



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 19, 2024 – 02:08 PM JST

PDB ID : 8K9W
Title : Crystal structure of plasmodium LysRS complexing with ASP3026 derived LysRS inhibitor 4 (ADKI4)
Authors : Zhou, J.; Xia, M.; Yang, G.; Li, P.; Fang, P.
Deposited on : 2023-08-01
Resolution : 2.30 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 : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.37.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

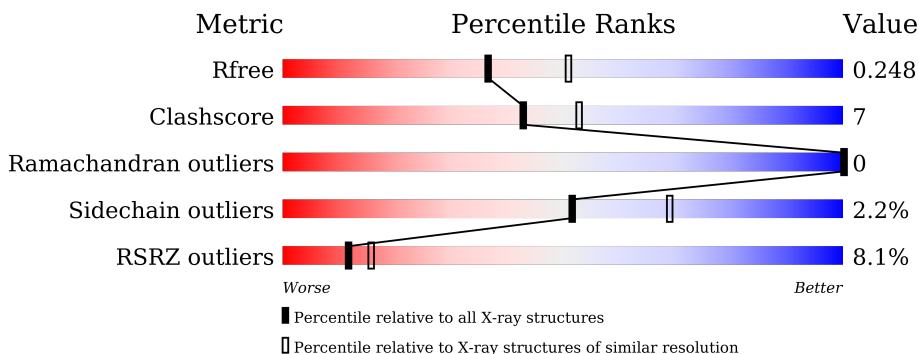
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	516	
1	B	516	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 7727 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.

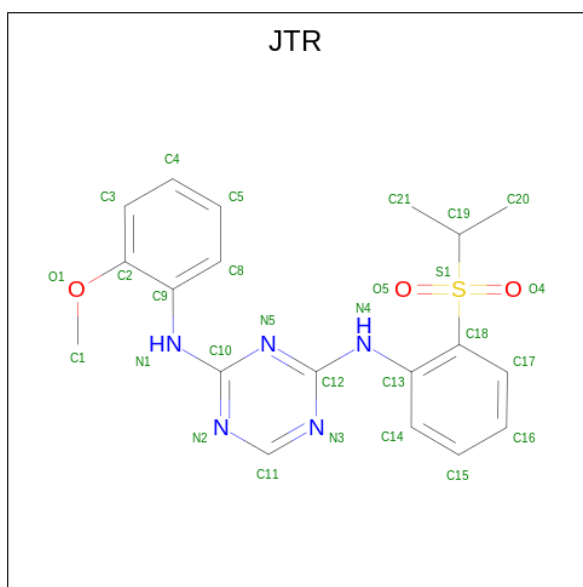
- Molecule 1 is a protein called Lysine-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	476	3741	2419	620	687	15	0	0	0
1	B	482	3770	2436	626	691	17	0	0	0

There are 18 discrepancies between the modelled and reference sequences:

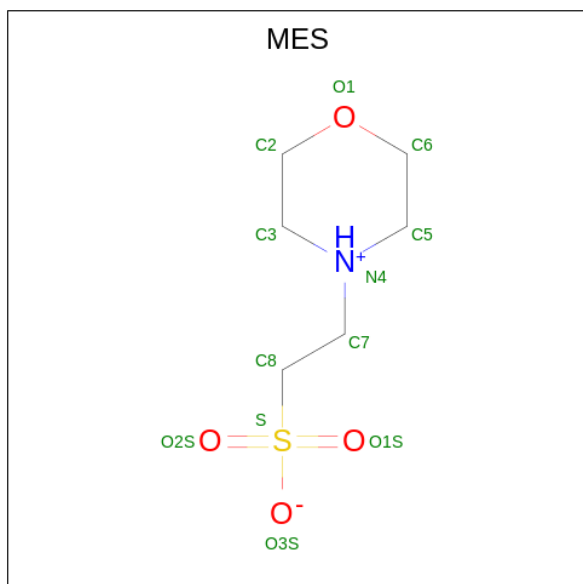
Chain	Residue	Modelled	Actual	Comment	Reference
A	76	MET	-	initiating methionine	UNP A0A024X378
A	584	GLY	-	expression tag	UNP A0A024X378
A	585	GLY	-	expression tag	UNP A0A024X378
A	586	HIS	-	expression tag	UNP A0A024X378
A	587	HIS	-	expression tag	UNP A0A024X378
A	588	HIS	-	expression tag	UNP A0A024X378
A	589	HIS	-	expression tag	UNP A0A024X378
A	590	HIS	-	expression tag	UNP A0A024X378
A	591	HIS	-	expression tag	UNP A0A024X378
B	76	MET	-	initiating methionine	UNP A0A024X378
B	584	GLY	-	expression tag	UNP A0A024X378
B	585	GLY	-	expression tag	UNP A0A024X378
B	586	HIS	-	expression tag	UNP A0A024X378
B	587	HIS	-	expression tag	UNP A0A024X378
B	588	HIS	-	expression tag	UNP A0A024X378
B	589	HIS	-	expression tag	UNP A0A024X378
B	590	HIS	-	expression tag	UNP A0A024X378
B	591	HIS	-	expression tag	UNP A0A024X378

- Molecule 2 is {N}2-(2-methoxyphenyl)- {N}4-(2-propan-2-ylsulfonylphenyl)-1,3,5-triazine-2,4-diamine (three-letter code: JTR) (formula: C₁₉H₂₁N₅O₃S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	28	19	5	3	1	0	0
2	B	1	28	19	5	3	1	0	0

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	12	6	1	4	1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	12	6	1	4	1	0	0

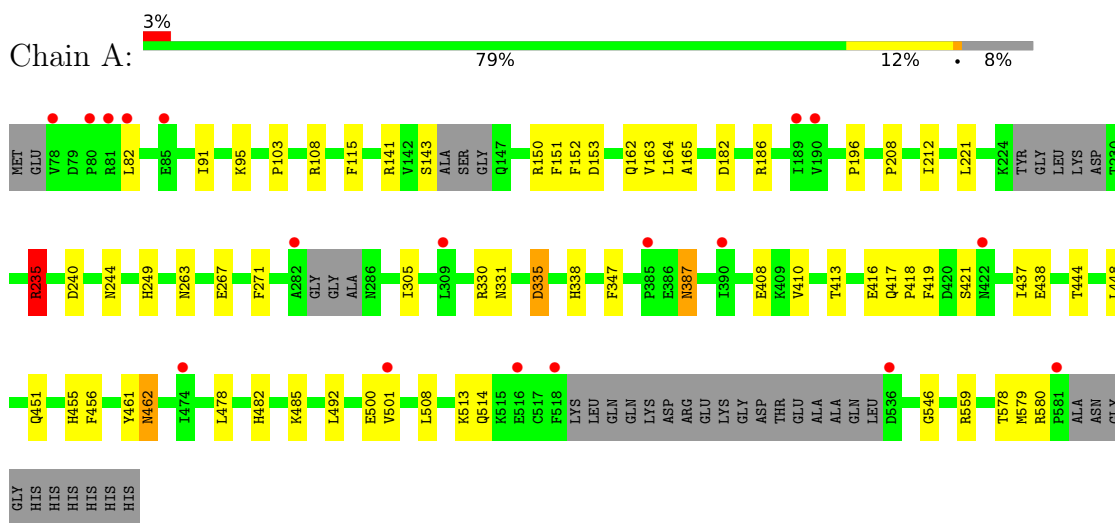
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	61	Total	O	0	0
			61	61		
4	B	75	Total	O	0	0
			75	75		

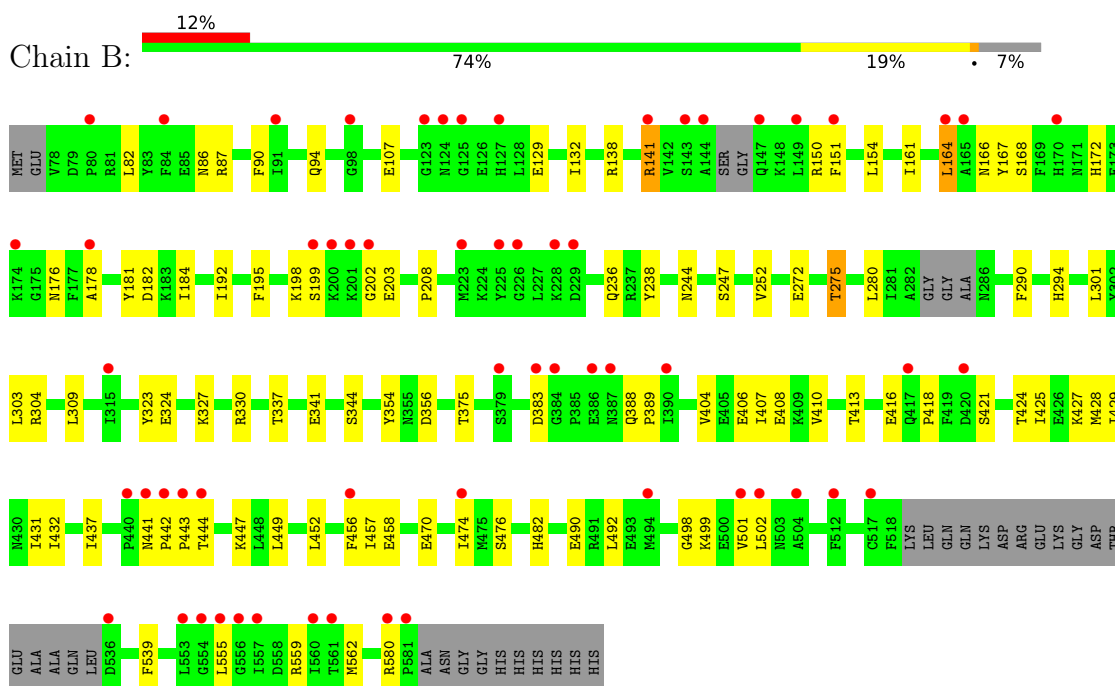
3 Residue-property plots i

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: Lysine-tRNA ligase



- Molecule 1: Lysine-tRNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.92Å 95.55Å 168.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.45 – 2.30 32.45 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (32.45-2.30) 99.9 (32.45-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.29Å)	Xtrriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.214 , 0.251 0.214 , 0.248	Depositor DCC
R_{free} test set	2000 reflections (3.83%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtrriage
Anisotropy	0.362	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 52.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7727	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.98% 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: MES, JTR

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	0.34	0/3835	0.54	1/5211 (0.0%)
1	B	0.35	0/3865	0.56	1/5250 (0.0%)
All	All	0.35	0/7700	0.55	2/10461 (0.0%)

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

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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	B	164	LEU	CB-CG-CD2	5.15	119.76	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	235	ARG	Sidechain

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	3741	0	3550	40	0
1	B	3770	0	3556	63	0
2	A	28	0	0	0	0
2	B	28	0	0	2	0
3	A	24	0	24	3	0
4	A	61	0	0	2	0
4	B	75	0	0	2	0
All	All	7727	0	7130	100	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 7.

All (100) 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:249:HIS:CD2	4:A:706:HOH:O	2.35	0.79
1:B:441:ASN:HA	1:B:443:PRO:N	1.99	0.77
1:B:181:TYR:HA	1:B:184:ILE:HD12	1.69	0.73
1:B:354:TYR:HB2	1:B:490:GLU:HG2	1.75	0.69
1:B:444:THR:HG23	1:B:447:LYS:H	1.56	0.69
1:B:199:SER:OG	1:B:203:GLU:N	2.26	0.69
1:A:235:ARG:NH2	1:A:580:ARG:O	2.26	0.68
1:B:421:SER:O	1:B:425:ILE:HG13	1.94	0.68
1:B:425:ILE:O	1:B:429:ILE:HG13	1.95	0.67
1:B:275:THR:HB	1:B:324:GLU:OE1	1.96	0.66
1:B:407:ILE:HG13	1:B:457:ILE:HD11	1.78	0.65
1:B:441:ASN:HA	1:B:442:PRO:C	2.16	0.65
1:A:482:HIS:CE1	1:A:485:LYS:HG3	2.30	0.65
1:B:337:THR:HG21	1:B:499:LYS:HD2	1.79	0.65
1:B:172:HIS:HD2	4:B:747:HOH:O	1.78	0.64
1:A:152:PHE:HB2	1:A:163:VAL:HB	1.81	0.63
1:B:176:ASN:HA	4:B:747:HOH:O	2.00	0.61
1:B:164:LEU:O	1:B:208:PRO:HD2	2.03	0.58
1:A:263:ASN:O	1:A:267:GLU:HG3	2.05	0.56
1:B:424:THR:O	1:B:428:MET:HG3	2.05	0.56
1:B:166:ASN:OD1	1:B:168:SER:OG	2.23	0.54
1:B:427:LYS:O	1:B:431:ILE:HG13	2.08	0.54
1:A:335:ASP:HB3	1:A:338:HIS:HB2	1.90	0.53
1:B:141:ARG:HH22	1:B:151:PHE:HB2	1.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:559:ARG:HA	1:B:562:MET:HE3	1.91	0.53
1:A:186:ARG:NH1	1:A:221:LEU:O	2.43	0.52
1:B:129:GLU:HG2	1:B:195:PHE:CE2	2.44	0.52
1:B:458:GLU:O	1:B:498:GLY:HA2	2.10	0.51
1:B:82:LEU:O	1:B:86:ASN:ND2	2.42	0.51
1:B:272:GLU:HB2	1:B:323:TYR:CZ	2.45	0.51
1:B:164:LEU:CD2	1:B:166:ASN:HB2	2.41	0.50
1:B:244:ASN:O	1:B:247:SER:OG	2.29	0.50
1:B:441:ASN:HA	1:B:443:PRO:CD	2.42	0.50
1:A:410:VAL:HG21	1:A:456:PHE:HB3	1.94	0.50
1:A:508:LEU:HD12	1:A:514:GLN:HG2	1.94	0.49
1:B:236:GLN:HG2	1:B:238:TYR:CZ	2.47	0.49
1:A:417:GLN:O	1:A:419:PHE:N	2.46	0.49
1:A:271:PHE:CZ	1:B:252:VAL:HA	2.48	0.48
1:A:500:GLU:HG2	1:A:559:ARG:NH1	2.28	0.48
1:B:192:ILE:HG23	1:B:208:PRO:HB3	1.94	0.48
1:B:418:PRO:HB2	1:B:421:SER:HB3	1.94	0.48
1:B:452:LEU:HD22	1:B:456:PHE:HE2	1.78	0.48
1:B:502:LEU:HD12	1:B:555:LEU:HD21	1.96	0.48
1:A:437:ILE:HD13	1:A:455:HIS:HD2	1.80	0.47
1:B:167:TYR:HE1	1:B:172:HIS:HE2	1.63	0.47
1:A:408:GLU:HG2	1:A:413:THR:O	2.15	0.46
1:B:388:GLN:HG2	1:B:389:PRO:HD2	1.97	0.46
1:B:275:THR:HG22	1:B:304:ARG:HH11	1.81	0.46
1:B:432:ILE:HG22	1:B:437:ILE:O	2.16	0.46
1:B:404:VAL:O	1:B:408:GLU:HG3	2.16	0.45
3:A:603:MES:H51	1:B:356:ASP:OD1	2.16	0.45
1:B:280:LEU:HD23	1:B:301:LEU:CD2	2.46	0.45
1:B:176:ASN:OD1	1:B:178:ALA:N	2.49	0.45
1:B:330:ARG:HD3	2:B:601:JTR:C5	2.46	0.45
1:B:164:LEU:HD21	1:B:166:ASN:HB2	1.99	0.44
1:B:449:LEU:CD1	1:B:474:ILE:HD11	2.46	0.44
1:B:327:LYS:NZ	1:B:341:GLU:OE2	2.47	0.44
1:A:478:LEU:HA	1:A:514:GLN:HE22	1.83	0.44
1:B:428:MET:HB3	1:B:428:MET:HE2	1.84	0.44
1:A:387:ASN:OD1	1:A:387:ASN:N	2.51	0.44
1:B:154:LEU:HD23	1:B:161:ILE:HD11	1.98	0.44
1:A:141:ARG:HB3	1:A:153:ASP:HB2	1.99	0.44
1:A:150:ARG:NH1	1:A:182:ASP:OD1	2.48	0.44
1:A:305:ILE:O	1:A:330:ARG:NH1	2.50	0.44
1:A:578:THR:O	1:A:579:MET:HG3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:601:JTR:N5	2:B:601:JTR:C14	2.80	0.44
1:A:115:PHE:CE2	1:A:196:PRO:HB3	2.52	0.43
1:A:578:THR:C	1:A:579:MET:HG3	2.39	0.43
1:A:438:GLU:O	1:A:451:GLN:NE2	2.51	0.43
1:B:309:LEU:HD23	1:B:539:PHE:HB2	2.00	0.43
1:A:103:PRO:HG3	1:A:212:ILE:HD11	2.01	0.43
1:B:150:ARG:HH12	1:B:182:ASP:CG	2.23	0.43
1:B:198:LYS:HD2	1:B:202:GLY:HA2	2.00	0.43
1:A:508:LEU:HD11	1:A:513:LYS:HB2	2.01	0.42
1:A:461:TYR:C	1:A:462:ASN:HD22	2.21	0.42
1:B:290:PHE:HB2	1:B:303:LEU:HB2	2.01	0.42
1:B:90:PHE:O	1:B:94:GLN:HG2	2.19	0.42
1:A:150:ARG:HB2	1:A:165:ALA:HB3	2.02	0.42
1:B:132:ILE:O	1:B:132:ILE:HD12	2.19	0.42
1:B:470:GLU:HG2	1:B:482:HIS:CD2	2.55	0.42
1:B:87:ARG:HD2	1:B:87:ARG:HA	1.80	0.42
1:A:164:LEU:O	1:A:208:PRO:HD2	2.20	0.41
1:A:418:PRO:HG2	1:A:421:SER:HB3	2.02	0.41
1:A:91:ILE:O	1:A:95:LYS:HG3	2.20	0.41
1:B:406:GLU:O	1:B:410:VAL:HG22	2.20	0.41
1:A:143:SER:OG	1:A:151:PHE:HB2	2.20	0.41
1:A:416:GLU:O	1:A:417:GLN:C	2.60	0.41
3:A:602:MES:H32	1:B:107:GLU:HA	2.02	0.41
1:A:108:ARG:HD2	3:A:603:MES:O3S	2.21	0.41
1:A:240:ASP:OD1	1:A:244:ASN:ND2	2.38	0.41
1:B:432:ILE:H	1:B:432:ILE:HG13	1.65	0.41
1:A:249:HIS:HD2	4:A:706:HOH:O	1.93	0.40
1:A:444:THR:O	1:A:448:LEU:HG	2.21	0.40
1:A:271:PHE:HZ	1:B:252:VAL:HA	1.87	0.40
1:A:331:ASN:OD1	1:B:294:HIS:HD2	2.04	0.40
1:A:546:GLY:O	1:B:138:ARG:NH2	2.51	0.40
1:B:416:GLU:H	1:B:424:THR:HG23	1.86	0.40
1:A:330:ARG:HG2	1:A:330:ARG:HH11	1.87	0.40
1:B:141:ARG:HH22	1:B:151:PHE:CB	2.32	0.40
1:B:408:GLU:HB3	1:B:413:THR:O	2.22	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	466/516 (90%)	457 (98%)	9 (2%)	0	100	100
1	B	474/516 (92%)	457 (96%)	17 (4%)	0	100	100
All	All	940/1032 (91%)	914 (97%)	26 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	392/464 (84%)	384 (98%)	8 (2%)	55	72
1	B	388/464 (84%)	379 (98%)	9 (2%)	50	67
All	All	780/928 (84%)	763 (98%)	17 (2%)	52	69

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	LEU
1	A	162	GLN
1	A	335	ASP
1	A	347	PHE
1	A	387	ASN
1	A	462	ASN
1	A	492	LEU
1	A	501	VAL

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Mol	Chain	Res	Type
1	B	141	ARG
1	B	275	THR
1	B	344	SER
1	B	375	THR
1	B	383	ASP
1	B	476	SER
1	B	492	LEU
1	B	501	VAL
1	B	580	ARG

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	176	ASN
1	A	455	HIS
1	A	462	ASN
1	A	514	GLN
1	B	86	ASN
1	B	294	HIS
1	B	295	ASN
1	B	388	GLN
1	B	514	GLN

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 monosaccharides 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	JTR	A	601	-	29,30,30	6.22	4 (13%)	37,42,42	2.75	10 (27%)
3	MES	A	602	-	12,12,12	2.13	1 (8%)	14,16,16	2.36	7 (50%)
2	JTR	B	601	-	29,30,30	6.14	5 (17%)	37,42,42	2.89	9 (24%)
3	MES	A	603	-	12,12,12	2.14	1 (8%)	14,16,16	2.44	5 (35%)

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	JTR	A	601	-	-	8/22/22/22	0/3/3/3
3	MES	A	602	-	-	1/6/14/14	0/1/1/1
2	JTR	B	601	-	-	7/22/22/22	0/3/3/3
3	MES	A	603	-	-	1/6/14/14	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	JTR	O4-S1	23.01	1.65	1.44
2	A	601	JTR	O5-S1	22.91	1.65	1.44
2	B	601	JTR	O4-S1	22.70	1.64	1.44
2	B	601	JTR	O5-S1	22.64	1.64	1.44
3	A	603	MES	C8-S	-7.19	1.67	1.77
3	A	602	MES	C8-S	-7.06	1.67	1.77
2	B	601	JTR	C12-N4	4.64	1.46	1.36
2	A	601	JTR	C12-N4	4.59	1.45	1.36
2	A	601	JTR	C10-N1	4.41	1.45	1.36
2	B	601	JTR	C10-N1	4.23	1.45	1.36
2	B	601	JTR	C10-N2	-2.13	1.31	1.34

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	JTR	O5-S1-O4	-10.68	108.92	118.71
2	A	601	JTR	O5-S1-O4	-9.13	110.33	118.71
2	A	601	JTR	N3-C12-N5	-6.96	119.96	126.55
2	B	601	JTR	N3-C12-N5	-6.87	120.04	126.55
2	A	601	JTR	N2-C10-N5	-6.64	120.26	126.55
2	B	601	JTR	N2-C10-N5	-6.28	120.59	126.55
2	B	601	JTR	N2-C11-N3	-5.89	119.39	128.60
2	A	601	JTR	N2-C11-N3	-5.82	119.50	128.60
3	A	603	MES	C5-N4-C3	5.50	121.21	108.83
3	A	602	MES	C5-N4-C3	5.06	120.21	108.83
2	B	601	JTR	O1-C2-C9	3.85	119.53	114.80
3	A	603	MES	C7-N4-C3	3.83	121.03	111.23
2	A	601	JTR	O1-C2-C9	3.67	119.30	114.80
3	A	603	MES	O1S-S-C8	3.35	110.95	106.92
3	A	602	MES	C7-N4-C3	3.27	119.58	111.23
2	A	601	JTR	C12-N5-C10	3.15	119.23	113.89
2	B	601	JTR	C12-N5-C10	2.91	118.83	113.89
3	A	602	MES	O3S-S-C8	2.84	110.37	105.77
3	A	602	MES	C7-N4-C5	2.82	118.44	111.23
3	A	603	MES	O3S-S-C8	2.81	110.32	105.77
3	A	602	MES	C2-C3-N4	-2.59	106.17	110.10
2	B	601	JTR	O1-C2-C3	-2.46	120.16	124.37
2	B	601	JTR	C1-O1-C2	-2.43	113.86	117.53
2	A	601	JTR	C13-N4-C12	-2.40	122.36	129.60
2	A	601	JTR	C18-S1-C19	2.31	110.17	105.55
3	A	602	MES	O2S-S-C8	2.31	109.69	106.92
2	A	601	JTR	O1-C2-C3	-2.23	120.55	124.37
3	A	603	MES	C7-N4-C5	2.22	116.92	111.23
3	A	602	MES	O1S-S-C8	2.20	109.57	106.92
2	B	601	JTR	C13-N4-C12	-2.14	123.14	129.60
2	A	601	JTR	C1-O1-C2	-2.08	114.39	117.53

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	JTR	N3-C12-N4-C13
2	A	601	JTR	N5-C12-N4-C13
2	A	601	JTR	C20-C19-S1-C18
2	A	601	JTR	C20-C19-S1-O4
2	A	601	JTR	C20-C19-S1-O5
2	A	601	JTR	C21-C19-S1-C18
2	A	601	JTR	C21-C19-S1-O4

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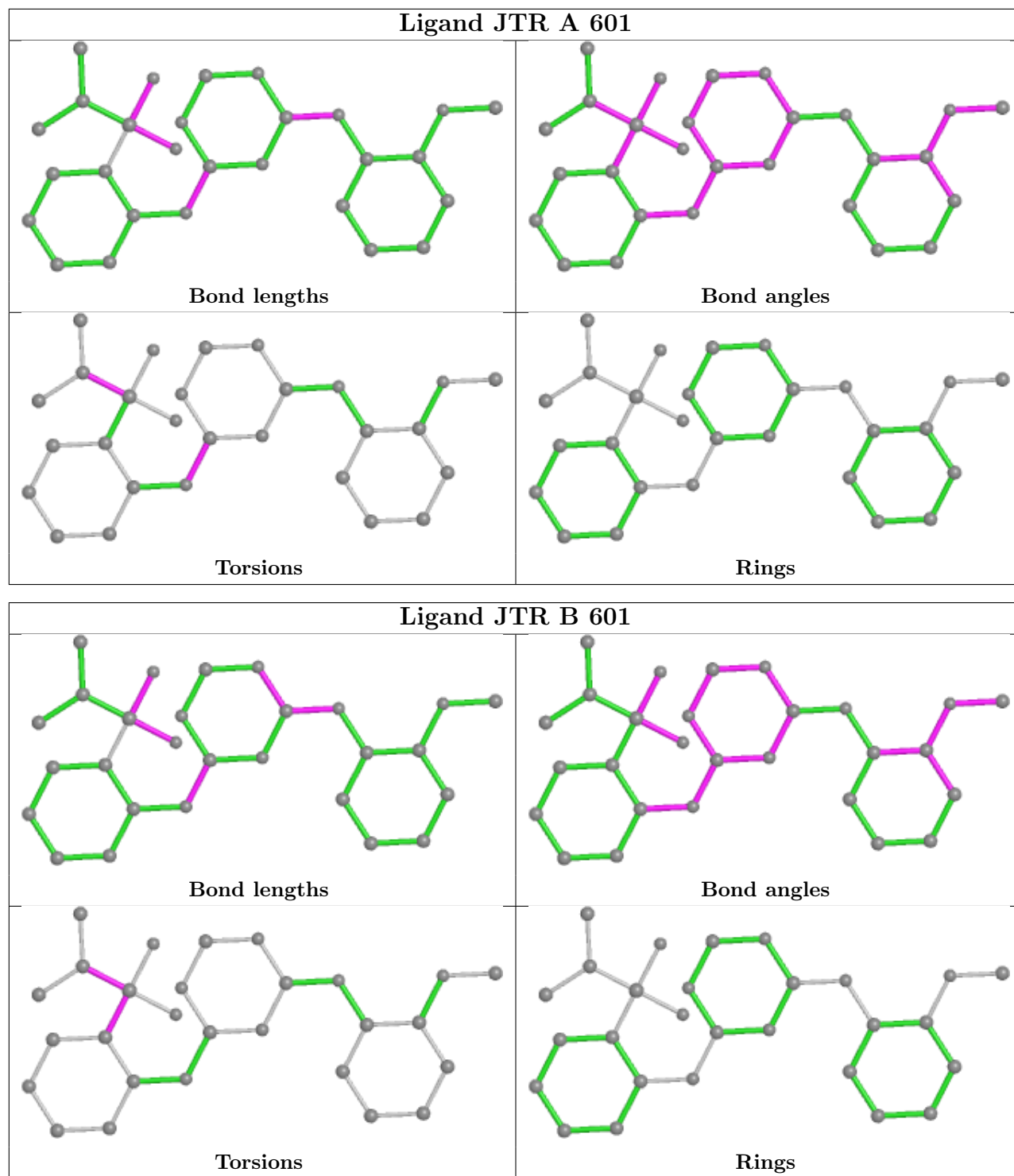
Mol	Chain	Res	Type	Atoms
2	A	601	JTR	C21-C19-S1-O5
2	B	601	JTR	C20-C19-S1-C18
2	B	601	JTR	C20-C19-S1-O4
2	B	601	JTR	C20-C19-S1-O5
2	B	601	JTR	C21-C19-S1-C18
2	B	601	JTR	C21-C19-S1-O4
2	B	601	JTR	C21-C19-S1-O5
3	A	602	MES	C8-C7-N4-C5
3	A	603	MES	C8-C7-N4-C5
2	B	601	JTR	C17-C18-S1-O4

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	MES	1	0
2	B	601	JTR	2	0
3	A	603	MES	2	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

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	476/516 (92%)	0.23	18 (3%) 40 47	45, 65, 101, 155	0
1	B	482/516 (93%)	0.69	60 (12%) 4 5	42, 76, 131, 186	0
All	All	958/1032 (92%)	0.46	78 (8%) 12 16	42, 70, 120, 186	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	225	TYR	15.2
1	B	144	ALA	8.4
1	B	202	GLY	7.2
1	B	125	GLY	5.9
1	B	226	GLY	5.9
1	A	518	PHE	5.4
1	B	443	PRO	5.0
1	B	80	PRO	4.7
1	A	536	ASP	4.6
1	B	200	LYS	4.6
1	B	536	ASP	4.6
1	B	229	ASP	4.5
1	B	123	GLY	4.5
1	B	390	ILE	4.3
1	B	501	VAL	4.3
1	B	581	PRO	4.1
1	B	201	LYS	4.0
1	B	124	ASN	3.9
1	A	82	LEU	3.9
1	B	199	SER	3.8
1	B	164	LEU	3.7
1	B	580	ARG	3.7
1	B	502	LEU	3.6
1	B	149	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	98	GLY	3.2
1	B	441	ASN	3.2
1	B	417	GLN	3.1
1	B	442	PRO	3.1
1	B	456	PHE	3.1
1	B	143	SER	3.1
1	B	387	ASN	3.0
1	B	512	PHE	3.0
1	B	151	PHE	3.0
1	B	560	ILE	2.9
1	B	440	PRO	2.9
1	B	556	GLY	2.9
1	A	385	PRO	2.9
1	B	557	ILE	2.9
1	A	189	ILE	2.8
1	B	553	LEU	2.8
1	A	501	VAL	2.8
1	A	581	PRO	2.7
1	A	309	LEU	2.6
1	B	165	ALA	2.6
1	B	554	GLY	2.6
1	A	390	ILE	2.6
1	A	516	GLU	2.6
1	A	78	VAL	2.6
1	B	379	SER	2.5
1	B	84	PHE	2.5
1	B	178	ALA	2.5
1	B	147	GLN	2.5
1	A	422	ASN	2.5
1	A	282	ALA	2.5
1	A	474	ILE	2.4
1	B	384	GLY	2.4
1	B	494	MET	2.4
1	B	386	GLU	2.4
1	A	80	PRO	2.3
1	B	420	ASP	2.3
1	B	141	ARG	2.3
1	B	383	ASP	2.3
1	B	223	MET	2.2
1	B	555	LEU	2.2
1	B	504	ALA	2.2
1	B	127	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	228	LYS	2.1
1	A	81	ARG	2.1
1	B	444	THR	2.1
1	B	474	ILE	2.1
1	B	174	LYS	2.1
1	A	190	VAL	2.1
1	B	517	CYS	2.1
1	A	85	GLU	2.0
1	B	170	HIS	2.0
1	B	91	ILE	2.0
1	B	315	ILE	2.0
1	B	561	THR	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](#)

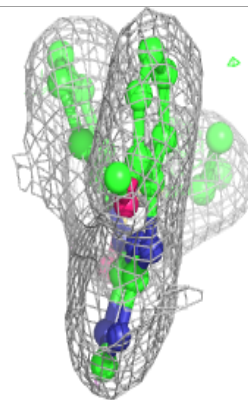
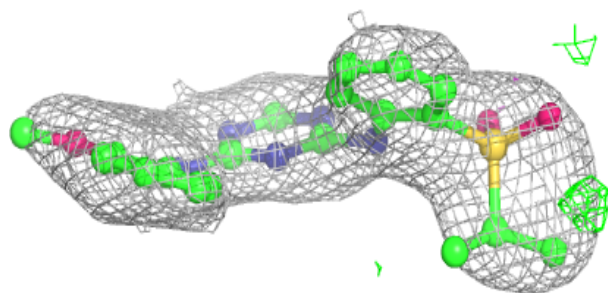
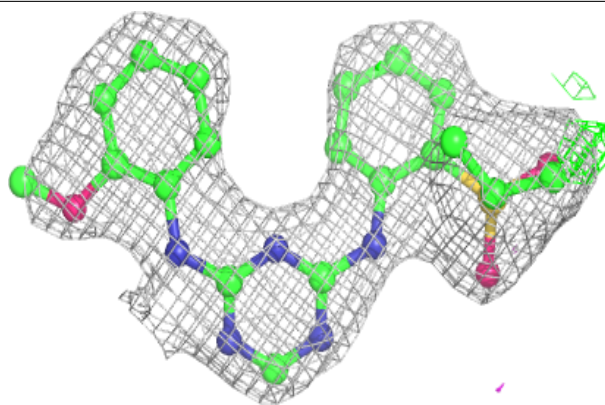
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	JTR	A	601	28/28	0.90	0.16	60,67,74,76	0
2	JTR	B	601	28/28	0.92	0.17	61,70,75,76	0
3	MES	A	602	12/12	0.97	0.09	66,71,78,80	0
3	MES	A	603	12/12	0.97	0.11	54,57,60,62	0

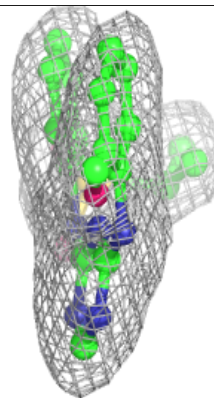
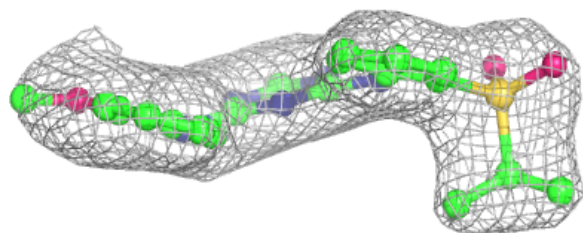
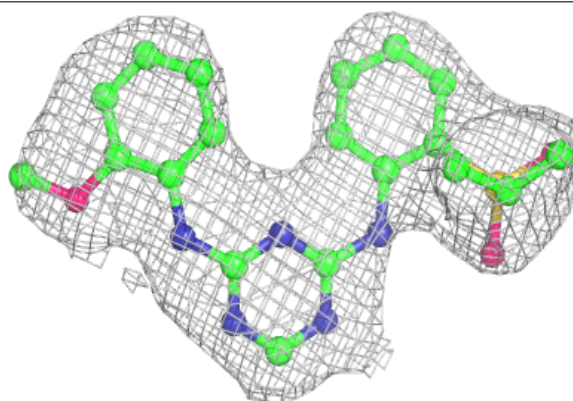
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around JTR A 601:

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

**Electron density around JTR B 601:**

$2mF_o-DF_c$ (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.