



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:28 AM UTC

PDB ID : 6S43 / pdb_00006s43
Title : Fumarate hydratase of Mycobacterium tuberculosis in complex with formate and allosteric modulator N-(5-(Azocan-1-ylsulfonyl)-2-methoxyphenyl)-2-(4-oxo-3,4-dihydrophthalazin-1-yl)acetamide
Authors : Whitehouse, A.J.; Libardo, M.D.; Kasbekar, M.; Brear, P.; Fischer, G.; Thomas, C.J.; Barry, C.E.; Boshoff, H.I.; Coyne, A.G.; Abell, C.
Deposited on : 2019-06-26
Resolution : 1.42 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](#) ①) 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

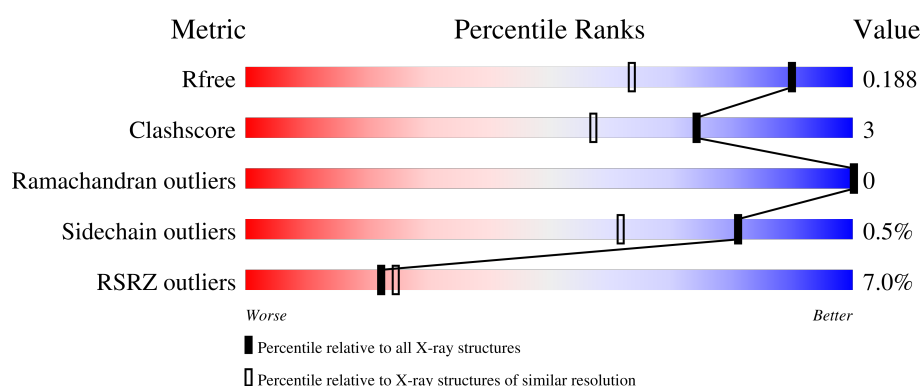
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.42 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	4041 (1.44-1.40)
Clashscore	190562	4154 (1.44-1.40)
Ramachandran outliers	187476	4083 (1.44-1.40)
Sidechain outliers	187428	4082 (1.44-1.40)
RSRZ outliers	180081	4039 (1.44-1.40)

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	474	6% 88% 7% 5%
1	B	474	8% 89% 7% •
1	C	474	5% 89% 5% 5%
1	D	474	7% 87% 7% 5%

2 Entry composition [i](#)

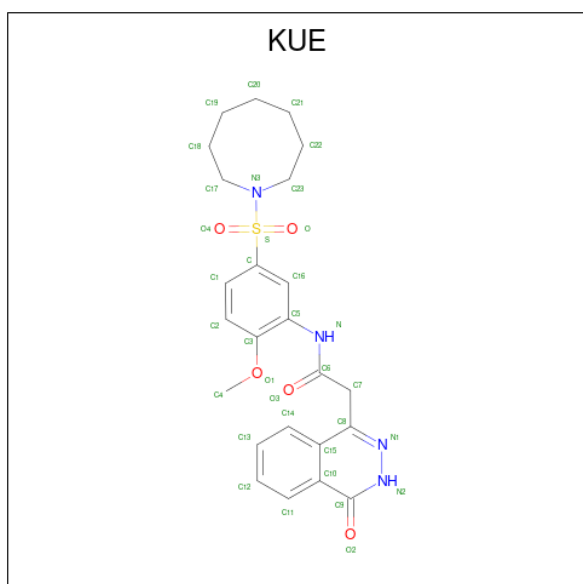
There are 5 unique types of molecules in this entry. The entry contains 15622 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fumarate hydratase class II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	13	0
			3471	2160	635	663	13			
1	B	458	Total	C	N	O	S	0	13	0
			3515	2187	641	673	14			
1	C	449	Total	C	N	O	S	0	19	0
			3493	2177	634	668	14			
1	D	448	Total	C	N	O	S	0	9	0
			3409	2122	621	654	12			

- Molecule 2 is {N}-[5-(azocan-1-ylsulfonyl)-2-methoxy-phenyl]-2-(4-oxidanylidene-3 {H}-phthalazin-1-yl)ethanamide (CCD ID: KUE) (formula: C₂₄H₂₈N₄O₅S) (labeled as "Ligand of Interest" by depositor).



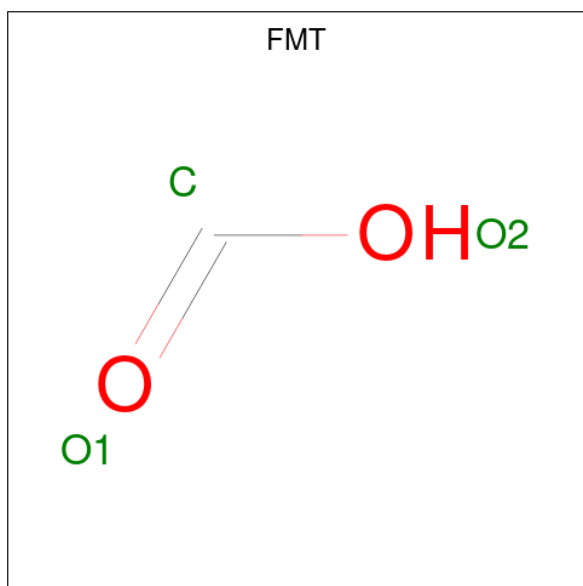
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			34	24	4	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			34	24	4	5	1		

- Molecule 3 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	446	Total	O	0	0
			446	446		
5	B	415	Total	O	0	0
			415	415		
5	C	405	Total	O	0	0
			405	405		

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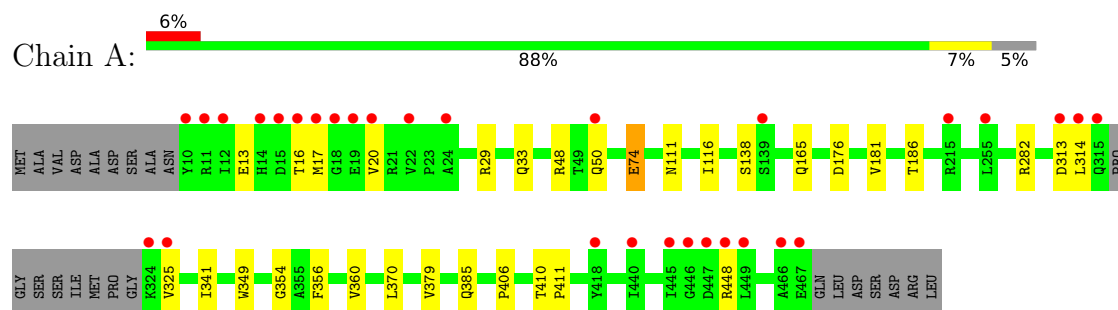
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	395	Total 395	O 395	0	0

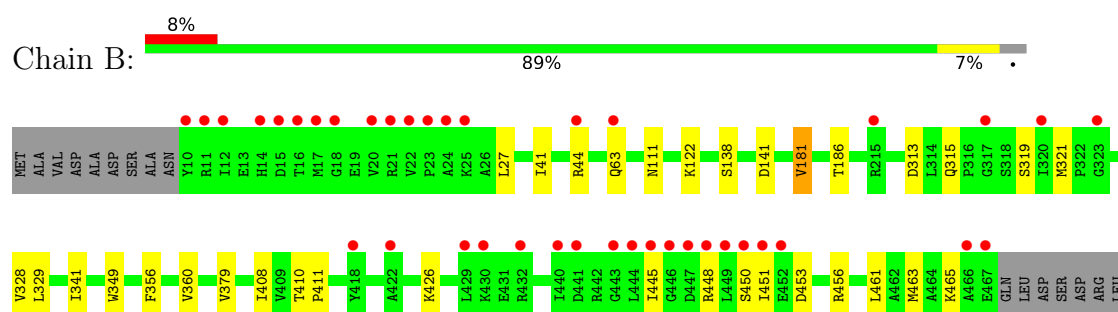
3 Residue-property plots [i](#)

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.

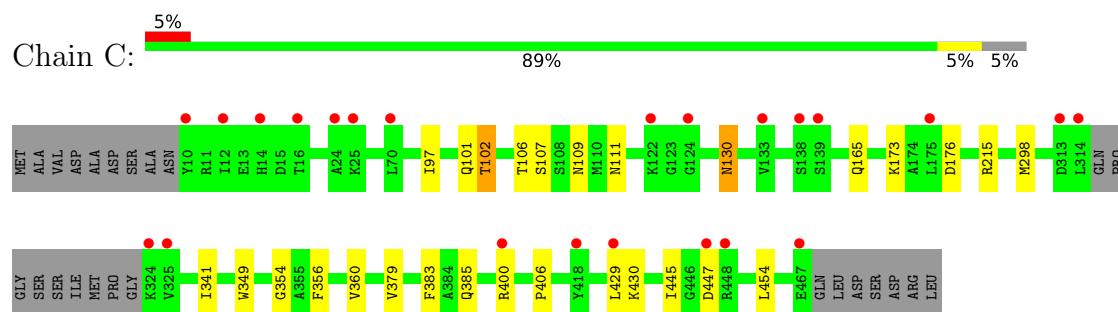
• Molecule 1: Fumarate hydratase class II



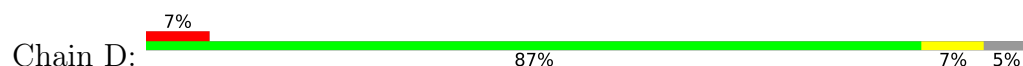
• Molecule 1: Fumarate hydratase class II

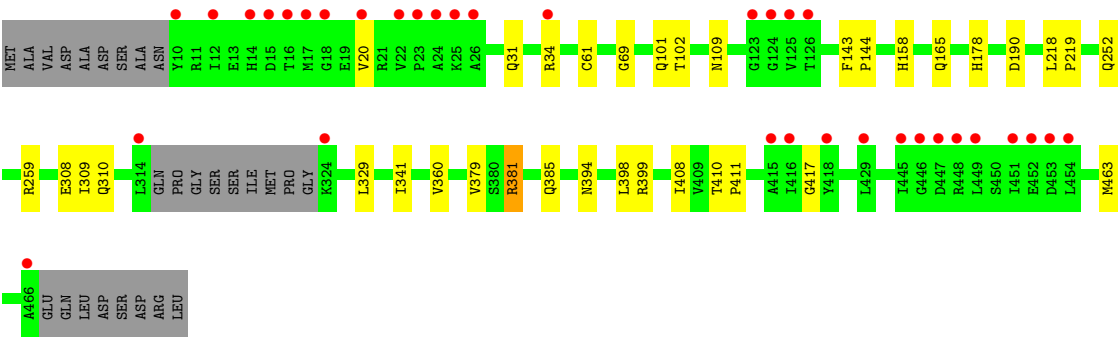


• Molecule 1: Fumarate hydratase class II



• Molecule 1: Fumarate hydratase class II





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	175.31Å 96.57Å 124.35Å 90.00° 102.70° 90.00°	Depositor
Resolution (Å)	84.09 – 1.42 84.09 – 1.42	Depositor EDS
% Data completeness (in resolution range)	95.5 (84.09-1.42) 95.4 (84.09-1.42)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 1.42Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.170 , 0.190 (Not available) , 0.188	Depositor DCC
R_{free} test set	17874 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	16.6	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.2	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	0.97	EDS
Total number of atoms	15622	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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: FMT, KUE, MG

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	1.13	3/3517 (0.1%)	1.24	4/4776 (0.1%)
1	B	1.16	3/3564 (0.1%)	1.30	9/4841 (0.2%)
1	C	1.18	2/3542 (0.1%)	1.30	4/4811 (0.1%)
1	D	1.14	3/3455 (0.1%)	1.30	7/4694 (0.1%)
All	All	1.15	11/14078 (0.1%)	1.28	24/19122 (0.1%)

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
1	B	0	2
1	C	0	5
All	All	0	8

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	321	MET	C-N	12.93	1.49	1.33
1	A	356	PHE	C-O	-6.41	1.18	1.24
1	D	399	ARG	C-O	-6.41	1.16	1.24
1	C	109	ASN	C-O	6.09	1.31	1.24
1	D	309	ILE	C-N	-5.98	1.25	1.33
1	B	138	SER	CA-CB	-5.71	1.45	1.53
1	B	356	PHE	C-O	-5.69	1.19	1.24
1	D	109	ASN	C-O	5.67	1.30	1.24
1	A	29	ARG	C-O	5.31	1.29	1.24
1	A	48	ARG	C-O	-5.29	1.17	1.24
1	C	356	PHE	C-O	-5.13	1.20	1.24

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	321	MET	CA-C-N	9.83	129.59	119.56
1	B	321	MET	C-N-CA	9.83	129.59	119.56
1	A	111	ASN	CA-CB-CG	7.07	119.67	112.60
1	B	321	MET	O-C-N	-6.71	115.26	121.36
1	B	313	ASP	CA-CB-CG	6.10	118.70	112.60
1	C	176	ASP	CA-CB-CG	5.92	118.52	112.60
1	D	178	HIS	CA-C-O	-5.79	114.37	120.63
1	B	111	ASN	CA-CB-CG	5.68	118.28	112.60
1	D	381	ARG	CG-CD-NE	-5.43	100.06	112.00
1	B	410	THR	CA-CB-OG1	-5.38	101.53	109.60
1	B	41	ILE	CA-C-O	-5.33	116.17	120.70
1	C	111	ASN	CA-CB-CG	5.32	117.92	112.60
1	D	190	ASP	CA-CB-CG	5.25	117.85	112.60
1	C	383	PHE	CA-CB-CG	5.20	119.00	113.80
1	A	186	THR	CA-C-O	-5.18	114.57	120.32
1	B	186	THR	CA-C-O	-5.14	114.61	120.32
1	B	141	ASP	CA-CB-CG	5.10	117.70	112.60
1	A	116	ILE	CA-C-O	-5.10	115.65	120.95
1	A	176	ASP	CA-CB-CG	5.07	117.67	112.60
1	C	447	ASP	CA-CB-CG	5.06	117.66	112.60
1	D	417	GLY	CA-C-N	5.05	127.76	120.38
1	D	417	GLY	C-N-CA	5.05	127.76	120.38
1	D	61	CYS	O-C-N	5.05	127.47	122.12
1	D	309	ILE	CA-C-O	-5.04	116.53	121.67

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	181	VAL	Mainchain
1	B	181[A]	VAL	Mainchain
1	B	181[B]	VAL	Mainchain
1	C	102[B]	THR	Mainchain
1	C	130[A]	ASN	Mainchain
1	C	130[B]	ASN	Mainchain
1	C	298[A]	MET	Mainchain
1	C	298[B]	MET	Mainchain

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	3471	0	3511	23	0
1	B	3515	0	3556	26	0
1	C	3493	0	3537	27	0
1	D	3409	0	3441	24	0
2	A	34	0	0	0	0
2	C	34	0	0	0	0
3	B	3	0	1	0	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
5	A	446	0	0	8	0
5	B	415	0	0	5	0
5	C	405	0	0	7	0
5	D	395	0	0	9	0
All	All	15622	0	14046	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 3.

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:C:102[A]:THR:HG21	5:C:878:HOH:O	1.56	1.04
1:B:445:ILE:HD13	1:B:451:ILE:HG12	1.58	0.86
1:B:319:SER:HB2	1:C:130[B]:ASN:OD1	1.91	0.70
1:D:360:VAL:HG22	5:D:847:HOH:O	1.95	0.66
1:B:319:SER:CB	1:C:130[B]:ASN:OD1	2.45	0.65
1:D:408:ILE:HG12	1:D:463:MET:HE1	1.81	0.62
1:D:310[A]:GLN:HE21	1:D:394:ASN:HD21	1.49	0.61
1:A:360:VAL:HG22	5:A:879:HOH:O	2.00	0.60
1:A:50[C]:GLN:HG2	1:A:370:LEU:CD1	2.32	0.60
1:A:74:GLU:H	1:A:74:GLU:CD	2.09	0.60
1:D:165:GLN:NE2	1:D:385:GLN:OE1	2.35	0.60
1:C:165:GLN:HE22	1:C:385:GLN:CG	2.15	0.59
1:B:360:VAL:HG22	5:B:865:HOH:O	2.03	0.59
1:C:165:GLN:NE2	1:C:385:GLN:HG3	2.19	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:360:VAL:HG22	5:C:830:HOH:O	2.05	0.57
1:B:445:ILE:HD13	1:B:451:ILE:CG1	2.33	0.56
1:D:69:GLY:HA2	5:D:766:HOH:O	2.04	0.56
1:D:259:ARG:NH2	5:D:606:HOH:O	2.39	0.56
1:D:381:ARG:HD3	5:D:866:HOH:O	2.04	0.56
1:A:448:ARG:HG2	5:A:685:HOH:O	2.04	0.56
1:B:44:ARG:HD3	5:B:826:HOH:O	2.08	0.53
1:B:181[B]:VAL:HG21	1:B:461:LEU:HB2	1.91	0.53
1:A:165:GLN:NE2	1:A:385:GLN:HG3	2.26	0.51
1:B:63:GLN:NE2	5:B:601:HOH:O	2.27	0.51
1:D:158:HIS:HE1	5:D:900:HOH:O	1.94	0.51
1:A:16:THR:HG23	1:A:17:MET:HG2	1.94	0.50
1:C:165:GLN:NE2	1:C:385:GLN:CG	2.73	0.50
1:B:341[A]:ILE:HD13	1:C:349:TRP:CH2	2.47	0.50
1:A:313:ASP:HA	1:A:325:VAL:CG1	2.42	0.50
1:B:411:PRO:HG2	1:B:463[B]:MET:HE2	1.94	0.50
1:D:165:GLN:HE21	1:D:385:GLN:HG3	1.77	0.49
1:B:450:SER:OG	1:B:453:ASP:HB2	2.13	0.49
1:B:328:VAL:CG1	1:C:102[B]:THR:HG22	2.43	0.49
1:C:406:PRO:HD2	5:C:723:HOH:O	2.13	0.48
1:A:20:VAL:HG13	1:A:20:VAL:O	2.12	0.48
1:B:453:ASP:OD1	1:B:456:ARG:NH2	2.46	0.48
5:A:752:HOH:O	1:B:465:LYS:HD3	2.13	0.48
1:B:319:SER:HB2	1:C:130[B]:ASN:CG	2.38	0.48
1:A:282:ARG:HG2	1:A:341[B]:ILE:HD12	1.95	0.48
1:C:97:ILE:HD13	1:C:107:SER:HB3	1.95	0.48
1:C:445:ILE:HD11	1:C:454:LEU:HD22	1.96	0.47
1:A:13:GLU:HB2	1:A:20:VAL:CG1	2.45	0.47
1:C:406:PRO:CG	1:C:429[B]:LEU:HD23	2.45	0.46
1:B:349:TRP:CZ2	1:C:341:ILE:HD13	2.51	0.46
1:D:252:GLN:OE1	5:D:601:HOH:O	2.21	0.45
1:C:106[B]:THR:HG21	5:C:878:HOH:O	2.16	0.45
1:A:410:THR:N	1:A:411:PRO:HD2	2.32	0.45
1:B:181[B]:VAL:CG2	1:B:461:LEU:HD13	2.46	0.45
1:B:448:ARG:NH1	5:B:615:HOH:O	2.49	0.45
1:A:406:PRO:HD2	5:A:725:HOH:O	2.15	0.45
1:B:315:GLN:NE2	1:C:102[B]:THR:HG21	2.32	0.45
1:C:215[A]:ARG:HD2	5:C:614:HOH:O	2.17	0.45
1:D:385:GLN:NE2	5:D:616:HOH:O	2.48	0.45
1:D:410:THR:HB	1:D:411:PRO:HD3	1.99	0.45
1:A:20:VAL:HG22	1:A:33:GLN:HG2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:400[A]:ARG:NE	5:C:610:HOH:O	2.43	0.44
1:D:20:VAL:HG22	1:D:34:ARG:CZ	2.47	0.44
1:A:165:GLN:HE21	1:A:385:GLN:HG3	1.82	0.44
1:C:354:GLY:HA2	5:C:830:HOH:O	2.17	0.43
1:B:341[A]:ILE:HD13	1:C:349:TRP:CZ2	2.53	0.43
1:C:101:GLN:O	1:C:102[B]:THR:C	2.61	0.43
1:C:165:GLN:HE22	1:C:385:GLN:HG3	1.77	0.43
1:C:173[A]:LYS:HA	1:C:173[A]:LYS:HD3	1.84	0.43
1:D:31:GLN:HE21	1:D:31:GLN:HA	1.84	0.43
1:D:308:GLU:HB3	1:D:398:LEU:HD21	2.00	0.43
1:B:328:VAL:HB	1:C:102[B]:THR:HG22	2.01	0.42
1:A:349:TRP:CH2	1:D:341:ILE:HD13	2.55	0.42
1:A:138:SER:HB2	5:A:660:HOH:O	2.18	0.42
1:B:408:ILE:O	1:B:411:PRO:HD2	2.20	0.42
1:A:50[C]:GLN:HG2	1:A:370:LEU:HD12	2.01	0.42
1:D:143:PHE:N	1:D:144:PRO:HD2	2.34	0.42
1:B:448:ARG:NH2	5:B:622:HOH:O	2.52	0.42
1:B:27:LEU:HD21	1:B:122:LYS:HE2	2.02	0.42
1:A:354:GLY:HA2	5:A:879:HOH:O	2.19	0.41
1:D:218:LEU:N	1:D:219:PRO:CD	2.83	0.41
1:A:20:VAL:HG21	1:A:33:GLN:HB3	2.01	0.41
1:D:158:HIS:CE1	5:D:900:HOH:O	2.70	0.41
1:A:385:GLN:NE2	5:A:618:HOH:O	2.52	0.41
1:C:430:LYS:HB2	1:C:430:LYS:HE3	1.81	0.41
1:D:310[A]:GLN:NE2	1:D:394:ASN:HD21	2.15	0.41
1:D:329:LEU:HD23	1:D:329:LEU:HA	1.97	0.41
1:D:259:ARG:CZ	5:D:606:HOH:O	2.68	0.41
1:B:315:GLN:HE21	1:C:102[B]:THR:HG21	1.86	0.41
1:A:314:LEU:HD23	5:A:607:HOH:O	2.21	0.40
1:D:101:GLN:O	1:D:102:THR:C	2.65	0.40
1:A:313:ASP:HA	1:A:325:VAL:HG11	2.02	0.40
1:A:349:TRP:CZ2	1:D:341:ILE:HD13	2.56	0.40
1:B:408:ILE:HA	1:B:463[A]:MET:HE1	2.04	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	460/474 (97%)	451 (98%)	9 (2%)	0	100	100
1	B	470/474 (99%)	462 (98%)	8 (2%)	0	100	100
1	C	464/474 (98%)	451 (97%)	13 (3%)	0	100	100
1	D	453/474 (96%)	442 (98%)	11 (2%)	0	100	100
All	All	1847/1896 (97%)	1806 (98%)	41 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	359/364 (99%)	357 (99%)	2 (1%)	78	56
1	B	365/364 (100%)	362 (99%)	3 (1%)	73	47
1	C	363/364 (100%)	362 (100%)	1 (0%)	86	71
1	D	352/364 (97%)	351 (100%)	1 (0%)	86	71
All	All	1439/1456 (99%)	1432 (100%)	7 (0%)	81	62

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	GLU
1	A	379	VAL

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Mol	Chain	Res	Type
1	B	329	LEU
1	B	379	VAL
1	B	426	LYS
1	C	379	VAL
1	D	379	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	164	GLN
1	A	165	GLN
1	A	268	GLN
1	A	310	GLN
1	A	339	GLN
1	A	343	ASN
1	A	368	ASN
1	A	385	GLN
1	B	90	GLN
1	B	164	GLN
1	B	165	GLN
1	B	236	ASN
1	B	268	GLN
1	B	339	GLN
1	B	343	ASN
1	B	368	ASN
1	B	385	GLN
1	C	63	GLN
1	C	164	GLN
1	C	339	GLN
1	C	343	ASN
1	C	368	ASN
1	D	14	HIS
1	D	31	GLN
1	D	158	HIS
1	D	165	GLN
1	D	339	GLN
1	D	343	ASN
1	D	368	ASN
1	D	385	GLN
1	D	394	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 ⓘ

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	FMT	B	501	-	2,2,2	0.17	0	1,1,1	0.43	0
2	KUE	A	501	-	37,37,37	0.50	0	44,52,52	0.51	0
2	KUE	C	501	-	37,37,37	0.52	0	44,52,52	0.61	0

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	KUE	A	501	-	-	4/22/32/32	0/3/4/4
2	KUE	C	501	-	-	6/22/32/32	0/3/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

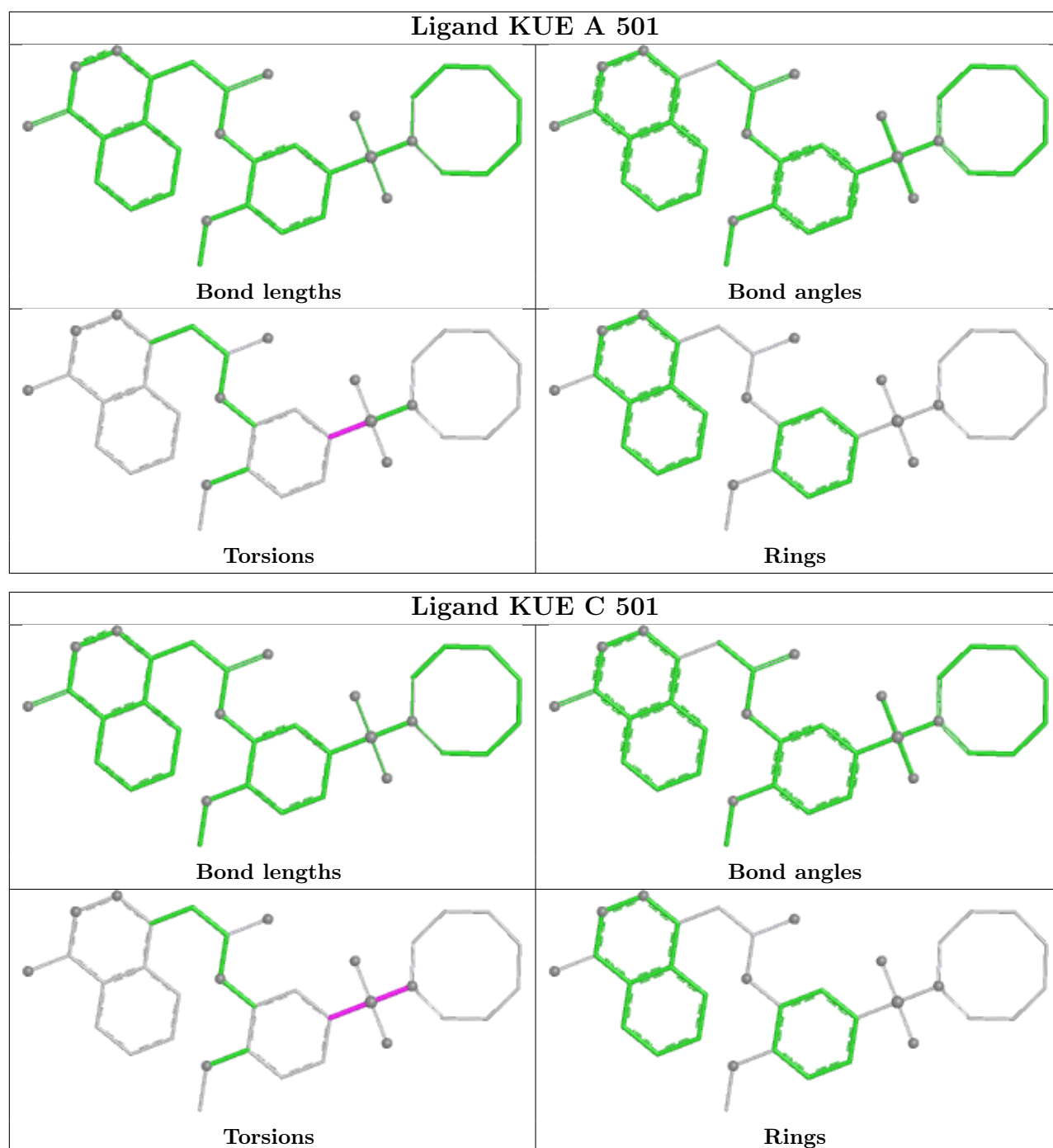
All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	501	KUE	C23-N3-S-C
2	C	501	KUE	C23-N3-S-O4
2	A	501	KUE	C16-C-S-O4
2	A	501	KUE	C1-C-S-O4
2	A	501	KUE	C16-C-S-N3
2	A	501	KUE	C1-C-S-N3
2	C	501	KUE	C1-C-S-O
2	C	501	KUE	C16-C-S-O
2	C	501	KUE	C1-C-S-N3
2	C	501	KUE	C16-C-S-N3

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	450/474 (94%)	0.10	30 (6%)	24	26	5, 17, 45, 111	13 (2%)
1	B	458/474 (96%)	0.25	39 (8%)	16	18	5, 18, 56, 123	13 (2%)
1	C	449/474 (94%)	0.16	23 (5%)	33	36	6, 18, 40, 74	19 (4%)
1	D	448/474 (94%)	0.26	34 (7%)	20	21	6, 17, 53, 120	9 (2%)
All	All	1805/1896 (95%)	0.20	126 (6%)	22	25	5, 17, 49, 123	54 (2%)

All (126) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	215[B]	ARG	12.1
1	A	215[A]	ARG	10.7
1	B	12	ILE	7.3
1	B	451	ILE	6.7
1	B	10	TYR	6.5
1	D	12	ILE	6.2
1	D	10	TYR	6.1
1	A	10	TYR	6.1
1	A	12	ILE	5.7
1	C	10	TYR	5.3
1	A	14	HIS	5.2
1	B	445	ILE	5.1
1	D	418	TYR	5.1
1	B	16	THR	5.1
1	A	20	VAL	4.9
1	C	447	ASP	4.7
1	B	17	MET	4.3
1	A	16	THR	4.3
1	A	17	MET	4.1
1	D	20	VAL	4.1
1	D	314	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
1	D	24	ALA	4.0
1	D	14	HIS	3.9
1	A	315	GLN	3.9
1	B	15	ASP	3.9
1	A	314	LEU	3.9
1	A	418	TYR	3.9
1	D	15	ASP	3.8
1	B	24	ALA	3.8
1	B	14	HIS	3.8
1	C	314	LEU	3.7
1	A	15	ASP	3.7
1	C	12	ILE	3.6
1	A	139	SER	3.6
1	D	416	ILE	3.6
1	C	313	ASP	3.5
1	C	14	HIS	3.5
1	B	323	GLY	3.4
1	D	16	THR	3.4
1	D	22	VAL	3.4
1	B	18	GLY	3.4
1	D	17	MET	3.4
1	B	450	SER	3.4
1	A	324	LYS	3.2
1	B	448	ARG	3.2
1	B	432	ARG	3.2
1	D	448	ARG	3.2
1	D	26	ALA	3.2
1	B	20	VAL	3.1
1	D	18	GLY	3.1
1	A	18	GLY	3.0
1	D	125	VAL	3.0
1	C	25	LYS	3.0
1	A	440	ILE	3.0
1	C	24	ALA	2.9
1	C	16	THR	2.9
1	D	445	ILE	2.9
1	B	452	GLU	2.9
1	B	418	TYR	2.9
1	B	449	LEU	2.9
1	C	324	LYS	2.9
1	B	446	GLY	2.8
1	D	415	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	22	VAL	2.8
1	B	422	ALA	2.8
1	D	123	GLY	2.8
1	D	23	PRO	2.8
1	D	451	ILE	2.7
1	B	11	ARG	2.7
1	B	443	GLY	2.6
1	B	466	ALA	2.6
1	A	11	ARG	2.6
1	A	446	GLY	2.6
1	D	447	ASP	2.6
1	B	429	LEU	2.6
1	D	452	GLU	2.6
1	C	139	SER	2.6
1	A	467	GLU	2.5
1	C	448	ARG	2.5
1	D	124	GLY	2.4
1	D	446	GLY	2.4
1	A	325	VAL	2.4
1	A	445	ILE	2.4
1	C	418	TYR	2.4
1	C	467	GLU	2.4
1	A	24	ALA	2.3
1	A	466	ALA	2.3
1	B	21	ARG	2.3
1	B	44	ARG	2.3
1	A	255[A]	LEU	2.3
1	B	320	ILE	2.3
1	C	429[A]	LEU	2.3
1	D	449	LEU	2.3
1	B	25	LYS	2.3
1	B	430	LYS	2.3
1	B	467	GLU	2.3
1	A	22	VAL	2.3
1	C	325	VAL	2.3
1	B	63	GLN	2.3
1	C	138	SER	2.2
1	B	440	ILE	2.2
1	D	466	ALA	2.2
1	C	400[A]	ARG	2.2
1	D	454	LEU	2.2
1	B	317	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	313	ASP	2.2
1	A	449	LEU	2.1
1	B	444	LEU	2.1
1	C	124	GLY	2.1
1	B	447	ASP	2.1
1	D	25	LYS	2.1
1	A	448	ARG	2.1
1	D	34	ARG	2.1
1	A	50[B]	GLN	2.1
1	A	19	GLU	2.1
1	D	453	ASP	2.1
1	A	447	ASP	2.1
1	B	23	PRO	2.1
1	C	133	VAL	2.0
1	C	175[A]	LEU	2.0
1	D	429	LEU	2.0
1	B	441	ASP	2.0
1	C	122	LYS	2.0
1	D	324	LYS	2.0
1	D	126	THR	2.0
1	C	70	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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
4	MG	D	501	1/1	0.95	0.34	30,30,30,30	0
4	MG	B	502	1/1	0.96	0.21	30,30,30,30	0

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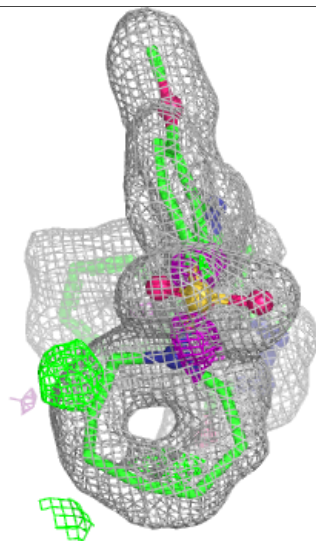
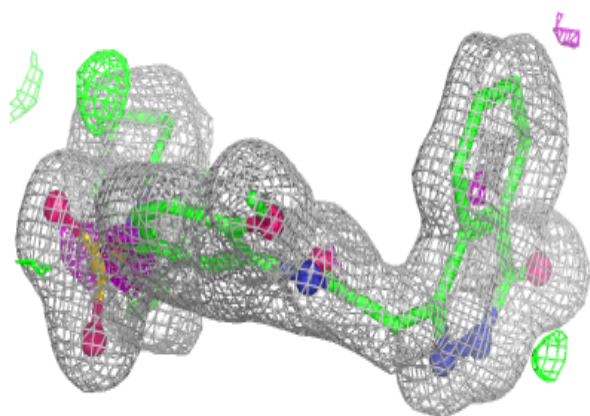
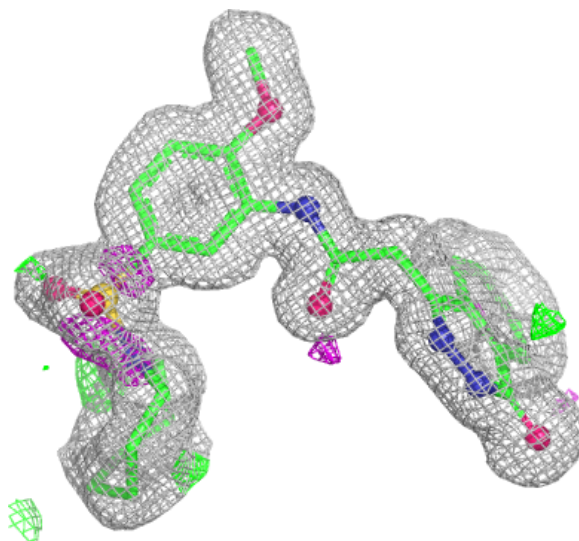
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	FMT	B	501	3/3	0.97	0.09	20,20,21,22	0
2	KUE	A	501	34/34	0.97	0.08	17,20,38,40	0
2	KUE	C	501	34/34	0.97	0.09	17,20,45,50	0

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.

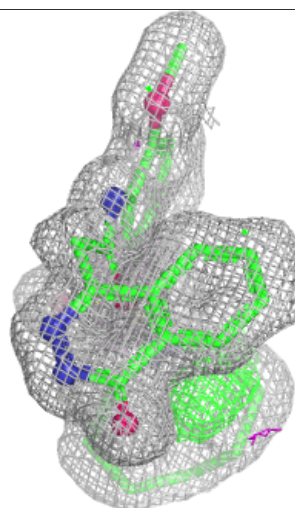
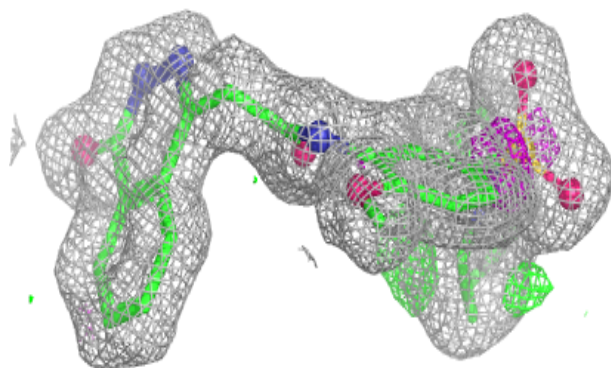
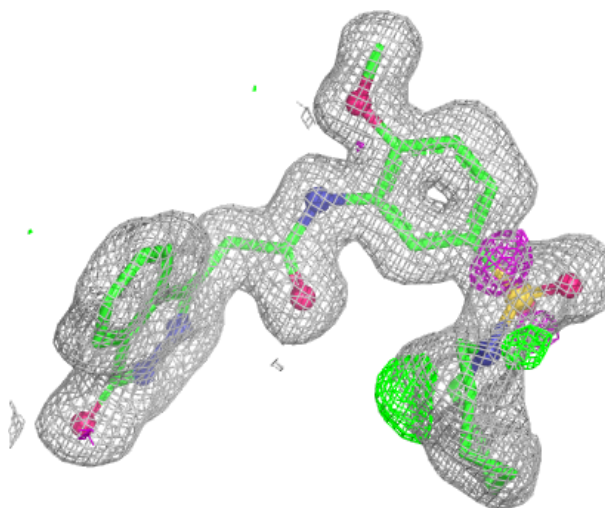
Electron density around KUE A 501:

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 KUE C 501:

$2mF_o-DF_c$ (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.