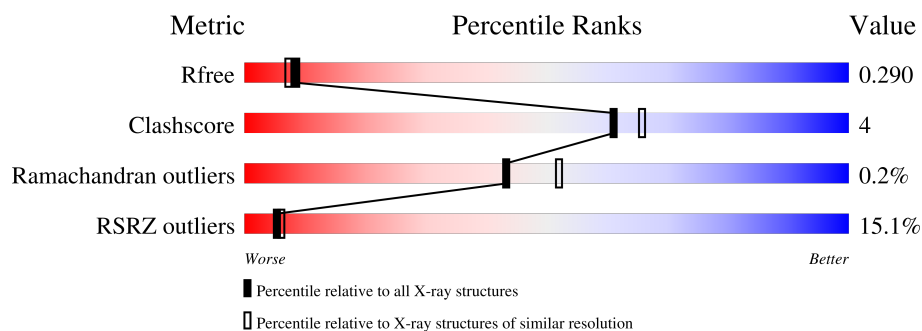
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	236	11% 81% 9% 10%
1	B	236	10% 78% 12% 10%
1	C	236	19% 76% 11% 13%
1	D	236	13% 80% 7% 14%

2 Entry composition [i](#)

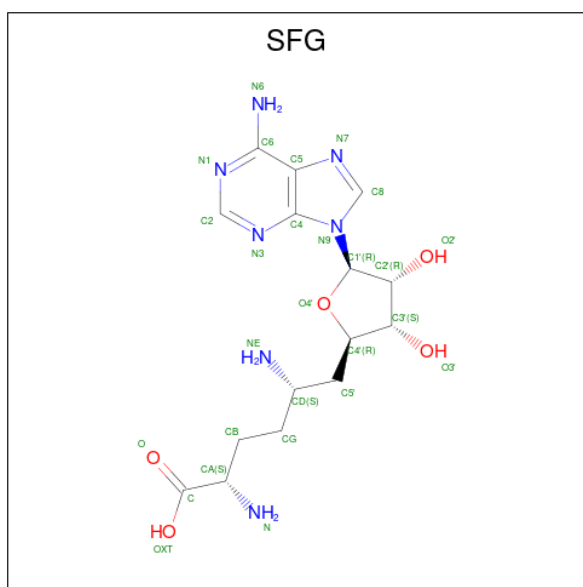
There are 3 unique types of molecules in this entry. The entry contains 13814 atoms, of which 6825 are hydrogens and 0 are deuteriums.

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 sensor upstream of mTORC1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	213	Total	C	H	N	O	S	0	0	0
			3433	1096	1722	306	298	11			
1	A	213	Total	C	H	N	O	S	0	0	0
			3408	1089	1707	301	300	11			
1	C	205	Total	C	H	N	O	S	0	0	0
			3313	1060	1662	289	291	11			
1	D	204	Total	C	H	N	O	S	0	0	0
			3292	1059	1646	283	293	11			

- Molecule 2 is SINEFUNGIN (CCD ID: SFG) (formula: C₁₅H₂₃N₇O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	H	N	O	0	0
			49	15	22	7	5		
2	A	1	Total	C	H	N	O	0	0
			49	15	22	7	5		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	H	N	O	0	0
			49	15	22	7	5		
2	D	1	Total	C	H	N	O	0	0
			49	15	22	7	5		

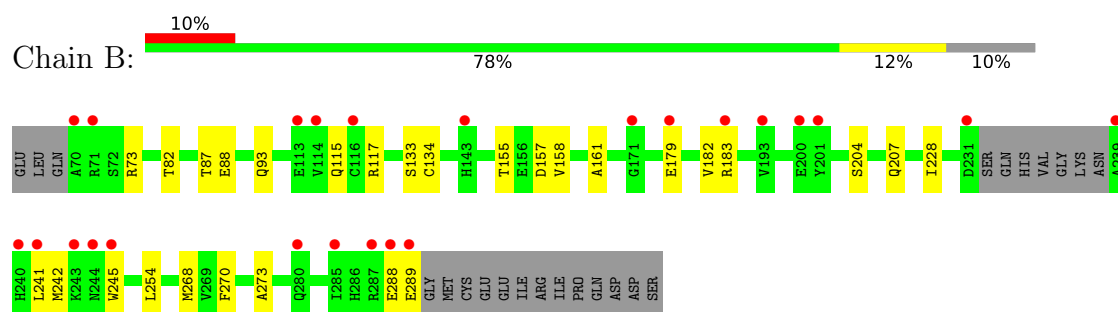
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	52	Total	O	0	0
			52	52		
3	A	46	Total	O	0	0
			46	46		
3	C	33	Total	O	0	0
			33	33		
3	D	41	Total	O	0	0
			41	41		

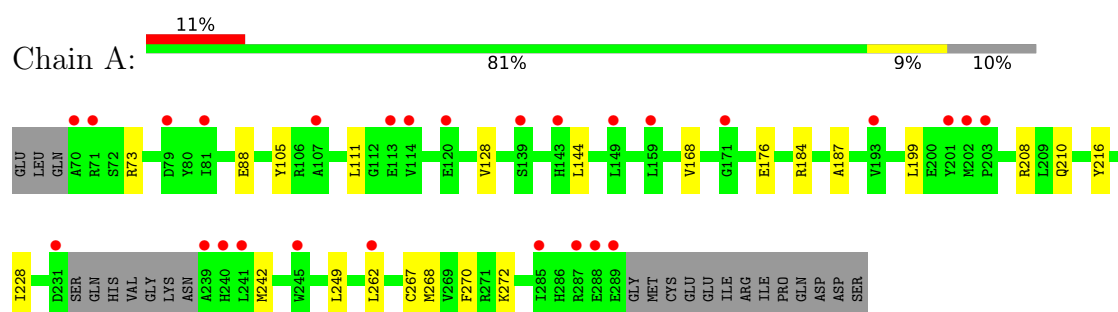
3 Residue-property plots

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.

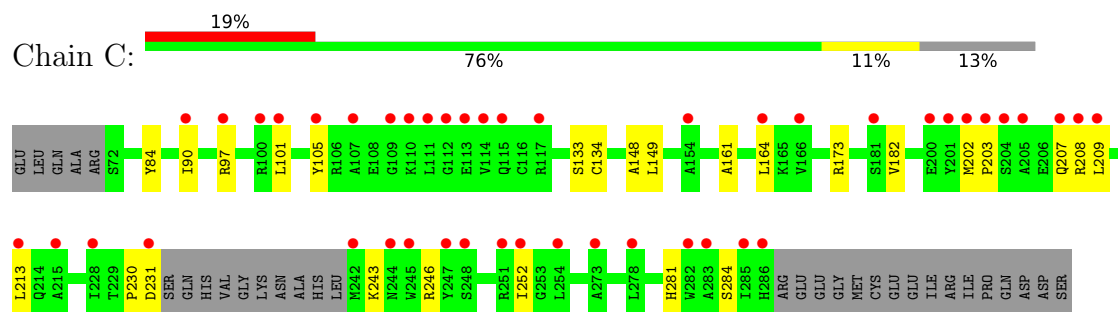
- Molecule 1: S-adenosylmethionine sensor upstream of mTORC1



- Molecule 1: S-adenosylmethionine sensor upstream of mTORC1

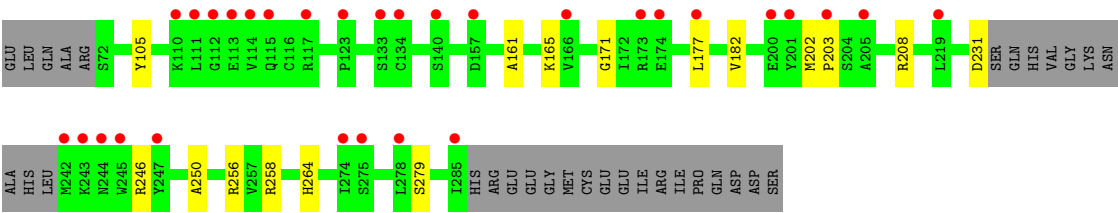


- Molecule 1: S-adenosylmethionine sensor upstream of mTORC1



- Molecule 1: S-adenosylmethionine sensor upstream of mTORC1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.33Å 78.90Å 100.10Å 90.00° 96.40° 90.00°	Depositor
Resolution (Å)	63.93 – 2.35 63.93 – 2.35	Depositor EDS
% Data completeness (in resolution range)	97.4 (63.93-2.35) 97.4 (63.93-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.240 , 0.291 (Not available) , 0.290	Depositor DCC
R_{free} test set	2055 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.792	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13814	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 56.92 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.5529e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: SFG

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	0.11	0/1737	0.28	0/2352
1	B	0.11	0/1749	0.29	0/2367
1	C	0.11	0/1688	0.31	0/2285
1	D	0.11	0/1683	0.28	0/2279
All	All	0.11	0/6857	0.29	0/9283

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	1701	1707	1705	13	0
1	B	1711	1722	1722	18	0
1	C	1651	1662	1662	16	0
1	D	1646	1646	1655	10	0
2	A	27	22	22	0	0
2	B	27	22	22	0	0
2	C	27	22	22	0	0
2	D	27	22	22	0	0
3	A	46	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	52	0	0	6	0
3	C	33	0	0	0	0
3	D	41	0	0	4	0
All	All	6989	6825	6832	57	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 (57) 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:C:213:LEU:HD11	1:C:252:ILE:HG21	1.64	0.80
1:B:134:CYS:N	3:B:505:HOH:O	2.16	0.79
1:A:128:VAL:HG21	1:A:144:LEU:HG	1.71	0.72
1:D:258:ARG:NE	3:D:504:HOH:O	2.25	0.69
1:A:88:GLU:OE1	3:A:501:HOH:O	2.13	0.67
1:B:242:MET:SD	1:B:268:MET:HE1	2.35	0.66
1:B:88:GLU:OE1	3:B:502:HOH:O	2.14	0.66
1:D:105:TYR:OH	3:D:501:HOH:O	2.14	0.65
1:B:179:GLU:OE1	3:B:503:HOH:O	2.15	0.65
1:C:97:ARG:HG2	1:C:97:ARG:HH11	1.61	0.65
1:B:82:THR:HG22	1:B:87:THR:HG23	1.79	0.64
1:C:281:HIS:O	1:C:284:SER:OG	2.14	0.63
1:D:165:LYS:NZ	3:D:506:HOH:O	2.32	0.62
1:B:93:GLN:OE1	3:B:504:HOH:O	2.16	0.62
1:D:250:ALA:O	1:D:279:SER:OG	2.21	0.58
1:A:168:VAL:HG12	1:A:187:ALA:HB2	1.88	0.55
1:B:161:ALA:HB2	1:B:182:VAL:HB	1.89	0.55
1:C:164:LEU:HD22	1:C:207:GLN:HG2	1.88	0.54
1:C:230:PRO:O	1:C:231:ASP:HB2	2.09	0.53
1:C:202:MET:HB2	1:C:208:ARG:HG3	1.90	0.52
1:C:161:ALA:HB2	1:C:182:VAL:HB	1.92	0.52
1:A:105:TYR:CB	1:A:111:LEU:HD12	2.40	0.51
1:C:148:ALA:C	1:C:149:LEU:HD12	2.35	0.51
1:C:209:LEU:HD12	1:C:213:LEU:HD13	1.92	0.51
1:B:73:ARG:HB2	1:B:228:ILE:HD13	1.94	0.50
1:A:210:GLN:OE1	3:A:502:HOH:O	2.19	0.49
1:A:73:ARG:HB2	1:A:228:ILE:HD13	1.94	0.48
1:C:101:LEU:HD11	1:C:105:TYR:CZ	2.48	0.47
1:A:249:LEU:HD12	1:A:270:PHE:CZ	2.50	0.47
1:B:288:GLU:O	1:B:289:GLU:C	2.58	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:246:ARG:HD2	1:D:256:ARG:CZ	2.46	0.46
1:D:171:GLY:O	3:D:502:HOH:O	2.21	0.46
1:B:241:LEU:HD22	1:B:245:TRP:NE1	2.32	0.45
1:A:105:TYR:HB3	1:A:111:LEU:HD12	1.98	0.45
1:C:97:ARG:HG2	1:C:97:ARG:NH1	2.28	0.45
1:D:161:ALA:HB2	1:D:182:VAL:HB	1.98	0.45
1:B:183:ARG:NH1	3:B:503:HOH:O	2.45	0.45
1:B:242:MET:SD	1:B:268:MET:CE	3.04	0.44
1:B:133:SER:O	1:B:158:VAL:HG11	2.18	0.43
1:A:262:LEU:HD22	1:A:267:CYS:SG	2.59	0.43
1:A:199:LEU:HB3	1:A:208:ARG:HG2	2.02	0.42
1:D:202:MET:HB2	1:D:208:ARG:HG3	2.01	0.42
1:C:133:SER:O	1:C:134:CYS:HB3	2.19	0.42
1:B:115:GLN:HG3	1:B:117:ARG:HH11	1.84	0.41
1:B:273:ALA:O	3:B:506:HOH:O	2.22	0.41
1:C:173:ARG:O	1:C:173:ARG:HG2	2.20	0.41
1:B:204:SER:OG	1:B:207:GLN:HG3	2.21	0.41
1:A:242:MET:SD	1:A:268:MET:CE	3.08	0.41
1:C:84:TYR:HA	1:C:90:ILE:HB	2.03	0.41
1:D:231:ASP:OD1	1:D:264:HIS:HA	2.21	0.41
1:B:155:THR:OG1	1:B:157:ASP:OD1	2.37	0.40
1:B:254:LEU:HB3	1:B:270:PHE:HB3	2.03	0.40
1:A:176:GLU:HB2	1:A:184:ARG:HB2	2.03	0.40
1:C:230:PRO:O	1:C:231:ASP:CB	2.70	0.40
1:C:243:LYS:HD3	1:C:246:ARG:HD3	2.03	0.40
1:D:177:LEU:N	1:D:177:LEU:HD12	2.36	0.40
1:A:216:TYR:CE2	1:A:272:LYS:HG3	2.57	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	209/236 (89%)	203 (97%)	6 (3%)	0	100	100
1	B	209/236 (89%)	201 (96%)	8 (4%)	0	100	100
1	C	201/236 (85%)	191 (95%)	9 (4%)	1 (0%)	24	27
1	D	200/236 (85%)	193 (96%)	6 (3%)	1 (0%)	24	27
All	All	819/944 (87%)	788 (96%)	29 (4%)	2 (0%)	43	52

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	203	PRO
1	C	203	PRO

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](#)

4 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	SFG	A	401	-	28,29,29	2.80	10 (35%)	34,42,42	3.67	13 (38%)
2	SFG	B	401	-	28,29,29	2.79	10 (35%)	34,42,42	3.68	13 (38%)
2	SFG	D	401	-	28,29,29	2.78	10 (35%)	34,42,42	3.63	13 (38%)
2	SFG	C	401	-	28,29,29	2.79	11 (39%)	34,42,42	3.66	13 (38%)

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	SFG	A	401	-	-	2/17/33/33	0/3/3/3
2	SFG	B	401	-	-	3/17/33/33	0/3/3/3
2	SFG	D	401	-	-	3/17/33/33	0/3/3/3
2	SFG	C	401	-	-	3/17/33/33	0/3/3/3

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	SFG	C5-N7	8.72	1.55	1.39
2	C	401	SFG	C5-N7	8.70	1.55	1.39
2	B	401	SFG	C5-N7	8.68	1.54	1.39
2	D	401	SFG	C5-N7	8.64	1.54	1.39
2	B	401	SFG	C6-N6	5.51	1.48	1.34
2	D	401	SFG	C6-N6	5.46	1.48	1.34
2	A	401	SFG	C6-N6	5.41	1.48	1.34
2	C	401	SFG	C6-N6	5.41	1.48	1.34
2	A	401	SFG	C8-N9	-5.06	1.28	1.37
2	B	401	SFG	C8-N9	-5.01	1.29	1.37
2	D	401	SFG	C8-N9	-5.00	1.29	1.37
2	C	401	SFG	C8-N9	-4.98	1.29	1.37
2	B	401	SFG	C5-C4	4.72	1.47	1.39
2	A	401	SFG	C5-C4	4.58	1.47	1.39
2	D	401	SFG	C5-C4	4.56	1.47	1.39
2	C	401	SFG	C5-C4	4.51	1.47	1.39
2	C	401	SFG	C4-N9	-3.89	1.29	1.37
2	A	401	SFG	C4-N9	-3.76	1.29	1.37
2	D	401	SFG	C4-N9	-3.73	1.29	1.37
2	A	401	SFG	O4'-C1'	3.71	1.50	1.42
2	D	401	SFG	O4'-C1'	3.61	1.50	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	SFG	O4'-C1'	3.61	1.50	1.42
2	B	401	SFG	C4-N9	-3.49	1.30	1.37
2	C	401	SFG	O4'-C1'	3.45	1.50	1.42
2	C	401	SFG	C8-N7	3.08	1.37	1.31
2	D	401	SFG	C8-N7	3.05	1.37	1.31
2	A	401	SFG	C8-N7	3.01	1.37	1.31
2	B	401	SFG	C8-N7	2.94	1.37	1.31
2	C	401	SFG	CG-CB	2.43	1.59	1.53
2	A	401	SFG	CG-CB	2.26	1.59	1.53
2	D	401	SFG	CG-CB	2.24	1.59	1.53
2	B	401	SFG	CG-CB	2.20	1.58	1.53
2	A	401	SFG	C2'-C1'	-2.19	1.46	1.53
2	D	401	SFG	C2'-C1'	-2.15	1.46	1.53
2	C	401	SFG	C2'-C1'	-2.11	1.46	1.53
2	B	401	SFG	C3'-C4'	-2.11	1.47	1.53
2	B	401	SFG	C6-N1	-2.11	1.29	1.35
2	C	401	SFG	C6-N1	-2.07	1.30	1.35
2	D	401	SFG	C6-N1	-2.05	1.30	1.35
2	A	401	SFG	C6-N1	-2.05	1.30	1.35
2	C	401	SFG	C3'-C4'	-2.03	1.47	1.53

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	SFG	C4-N9-C8	15.00	121.49	105.74
2	C	401	SFG	C4-N9-C8	14.95	121.43	105.74
2	D	401	SFG	C4-N9-C8	14.80	121.28	105.74
2	B	401	SFG	C4-N9-C8	14.73	121.21	105.74
2	B	401	SFG	N3-C4-N9	7.28	139.54	127.17
2	A	401	SFG	N3-C4-N9	6.85	138.82	127.17
2	D	401	SFG	N3-C4-N9	6.77	138.69	127.17
2	C	401	SFG	N3-C4-N9	6.63	138.44	127.17
2	B	401	SFG	C5-C4-N3	-6.25	118.11	126.72
2	D	401	SFG	C5-C4-N3	-5.78	118.76	126.72
2	A	401	SFG	C5-C4-N3	-5.74	118.81	126.72
2	C	401	SFG	C5-C4-N3	-5.63	118.97	126.72
2	C	401	SFG	C4-C5-N7	-5.54	104.25	110.58
2	D	401	SFG	C4-C5-N7	-5.51	104.28	110.58
2	A	401	SFG	C4-C5-N7	-5.44	104.36	110.58
2	B	401	SFG	C4-C5-N7	-5.44	104.36	110.58
2	C	401	SFG	N9-C8-N7	-4.64	107.36	113.94
2	A	401	SFG	N9-C8-N7	-4.56	107.47	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	SFG	N9-C8-N7	-4.56	107.47	113.94
2	B	401	SFG	N9-C8-N7	-4.42	107.67	113.94
2	C	401	SFG	C6-C5-N7	4.02	139.85	132.09
2	D	401	SFG	C6-C5-N7	3.92	139.65	132.09
2	A	401	SFG	N3-C2-N1	-3.92	122.65	128.58
2	A	401	SFG	C6-C5-N7	3.92	139.65	132.09
2	C	401	SFG	N3-C2-N1	-3.91	122.67	128.58
2	C	401	SFG	C4-N9-C1'	-3.89	117.53	126.63
2	B	401	SFG	N3-C2-N1	-3.87	122.72	128.58
2	D	401	SFG	N3-C2-N1	-3.85	122.75	128.58
2	B	401	SFG	C1'-N9-C8	-3.72	118.84	127.09
2	B	401	SFG	C6-C5-N7	3.67	139.17	132.09
2	D	401	SFG	C4-N9-C1'	-3.53	118.37	126.63
2	B	401	SFG	C2-N3-C4	3.46	120.29	111.83
2	A	401	SFG	C4-N9-C1'	-3.45	118.56	126.63
2	A	401	SFG	C2-N3-C4	3.34	119.98	111.83
2	D	401	SFG	C2-N3-C4	3.33	119.96	111.83
2	C	401	SFG	C2-N3-C4	3.32	119.93	111.83
2	A	401	SFG	C1'-N9-C8	-3.23	119.92	127.09
2	B	401	SFG	C5-C4-N9	-3.18	102.34	105.81
2	A	401	SFG	C5-C4-N9	-3.16	102.37	105.81
2	D	401	SFG	C1'-N9-C8	-3.05	120.33	127.09
2	D	401	SFG	C5-C4-N9	-2.99	102.56	105.81
2	C	401	SFG	C5-C4-N9	-2.95	102.60	105.81
2	B	401	SFG	C4-N9-C1'	-2.86	119.95	126.63
2	C	401	SFG	C1'-N9-C8	-2.74	121.02	127.09
2	B	401	SFG	C2'-C3'-C4'	2.53	107.49	102.61
2	A	401	SFG	OXT-C-O	-2.46	118.50	124.08
2	B	401	SFG	C3'-C2'-C1'	2.37	105.95	101.46
2	D	401	SFG	C2'-C3'-C4'	2.29	107.03	102.61
2	C	401	SFG	C2'-C3'-C4'	2.22	106.91	102.61
2	A	401	SFG	C2'-C3'-C4'	2.22	106.90	102.61
2	D	401	SFG	C3'-C2'-C1'	2.18	105.59	101.46
2	C	401	SFG	C3'-C2'-C1'	2.02	105.29	101.46

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	SFG	NE-CD-CG-CB
2	B	401	SFG	C5'-CD-CG-CB
2	A	401	SFG	NE-CD-CG-CB

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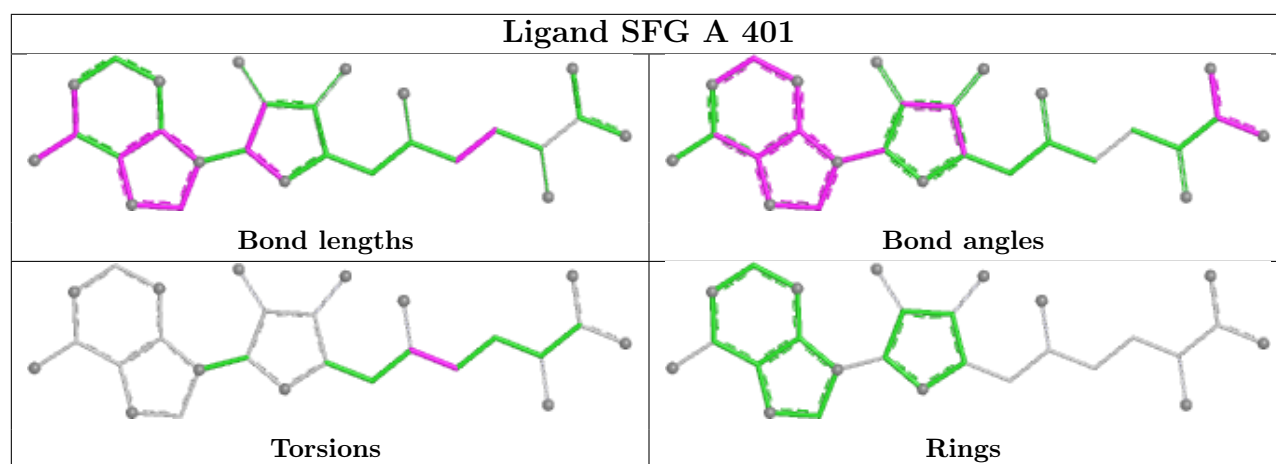
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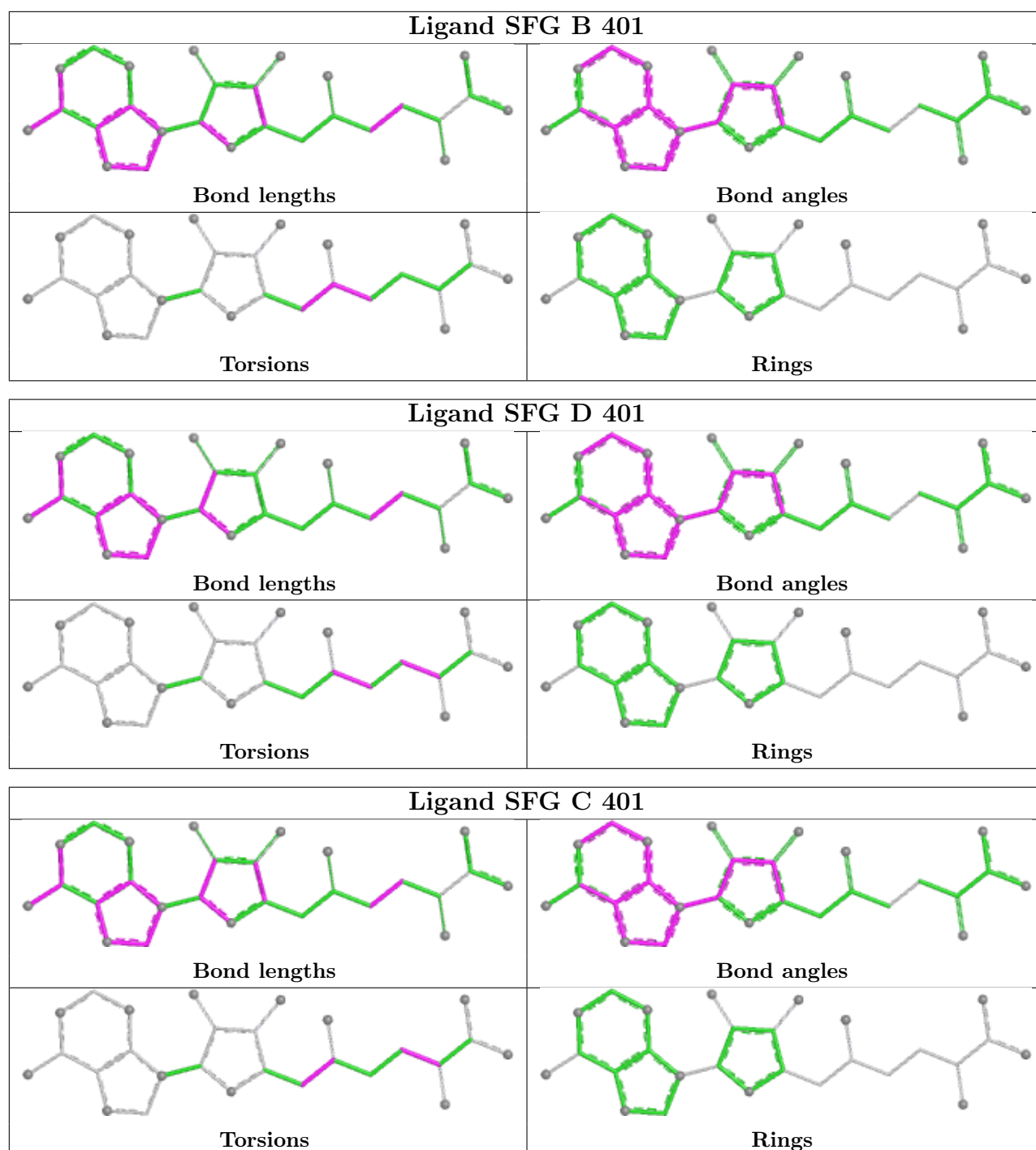
Mol	Chain	Res	Type	Atoms
2	A	401	SFG	C5'-CD-CG-CB
2	C	401	SFG	N-CA-CB-CG
2	D	401	SFG	NE-CD-CG-CB
2	D	401	SFG	C5'-CD-CG-CB
2	C	401	SFG	C-CA-CB-CG
2	C	401	SFG	C4'-C5'-CD-CG
2	B	401	SFG	C4'-C5'-CD-NE
2	D	401	SFG	N-CA-CB-CG

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	213/236 (90%)	1.10	27 (12%) 8 9	29, 42, 70, 112	0
1	B	213/236 (90%)	1.05	24 (11%) 10 11	28, 41, 75, 107	0
1	C	205/236 (86%)	1.27	45 (21%) 2 2	31, 45, 80, 112	0
1	D	204/236 (86%)	1.13	30 (14%) 6 7	30, 44, 74, 113	0
All	All	835/944 (88%)	1.14	126 (15%) 5 6	28, 43, 76, 113	0

All (126) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	70	ALA	5.3
1	D	242	MET	5.2
1	C	245	TRP	5.1
1	D	245	TRP	4.8
1	B	288	GLU	4.7
1	A	70	ALA	4.7
1	D	113	GLU	4.5
1	C	242	MET	4.4
1	C	113	GLU	4.2
1	C	201	TYR	4.2
1	D	201	TYR	4.1
1	B	240	HIS	4.1
1	B	239	ALA	4.1
1	A	201	TYR	4.0
1	C	111	LEU	4.0
1	C	200	GLU	3.9
1	B	231	ASP	3.8
1	B	113	GLU	3.8
1	A	289	GLU	3.8
1	C	282	TRP	3.7
1	C	115	GLN	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	285	ILE	3.7
1	A	285	ILE	3.7
1	A	245	TRP	3.5
1	C	109	GLY	3.5
1	B	285	ILE	3.5
1	D	133	SER	3.4
1	C	247	TYR	3.4
1	C	202	MET	3.3
1	D	110	LYS	3.2
1	B	114	VAL	3.2
1	A	239	ALA	3.2
1	D	111	LEU	3.2
1	C	231	ASP	3.1
1	D	173	ARG	3.1
1	D	112	GLY	3.0
1	C	107	ALA	3.0
1	A	203	PRO	3.0
1	C	205	ALA	2.9
1	C	117	ARG	2.9
1	C	203	PRO	2.9
1	C	101	LEU	2.9
1	D	114	VAL	2.9
1	C	110	LYS	2.8
1	B	245	TRP	2.8
1	D	134	CYS	2.8
1	A	79	ASP	2.7
1	C	213	LEU	2.7
1	B	171	GLY	2.7
1	A	107	ALA	2.7
1	A	241	LEU	2.6
1	D	177	LEU	2.6
1	A	231	ASP	2.6
1	C	105	TYR	2.6
1	B	143	HIS	2.6
1	C	209	LEU	2.6
1	B	193	VAL	2.6
1	D	140	SER	2.6
1	A	143	HIS	2.6
1	B	116	CYS	2.5
1	C	97	ARG	2.5
1	C	208	ARG	2.5
1	B	241	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	215	ALA	2.5
1	B	287	ARG	2.5
1	C	285	ILE	2.5
1	C	286	HIS	2.5
1	D	115	GLN	2.5
1	C	164	LEU	2.4
1	C	244	ASN	2.4
1	B	201	TYR	2.4
1	D	203	PRO	2.4
1	A	139	SER	2.4
1	D	275	SER	2.4
1	A	149	LEU	2.4
1	C	207	GLN	2.4
1	D	174	GLU	2.4
1	D	200	GLU	2.4
1	B	71	ARG	2.4
1	A	288	GLU	2.4
1	D	244	ASN	2.3
1	D	205	ALA	2.3
1	C	181	SER	2.3
1	C	204	SER	2.3
1	A	120	GLU	2.3
1	C	112	GLY	2.3
1	C	254	LEU	2.3
1	B	243	LYS	2.3
1	D	243	LYS	2.3
1	A	71	ARG	2.3
1	D	117	ARG	2.3
1	B	289	GLU	2.3
1	C	228	ILE	2.3
1	A	114	VAL	2.3
1	D	166	VAL	2.3
1	A	202	MET	2.3
1	C	251	ARG	2.3
1	A	113	GLU	2.3
1	A	262	LEU	2.2
1	C	100	ARG	2.2
1	C	114	VAL	2.2
1	D	219	LEU	2.2
1	C	166	VAL	2.2
1	B	280	GLN	2.2
1	A	240	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	273	ALA	2.2
1	D	123	PRO	2.1
1	C	283	ALA	2.1
1	D	247	TYR	2.1
1	A	159	LEU	2.1
1	C	252	ILE	2.1
1	C	278	LEU	2.1
1	A	171	GLY	2.1
1	B	183	ARG	2.1
1	B	179	GLU	2.1
1	B	200	GLU	2.1
1	B	244	ASN	2.1
1	D	157	ASP	2.1
1	A	287	ARG	2.1
1	C	248	SER	2.0
1	C	154	ALA	2.0
1	D	278	LEU	2.0
1	D	274	ILE	2.0
1	A	193	VAL	2.0
1	A	81	ILE	2.0
1	C	90	ILE	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	SFG	B	401	27/27	0.86	0.11	26,34,41,42	0
2	SFG	C	401	27/27	0.87	0.13	30,41,51,53	0

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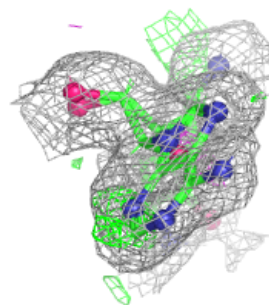
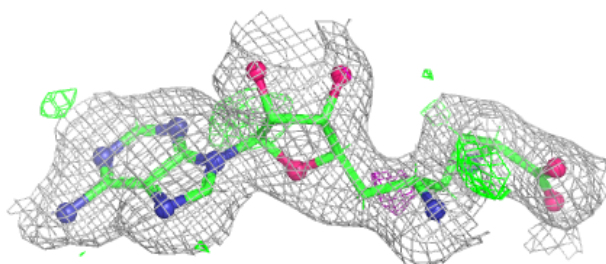
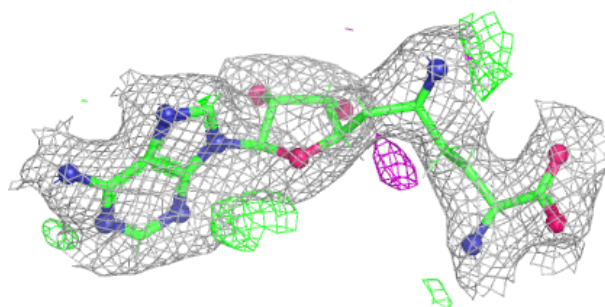
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SFG	D	401	27/27	0.91	0.10	27,37,48,49	0
2	SFG	A	401	27/27	0.92	0.09	27,36,45,47	0

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.

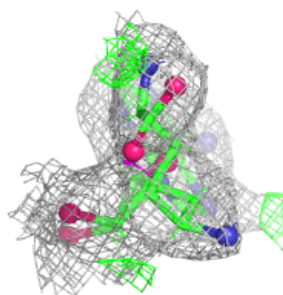
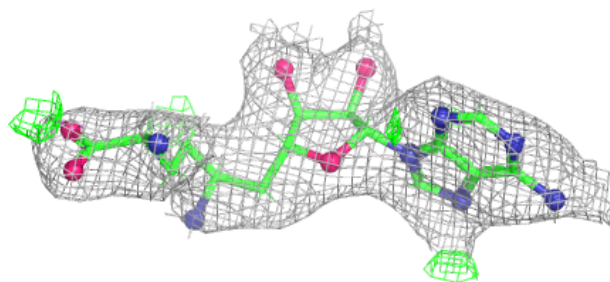
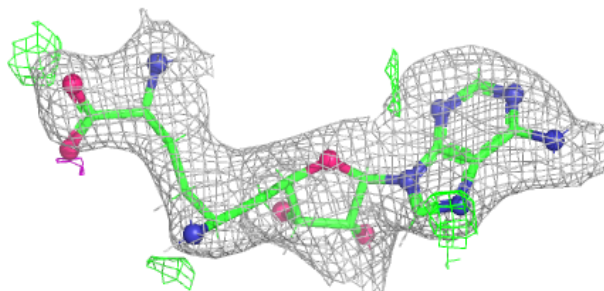
Electron density around SFG B 401:

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

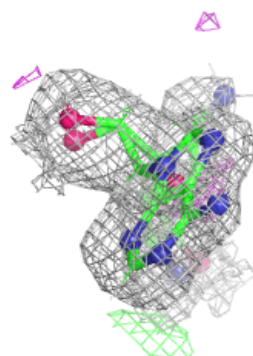
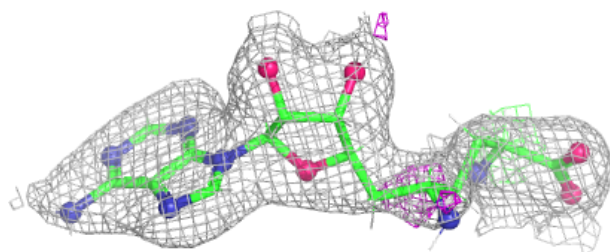
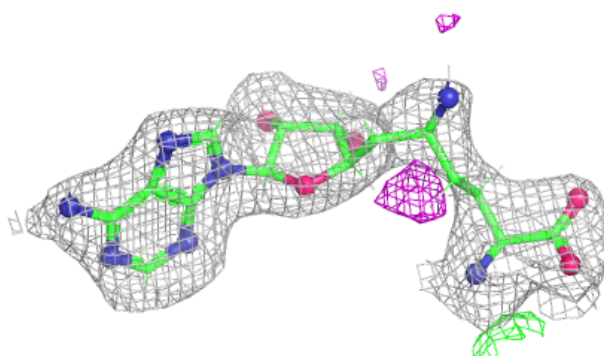


Electron density around SFG C 401:

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

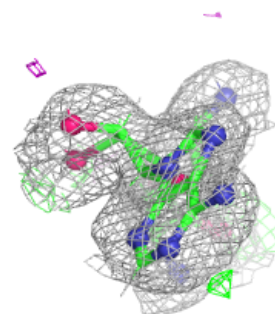
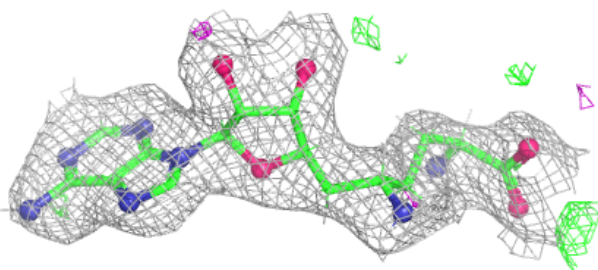
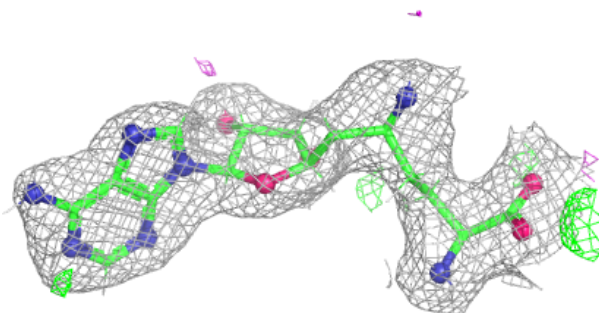
**Electron density around SFG D 401:**

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



Electron density around SFG A 401:

$2mF_o-DF_c$ (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.