



Full wwPDB X-ray Structure Validation Report

May 25, 2020 – 08:23 am BST

PDB ID : 4W1O
Title : PDE4D complexed with inhibitor
Authors : Sorensen, M.D.
Deposited on : 2014-08-14
Resolution : 2.20 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 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 : 2.11
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) : 2.11

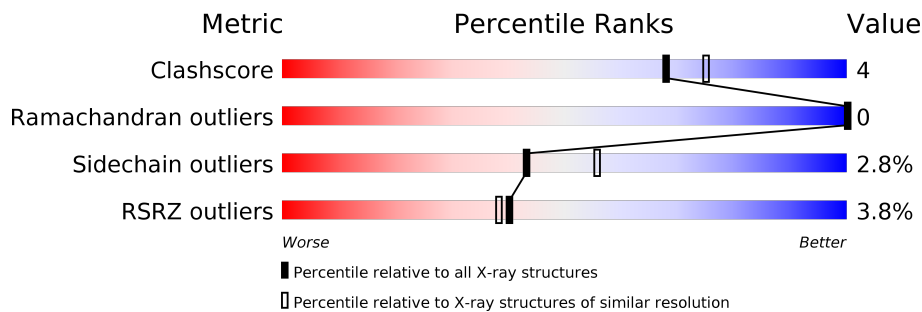
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 on 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	359	<div style="display: flex; align-items: center;"> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 4%, orange 4%, yellow 9%, green 81%, grey 100%);"></div> <div style="margin-left: 10px;"> <p>4%</p> <p>81%</p> <p>9%</p> <p>10%</p> </div> </div>
1	B	359	<div style="display: flex; align-items: center;"> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 4%, orange 4%, yellow 10%, green 80%, grey 100%);"></div> <div style="margin-left: 10px;"> <p>4%</p> <p>80%</p> <p>10%</p> <p>9%</p> </div> </div>
1	C	359	<div style="display: flex; align-items: center;"> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 4%, orange 4%, yellow 11%, green 78%, grey 100%);"></div> <div style="margin-left: 10px;"> <p>4%</p> <p>78%</p> <p>11%</p> <p>11%</p> </div> </div>
1	D	359	<div style="display: flex; align-items: center;"> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 4%, orange 4%, yellow 11%, green 78%, grey 100%);"></div> <div style="margin-left: 10px;"> <p>4%</p> <p>78%</p> <p>11%</p> <p>12%</p> </div> </div>

2 Entry composition i

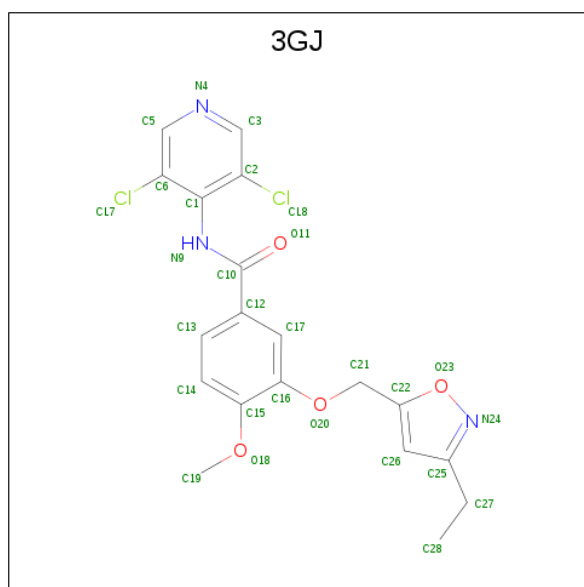
There are 4 unique types of molecules in this entry. The entry contains 11010 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 cAMP-specific 3',5'-cyclic phosphodiesterase 4D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	322	Total 2604	C 1649	N 445	O 496	S 14	0	0	0
1	B	325	Total 2635	C 1669	N 450	O 502	S 14	0	0	0
1	C	321	Total 2608	C 1652	N 445	O 497	S 14	0	0	0
1	D	317	Total 2570	C 1630	N 440	O 486	S 14	0	0	0

- Molecule 2 is N-(3,5-dichloropyridin-4-yl)-3-[(3-ethyl-1,2-oxazol-5-yl)methoxy]-4-methoxybenzamide (three-letter code: 3GJ) (formula: C₁₉H₁₇Cl₂N₃O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Cl	N			O
2	A	1	Total 28	C 19	Cl 2	N 3	O 4	0	0
2	B	1	Total 28	C 19	Cl 2	N 3	O 4	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	C	1	Total	C	Cl	N	O	0	0
			28	19	2	3	4		
2	D	1	Total	C	Cl	N	O	0	0
			28	19	2	3	4		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Zn	0	0
			2	2		
3	A	2	Total	Zn	0	0
			2	2		
3	D	2	Total	Zn	0	0
			2	2		
3	C	2	Total	Zn	0	0
			2	2		

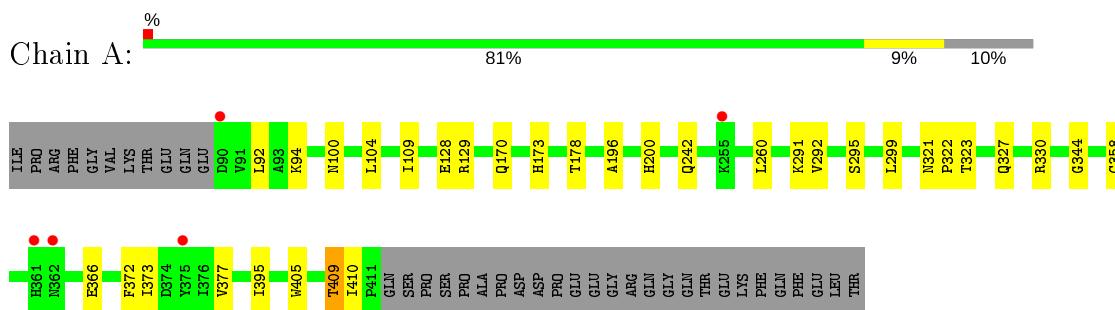
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	144	Total	O	0	0
			144	144		
4	B	121	Total	O	0	0
			121	121		
4	C	103	Total	O	0	0
			103	103		
4	D	105	Total	O	0	0
			105	105		

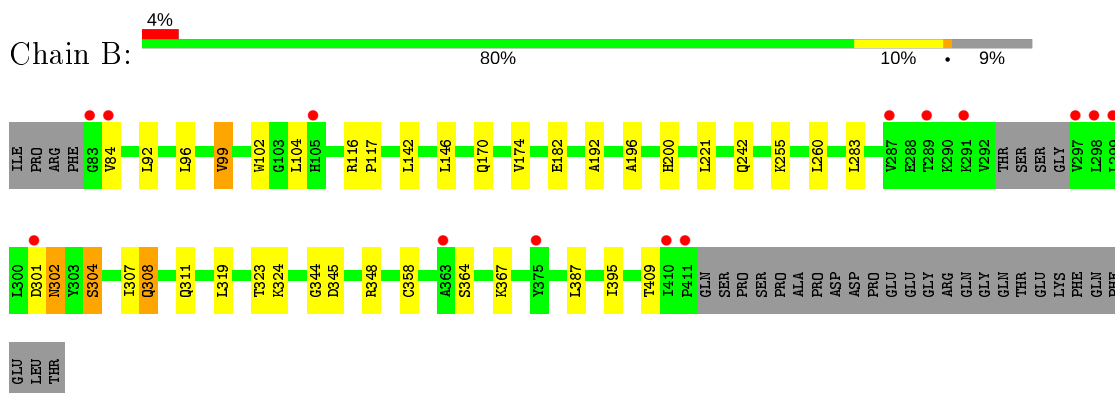
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

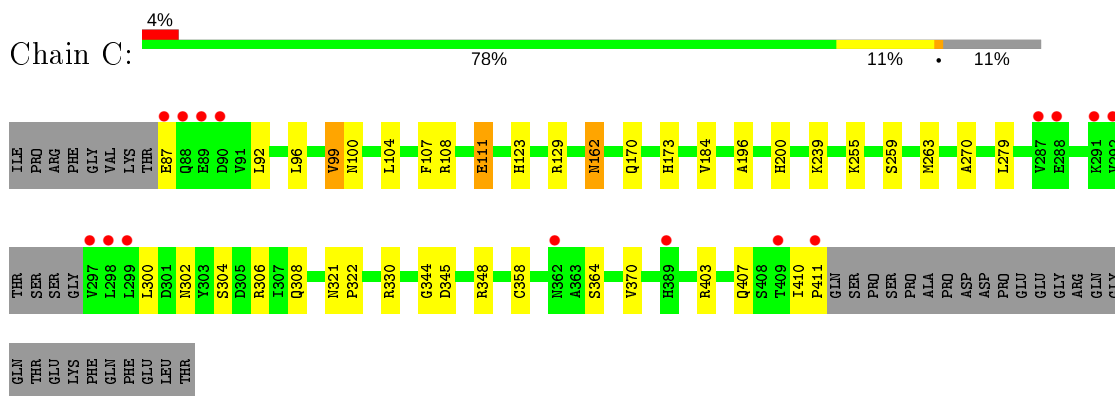
- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



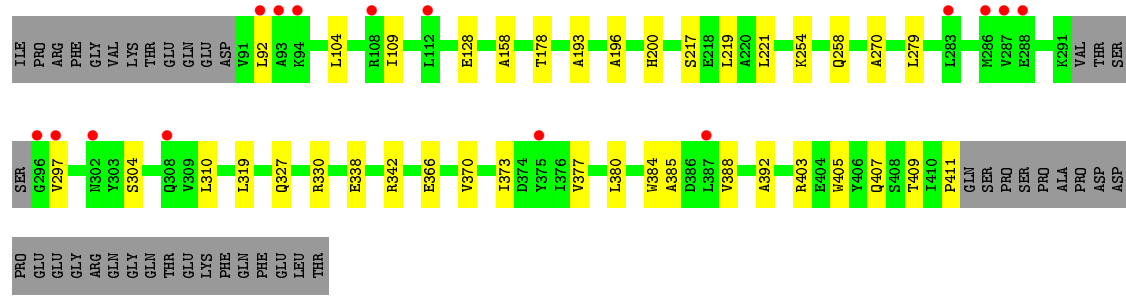
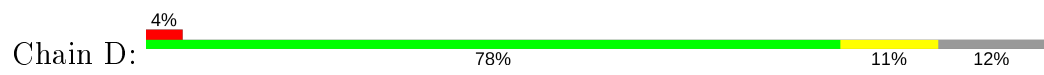
- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.43Å 112.67Å 161.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.45 – 2.20 41.87 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.4 (92.45-2.20) 96.0 (41.87-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.237 , 0.281 0.227 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtrriage
Anisotropy	0.209	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11010	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: ZN, 3GJ

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	0.56	0/2658	0.61	0/3612
1	B	0.53	0/2688	0.63	0/3650
1	C	0.49	0/2661	0.62	0/3614
1	D	0.50	0/2623	0.62	0/3562
All	All	0.52	0/10630	0.62	0/14438

There are no bond length outliers.

There are no bond angle outliers.

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	2604	0	2564	17	1
1	B	2635	0	2595	25	1
1	C	2608	0	2563	24	0
1	D	2570	0	2533	18	0
2	A	28	0	17	1	0
2	B	28	0	17	1	0
2	C	28	0	17	0	0
2	D	28	0	17	0	0
3	A	2	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	144	0	0	0	0
4	B	121	0	0	1	0
4	C	103	0	0	0	0
4	D	105	0	0	0	0
All	All	11010	0	10323	80	1

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 (80) 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:366:GLU:HG2	1:A:409:THR:HG23	1.52	0.90
1:A:100:ASN:HD22	1:A:173:HIS:HE1	1.25	0.84
1:A:129:ARG:HE	1:A:173:HIS:HD2	1.27	0.82
1:B:96:LEU:O	1:B:99:VAL:HG22	1.86	0.76
1:A:366:GLU:HG2	1:A:409:THR:CG2	2.15	0.75
1:B:409:THR:HG23	1:D:411:PRO:HB3	1.67	0.74
1:C:302:ASN:O	1:C:306:ARG:HG3	1.90	0.71
1:A:295:SER:HB2	1:B:255:LYS:HG3	1.73	0.70
1:A:100:ASN:ND2	1:A:173:HIS:HE1	1.90	0.69
1:C:129:ARG:HE	1:C:173:HIS:HD1	1.42	0.68
1:C:162:ASN:H	1:C:162:ASN:HD22	1.40	0.67
1:B:307:ILE:O	1:B:311:GLN:HG3	1.96	0.65
1:C:345:ASP:OD1	1:C:348:ARG:NH2	2.29	0.65
1:B:84:VAL:HG11	1:C:123:HIS:CD2	2.31	0.65
1:D:104:LEU:HD11	1:D:109:ILE:HD11	1.79	0.64
1:D:366:GLU:HG2	1:D:409:THR:HG22	1.82	0.62
1:D:330:ARG:HD3	1:D:405:TRP:CH2	2.35	0.62
1:A:129:ARG:HE	1:A:173:HIS:CD2	2.16	0.62
1:D:254:LYS:HE3	1:D:258:GLN:OE1	1.99	0.61
1:B:364:SER:HB2	1:B:367:LYS:HB2	1.82	0.61
1:B:301:ASP:HB3	1:B:302:ASN:HD22	1.65	0.60
1:C:162:ASN:N	1:C:162:ASN:HD22	1.99	0.60
1:B:283:LEU:HD11	1:B:387:LEU:HD22	1.84	0.60
1:C:370:VAL:HG11	1:C:407:GLN:HE21	1.66	0.59
1:C:403:ARG:HH12	1:C:407:GLN:HG3	1.69	0.58
1:A:321:ASN:HB2	1:A:322:PRO:HD3	1.86	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:GLY:HA3	1:B:358:CYS:O	2.04	0.57
1:B:304:SER:O	1:B:308:GLN:HB2	2.05	0.55
1:C:100:ASN:ND2	1:C:173:HIS:NE2	2.54	0.55
1:A:196:ALA:O	1:A:200:HIS:HB3	2.08	0.54
1:D:403:ARG:NH1	1:D:407:GLN:HG3	2.24	0.53
1:D:338:GLU:O	1:D:342:ARG:HD2	2.09	0.52
1:B:302:ASN:H	1:B:302:ASN:ND2	2.09	0.51
1:D:193:ALA:HB2	1:D:310:LEU:HD22	1.93	0.51
1:C:344:GLY:HA3	1:C:358:CYS:O	2.11	0.51
1:C:96:LEU:O	1:C:99:VAL:HB	2.10	0.51
1:D:270:ALA:HB1	1:D:279:LEU:HD11	1.93	0.51
1:D:254:LYS:O	1:D:258:GLN:HG3	2.10	0.50
1:C:304:SER:O	1:C:308:GLN:HB2	2.12	0.50
1:A:100:ASN:HD22	1:A:173:HIS:CE1	2.17	0.49
1:A:323:THR:HB	1:A:395:ILE:HG23	1.95	0.48
1:D:373:ILE:HA	1:D:377:VAL:HB	1.95	0.48
1:D:158:ALA:H	1:D:342:ARG:HH12	1.62	0.48
1:A:344:GLY:HA3	1:A:358:CYS:O	2.14	0.47
1:B:102:TRP:CE2	1:B:324:LYS:HE2	2.49	0.47
1:C:107:PHE:O	1:C:111:GLU:HG2	2.15	0.47
1:B:192:ALA:HB2	1:B:260:LEU:HD12	1.97	0.47
1:D:385:ALA:HA	1:D:392:ALA:HB3	1.97	0.46
1:B:302:ASN:OD1	1:B:304:SER:HB2	2.16	0.46
1:B:196:ALA:O	1:B:200:HIS:HB3	2.16	0.45
1:B:221:LEU:O	1:B:221:LEU:HD23	2.15	0.45
1:B:323:THR:HB	1:B:395:ILE:HG23	1.98	0.45
1:B:409:THR:HG21	4:B:704:HOH:O	2.16	0.45
1:A:104:LEU:HD22	1:A:170:GLN:HG3	1.98	0.45
1:C:104:LEU:HD22	1:C:170:GLN:HG3	1.99	0.45
1:C:403:ARG:NH1	1:C:407:GLN:HG3	2.32	0.45
1:D:196:ALA:O	1:D:200:HIS:HB3	2.17	0.44
1:B:116:ARG:N	1:B:117:PRO:CD	2.80	0.44
1:C:184:VAL:O	1:C:306:ARG:NH2	2.49	0.44
1:D:217:SER:O	1:D:221:LEU:HG	2.18	0.43
1:B:104:LEU:HD22	1:B:170:GLN:HG3	2.01	0.43
1:B:345:ASP:OD1	1:B:348:ARG:NH2	2.51	0.43
1:D:370:VAL:HG21	1:D:407:GLN:HG2	2.01	0.43
1:A:410:ILE:O	1:C:330:ARG:NH2	2.51	0.43
1:B:170:GLN:O	1:B:174:VAL:HG23	2.20	0.42
1:C:321:ASN:HB2	1:C:322:PRO:HD3	2.02	0.42
1:A:330:ARG:HD3	1:A:405:TRP:CH2	2.55	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:ASN:H	1:C:162:ASN:ND2	2.13	0.42
1:B:319:LEU:CD2	2:B:501:3GJ:CL7	3.05	0.41
1:C:270:ALA:HB1	1:C:279:LEU:HD11	2.02	0.41
1:D:384:TRP:O	1:D:388:VAL:HG22	2.21	0.41
1:A:372:PHE:HB2	2:A:501:3GJ:C26	2.51	0.41
1:C:196:ALA:O	1:C:200:HIS:HB3	2.21	0.41
1:C:300:LEU:HD13	1:C:306:ARG:HA	2.01	0.41
1:D:319:LEU:HD12	1:D:380:LEU:HD22	2.02	0.41
1:A:373:ILE:HA	1:A:377:VAL:HB	2.02	0.40
1:B:302:ASN:N	1:B:302:ASN:ND2	2.69	0.40
1:C:259:SER:O	1:C:263:MET:HG3	2.22	0.40
1:B:142:LEU:O	1:B:146:LEU:HG	2.22	0.40
1:C:410:ILE:HA	1:C:411:PRO:HD3	1.90	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:GLN:NE2	1:B:242:GLN:OE1[4_555]	2.10	0.10

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	320/359 (89%)	313 (98%)	7 (2%)	0	100	100
1	B	321/359 (89%)	317 (99%)	4 (1%)	0	100	100
1	C	317/359 (88%)	309 (98%)	8 (2%)	0	100	100
1	D	313/359 (87%)	306 (98%)	7 (2%)	0	100	100
All	All	1271/1436 (88%)	1245 (98%)	26 (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	294/327 (90%)	283 (96%)	11 (4%)	34	43
1	B	297/327 (91%)	291 (98%)	6 (2%)	55	69
1	C	294/327 (90%)	285 (97%)	9 (3%)	40	51
1	D	289/327 (88%)	282 (98%)	7 (2%)	49	62
All	All	1174/1308 (90%)	1141 (97%)	33 (3%)	43	56

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	LEU
1	A	94	LYS
1	A	109	ILE
1	A	128	GLU
1	A	178	THR
1	A	260	LEU
1	A	291	LYS
1	A	292	VAL
1	A	299	LEU
1	A	327	GLN
1	A	409	THR
1	B	92	LEU
1	B	99	VAL
1	B	182	GLU
1	B	302	ASN
1	B	304	SER
1	B	308	GLN
1	C	87	GLU
1	C	92	LEU
1	C	99	VAL
1	C	108	ARG
1	C	111	GLU
1	C	162	ASN
1	C	239	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	255	LYS
1	C	364	SER
1	D	92	LEU
1	D	128	GLU
1	D	178	THR
1	D	219	LEU
1	D	297	VAL
1	D	304	SER
1	D	327	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	ASN
1	A	173	HIS
1	A	242	GLN
1	A	278	ASN
1	B	123	HIS
1	B	127	GLN
1	B	302	ASN
1	B	407	GLN
1	C	100	ASN
1	C	123	HIS
1	C	162	ASN
1	C	258	GLN
1	C	407	GLN
1	D	123	HIS
1	D	245	ASN
1	D	250	GLN
1	D	308	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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
2	3GJ	D	501	-	28,30,30	2.02	6 (21%)	35,41,41	1.96	11 (31%)
2	3GJ	C	501	-	28,30,30	2.08	6 (21%)	35,41,41	2.05	12 (34%)
2	3GJ	B	501	-	28,30,30	2.10	5 (17%)	35,41,41	2.06	8 (22%)
2	3GJ	A	501	-	28,30,30	1.97	5 (17%)	35,41,41	2.02	9 (25%)

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	3GJ	D	501	-	-	3/14/17/17	0/3/3/3
2	3GJ	C	501	-	-	5/14/17/17	0/3/3/3
2	3GJ	B	501	-	-	5/14/17/17	0/3/3/3
2	3GJ	A	501	-	-	3/14/17/17	0/3/3/3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	3GJ	C1-C2	6.02	1.49	1.40
2	D	501	3GJ	C1-C6	5.87	1.49	1.40
2	B	501	3GJ	C1-C6	5.78	1.48	1.40
2	C	501	3GJ	C1-C2	5.67	1.48	1.40
2	A	501	3GJ	C1-C2	5.62	1.48	1.40
2	D	501	3GJ	C1-C2	5.54	1.48	1.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	3GJ	C1-C6	5.50	1.48	1.40
2	A	501	3GJ	C1-C6	5.24	1.48	1.40
2	B	501	3GJ	C15-C16	4.87	1.51	1.40
2	C	501	3GJ	C15-C16	4.66	1.50	1.40
2	A	501	3GJ	C15-C16	4.50	1.50	1.40
2	D	501	3GJ	C15-C16	4.13	1.49	1.40
2	C	501	3GJ	C1-N9	-3.57	1.36	1.43
2	D	501	3GJ	C1-N9	-3.25	1.37	1.43
2	A	501	3GJ	C1-N9	-3.16	1.37	1.43
2	B	501	3GJ	C1-N9	-3.13	1.37	1.43
2	D	501	3GJ	C2-CL8	2.55	1.79	1.73
2	B	501	3GJ	C2-CL8	2.51	1.79	1.73
2	C	501	3GJ	C6-CL7	2.36	1.79	1.73
2	C	501	3GJ	C2-CL8	2.33	1.79	1.73
2	A	501	3GJ	C2-CL8	2.23	1.78	1.73
2	D	501	3GJ	C6-CL7	2.01	1.78	1.73

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	3GJ	C22-C26-C25	5.66	110.74	104.63
2	C	501	3GJ	C22-C26-C25	5.56	110.63	104.63
2	A	501	3GJ	C22-C26-C25	5.06	110.09	104.63
2	D	501	3GJ	C22-C26-C25	4.97	110.00	104.63
2	A	501	3GJ	C19-O18-C15	4.72	124.66	117.53
2	B	501	3GJ	C19-O18-C15	4.25	123.94	117.53
2	D	501	3GJ	C19-O18-C15	4.06	123.66	117.53
2	C	501	3GJ	O18-C15-C16	3.96	120.92	115.41
2	B	501	3GJ	C27-C25-C26	-3.90	124.84	128.68
2	B	501	3GJ	O18-C15-C16	3.88	120.82	115.41
2	D	501	3GJ	C3-N4-C5	3.81	122.68	117.48
2	A	501	3GJ	C27-C25-C26	-3.76	124.98	128.68
2	B	501	3GJ	C3-N4-C5	3.75	122.60	117.48
2	C	501	3GJ	C3-N4-C5	3.73	122.57	117.48
2	C	501	3GJ	C27-C25-C26	-3.71	125.03	128.68
2	D	501	3GJ	C27-C25-C26	-3.47	125.26	128.68
2	A	501	3GJ	C3-N4-C5	3.44	122.18	117.48
2	A	501	3GJ	C21-O20-C16	3.19	124.03	117.76
2	A	501	3GJ	O18-C15-C16	3.10	119.72	115.41
2	C	501	3GJ	O20-C16-C15	2.88	121.68	115.73
2	D	501	3GJ	C21-O20-C16	2.86	123.40	117.76
2	B	501	3GJ	O20-C16-C15	2.85	121.63	115.73

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	3GJ	C3-C2-CL8	2.85	122.47	118.60
2	D	501	3GJ	O18-C15-C16	2.79	119.30	115.41
2	C	501	3GJ	C21-O20-C16	2.69	123.05	117.76
2	C	501	3GJ	O18-C15-C14	-2.55	120.00	124.37
2	C	501	3GJ	C19-O18-C15	2.50	121.30	117.53
2	C	501	3GJ	C3-C2-CL8	2.46	121.94	118.60
2	A	501	3GJ	O20-C16-C15	2.43	120.75	115.73
2	D	501	3GJ	C12-C10-N9	2.39	121.19	115.92
2	B	501	3GJ	C26-C25-N24	-2.34	108.06	111.51
2	D	501	3GJ	C3-C2-C1	-2.33	118.88	120.16
2	A	501	3GJ	C3-C2-CL8	2.27	121.68	118.60
2	D	501	3GJ	C6-C5-N4	-2.18	120.95	122.85
2	C	501	3GJ	C26-C25-N24	-2.17	108.31	111.51
2	C	501	3GJ	C12-C10-N9	2.16	120.68	115.92
2	A	501	3GJ	C12-C10-N9	2.14	120.63	115.92
2	B	501	3GJ	C3-C2-C1	-2.14	118.98	120.16
2	D	501	3GJ	C26-C25-N24	-2.10	108.41	111.51
2	C	501	3GJ	C5-C6-CL7	2.05	121.38	118.60

There are no chirality outliers.

All (16) torsion outliers are listed below:

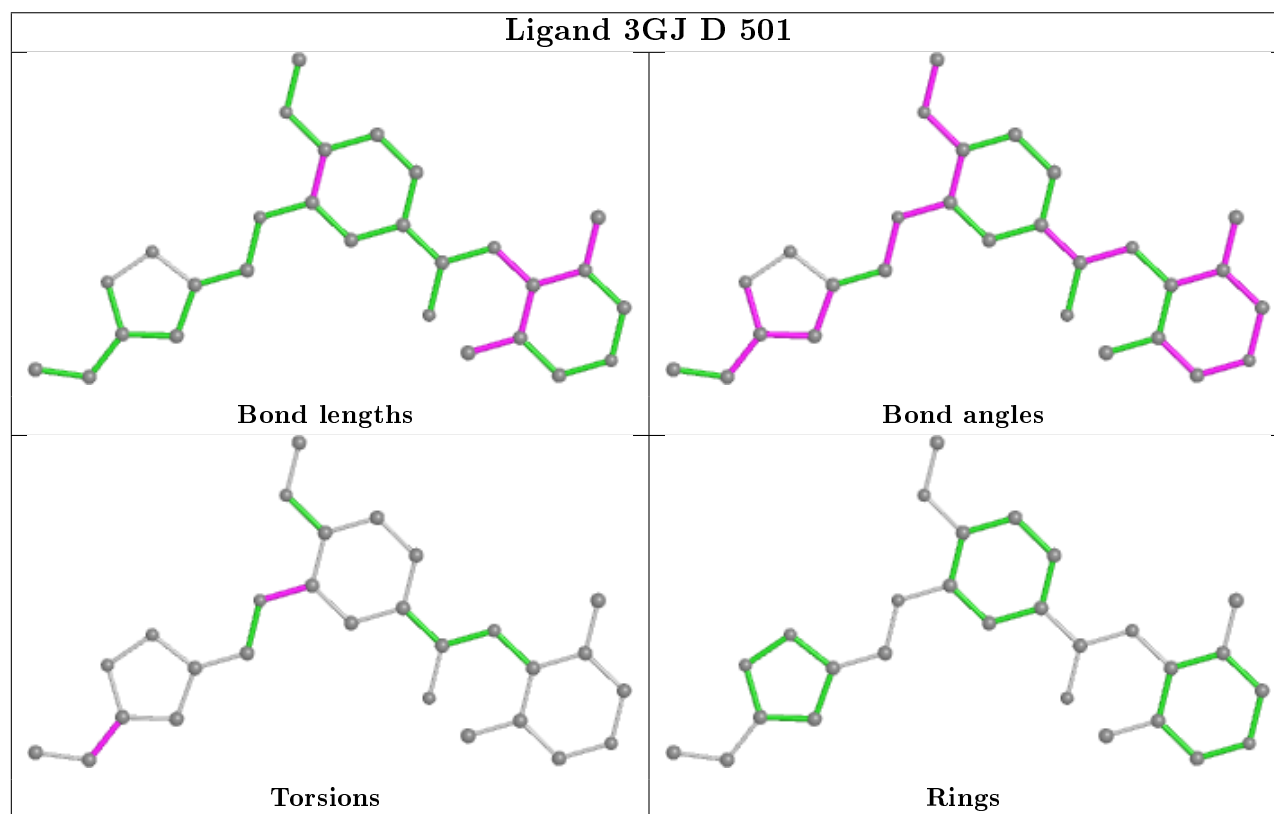
Mol	Chain	Res	Type	Atoms
2	B	501	3GJ	C26-C25-C27-C28
2	D	501	3GJ	C26-C25-C27-C28
2	C	501	3GJ	C26-C25-C27-C28
2	A	501	3GJ	C26-C25-C27-C28
2	C	501	3GJ	C16-C15-O18-C19
2	C	501	3GJ	C14-C15-O18-C19
2	B	501	3GJ	C16-C15-O18-C19
2	D	501	3GJ	C15-C16-O20-C21
2	B	501	3GJ	C14-C15-O18-C19
2	B	501	3GJ	C15-C16-O20-C21
2	C	501	3GJ	C15-C16-O20-C21
2	A	501	3GJ	C15-C16-O20-C21
2	D	501	3GJ	C17-C16-O20-C21
2	B	501	3GJ	C17-C16-O20-C21
2	C	501	3GJ	C17-C16-O20-C21
2	A	501	3GJ	C17-C16-O20-C21

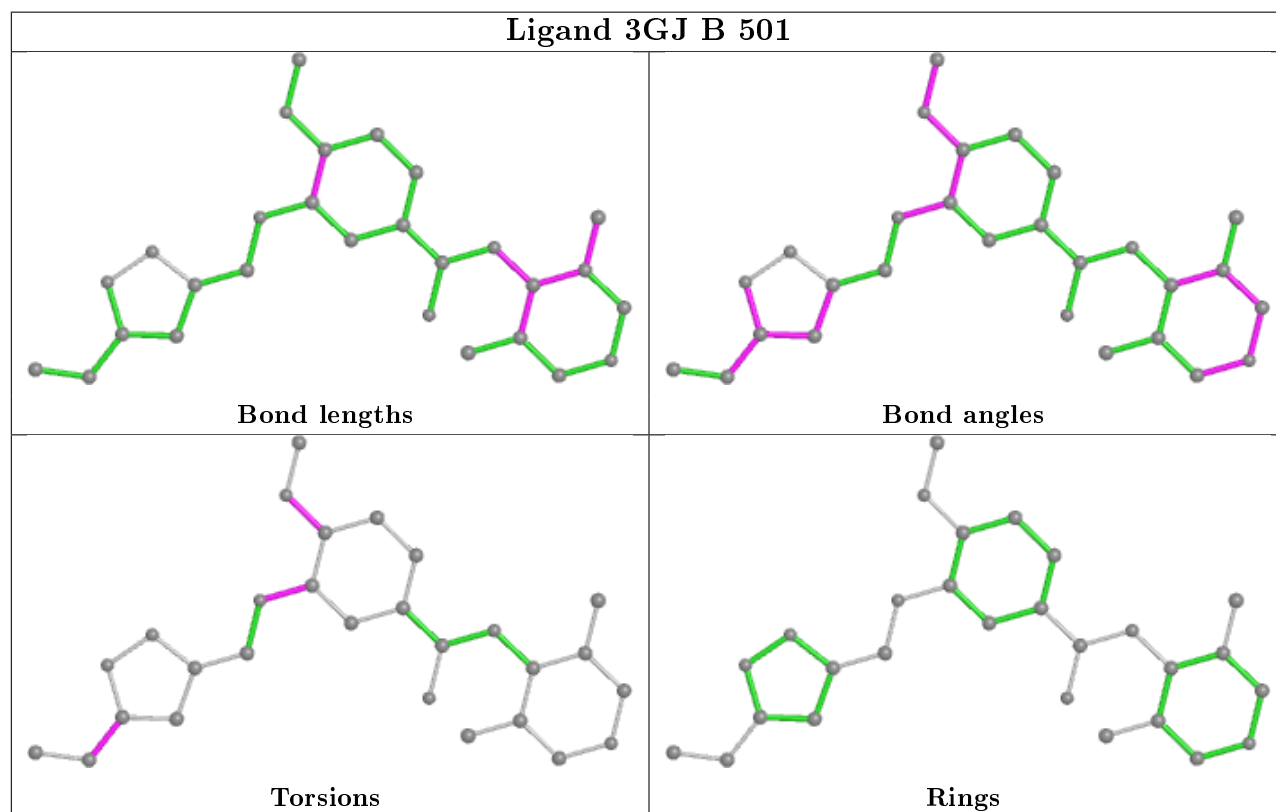
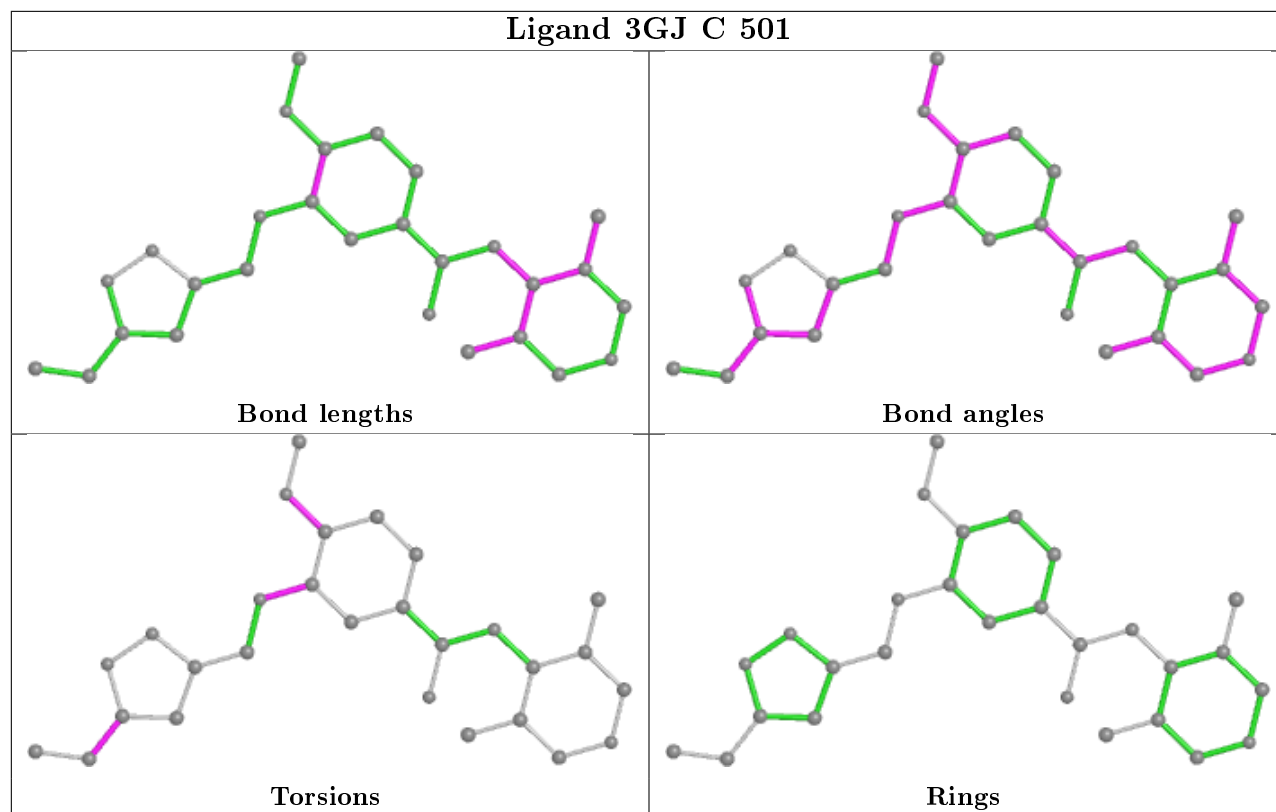
There are no ring outliers.

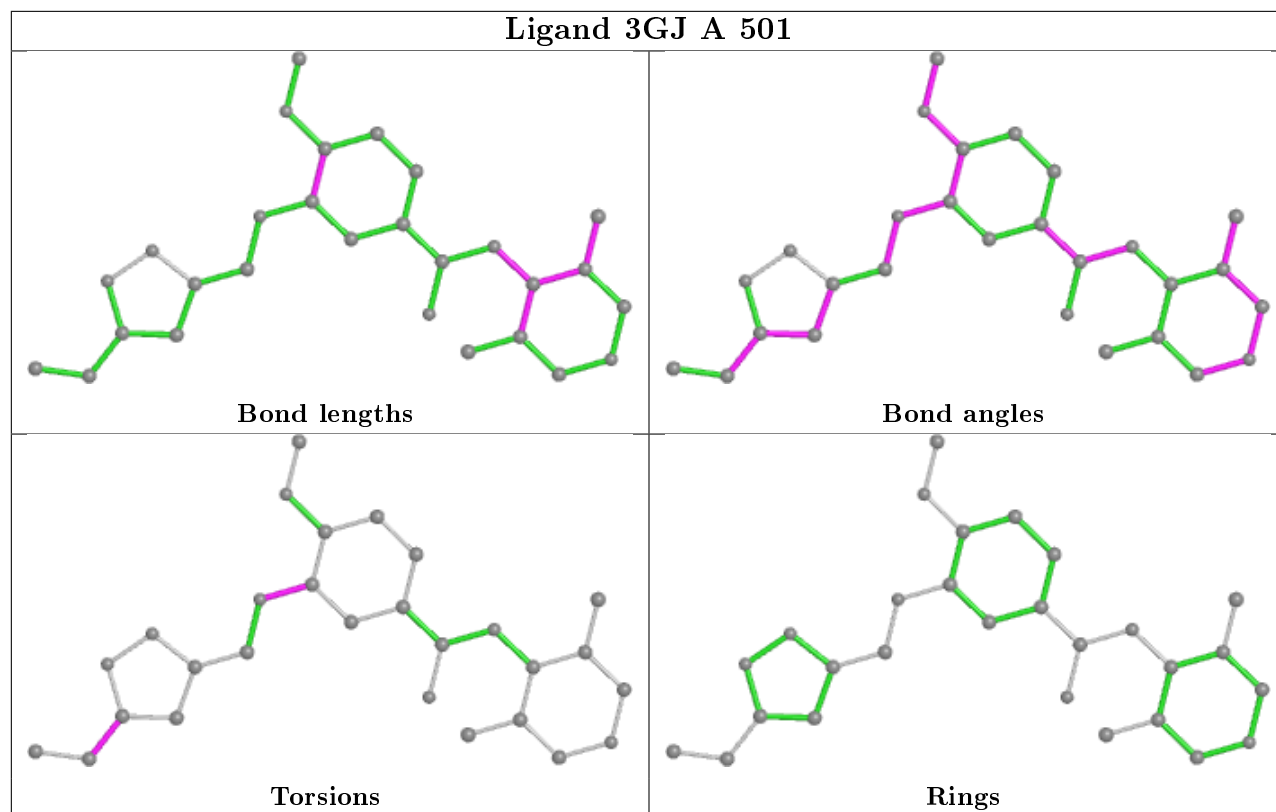
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	3GJ	1	0
2	A	501	3GJ	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	322/359 (89%)	-0.03	5 (1%) 72 70	24, 35, 50, 62	0
1	B	325/359 (90%)	0.03	14 (4%) 35 33	24, 36, 57, 70	0
1	C	321/359 (89%)	0.25	15 (4%) 31 30	25, 43, 69, 84	0
1	D	317/359 (88%)	0.28	15 (4%) 31 30	28, 45, 64, 71	0
All	All	1285/1436 (89%)	0.13	49 (3%) 40 38	24, 40, 62, 84	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	411	PRO	7.1
1	C	292	VAL	5.9
1	B	83	GLY	4.7
1	C	87	GLU	4.6
1	C	287	VAL	4.4
1	D	296	GLY	4.4
1	C	291	LYS	4.1
1	C	297	VAL	3.8
1	C	409	THR	3.7
1	B	84	VAL	3.6
1	D	287	VAL	3.6
1	D	387	LEU	3.4
1	C	89	GLU	3.3
1	C	298	LEU	3.3
1	B	297	VAL	3.3
1	B	287	VAL	3.2
1	D	94	LYS	3.2
1	D	288	GLU	3.1
1	B	289	THR	3.1
1	C	299	LEU	3.1
1	A	90	ASP	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	362	ASN	2.8
1	B	410	ILE	2.8
1	A	375	TYR	2.7
1	D	108	ARG	2.7
1	B	298	LEU	2.6
1	D	112	LEU	2.6
1	B	299	LEU	2.5
1	D	92	LEU	2.5
1	C	90	ASP	2.5
1	B	105	HIS	2.5
1	D	286	MET	2.5
1	B	363	ALA	2.4
1	D	297	VAL	2.4
1	B	375	TYR	2.4
1	C	389	HIS	2.4
1	B	301	ASP	2.4
1	A	255	LYS	2.4
1	D	93	ALA	2.3
1	D	375	TYR	2.3
1	C	288	GLU	2.3
1	C	411	PRO	2.3
1	B	291	LYS	2.2
1	D	283	LEU	2.2
1	C	88	GLN	2.2
1	A	362	ASN	2.2
1	D	302	ASN	2.1
1	D	308	GLN	2.0
1	A	361	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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,

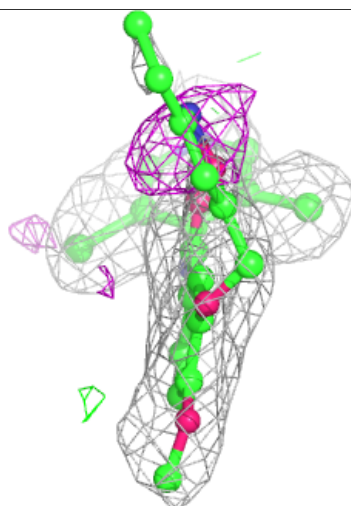
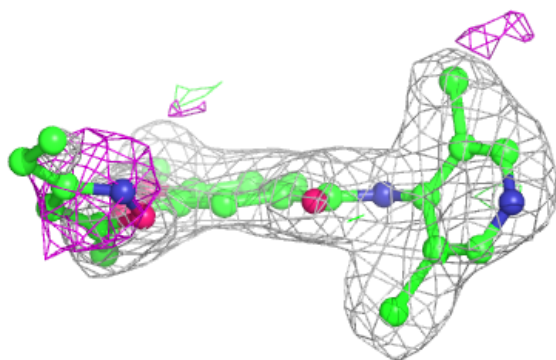
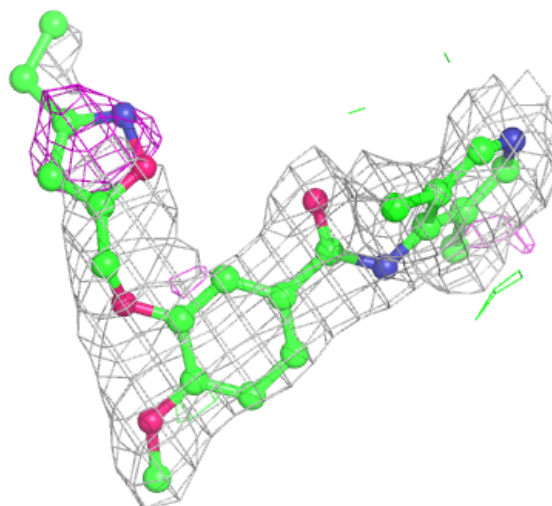
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	3GJ	D	501	28/28	0.66	0.24	59,63,67,68	0
2	3GJ	A	501	28/28	0.76	0.25	50,54,58,59	0
2	3GJ	B	501	28/28	0.78	0.23	54,57,66,67	0
2	3GJ	C	501	28/28	0.81	0.20	58,61,63,64	0
3	ZN	C	503	1/1	0.92	0.05	78,78,78,78	0
3	ZN	B	503	1/1	0.93	0.05	82,82,82,82	0
3	ZN	A	503	1/1	0.94	0.04	75,75,75,75	0
3	ZN	C	502	1/1	0.94	0.13	49,49,49,49	0
3	ZN	D	503	1/1	0.95	0.04	87,87,87,87	0
3	ZN	B	502	1/1	0.95	0.08	46,46,46,46	0
3	ZN	D	502	1/1	0.96	0.10	50,50,50,50	0
3	ZN	A	502	1/1	0.97	0.09	44,44,44,44	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.

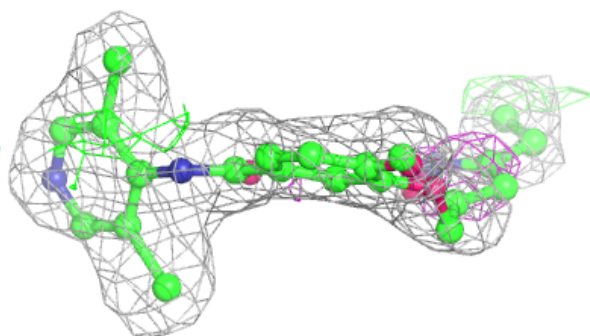
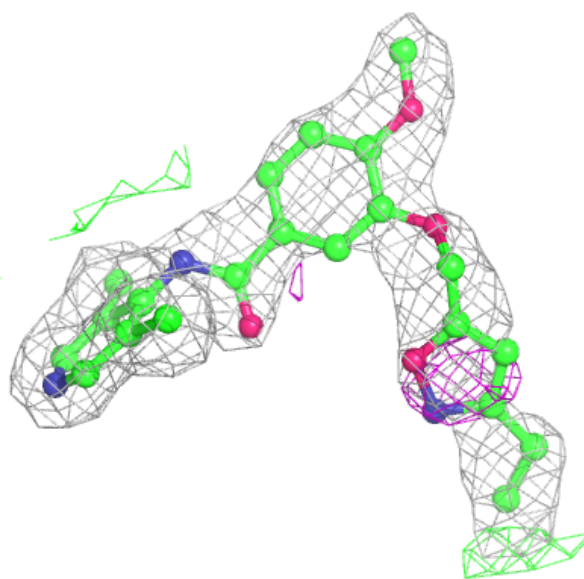
Electron density around 3GJ D 501:

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



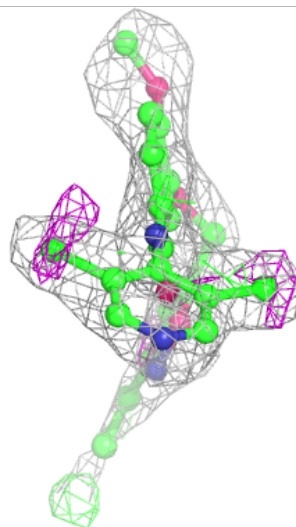
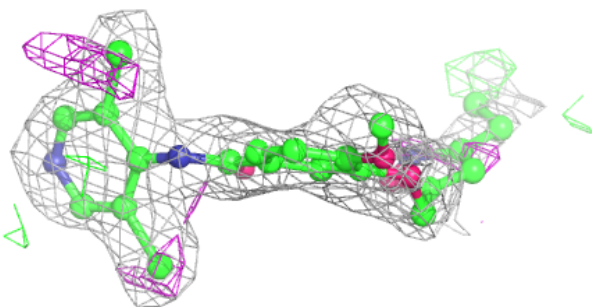
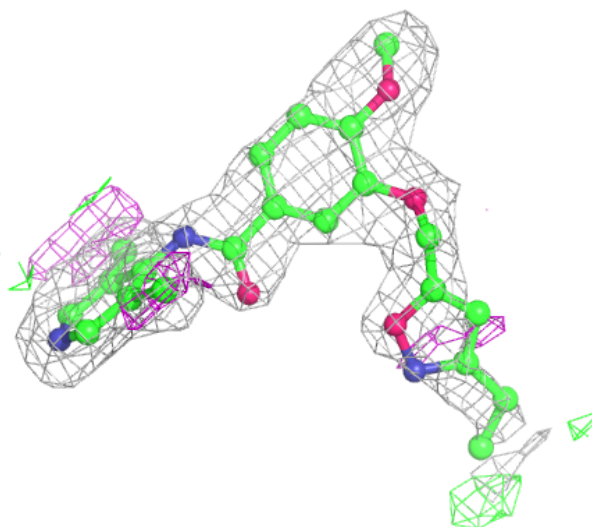
Electron density around 3GJ A 501:

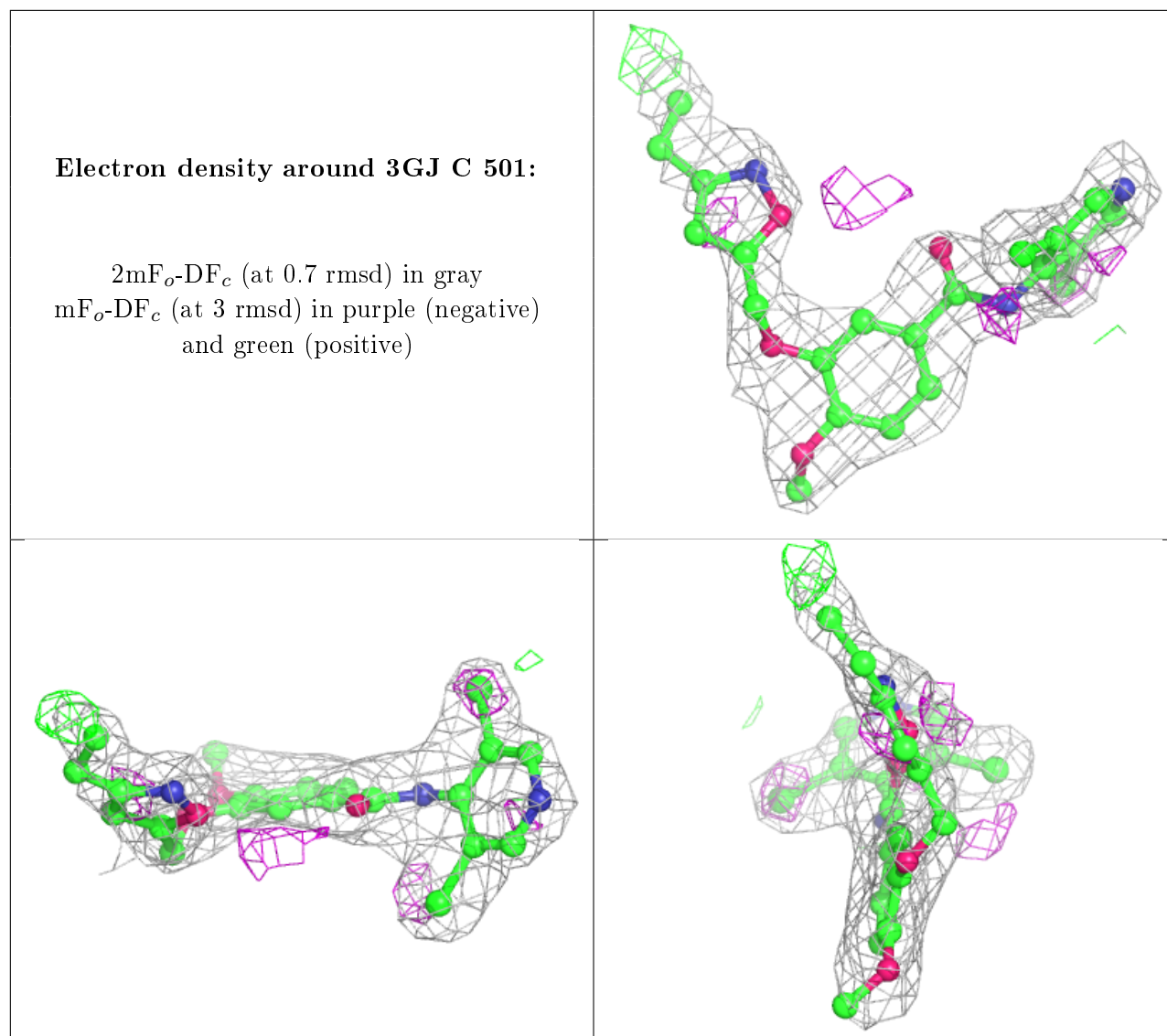
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 3GJ B 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.