



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 15, 2023 – 11:59 PM JST

PDB ID : 6KBF
Title : Crystal structure of plasmodium lysyl-tRNA synthetase in complex with a cladosporin derivative 3
Authors : Zhou, J.; Fang, P.
Deposited on : 2019-06-24
Resolution : 1.92 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 : 2.36
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) : 2.36

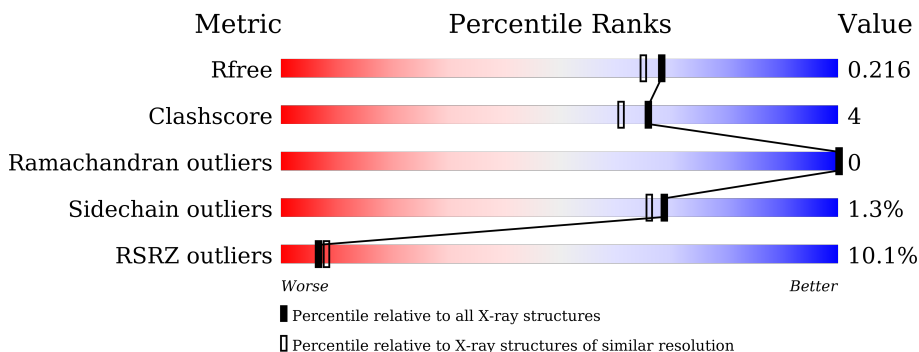
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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	

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
5	DMS	B	605	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8696 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.

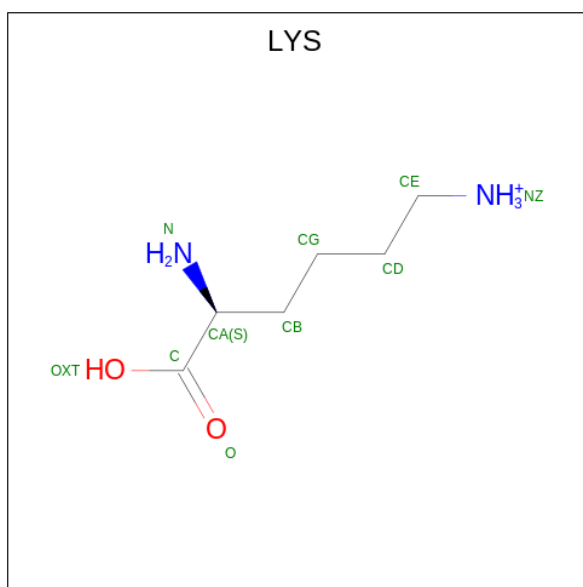
- Molecule 1 is a protein called Lysine-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	504	4037	2597	674	750	16	0	3	0
1	B	503	3903	2495	662	731	15	0	2	0

There are 18 discrepancies between the modelled and reference sequences:

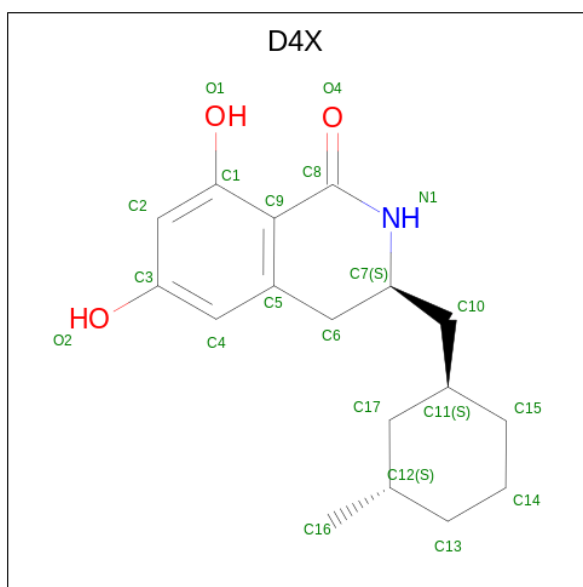
Chain	Residue	Modelled	Actual	Comment	Reference
A	76	MET	-	initiating methionine	UNP W7JP72
A	584	GLY	-	expression tag	UNP W7JP72
A	585	GLY	-	expression tag	UNP W7JP72
A	586	HIS	-	expression tag	UNP W7JP72
A	587	HIS	-	expression tag	UNP W7JP72
A	588	HIS	-	expression tag	UNP W7JP72
A	589	HIS	-	expression tag	UNP W7JP72
A	590	HIS	-	expression tag	UNP W7JP72
A	591	HIS	-	expression tag	UNP W7JP72
B	76	MET	-	initiating methionine	UNP W7JP72
B	584	GLY	-	expression tag	UNP W7JP72
B	585	GLY	-	expression tag	UNP W7JP72
B	586	HIS	-	expression tag	UNP W7JP72
B	587	HIS	-	expression tag	UNP W7JP72
B	588	HIS	-	expression tag	UNP W7JP72
B	589	HIS	-	expression tag	UNP W7JP72
B	590	HIS	-	expression tag	UNP W7JP72
B	591	HIS	-	expression tag	UNP W7JP72

- Molecule 2 is LYSINE (three-letter code: LYS) (formula: $C_6H_{15}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			10	6	2	2		
2	B	1	Total	C	N	O	0	0
			10	6	2	2		

- Molecule 3 is (3 {S})-3-[[[(1 {S}),3 {S})-3-methylcyclohexyl]methyl]-6,8-bis(oxidanyl)-3,4-dihydro-2 {H}-isoquinolin-1-one (three-letter code: D4X) (formula: C₁₇H₂₃NO₃) (labeled as "Ligand of Interest" by depositor).



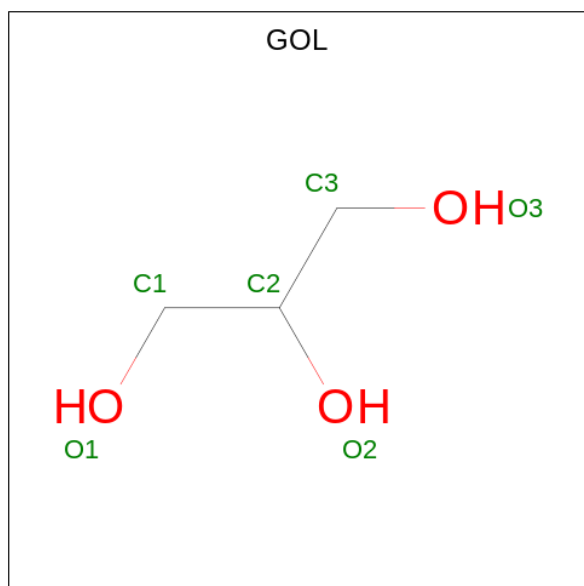
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			21	17	1	3		

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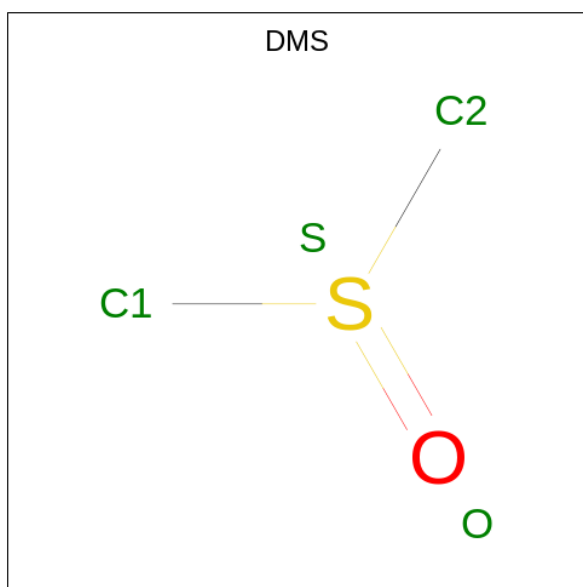
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	B	1	21	17	1	3	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	6	3	3	0	0
4	A	1	6	3	3	0	0
4	A	1	6	3	3	0	0
4	B	1	6	3	3	0	0
4	B	1	6	3	3	0	0

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
5	B	1	4	2	1	1	0	0

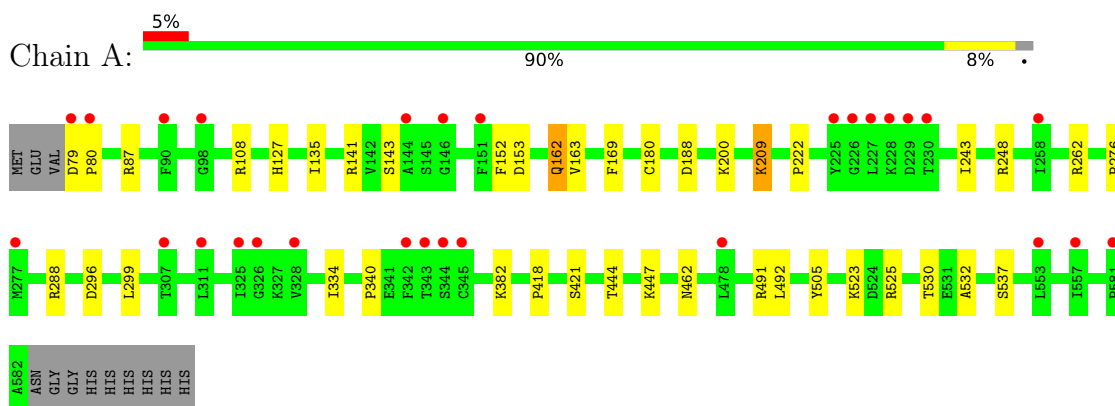
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	418	418	418	0	0
6	B	242	242	242	0	0

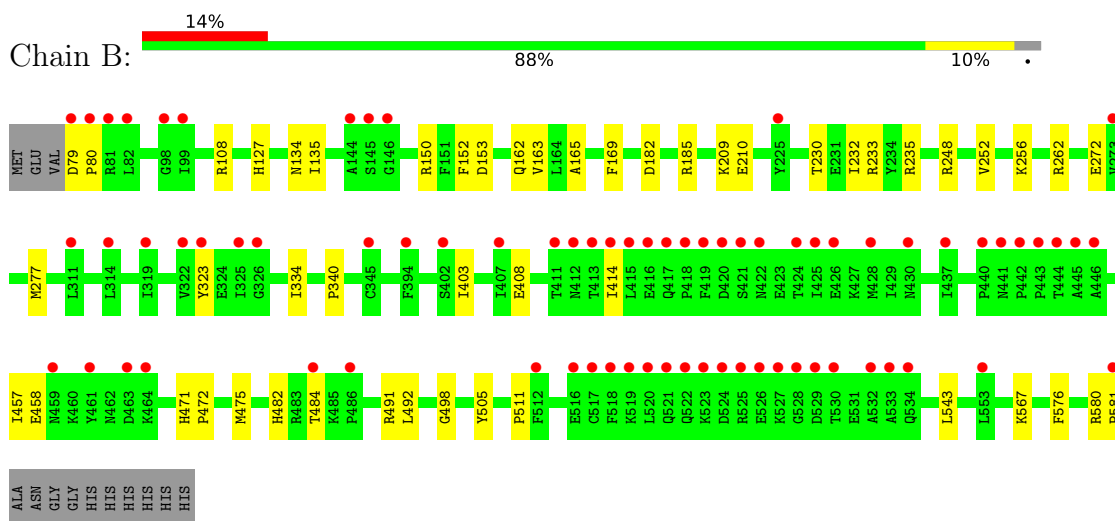
3 Residue-property plots [i](#)

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: Lysine-tRNA ligase



- Molecule 1: Lysine-tRNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.93Å 70.82Å 101.12Å 90.00° 111.45° 90.00°	Depositor
Resolution (Å)	43.84 – 1.92 47.06 – 1.92	Depositor EDS
% Data completeness (in resolution range)	99.0 (43.84-1.92) 99.0 (47.06-1.92)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 1.92Å)	Xtrriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.180 , 0.216 0.180 , 0.216	Depositor DCC
R_{free} test set	4257 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	29.8	Xtrriage
Anisotropy	0.290	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8696	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% 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: D4X, DMS, GOL

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	0.50	0/4136	0.62	2/5604 (0.0%)
1	B	0.43	0/3998	0.57	0/5428
All	All	0.46	0/8134	0.60	2/11032 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	ARG	NE-CZ-NH2	5.65	123.12	120.30
1	A	248	ARG	NE-CZ-NH1	-5.56	117.52	120.30

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	4037	0	3911	25	0
1	B	3903	0	3643	39	0
2	A	10	0	12	0	0
2	B	10	0	12	0	0
3	A	21	0	0	0	0
3	B	21	0	0	0	0
4	A	18	0	24	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	12	0	14	2	0
5	B	4	0	6	9	0
6	A	418	0	0	6	0
6	B	242	0	0	2	0
All	All	8696	0	7622	62	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 (62) 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:262:ARG:HH12	4:A:604:GOL:H11	1.38	0.89
1:B:248:ARG:HH21	5:B:605:DMS:C2	1.91	0.82
1:B:248:ARG:HH21	5:B:605:DMS:H23	1.47	0.80
1:A:525:ARG:HG3	1:A:532:ALA:HB3	1.62	0.79
1:A:169:PHE:O	1:A:209:LYS:HE2	1.89	0.71
1:B:230:THR:HB	1:B:232:ILE:HD13	1.73	0.69
1:B:248:ARG:HE	5:B:605:DMS:C2	2.05	0.69
1:B:169:PHE:O	1:B:209:LYS:HE2	1.93	0.68
1:B:182:ASP:O	1:B:185:ARG:NH1	2.24	0.68
1:B:403:ILE:HD12	1:B:457:ILE:HD12	1.79	0.65
1:B:262:ARG:HH12	4:B:604:GOL:H32	1.63	0.64
1:A:200:LYS:NZ	6:A:701:HOH:O	2.32	0.63
1:A:141:ARG:NH1	1:A:143:SER:OG	2.32	0.63
1:B:262:ARG:HH12	4:B:604:GOL:C3	2.13	0.62
1:B:408:GLU:HG2	1:B:414:ILE:HA	1.81	0.62
1:B:134:ASN:C	1:B:135:ILE:HD12	2.21	0.61
1:B:248:ARG:NH2	5:B:605:DMS:C2	2.64	0.60
1:A:79:ASP:HB3	1:A:80:PRO:HD3	1.84	0.59
1:A:153:ASP:OD1	1:A:162[A]:GLN:NE2	2.35	0.59
1:A:127:HIS:HE1	6:A:951:HOH:O	1.84	0.58
1:B:127:HIS:NE2	6:B:704:HOH:O	2.32	0.58
1:B:511:PRO:HB3	1:B:543:LEU:HB3	1.85	0.56
1:B:458:GLU:O	1:B:498:GLY:HA2	2.07	0.54
1:B:403:ILE:HG21	1:B:471:HIS:HA	1.89	0.53
1:B:248:ARG:NH2	5:B:605:DMS:H22	2.23	0.53
1:B:248:ARG:NH2	5:B:605:DMS:H23	2.22	0.52
1:B:153:ASP:OD1	1:B:162:GLN:NE2	2.43	0.51
1:A:108:ARG:HD2	6:A:830:HOH:O	2.11	0.51
1:B:248:ARG:NE	5:B:605:DMS:C2	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:LEU:HD13	1:B:581:PRO:HG3	1.95	0.48
1:B:230:THR:O	1:B:233:ARG:NH1	2.46	0.48
1:B:403:ILE:CD1	1:B:457:ILE:HD12	2.44	0.48
1:B:334:ILE:HG12	1:B:340:PRO:HD3	1.96	0.48
1:B:152:PHE:HB2	1:B:163:VAL:HB	1.95	0.47
1:A:444:THR:OG1	1:A:447:LYS:HG3	2.15	0.47
1:B:567:LYS:HD3	1:B:567:LYS:HA	1.78	0.47
1:A:288:ARG:HH21	1:A:530:THR:HG23	1.79	0.46
1:A:530:THR:HG22	6:A:1007:HOH:O	2.15	0.46
1:A:87:ARG:NH2	1:A:188:ASP:OD1	2.48	0.46
1:A:288:ARG:HH21	1:A:530:THR:CG2	2.28	0.46
1:B:248:ARG:HE	5:B:605:DMS:H23	1.79	0.46
1:A:276:PRO:HG3	1:B:576:PHE:HB2	1.98	0.45
1:A:262:ARG:NH1	4:A:604:GOL:H11	2.18	0.45
1:A:382:LYS:HD2	6:A:733:HOH:O	2.15	0.45
1:A:222:PRO:HG2	1:A:243:ILE:HG13	1.99	0.45
1:A:418:PRO:HG2	1:A:421:SER:HB3	1.98	0.45
1:B:252:VAL:HG12	1:B:256:LYS:HE2	1.98	0.45
1:B:482:HIS:CE1	1:B:484:THR:HG22	2.51	0.45
1:A:334:ILE:HG12	1:A:340:PRO:HD3	1.99	0.44
1:B:162:GLN:HG2	6:B:877:HOH:O	2.18	0.44
1:B:248:ARG:HH21	5:B:605:DMS:H22	1.73	0.43
1:B:491:ARG:HA	1:B:505:TYR:HB3	2.01	0.43
1:B:134:ASN:O	1:B:135:ILE:HD12	2.18	0.43
1:B:150:ARG:HB2	1:B:165:ALA:HB3	2.02	0.42
1:B:472:PRO:HD2	1:B:475:MET:SD	2.60	0.42
1:A:523:LYS:HB3	1:A:523:LYS:HE3	1.73	0.42
1:A:491:ARG:HA	1:A:505:TYR:HB3	2.02	0.42
1:B:79:ASP:HB3	1:B:80:PRO:HD3	2.02	0.42
1:A:152:PHE:HB2	1:A:163:VAL:HB	2.02	0.41
1:A:135:ILE:HG13	6:A:830:HOH:O	2.21	0.40
1:B:235:ARG:NH2	1:B:580:ARG:O	2.39	0.40
1:B:272:GLU:HB2	1:B:323:TYR:CZ	2.56	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	505/516 (98%)	494 (98%)	11 (2%)	0	100	100
1	B	503/516 (98%)	495 (98%)	8 (2%)	0	100	100
All	All	1008/1032 (98%)	989 (98%)	19 (2%)	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	431/464 (93%)	423 (98%)	8 (2%)	57	51
1	B	396/464 (85%)	392 (99%)	4 (1%)	76	75
All	All	827/928 (89%)	815 (98%)	12 (2%)	69	61

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162[A]	GLN
1	A	162[B]	GLN
1	A	180	CYS
1	A	209	LYS
1	A	296	ASP
1	A	462	ASN
1	A	492	LEU
1	A	537	SER

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Mol	Chain	Res	Type
1	B	108	ARG
1	B	210	GLU
1	B	277	MET
1	B	492	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

10 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	$\# Z > 2$
4	GOL	A	605	-	5,5,5	0.95	0	5,5,5	0.98	0
3	D4X	A	602	-	23,23,23	0.79	0	25,33,33	1.39	4 (16%)
4	GOL	B	603	-	5,5,5	1.25	1 (20%)	5,5,5	0.90	0
4	GOL	B	604	-	5,5,5	1.02	0	5,5,5	1.38	1 (20%)
4	GOL	A	604	-	5,5,5	0.60	0	5,5,5	1.08	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	D4X	B	602	-	23,23,23	0.84	1 (4%)	25,33,33	1.07	2 (8%)
2	LYS	A	601	-	8,9,9	0.90	0	9,10,10	0.89	0
5	DMS	B	605	-	3,3,3	0.75	0	3,3,3	1.30	0
2	LYS	B	601	-	8,9,9	0.91	0	9,10,10	1.07	0
4	GOL	A	603	-	5,5,5	1.15	0	5,5,5	0.86	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
4	GOL	A	605	-	-	3/4/4/4	-
3	D4X	A	602	-	-	0/4/26/26	0/3/3/3
4	GOL	B	603	-	-	2/4/4/4	-
4	GOL	B	604	-	-	3/4/4/4	-
4	GOL	A	604	-	-	1/4/4/4	-
3	D4X	B	602	-	-	0/4/26/26	0/3/3/3
2	LYS	A	601	-	-	0/9/9/9	-
2	LYS	B	601	-	-	0/9/9/9	-
4	GOL	A	603	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	603	GOL	O2-C2	-2.25	1.36	1.43
3	B	602	D4X	C9-C8	2.04	1.51	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	D4X	C10-C11-C15	-3.77	103.53	111.73
3	A	602	D4X	C17-C12-C13	2.94	114.04	109.41
3	B	602	D4X	C17-C12-C13	2.76	113.77	109.41
3	A	602	D4X	O4-C8-C9	-2.56	118.61	123.98
4	B	604	GOL	C3-C2-C1	-2.43	102.26	111.70
3	B	602	D4X	C10-C11-C15	-2.37	106.59	111.73
3	A	602	D4X	C3-C4-C5	-2.06	118.54	120.83
4	A	604	GOL	C3-C2-C1	-2.03	103.82	111.70

There are no chirality outliers.

All (9) torsion outliers are listed below:

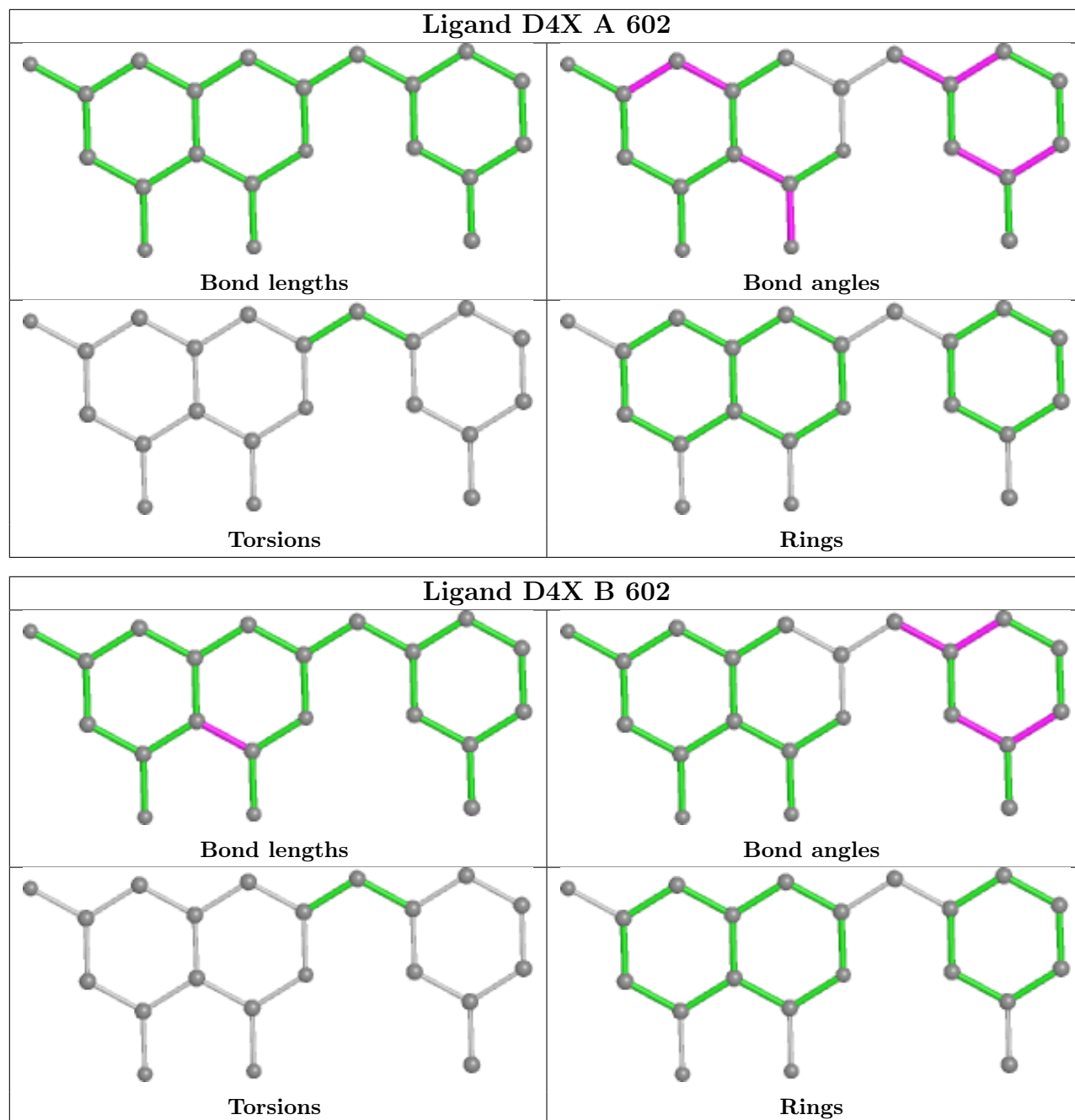
Mol	Chain	Res	Type	Atoms
4	A	605	GOL	C1-C2-C3-O3
4	B	603	GOL	O1-C1-C2-C3
4	B	604	GOL	O1-C1-C2-C3
4	A	604	GOL	C1-C2-C3-O3
4	A	605	GOL	O2-C2-C3-O3
4	B	603	GOL	O1-C1-C2-O2
4	A	605	GOL	O1-C1-C2-C3
4	B	604	GOL	O1-C1-C2-O2
4	B	604	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	604	GOL	2	0
4	A	604	GOL	2	0
5	B	605	DMS	9	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	504/516 (97%)	0.12	28 (5%) 24 27	20, 36, 67, 121	0
1	B	503/516 (97%)	0.56	74 (14%) 2 2	22, 48, 89, 105	0
All	All	1007/1032 (97%)	0.34	102 (10%) 7 8	20, 42, 85, 121	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	229	ASP	9.7
1	B	415	LEU	8.0
1	B	437	ILE	6.6
1	B	225	TYR	6.1
1	A	227	LEU	5.5
1	B	425	ILE	5.2
1	B	421	SER	4.9
1	B	530	THR	4.5
1	B	533	ALA	4.5
1	A	225	TYR	4.4
1	B	411	THR	4.4
1	B	426	GLU	4.3
1	B	523	LYS	4.1
1	B	443	PRO	4.0
1	B	82	LEU	4.0
1	B	517	CYS	3.8
1	B	424	THR	3.7
1	B	528	GLY	3.7
1	B	430	ASN	3.7
1	B	428	MET	3.6
1	B	524	ASP	3.6
1	B	526	GLU	3.6
1	B	98	GLY	3.6
1	B	520	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	80	PRO	3.5
1	B	412	ASN	3.5
1	B	446	ALA	3.5
1	A	146	GLY	3.5
1	B	518	PHE	3.5
1	B	419	PHE	3.4
1	A	151	PHE	3.3
1	B	445	ALA	3.3
1	B	81	ARG	3.3
1	B	522	GLN	3.2
1	B	146	GLY	3.2
1	A	79	ASP	3.2
1	B	322	VAL	3.2
1	A	228	LYS	3.2
1	B	311	LEU	3.1
1	B	407	ILE	3.1
1	B	414	ILE	3.1
1	B	521	GLN	3.1
1	B	417	GLN	3.1
1	B	444	THR	3.0
1	B	144	ALA	3.0
1	A	325	ILE	3.0
1	B	461	TYR	3.0
1	B	99	ILE	2.9
1	A	345	CYS	2.8
1	B	325	ILE	2.8
1	B	145	SER	2.8
1	B	418	PRO	2.8
1	A	328	VAL	2.8
1	B	525	ARG	2.7
1	A	342	PHE	2.7
1	B	442	PRO	2.7
1	A	343	THR	2.7
1	B	463	ASP	2.6
1	B	323	TYR	2.6
1	B	441	ASN	2.6
1	A	230	THR	2.6
1	B	527	LYS	2.6
1	A	90	PHE	2.5
1	B	416	GLU	2.5
1	B	516	GLU	2.5
1	B	532	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	307	THR	2.5
1	B	314	LEU	2.5
1	B	413	THR	2.4
1	B	440	PRO	2.4
1	B	422	ASN	2.4
1	B	420	ASP	2.4
1	A	326	GLY	2.4
1	B	529	ASP	2.3
1	B	519	LYS	2.3
1	B	345	CYS	2.3
1	A	98	GLY	2.3
1	A	478	LEU	2.3
1	A	144	ALA	2.3
1	A	581	PRO	2.3
1	B	326	GLY	2.2
1	B	581	PRO	2.2
1	A	557	ILE	2.2
1	A	226	GLY	2.2
1	A	311	LEU	2.2
1	B	394	PHE	2.2
1	B	79	ASP	2.1
1	B	464	LYS	2.1
1	B	534	GLN	2.1
1	A	277	MET	2.1
1	B	484	THR	2.1
1	A	80	PRO	2.1
1	B	459	ASN	2.1
1	B	273	VAL	2.1
1	B	402	SER	2.1
1	A	258	ILE	2.1
1	B	486	PRO	2.0
1	B	319	ILE	2.0
1	B	553	LEU	2.0
1	B	512	PHE	2.0
1	A	553	LEU	2.0
1	A	344	SER	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 [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

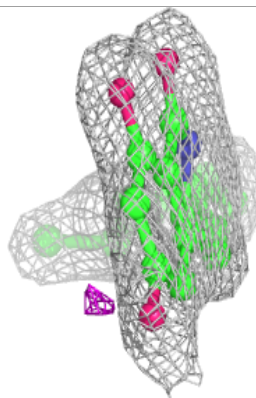
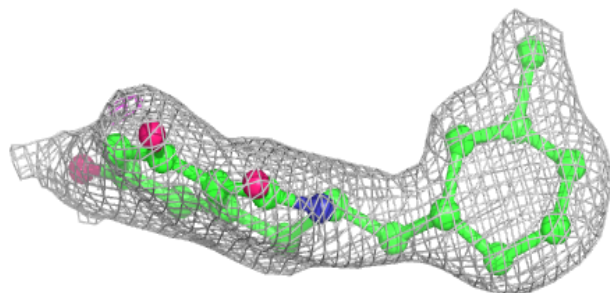
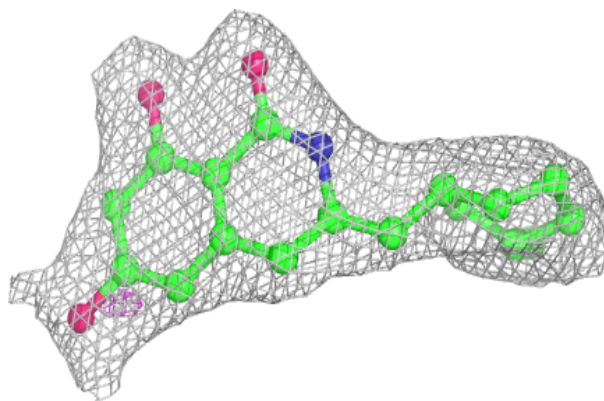
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	DMS	B	605	4/4	0.66	0.23	72,74,78,81	0
4	GOL	A	605	6/6	0.75	0.19	43,53,53,55	0
4	GOL	A	604	6/6	0.89	0.37	48,49,55,61	0
4	GOL	B	604	6/6	0.90	0.24	39,47,50,53	0
3	D4X	B	602	21/21	0.92	0.11	34,40,43,44	0
3	D4X	A	602	21/21	0.94	0.16	20,29,37,41	0
2	LYS	B	601	10/10	0.95	0.17	32,34,39,39	0
4	GOL	A	603	6/6	0.96	0.07	32,34,36,37	0
4	GOL	B	603	6/6	0.97	0.09	42,42,43,44	0
2	LYS	A	601	10/10	0.98	0.21	22,24,26,28	0

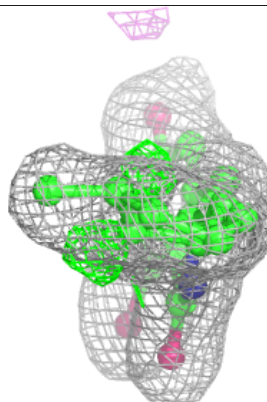
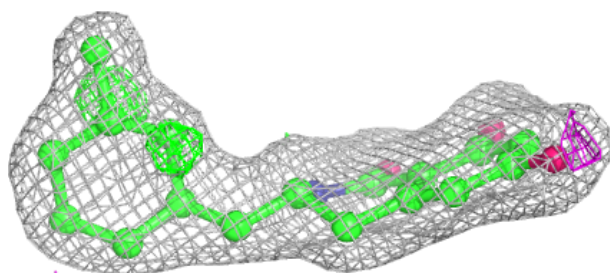
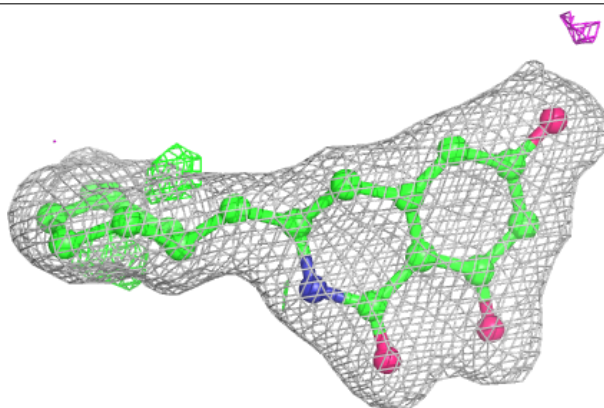
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around D4X B 602:

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

**Electron density around D4X A 602:**

$2mF_o-DF_c$ (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.