



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 20, 2024 – 04:54 AM EST

PDB ID : 4MUJ
Title : Crystal structure of pantothenate synthetase in complex with 2-(2-(benzylsulfonylcarbonyl)-5-methoxy-1H-indol-1-yl)acetic acid
Authors : Silvestre, H.L.; Blundell, T.L.
Deposited on : 2013-09-22
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)
Xtriage (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

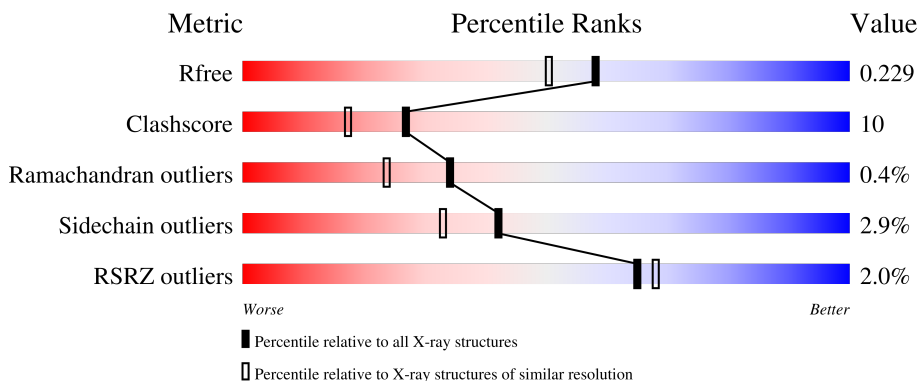
1 Overall quality at a glance

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	300	
1	B	300	

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
2	EOH	A	403	-	-	X	-
2	EOH	B	407	-	-	X	-
4	EDO	A	407	-	-	X	-
5	GOL	B	403	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4729 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.

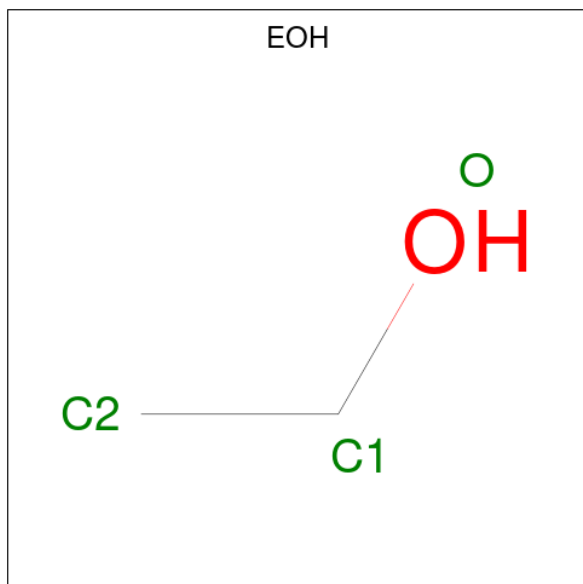
- Molecule 1 is a protein called Pantothenate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	288	Total 2185	C 1378	N 395	O 406	S 6	0	9	0
1	B	272	Total 2049	C 1295	N 369	O 379	S 6	0	6	0

There are 4 discrepancies between the modelled and reference sequences:

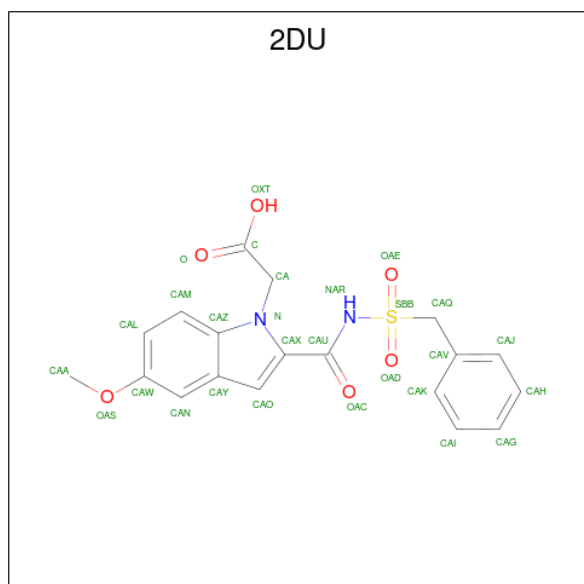
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ALA	THR	engineered mutation	UNP P0A5R0
A	77	GLY	GLU	engineered mutation	UNP P0A5R0
B	2	ALA	THR	engineered mutation	UNP P0A5R0
B	77	GLY	GLU	engineered mutation	UNP P0A5R0

- Molecule 2 is ETHANOL (three-letter code: EOH) (formula: C₂H₆O).



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

- Molecule 3 is {2-[(benzylsulfonyl)carbamoyl]-5-methoxy-1H-indol-1-yl}acetic acid (three-letter code: 2DU) (formula: C₁₉H₁₈N₂O₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			28	19	2	6	1		
3	B	1	Total	C	N	O	S	0	0
			28	19	2	6	1		
3	B	1	Total	C	N	O	S	0	0
			28	19	2	6	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

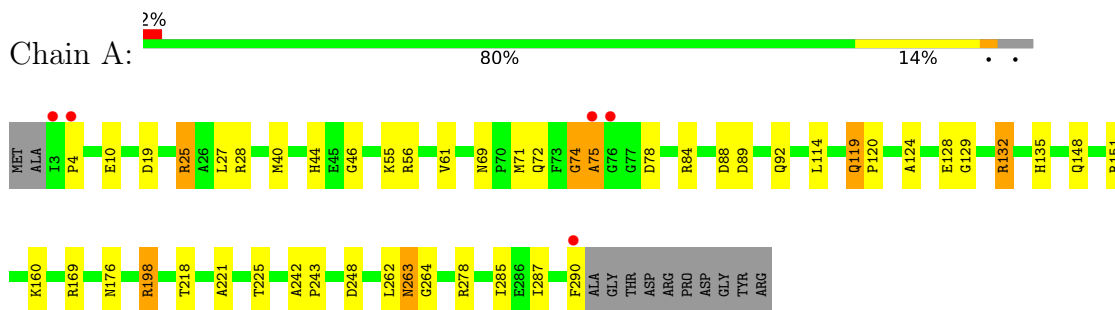
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	192	Total	O	0	0
			192	192		
6	B	175	Total	O	0	0
			175	175		

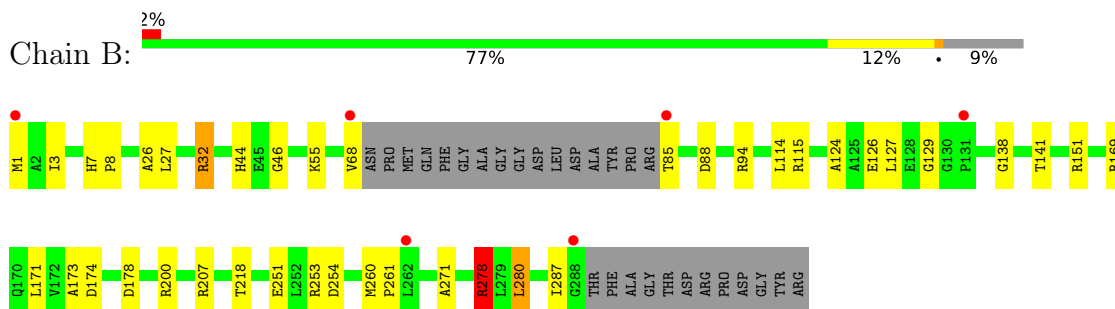
3 Residue-property plots [i](#)

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: Pantothenate synthetase



- Molecule 1: Pantothenate synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.73Å 70.77Å 81.82Å 90.00° 99.36° 90.00°	Depositor
Resolution (Å)	98.89 – 1.92 40.37 – 1.92	Depositor EDS
% Data completeness (in resolution range)	98.9 (98.89-1.92) 98.9 (40.37-1.92)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.26 (at 1.92Å)	Xtrriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.171 , 0.229 0.172 , 0.229	Depositor DCC
R_{free} test set	2093 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	19.5	Xtrriage
Anisotropy	0.088	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4729	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% 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: EOH, 2DU, EDO, 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	1.05	0/2224	0.97	6/3037 (0.2%)
1	B	1.08	1/2085 (0.0%)	0.96	7/2846 (0.2%)
All	All	1.07	1/4309 (0.0%)	0.97	13/5883 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	173	ALA	CA-CB	5.22	1.63	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	84	ARG	NE-CZ-NH1	6.34	123.47	120.30
1	B	207	ARG	NE-CZ-NH1	-6.15	117.22	120.30
1	B	278	ARG	NE-CZ-NH1	-6.01	117.29	120.30
1	A	40	MET	CG-SD-CE	-5.71	91.06	100.20
1	B	94	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	B	169	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	B	280	LEU	CA-CB-CG	5.21	127.29	115.30
1	B	171	LEU	CB-CG-CD1	-5.17	102.21	111.00
1	A	169	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	A	89	ASP	CB-CG-OD1	5.06	122.85	118.30
1	A	132[A]	ARG	NE-CZ-NH2	-5.03	117.78	120.30
1	A	132[B]	ARG	NE-CZ-NH2	-5.03	117.78	120.30
1	B	32	ARG	NE-CZ-NH1	5.02	122.81	120.30

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	2185	0	2215	51	0
1	B	2049	0	2105	36	0
2	A	15	0	30	5	0
2	B	3	0	6	5	0
3	A	28	0	17	2	0
3	B	56	0	34	3	0
4	A	8	0	12	5	0
4	B	12	0	18	1	0
5	B	6	0	8	5	0
6	A	192	0	0	6	0
6	B	175	0	0	5	0
All	All	4729	0	4445	90	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 10.

All (90) 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:B:126[B]:GLU:HA	1:B:126[B]:GLU:OE1	1.43	1.11
1:B:126[B]:GLU:OE1	1:B:126[B]:GLU:CA	1.96	1.10
1:A:148:GLN:HE22	2:B:407:EOH:H21	1.16	1.06
1:A:74:GLY:HA3	1:A:75:ALA:HB2	1.38	1.06
1:A:263:ASN:CG	1:A:264:GLY:H	1.59	1.05
1:B:251[B]:GLU:OE1	1:B:253[B]:ARG:NH1	1.96	0.98
1:A:148:GLN:NE2	2:B:407:EOH:H21	1.78	0.97
2:A:403:EOH:H22	6:A:690:HOH:O	1.66	0.95
1:B:32:ARG:HD3	3:B:402:2DU:OXT	1.76	0.86
1:A:25:ARG:HH11	1:A:25:ARG:HG2	1.40	0.85
1:A:71:MET:SD	1:A:114:LEU:HD11	2.17	0.84
1:A:128:GLU:O	1:A:132[B]:ARG:HD2	1.78	0.84
1:A:263:ASN:ND2	1:A:264:GLY:H	1.76	0.82
1:A:263:ASN:CG	1:A:264:GLY:N	2.34	0.80
1:B:218[B]:THR:HG21	6:B:504:HOH:O	1.80	0.80
1:A:74:GLY:HA3	1:A:75:ALA:CB	2.13	0.79

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:HIS:HD2	1:A:46:GLY:H	1.29	0.76
1:B:114:LEU:HD11	5:B:403:GOL:H2	1.67	0.76
1:A:176:ASN:HD22	2:A:403:EOH:H12	1.49	0.75
1:A:69:ASN:HD22	1:A:72:GLN:HE21	1.33	0.75
1:A:176:ASN:HD22	2:A:403:EOH:C1	2.01	0.72
1:A:44:HIS:HE1	3:A:404:2DU:OXT	1.72	0.72
1:B:85:THR:HB	1:B:88:ASP:OD1	1.91	0.71
1:A:135:HIS:N	4:A:407:EDO:H11	2.06	0.71
1:B:44:HIS:HD2	1:B:46:GLY:H	1.40	0.70
1:B:126[B]:GLU:OE1	1:B:126[B]:GLU:N	2.25	0.69
1:A:44:HIS:HD2	1:A:46:GLY:N	1.90	0.68
1:B:254:ASP:HB3	1:B:261:PRO:HD3	1.74	0.68
1:A:25:ARG:HG2	1:A:25:ARG:NH1	2.08	0.67
1:A:75:ALA:HB3	1:A:78:ASP:OD2	1.97	0.65
1:B:278:ARG:HG2	1:B:278:ARG:HH11	1.61	0.65
1:B:278:ARG:HG2	1:B:278:ARG:NH1	2.15	0.61
1:A:128:GLU:O	1:A:132[B]:ARG:CD	2.50	0.60
1:B:151:ARG:CZ	6:B:611:HOH:O	2.49	0.60
1:A:263:ASN:ND2	1:A:264:GLY:N	2.47	0.59
1:B:178[B]:ASP:H	2:B:407:EOH:H22	1.68	0.59
1:B:138:GLY:HA2	5:B:403:GOL:H31	1.85	0.58
1:A:221:ALA:HA	1:A:285:ILE:HD13	1.86	0.57
1:B:178[A]:ASP:OD1	1:B:178[A]:ASP:O	2.22	0.57
1:A:151:ARG:HD2	6:A:659:HOH:O	2.05	0.56
4:A:405:EDO:H12	1:B:174:ASP:HB3	1.87	0.56
1:A:218[B]:THR:HG23	6:A:670:HOH:O	2.07	0.54
1:A:44:HIS:CD2	1:A:46:GLY:H	2.19	0.53
1:A:176:ASN:HB3	2:A:403:EOH:H11	1.91	0.52
1:A:263:ASN:HB3	6:A:615:HOH:O	2.10	0.52
1:A:198:ARG:HH11	1:A:198:ARG:HB3	1.74	0.52
1:B:44:HIS:HD2	1:B:46:GLY:N	2.06	0.52
1:A:160[A]:LYS:HE3	6:A:532:HOH:O	2.10	0.51
1:B:114:LEU:HD11	5:B:403:GOL:H12	1.91	0.51
1:A:135:HIS:H	4:A:407:EDO:H11	1.77	0.50
1:B:260:MET:CE	1:B:287:ILE:HG12	2.41	0.50
1:B:114:LEU:HD11	5:B:403:GOL:C2	2.41	0.49
1:A:75:ALA:CB	1:A:78:ASP:OD2	2.61	0.49
1:A:151:ARG:HH12	2:B:407:EOH:C2	2.26	0.49
1:A:71:MET:SD	1:A:114:LEU:CD1	2.97	0.48
1:A:262:LEU:HD21	1:A:287:ILE:HG13	1.94	0.48
1:B:124:ALA:O	1:B:129:GLY:HA3	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:LEU:HD22	1:B:32:ARG:HG3	1.95	0.48
1:B:178[A]:ASP:H	2:B:407:EOH:H22	1.77	0.48
1:B:260:MET:SD	1:B:261:PRO:HD2	2.55	0.47
1:A:225:THR:HG22	1:A:290:PHE:CZ	2.50	0.47
1:A:119:GLN:HE21	1:A:119:GLN:HA	1.80	0.47
1:A:132[B]:ARG:NH1	1:A:135:HIS:CD2	2.83	0.46
1:A:132[A]:ARG:HB3	4:A:407:EDO:H12	1.97	0.46
1:A:25:ARG:NH1	6:A:517:HOH:O	2.48	0.46
1:A:198:ARG:HH11	1:A:198:ARG:CB	2.31	0.43
1:A:27:LEU:HD13	1:A:61:VAL:HG11	2.00	0.43
1:B:3:ILE:HD13	1:B:26:ALA:HB2	2.00	0.43
1:B:151:ARG:NH2	6:B:611:HOH:O	2.52	0.43
1:A:114:LEU:HD12	1:A:114:LEU:HA	1.68	0.42
1:A:120:PRO:HB2	1:A:124:ALA:HB2	2.01	0.42
1:B:141:THR:HG21	5:B:403:GOL:H11	2.00	0.42
1:A:248[B]:ASP:OD2	1:A:278:ARG:NH2	2.44	0.42
1:B:126[B]:GLU:HB3	1:B:127:LEU:HG	2.01	0.42
3:B:402:2DU:OAC	3:B:402:2DU:CA	2.68	0.42
1:B:7:HIS:HA	1:B:8:PRO:HD3	1.83	0.41
1:B:278:ARG:HH11	1:B:278:ARG:CG	2.26	0.41
1:A:242:ALA:HA	1:A:243:PRO:HD2	1.68	0.41
1:B:115:ARG:H	4:B:406:EDO:C2	2.33	0.41
1:B:280:LEU:HD12	6:B:665:HOH:O	2.19	0.41
3:B:402:2DU:OAC	3:B:402:2DU:H12	2.19	0.41
1:A:124:ALA:O	1:A:129:GLY:HA3	2.21	0.41
1:B:68:VAL:HG13	6:B:674:HOH:O	2.20	0.41
1:A:71:MET:HG3	2:A:408:EOH:H23	2.02	0.41
3:A:404:2DU:H4	3:A:404:2DU:H3	1.91	0.40
1:A:151:ARG:CZ	1:B:178[B]:ASP:OD2	2.70	0.40
1:A:88[A]:ASP:O	1:A:92:GLN:HG3	2.21	0.40
1:A:132[A]:ARG:CA	4:A:407:EDO:H12	2.51	0.40
1:A:4:PRO:HG3	1:A:19:ASP:HB3	2.04	0.40
1:B:271:ALA:HB2	1:B:280:LEU:HD22	2.04	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	293/300 (98%)	284 (97%)	7 (2%)	2 (1%)	22	11
1	B	274/300 (91%)	270 (98%)	4 (2%)	0	100	100
All	All	567/600 (94%)	554 (98%)	11 (2%)	2 (0%)	34	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	74	GLY
1	A	75	ALA

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	222/223 (100%)	214 (96%)	8 (4%)	35	24
1	B	209/223 (94%)	205 (98%)	4 (2%)	57	51
All	All	431/446 (97%)	419 (97%)	12 (3%)	42	34

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

Mol	Chain	Res	Type
1	A	10	GLU
1	A	25	ARG
1	A	28	ARG
1	A	55	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	56	ARG
1	A	119	GLN
1	A	198	ARG
1	A	263	ASN
1	B	1	MET
1	B	55	LYS
1	B	200	ARG
1	B	278	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS
1	A	72	GLN
1	A	119	GLN
1	A	148	GLN
1	A	222	HIS
1	B	44	HIS

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](#)

15 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	EOH	B	407	-	2,2,2	0.41	0	1,1,1	0.32	0
2	EOH	A	402	-	2,2,2	0.45	0	1,1,1	0.36	0
2	EOH	A	406	-	2,2,2	0.44	0	1,1,1	0.04	0
4	EDO	A	405	-	3,3,3	0.09	0	2,2,2	0.83	0
5	GOL	B	403	-	5,5,5	0.32	0	5,5,5	0.89	0
3	2DU	B	402	-	28,30,30	3.18	13 (46%)	34,43,43	1.96	5 (14%)
2	EOH	A	403	-	2,2,2	0.51	0	1,1,1	0.07	0
4	EDO	B	404	-	3,3,3	0.41	0	2,2,2	0.48	0
4	EDO	A	407	-	3,3,3	0.29	0	2,2,2	0.75	0
4	EDO	B	405	-	3,3,3	0.45	0	2,2,2	0.42	0
3	2DU	A	404	-	28,30,30	2.75	15 (53%)	34,43,43	2.48	6 (17%)
3	2DU	B	401	-	28,30,30	2.57	8 (28%)	34,43,43	1.65	7 (20%)
4	EDO	B	406	-	3,3,3	0.41	0	2,2,2	1.11	0
2	EOH	A	408	-	2,2,2	0.38	0	1,1,1	0.52	0
2	EOH	A	401	-	2,2,2	0.57	0	1,1,1	0.00	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	EDO	A	405	-	-	1/1/1/1	-
5	GOL	B	403	-	-	2/4/4/4	-
3	2DU	B	402	-	-	2/17/20/20	0/3/3/3
4	EDO	A	407	-	-	1/1/1/1	-
4	EDO	B	405	-	-	0/1/1/1	-
3	2DU	A	404	-	-	2/17/20/20	0/3/3/3
3	2DU	B	401	-	-	1/17/20/20	0/3/3/3
4	EDO	B	406	-	-	1/1/1/1	-
4	EDO	B	404	-	-	1/1/1/1	-

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	2DU	OAE-SBB	8.28	1.55	1.43
3	B	402	2DU	OAD-SBB	7.90	1.54	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	2DU	OAE-SBB	7.06	1.53	1.43
3	A	404	2DU	CAQ-CAV	-6.89	1.39	1.50
3	B	402	2DU	CA-C	6.47	1.59	1.51
3	B	401	2DU	CAQ-CAV	-6.27	1.40	1.50
3	A	404	2DU	CAU-NAR	-6.21	1.31	1.39
3	B	401	2DU	OAD-SBB	5.87	1.51	1.43
3	B	402	2DU	CAQ-CAV	-5.68	1.41	1.50
3	B	401	2DU	SBB-NAR	3.93	1.71	1.63
3	A	404	2DU	CAY-CAZ	-3.77	1.33	1.41
3	A	404	2DU	CA-N	-3.76	1.41	1.49
3	A	404	2DU	CAM-CAZ	-3.63	1.33	1.41
3	A	404	2DU	CAO-CAX	-3.61	1.34	1.39
3	A	404	2DU	SBB-NAR	3.21	1.70	1.63
3	B	402	2DU	CAM-CAZ	-3.15	1.34	1.41
3	B	402	2DU	SBB-NAR	3.14	1.69	1.63
3	B	401	2DU	CAO-CAX	-2.90	1.35	1.39
3	A	404	2DU	CAN-CAY	-2.83	1.35	1.42
3	B	402	2DU	CAN-CAW	2.81	1.42	1.37
3	A	404	2DU	OAS-CAA	-2.70	1.34	1.42
3	B	402	2DU	CAO-CAX	-2.69	1.35	1.39
3	B	402	2DU	CAN-CAY	-2.69	1.35	1.42
3	B	402	2DU	CAU-NAR	-2.64	1.36	1.39
3	B	402	2DU	CAY-CAZ	-2.58	1.35	1.41
3	A	404	2DU	OXT-C	-2.54	1.22	1.30
3	B	401	2DU	CAU-NAR	-2.49	1.36	1.39
3	B	401	2DU	CAN-CAY	-2.37	1.36	1.42
3	A	404	2DU	CAX-CAU	-2.30	1.36	1.50
3	A	404	2DU	OAC-CAU	-2.27	1.18	1.23
3	A	404	2DU	CAX-N	-2.24	1.32	1.36
3	A	404	2DU	CAN-CAW	2.19	1.41	1.37
3	A	404	2DU	OAS-CAW	-2.14	1.33	1.37
3	B	401	2DU	CAX-CAU	-2.09	1.37	1.50
3	B	402	2DU	CAX-CAU	-2.09	1.37	1.50
3	B	402	2DU	CA-N	2.01	1.54	1.49

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	404	2DU	C-CA-N	11.31	126.34	111.52
3	B	402	2DU	OAE-SBB-OAD	-6.71	109.63	119.35
3	B	402	2DU	CAU-NAR-SBB	-6.34	115.37	123.38
3	A	404	2DU	CAL-CAW-CAN	-4.57	114.67	120.81

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	2DU	OAE-SBB-OAD	-4.44	112.92	119.35
3	A	404	2DU	OAE-SBB-OAD	-4.41	112.95	119.35
3	B	401	2DU	CAA-OAS-CAW	3.92	126.01	117.51
3	B	402	2DU	CAV-CAQ-SBB	3.43	119.67	112.34
3	B	401	2DU	CAO-CAY-CAZ	3.19	109.05	106.27
3	B	401	2DU	CAU-NAR-SBB	-2.95	119.65	123.38
3	B	401	2DU	OAE-SBB-CAQ	2.89	112.70	108.30
3	A	404	2DU	CAX-CAO-CAY	2.69	109.91	106.55
3	B	402	2DU	CAQ-SBB-NAR	2.62	112.32	105.07
3	B	401	2DU	CAL-CAW-CAN	-2.35	117.66	120.81
3	B	401	2DU	CAL-CAM-CAZ	-2.34	116.11	119.70
3	A	404	2DU	OXT-C-CA	2.25	119.93	112.29
3	B	402	2DU	CAM-CAL-CAW	-2.24	117.16	120.17
3	A	404	2DU	O-C-CA	-2.06	114.28	121.69

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	404	2DU	CAU-NAR-SBB-OAD
3	A	404	2DU	CAU-NAR-SBB-CAQ
3	B	401	2DU	CAU-NAR-SBB-CAQ
3	B	402	2DU	CAU-NAR-SBB-OAE
3	B	402	2DU	CAU-NAR-SBB-CAQ
5	B	403	GOL	O1-C1-C2-C3
4	A	405	EDO	O1-C1-C2-O2
5	B	403	GOL	O1-C1-C2-O2
4	A	407	EDO	O1-C1-C2-O2
4	B	406	EDO	O1-C1-C2-O2
4	B	404	EDO	O1-C1-C2-O2

There are no ring outliers.

9 monomers are involved in 26 short contacts:

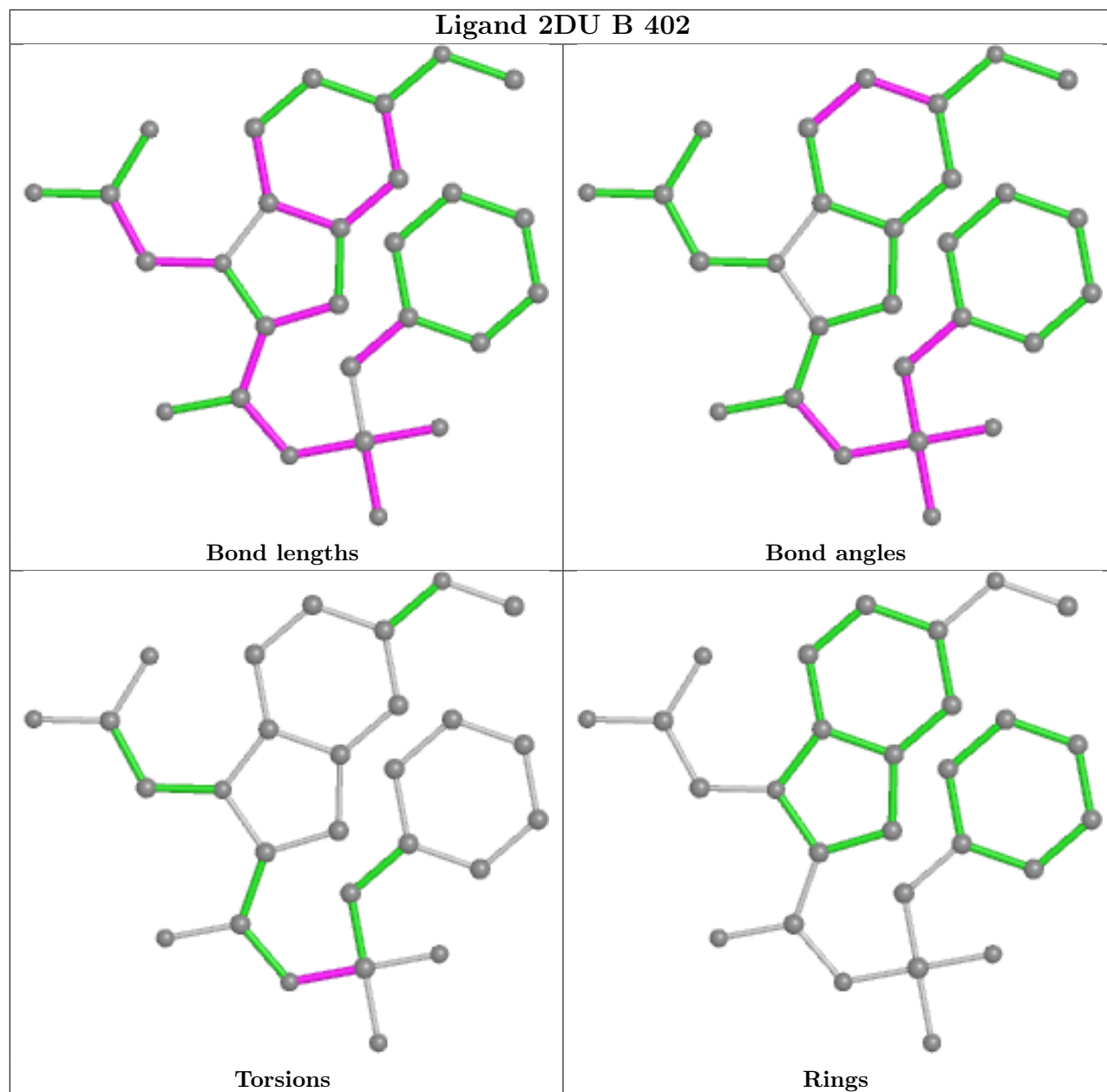
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	407	EOH	5	0
4	A	405	EDO	1	0
5	B	403	GOL	5	0
3	B	402	2DU	3	0
2	A	403	EOH	4	0
4	A	407	EDO	4	0

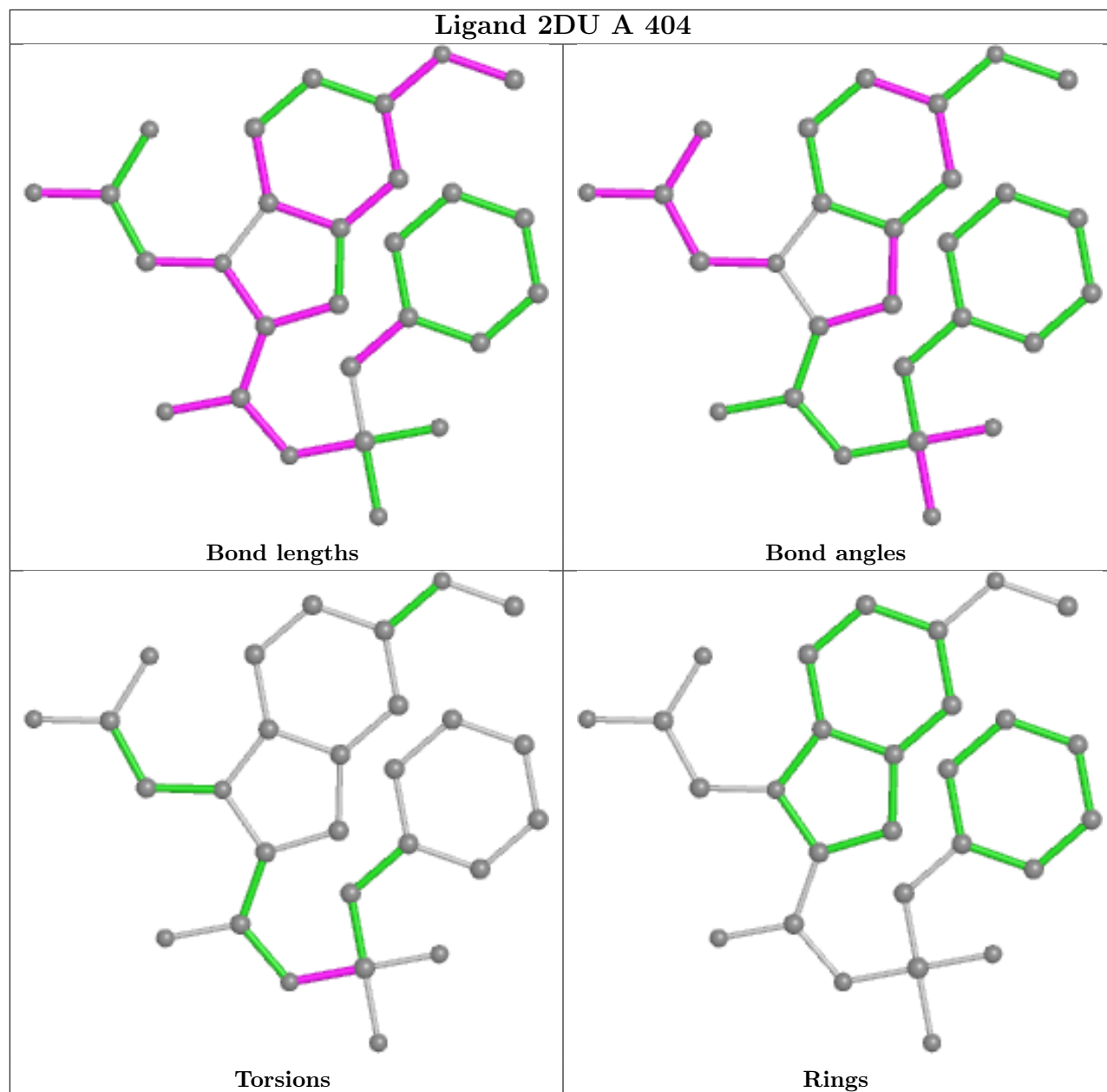
Continued on next page...

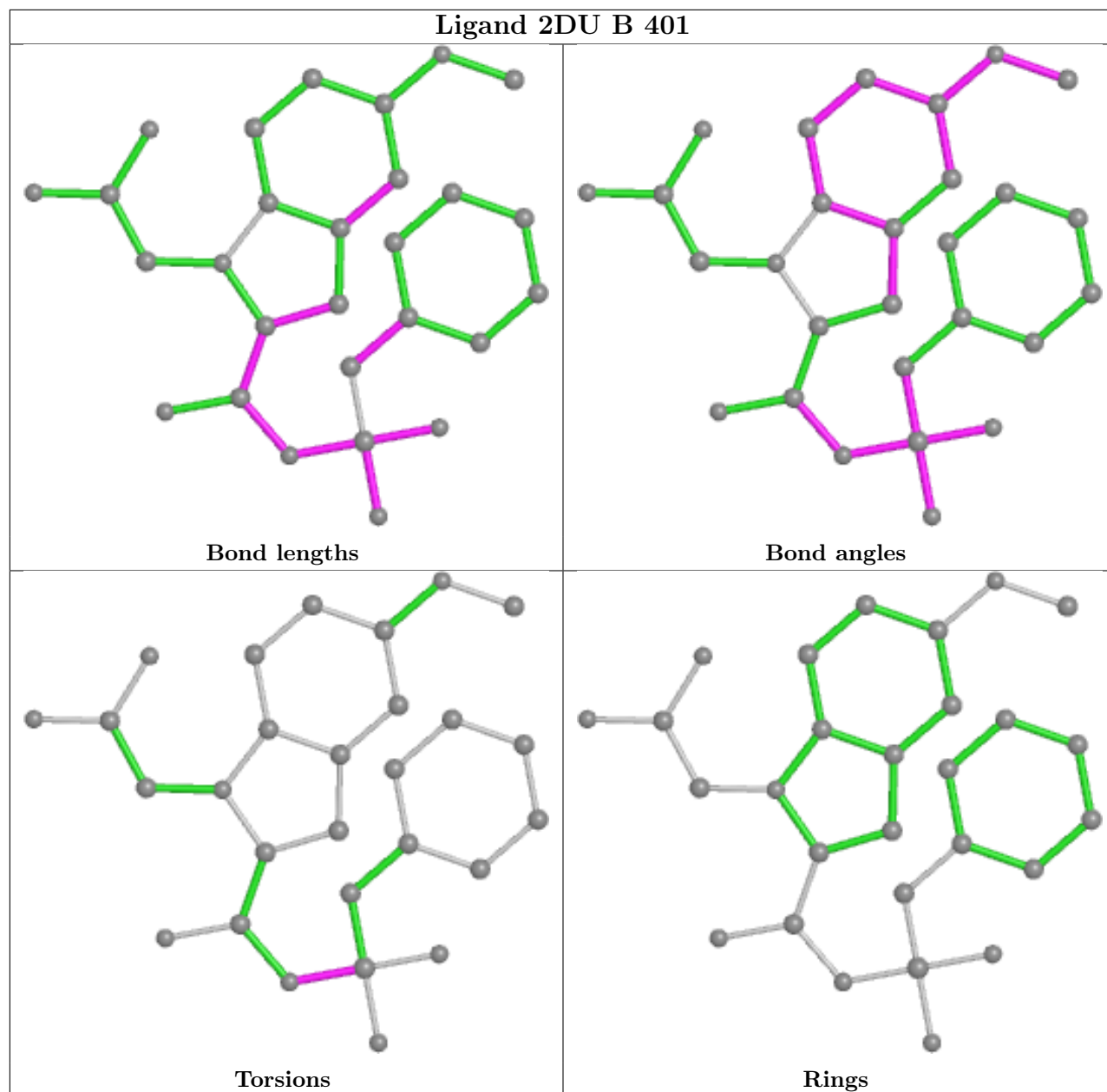
Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	404	2DU	2	0
4	B	406	EDO	1	0
2	A	408	EOH	1	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	288/300 (96%)	-0.28	5 (1%) 70 72	11, 22, 41, 53	3 (1%)
1	B	272/300 (90%)	-0.30	6 (2%) 62 65	11, 20, 42, 57	0
All	All	560/600 (93%)	-0.29	11 (1%) 65 68	11, 21, 42, 57	3 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	68	VAL	6.4
1	A	3	ILE	5.7
1	B	288	GLY	4.7
1	B	262	LEU	3.5
1	B	85	THR	3.3
1	B	131	PRO	3.2
1	A	290	PHE	3.1
1	A	76	GLY	3.1
1	B	1	MET	2.7
1	A	4	PRO	2.2
1	A	75	ALA	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

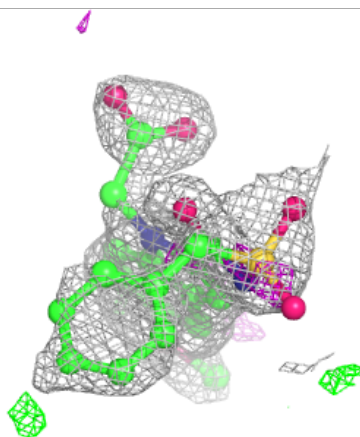
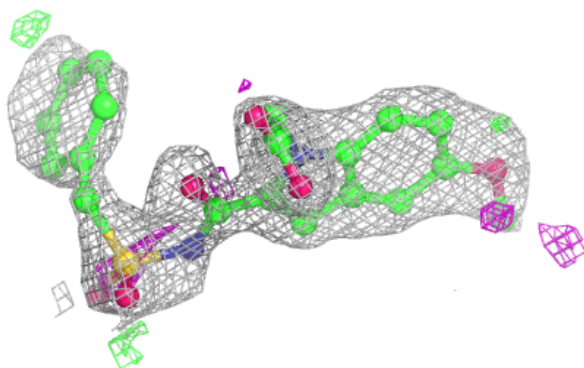
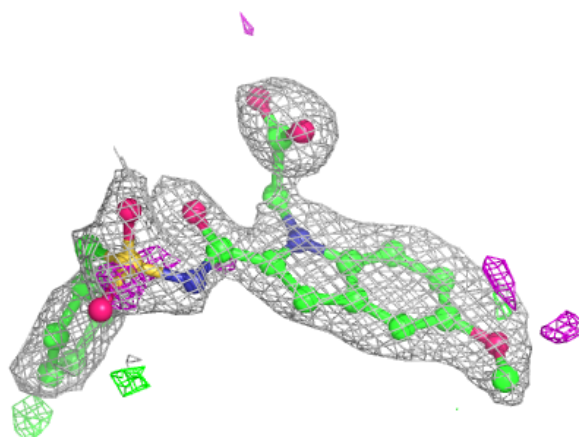
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	EOH	A	406	3/3	0.79	0.22	37,37,39,39	0
2	EOH	A	402	3/3	0.80	0.25	42,42,43,43	0
3	2DU	B	402	28/28	0.82	0.17	33,38,55,58	0
4	EDO	B	405	4/4	0.88	0.15	48,48,49,52	0
2	EOH	A	403	3/3	0.90	0.13	39,39,40,41	0
2	EOH	A	401	3/3	0.90	0.17	36,36,37,38	0
5	GOL	B	403	6/6	0.90	0.22	36,39,41,44	0
4	EDO	A	407	4/4	0.92	0.13	30,33,36,41	0
4	EDO	B	404	4/4	0.93	0.14	39,39,43,44	0
2	EOH	A	408	3/3	0.93	0.20	37,37,39,40	0
3	2DU	A	404	28/28	0.93	0.10	15,21,27,29	0
4	EDO	B	406	4/4	0.95	0.15	29,32,33,35	0
4	EDO	A	405	4/4	0.96	0.11	23,24,29,30	0
2	EOH	B	407	3/3	0.96	0.09	38,38,40,40	0
3	2DU	B	401	28/28	0.97	0.09	15,21,26,28	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.

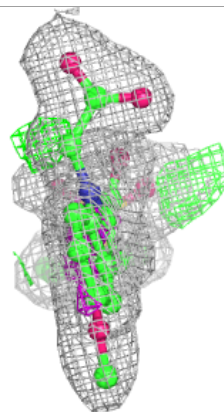
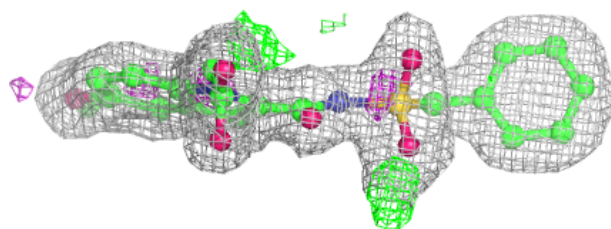
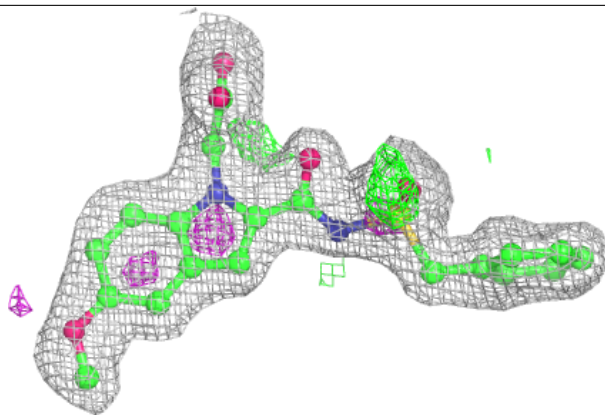
Electron density around 2DU B 402:

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

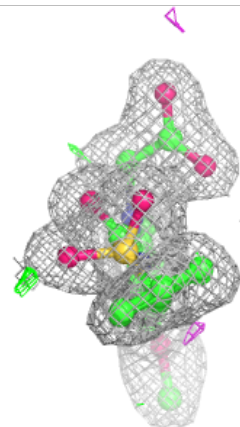
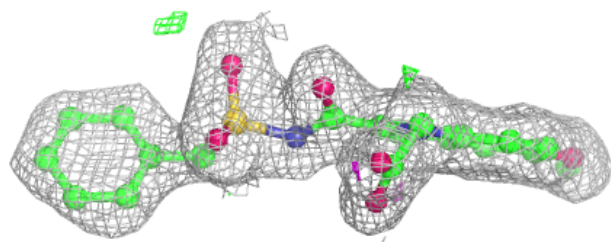
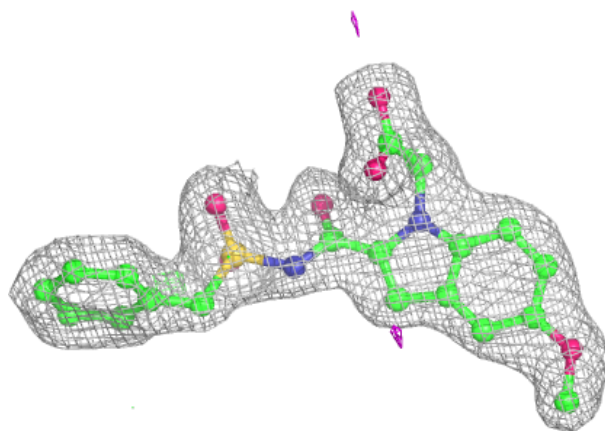


Electron density around 2DU A 404:

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

**Electron density around 2DU B 401:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.