



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:54 PM UTC

PDB ID : 5R0G / pdb_00005r0g
Title : PanDDA analysis group deposition – Aar2/RNaseH in complex with fragment F2X-Entry D07, DMSO-free
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Deposited on : 2020-02-12
Resolution : 1.73 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

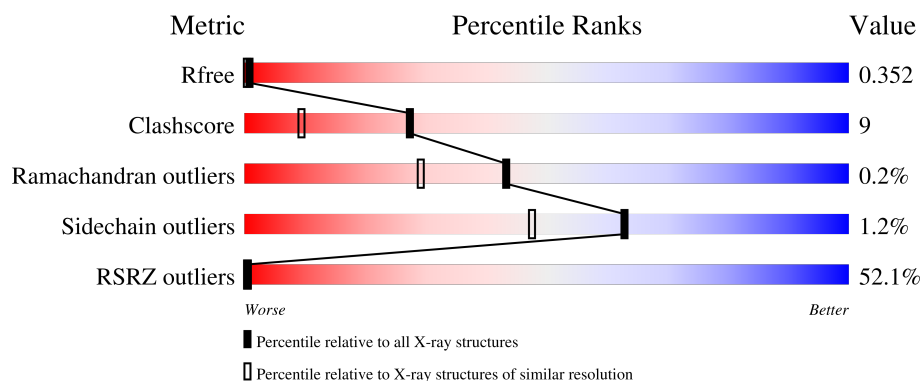
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.73 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	1187 (1.74-1.74)
Clashscore	190562	1207 (1.74-1.74)
Ramachandran outliers	187476	1200 (1.74-1.74)
Sidechain outliers	187428	1200 (1.74-1.74)
RSRZ outliers	180081	1188 (1.74-1.74)

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	258	48% 73% 16% • • 8%
2	B	308	51% 81% 17% •

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SYJ	A	2501	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4675 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 Pre-mRNA-splicing factor 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	237	1995	1278	334	371	12	0	12	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1833	GLY	-	expression tag	UNP P33334
A	1834	ALA	-	expression tag	UNP P33334
A	1835	MET	-	expression tag	UNP P33334

- Molecule 2 is a protein called A1 cistron-splicing factor AAR2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	300	2580	1654	421	485	20	0	9	0

There are 20 discrepancies between the modelled and reference sequences:

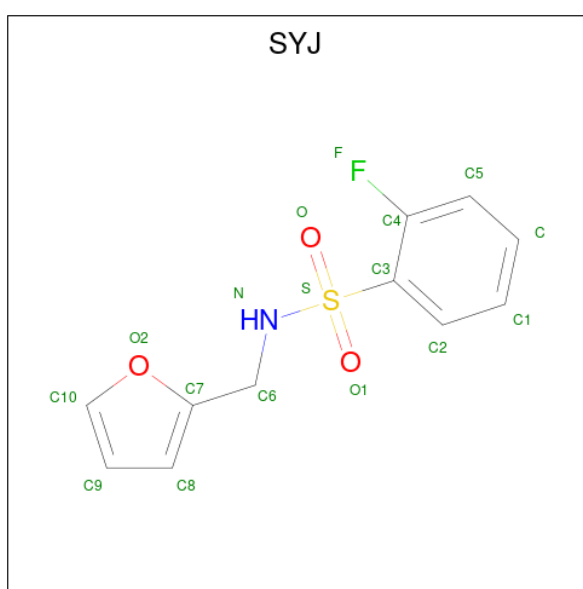
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P32357
B	-2	ALA	-	expression tag	UNP P32357
B	-1	MET	-	expression tag	UNP P32357
B	0	ALA	-	expression tag	UNP P32357
B	166	SER	LEU	conflict	UNP P32357
B	167	SER	LYS	conflict	UNP P32357
B	170	SER	LEU	conflict	UNP P32357
B	?	-	GLN	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	GLY	deletion	UNP P32357
B	?	-	SER	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	MET	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ASN	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	?	-	ASP	deletion	UNP P32357

- Molecule 3 is 2-fluoranyl- {N}-(furan-2-ylmethyl)benzenesulfonamide (CCD ID: SYJ) (formula: C₁₁H₁₀FNO₃S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	S	0	0
			17	11	1	1	3	1		

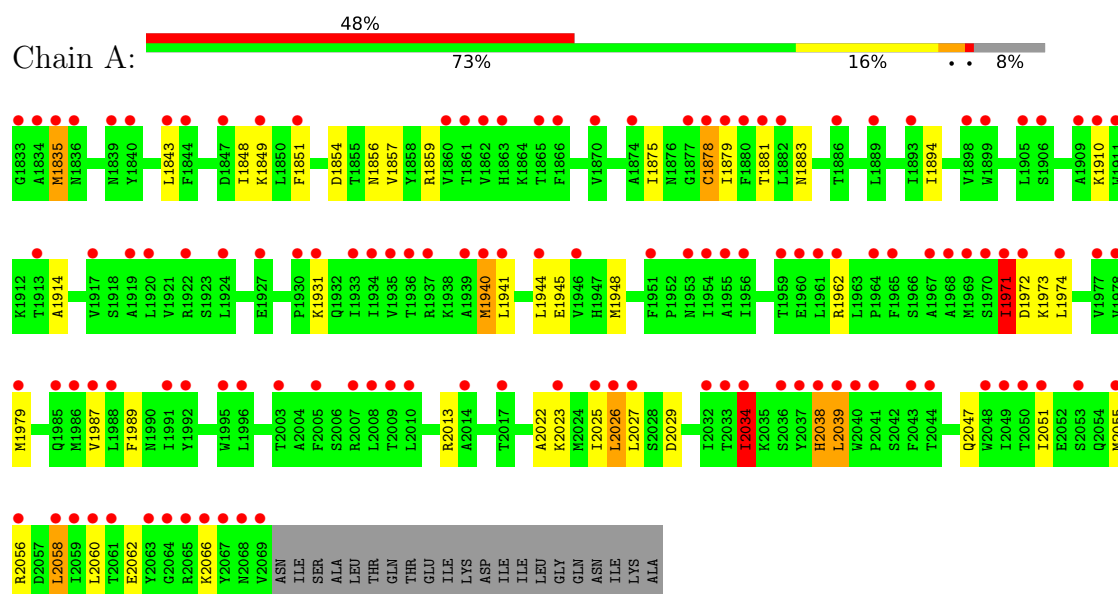
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	50	Total	O	0	0
			50	50		
4	B	33	Total	O	0	0
			33	33		

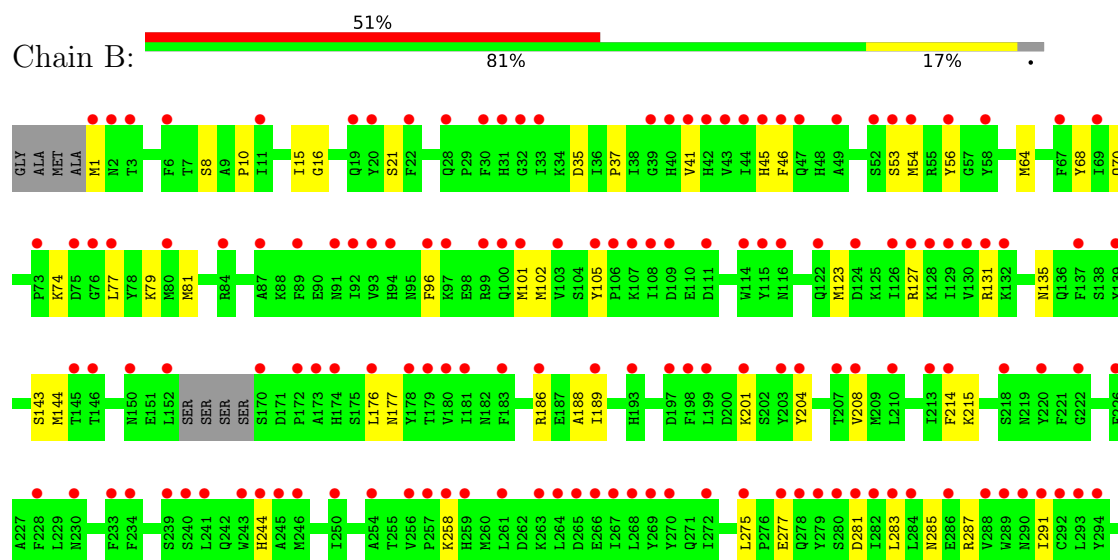
3 Residue-property plots

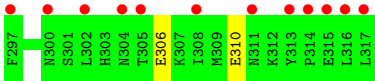
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Pre-mRNA-splicing factor 8



• Molecule 2: A1 cistron-splicing factor AAR2





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.20Å 82.30Å 93.11Å 90.00° 108.08° 90.00°	Depositor
Resolution (Å)	21.22 – 1.73 21.22 – 1.73	Depositor EDS
% Data completeness (in resolution range)	97.9 (21.22-1.73) 98.2 (21.22-1.73)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 1.73Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.289 , 0.340 0.304 , 0.352	Depositor DCC
R_{free} test set	2100 reflections (3.17%)	wwPDB-VP
Wilson B-factor (Å ²)	42.6	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4675	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% 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: SYJ

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	0.95	4/2042 (0.2%)	1.10	14/2765 (0.5%)
2	B	0.69	0/2651	0.82	0/3581
All	All	0.81	4/4693 (0.1%)	0.95	14/6346 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1972	ASP	C-O	-19.97	1.00	1.24
1	A	1971	ILE	C-O	-8.30	1.14	1.24
1	A	1883	ASN	CB-CG	5.52	1.65	1.52
1	A	2038	HIS	C-O	-5.29	1.17	1.23

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1940	MET	CG-SD-CE	-12.48	73.44	100.90
1	A	1883	ASN	CB-CG-ND2	-8.00	104.41	116.40
1	A	2034	ILE	CB-CA-C	6.78	120.86	110.96
1	A	2026	LEU	CB-CA-C	6.21	123.06	110.38
1	A	1972	ASP	CA-C-O	-6.17	114.00	120.92
1	A	2038	HIS	CB-CA-C	-6.17	97.14	109.35
1	A	1878	CYS	N-CA-C	5.97	118.19	108.76
1	A	1973	LYS	CA-C-O	-5.78	114.75	120.82
1	A	2034	ILE	N-CA-C	-5.66	99.60	107.80
1	A	2034	ILE	N-CA-CB	-5.63	104.36	111.46
1	A	2022	ALA	O-C-N	-5.55	116.35	122.07
1	A	2025	ILE	CA-C-N	-5.20	111.83	120.68
1	A	2025	ILE	C-N-CA	-5.20	111.83	120.68
1	A	1973	LYS	CB-CA-C	5.09	118.87	110.88

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	1995	0	2020	42	0
2	B	2580	0	2450	40	0
3	A	17	0	0	8	0
4	A	50	0	0	3	0
4	B	33	0	0	4	0
All	All	4675	0	4470	83	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 9.

All (83) 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:2026:LEU:HD11	1:A:2055:MET:CE	1.60	1.30
1:A:2023:LYS:HE2	3:A:2501:SYJ:C6	1.80	1.10
1:A:2026:LEU:HD11	1:A:2055:MET:HE2	1.11	1.08
1:A:2026:LEU:CD1	1:A:2055:MET:CE	2.31	1.07
1:A:2026:LEU:CD1	1:A:2055:MET:HE2	1.91	0.99
1:A:2023:LYS:CE	3:A:2501:SYJ:C6	2.55	0.84
2:B:96:PHE:HB2	2:B:102:MET:HE3	1.62	0.81
1:A:2026:LEU:CD1	1:A:2055:MET:HE1	2.15	0.76
1:A:1848:ILE:H	1:A:1931[A]:LYS:HZ2	1.36	0.71
2:B:96:PHE:CB	2:B:102:MET:HE3	2.19	0.70
2:B:70:GLN:HB3	2:B:81:MET:HE2	1.77	0.65
1:A:2038:HIS:CE1	4:A:2613:HOH:O	2.51	0.62
1:A:2062:GLU:O	1:A:2066:LYS:HG2	2.00	0.60
1:A:2038:HIS:HE1	4:A:2613:HOH:O	1.84	0.59
1:A:1962:ARG:O	1:A:2013:ARG:NH1	2.29	0.59
2:B:74:LYS:NZ	4:B:401:HOH:O	2.35	0.59
1:A:2026:LEU:HD11	1:A:2055:MET:HE3	1.76	0.59
2:B:54[B]:MET:HE1	2:B:143:SER:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1979[B]:MET:HA	1:A:1979[B]:MET:HE2	1.85	0.58
2:B:188:ALA:HA	2:B:204:TYR:CD1	2.39	0.58
1:A:1987:VAL:HG12	1:A:1989:PHE:CE1	2.39	0.58
2:B:306:GLU:O	2:B:310:GLU:HG3	2.05	0.56
3:A:2501:SYJ:C6	3:A:2501:SYJ:C2	2.83	0.56
1:A:1910:LYS:HG2	1:A:1940:MET:SD	2.46	0.55
2:B:144:MET:HE2	2:B:176:LEU:HG	1.88	0.54
1:A:2034:ILE:HG13	3:A:2501:SYJ:C5	2.38	0.54
1:A:1941:LEU:O	1:A:1945:GLU:HB2	2.08	0.53
2:B:186:ARG:NH1	2:B:189:ILE:O	2.41	0.52
2:B:1:MET:N	4:B:402:HOH:O	2.41	0.52
1:A:1843:LEU:HA	1:A:1849:LYS:HD2	1.92	0.52
1:A:2047:GLN:O	1:A:2051:ILE:HG12	2.10	0.51
1:A:2056[B]:ARG:O	1:A:2060:LEU:HG	2.11	0.51
2:B:201:LYS:NZ	4:B:403:HOH:O	2.42	0.51
2:B:70:GLN:HB3	2:B:81:MET:CE	2.40	0.51
2:B:277:GLU:CD	2:B:277:GLU:H	2.20	0.50
1:A:1875:ILE:HB	1:A:1979[C]:MET:HE1	1.94	0.49
2:B:68:TYR:CE2	2:B:81:MET:HE3	2.48	0.49
2:B:123:MET:O	2:B:127:ARG:HG3	2.13	0.49
1:A:1856:ASN:O	1:A:1878:CYS:N	2.37	0.48
1:A:2062:GLU:HB3	1:A:2066:LYS:HE3	1.95	0.48
2:B:214:PHE:O	2:B:215:LYS:HB2	2.12	0.48
1:A:1854:ASP:OD2	1:A:1940:MET:SD	2.71	0.48
2:B:15:ILE:O	2:B:21:SER:HA	2.12	0.48
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.38	0.48
1:A:1857:VAL:HG22	1:A:1879:ILE:HG13	1.96	0.47
1:A:1859:ARG:HH12	1:A:1979[A]:MET:CE	2.27	0.47
1:A:1971:ILE:HD12	1:A:1974:LEU:HD12	1.97	0.47
2:B:244:HIS:CD2	2:B:285:ASN:HB3	2.50	0.47
1:A:2023:LYS:CG	3:A:2501:SYJ:C8	2.93	0.47
1:A:2023:LYS:HG3	3:A:2501:SYJ:C8	2.46	0.46
2:B:77:LEU:HD21	2:B:79:LYS:HE3	1.96	0.46
2:B:144:MET:HE2	2:B:176:LEU:CD2	2.45	0.46
2:B:8:SER:C	2:B:10:PRO:HD3	2.41	0.46
1:A:2039:LEU:HD23	1:A:2039:LEU:N	2.30	0.46
2:B:37:PRO:HD3	2:B:105:TYR:CD1	2.51	0.46
2:B:204:TYR:O	2:B:208:VAL:HB	2.15	0.46
2:B:131:ARG:HH21	2:B:177[B]:ASN:ND2	2.14	0.46
1:A:2034:ILE:HG21	3:A:2501:SYJ:F	2.06	0.46
2:B:258:LYS:HD2	2:B:258:LYS:H	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1944:LEU:O	1:A:1948:MET:HG2	2.16	0.45
2:B:287:ARG:O	2:B:291:ILE:HD13	2.16	0.45
2:B:64[A]:MET:SD	2:B:123:MET:HG2	2.57	0.44
2:B:188:ALA:HB2	2:B:204:TYR:CZ	2.52	0.44
1:A:1848:ILE:H	1:A:1931[A]:LYS:NZ	2.12	0.44
1:A:1987:VAL:CG1	1:A:1989:PHE:CE1	3.02	0.43
2:B:46:PHE:HB2	2:B:56:TYR:CZ	2.54	0.43
1:A:1835:MET:HE3	1:A:1835:MET:HB3	1.94	0.42
2:B:16:GLY:HA3	2:B:45:HIS:CE1	2.54	0.42
1:A:1851:PHE:O	1:A:1881:THR:HA	2.19	0.42
2:B:41[A]:VAL:HG12	4:B:410:HOH:O	2.19	0.42
2:B:127:ARG:NH1	2:B:135:ASN:O	2.49	0.42
2:B:275:LEU:CD2	2:B:283:LEU:HD13	2.50	0.42
2:B:275:LEU:HD22	2:B:283:LEU:HD13	2.02	0.42
2:B:281:ASP:OD1	2:B:281:ASP:N	2.42	0.42
1:A:1914:ALA:HB1	1:A:1944:LEU:HA	2.02	0.41
2:B:1:MET:HB3	2:B:35:ASP:HA	2.02	0.41
2:B:96:PHE:HB3	2:B:102:MET:HE3	1.99	0.41
1:A:1894:ILE:HA	4:A:2623:HOH:O	2.21	0.41
1:A:2023:LYS:HG2	3:A:2501:SYJ:C8	2.50	0.41
1:A:2029:ASP:OD1	1:A:2029:ASP:C	2.64	0.41
2:B:54[B]:MET:HE1	2:B:143:SER:CB	2.51	0.41
2:B:244:HIS:HE1	2:B:283:LEU:O	2.04	0.40
2:B:53:SER:O	2:B:54[A]:MET:HB3	2.21	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	247/258 (96%)	239 (97%)	7 (3%)	1 (0%)	30 16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	306/308 (99%)	290 (95%)	16 (5%)	0	100	100
All	All	553/566 (98%)	529 (96%)	23 (4%)	1 (0%)	43	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1971	ILE

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	225/233 (97%)	220 (98%)	5 (2%)	45	22
2	B	287/284 (101%)	286 (100%)	1 (0%)	86	82
All	All	512/517 (99%)	506 (99%)	6 (1%)	63	47

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1835	MET
1	A	2027	LEU
1	A	2034	ILE
1	A	2039	LEU
1	A	2058	LEU
2	B	101	MET

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

Mol	Chain	Res	Type
2	B	2	ASN
2	B	47	GLN

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 ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SYJ	A	2501	-	18,18,18	0.68	0	25,25,25	1.62	2 (8%)

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	SYJ	A	2501	-	-	5/12/12/12	0/2/2/2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2501	SYJ	C4-C3-S	5.47	124.71	121.21
3	A	2501	SYJ	C7-C6-N	5.36	124.48	112.66

There are no chirality outliers.

All (5) torsion outliers are listed below:

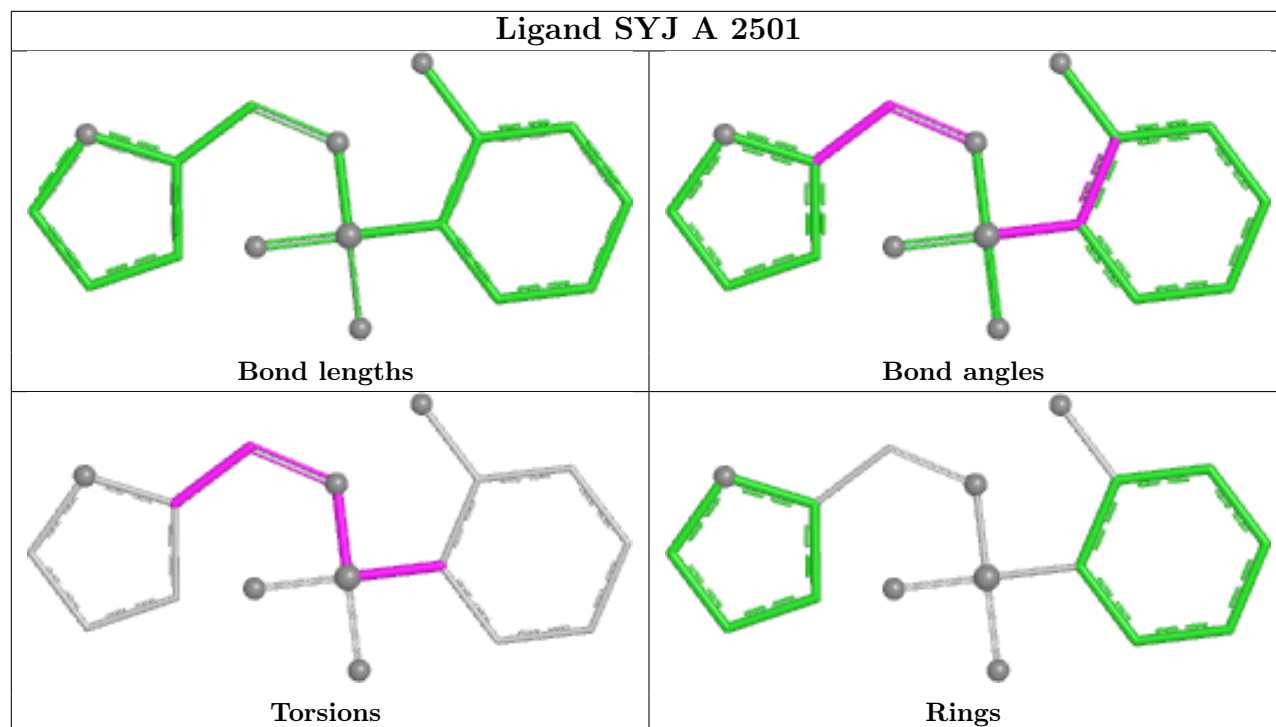
Mol	Chain	Res	Type	Atoms
3	A	2501	SYJ	C7-C6-N-S
3	A	2501	SYJ	C4-C3-S-N
3	A	2501	SYJ	C2-C3-S-N
3	A	2501	SYJ	N-C6-C7-C8
3	A	2501	SYJ	C6-N-S-O

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2501	SYJ	8	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	237/258 (91%)	2.50	124 (52%) 0 0	28, 74, 139, 198	11 (4%)
2	B	300/308 (97%)	2.32	156 (52%) 0 0	30, 78, 140, 220	9 (3%)
All	All	537/566 (94%)	2.40	280 (52%) 0 0	28, 76, 139, 220	20 (3%)

All (280) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	203[A]	TYR	13.1
1	A	2039	LEU	9.9
2	B	22	PHE	9.1
1	A	2034	ILE	8.5
1	A	1840	TYR	8.1
1	A	1877	GLY	8.0
1	A	2060	LEU	8.0
1	A	2026	LEU	7.8
1	A	1940	MET	7.6
2	B	54[A]	MET	7.1
1	A	1878	CYS	6.7
2	B	243	TRP	6.3
2	B	316	LEU	6.0
1	A	2040	TRP	6.0
1	A	1954	ILE	5.9
2	B	208	VAL	5.7
2	B	313	TYR	5.7
1	A	2038	HIS	5.6
1	A	1979[A]	MET	5.6
1	A	1969	MET	5.6
1	A	2058	LEU	5.5
2	B	267	ILE	5.4
2	B	268	LEU	5.2
1	A	1939	ALA	5.2

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Mol	Chain	Res	Type	RSRZ
2	B	31	HIS	5.2
2	B	283	LEU	5.2
2	B	89	PHE	5.1
1	A	2017[A]	THR	5.1
1	A	1927	GLU	5.0
1	A	1956	ILE	5.0
2	B	279	TYR	5.0
2	B	317	LEU	4.9
1	A	1880	PHE	4.8
1	A	1835	MET	4.8
1	A	1988	LEU	4.7
2	B	92	ILE	4.6
1	A	1934	ILE	4.6
2	B	174	HIS	4.6
2	B	52	SER	4.6
1	A	1889	LEU	4.5
2	B	264	LEU	4.5
1	A	1955	ALA	4.5
1	A	1866	PHE	4.5
1	A	1860	VAL	4.4
1	A	1935	VAL	4.4
2	B	275	LEU	4.4
2	B	53	SER	4.4
1	A	1977	VAL	4.3
1	A	1930	PRO	4.2
2	B	11	ILE	4.2
2	B	126	ILE	4.2
1	A	2010	LEU	4.2
1	A	1971	ILE	4.2
2	B	246	MET	4.1
2	B	130	VAL	4.1
1	A	1879	ILE	4.1
1	A	1944	LEU	4.1
2	B	129	ILE	4.1
1	A	2043	PHE	4.0
2	B	6	PHE	4.0
1	A	2059	ILE	4.0
1	A	2027	LEU	4.0
2	B	137	PHE	4.0
1	A	1834	ALA	4.0
1	A	2009	THR	3.8
1	A	1905	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
2	B	261	LEU	3.8
2	B	20	TYR	3.8
2	B	170	SER	3.8
2	B	189	ILE	3.8
2	B	67	PHE	3.8
2	B	294	TYR	3.8
2	B	291	ILE	3.7
1	A	1968	ALA	3.7
2	B	96	PHE	3.7
1	A	1862	VAL	3.7
1	A	2069	VAL	3.7
1	A	1886	THR	3.7
2	B	3	THR	3.7
2	B	1	MET	3.7
2	B	101	MET	3.7
1	A	1913	THR	3.6
2	B	250	ILE	3.6
1	A	1861	THR	3.6
2	B	288	VAL	3.6
2	B	259	HIS	3.6
1	A	2037	TYR	3.6
1	A	2068	ASN	3.5
2	B	284	LEU	3.5
2	B	228	PHE	3.5
2	B	193	HIS	3.5
1	A	1882	LEU	3.5
1	A	1924	LEU	3.5
2	B	281	ASP	3.5
2	B	282	ILE	3.5
2	B	226	PHE	3.5
2	B	234	PHE	3.4
1	A	1851	PHE	3.4
1	A	1961	LEU	3.4
1	A	1881	THR	3.4
2	B	32	GLY	3.4
1	A	1936	THR	3.4
1	A	2025	ILE	3.3
2	B	44[A]	ILE	3.3
2	B	77	LEU	3.3
2	B	199	LEU	3.3
1	A	1844	PHE	3.3
1	A	1898[A]	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	269	TYR	3.3
2	B	145	THR	3.3
1	A	1911	TRP	3.3
1	A	2066	LYS	3.3
1	A	2065	ARG	3.2
1	A	1899	TRP	3.2
2	B	49	ALA	3.2
1	A	1920	LEU	3.2
2	B	152	LEU	3.2
1	A	1995	TRP	3.2
1	A	1833	GLY	3.2
1	A	1965	PHE	3.2
2	B	30	PHE	3.2
2	B	100	GLN	3.1
1	A	1946	VAL	3.1
2	B	84	ARG	3.1
1	A	2067	TYR	3.1
2	B	204	TYR	3.1
1	A	1941	LEU	3.1
1	A	1870	VAL	3.1
1	A	1917	VAL	3.1
1	A	2063	TYR	3.1
1	A	2051	ILE	3.1
2	B	207	THR	3.1
2	B	256	VAL	3.0
2	B	76	GLY	3.0
2	B	308	ILE	3.0
1	A	1964	PRO	3.0
1	A	1978	VAL	3.0
2	B	91	ASN	3.0
2	B	114	TRP	3.0
1	A	2049	ILE	3.0
2	B	73	PRO	3.0
1	A	2005	PHE	3.0
2	B	214	PHE	3.0
1	A	2061	THR	2.9
2	B	210	LEU	2.9
2	B	302	LEU	2.9
2	B	56	TYR	2.9
2	B	222	GLY	2.9
1	A	1986	MET	2.9
2	B	244	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	241	LEU	2.9
1	A	2032	ILE	2.9
1	A	2041	PRO	2.9
2	B	109	ASP	2.9
2	B	257	PRO	2.9
1	A	1836	ASN	2.9
2	B	254	ALA	2.8
2	B	139	TYR	2.8
2	B	240	SER	2.8
2	B	314	PRO	2.8
2	B	289	TRP	2.8
2	B	198	PHE	2.8
2	B	293	LEU	2.8
1	A	1972	ASP	2.8
1	A	1863	HIS	2.8
2	B	183	PHE	2.8
2	B	300	ASN	2.8
2	B	115	TYR	2.8
1	A	2008	LEU	2.8
1	A	1933	ILE	2.8
1	A	1970	SER	2.8
2	B	131	ARG	2.7
1	A	1953	ASN	2.7
2	B	272	ILE	2.7
2	B	40	HIS	2.7
2	B	2	ASN	2.7
2	B	245	ALA	2.7
1	A	1910	LYS	2.7
1	A	1839	ASN	2.6
1	A	2023	LYS	2.6
1	A	1843	LEU	2.6
2	B	106	PRO	2.6
2	B	107	LYS	2.6
1	A	2055	MET	2.6
2	B	46	PHE	2.6
2	B	176	LEU	2.6
1	A	1991	ILE	2.6
1	A	2048	TRP	2.6
2	B	45	HIS	2.6
2	B	43	VAL	2.6
1	A	2064	GLY	2.5
2	B	80	MET	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	93	VAL	2.5
1	A	1959	THR	2.5
1	A	1974	LEU	2.5
2	B	103	VAL	2.5
2	B	181	ILE	2.5
1	A	2056[A]	ARG	2.5
2	B	178	TYR	2.5
2	B	270	TYR	2.5
2	B	258	LYS	2.5
1	A	2033	THR	2.5
1	A	2036	SER	2.5
2	B	99	ARG	2.4
2	B	41[A]	VAL	2.4
2	B	290	ASN	2.4
2	B	278	GLN	2.4
2	B	124	ASP	2.4
2	B	197	ASP	2.4
1	A	1922	ARG	2.4
2	B	230[A]	ASN	2.4
1	A	1865	THR	2.4
2	B	19	GLN	2.4
1	A	1987	VAL	2.4
2	B	69	ILE	2.4
2	B	263	LYS	2.4
2	B	47	GLN	2.4
2	B	218	SER	2.4
1	A	1960[A]	GLU	2.4
2	B	186	ARG	2.4
2	B	220	TYR	2.4
1	A	1906	SER	2.3
1	A	1909	ALA	2.3
2	B	87	ALA	2.3
2	B	146	THR	2.3
2	B	233	PHE	2.3
2	B	28	GLN	2.3
2	B	122[A]	GLN	2.3
2	B	280	SER	2.3
2	B	286	GLU	2.3
2	B	179	THR	2.3
1	A	1874	ALA	2.3
2	B	315	GLU	2.3
2	B	127	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	1847	ASP	2.3
1	A	2050	THR	2.3
2	B	94	HIS	2.3
2	B	173	ALA	2.3
2	B	277	GLU	2.3
1	A	1849	LYS	2.3
1	A	2003	THR	2.2
1	A	2044	THR	2.2
2	B	97	LYS	2.2
2	B	33	ILE	2.2
2	B	108	ILE	2.2
1	A	2014	ALA	2.2
1	A	2053[A]	SER	2.2
2	B	201	LYS	2.2
2	B	75	ASP	2.2
2	B	39	GLY	2.2
1	A	2007	ARG	2.2
2	B	105	TYR	2.2
2	B	304	ASN	2.2
1	A	1919	ALA	2.2
1	A	1985	GLN	2.2
1	A	1996	LEU	2.2
2	B	297	PHE	2.2
2	B	132	LYS	2.2
1	A	1893	ILE	2.2
2	B	172	PRO	2.2
2	B	305	THR	2.2
1	A	1937	ARG	2.1
2	B	311	ASN	2.1
2	B	111	ASP	2.1
2	B	239	SER	2.1
2	B	150	ASN	2.1
2	B	266	GLU	2.1
1	A	1951	PHE	2.1
2	B	58	TYR	2.1
1	A	1967	ALA	2.1
2	B	180	VAL	2.1
2	B	116	ASN	2.1
2	B	213	ILE	2.1
1	A	1931[A]	LYS	2.0
2	B	128	LYS	2.0
1	A	1962	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	1992	TYR	2.0
2	B	265	ASP	2.0
2	B	292	CYS	2.0
2	B	42	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](#)

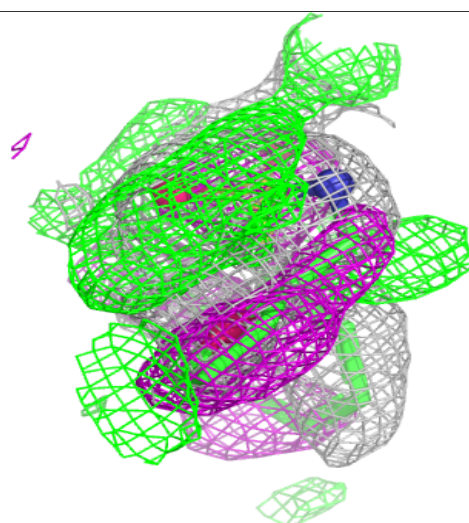
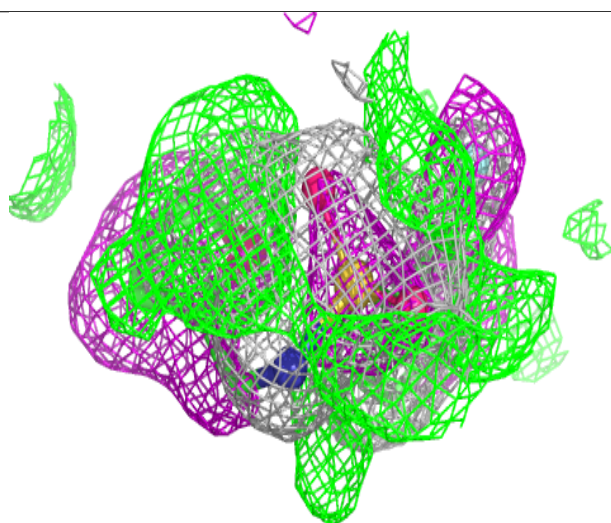
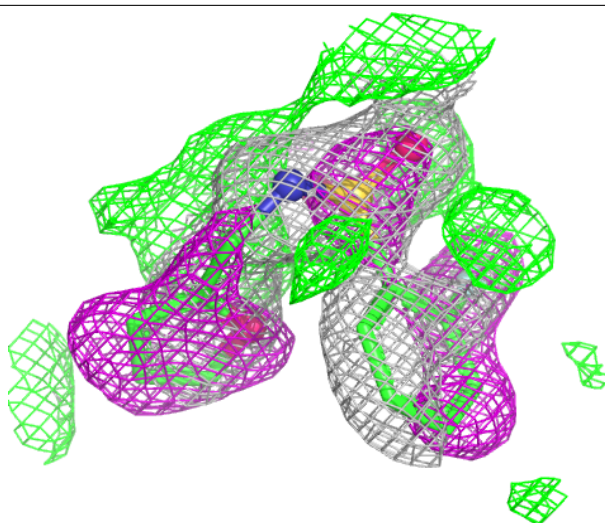
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	SYJ	A	2501	17/17	0.76	0.17	20,20,20,20	17

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 SYJ A 2501:

$2mF_o-DF_c$ (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.