



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 11:38 am BST

PDB ID : 1Q6T
Title : THE STRUCTURE OF PHOSPHOTYROSINE PHOSPHATASE 1B IN COMPLEX WITH COMPOUND 11
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Deposited on : 2003-08-13
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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

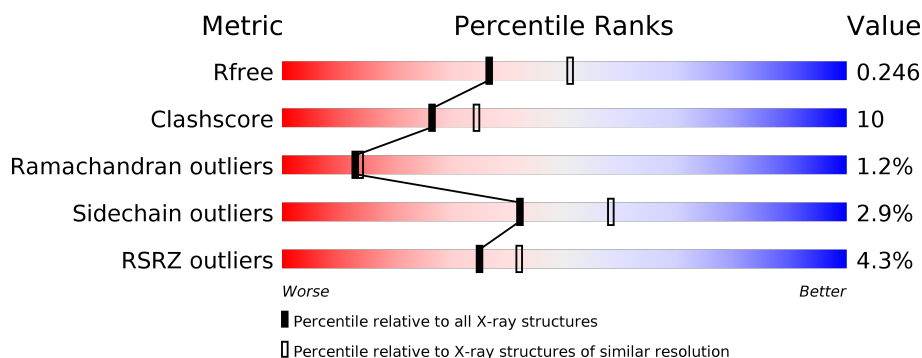
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	310	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>18%</div> <div>• 7%</div> </div> </div>
1	B	310	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>17%</div> <div>• 7%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5018 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 Protein-tyrosine phosphatase, non-receptor type 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2349	1490	404	439	16			
1	B	289	Total	C	N	O	S	0	0	0
			2346	1488	404	439	15			

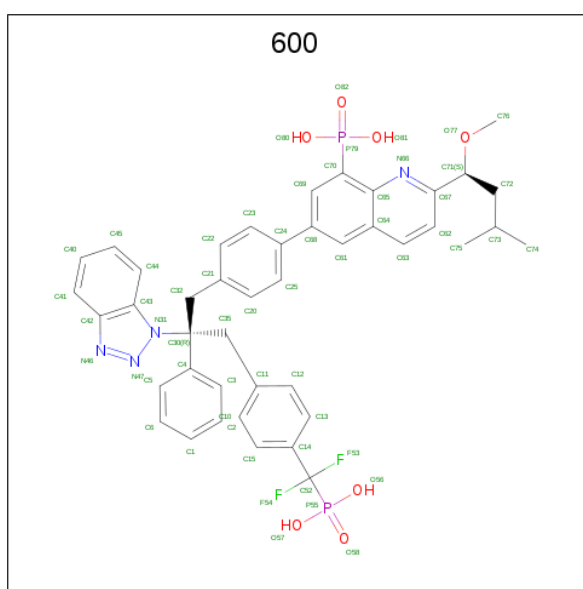
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	489	MET	-	CLONING ARTIFACT	UNP P18031
A	490	ASP	-	CLONING ARTIFACT	UNP P18031
A	491	TYR	-	CLONING ARTIFACT	UNP P18031
A	492	LYS	-	CLONING ARTIFACT	UNP P18031
A	493	ASP	-	CLONING ARTIFACT	UNP P18031
A	494	ASP	-	CLONING ARTIFACT	UNP P18031
A	495	ASP	-	CLONING ARTIFACT	UNP P18031
A	496	ASP	-	CLONING ARTIFACT	UNP P18031
A	497	LYS	-	CLONING ARTIFACT	UNP P18031
A	498	LEU	-	CLONING ARTIFACT	UNP P18031
A	499	GLU	-	CLONING ARTIFACT	UNP P18031
A	500	PHE	-	CLONING ARTIFACT	UNP P18031
B	989	MET	-	CLONING ARTIFACT	UNP P18031
B	990	ASP	-	CLONING ARTIFACT	UNP P18031
B	991	TYR	-	CLONING ARTIFACT	UNP P18031
B	992	LYS	-	CLONING ARTIFACT	UNP P18031
B	993	ASP	-	CLONING ARTIFACT	UNP P18031
B	994	ASP	-	CLONING ARTIFACT	UNP P18031
B	995	ASP	-	CLONING ARTIFACT	UNP P18031
B	996	ASP	-	CLONING ARTIFACT	UNP P18031
B	997	LYS	-	CLONING ARTIFACT	UNP P18031
B	998	LEU	-	CLONING ARTIFACT	UNP P18031
B	999	GLU	-	CLONING ARTIFACT	UNP P18031
B	1000	PHE	-	CLONING ARTIFACT	UNP P18031

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is 6-[4-((2S)-2-(1H-1,2,3-BENZOTRIAZOL-1-YL)-3-{4-[DIFLUORO(PHOSPHONO)METHYL]PHENYL}-2-PHENYLPROPYL)PHENYL]-2-[(1S)-1-METHOXY-3-METHYLBUTYL]QUINOLIN-8-YLPHOSPHONIC ACID (three-letter code: 600) (formula: C₄₃H₄₂F₂N₄O₇P₂).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	P	0	0
			58	43	2	4	7	2		
3	B	1	Total	C	F	N	O	P	0	0
			58	43	2	4	7	2		

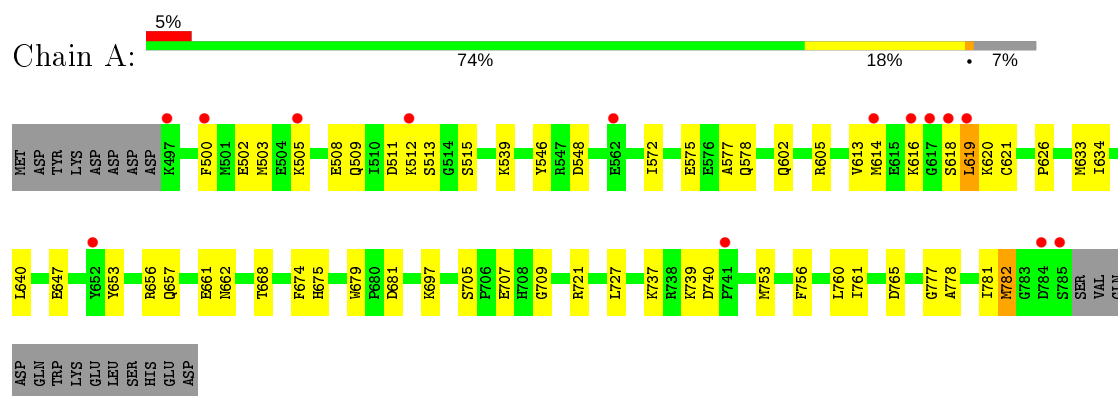
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	97	Total	O	0	0
			97	97		
4	B	108	Total	O	0	0
			108	108		

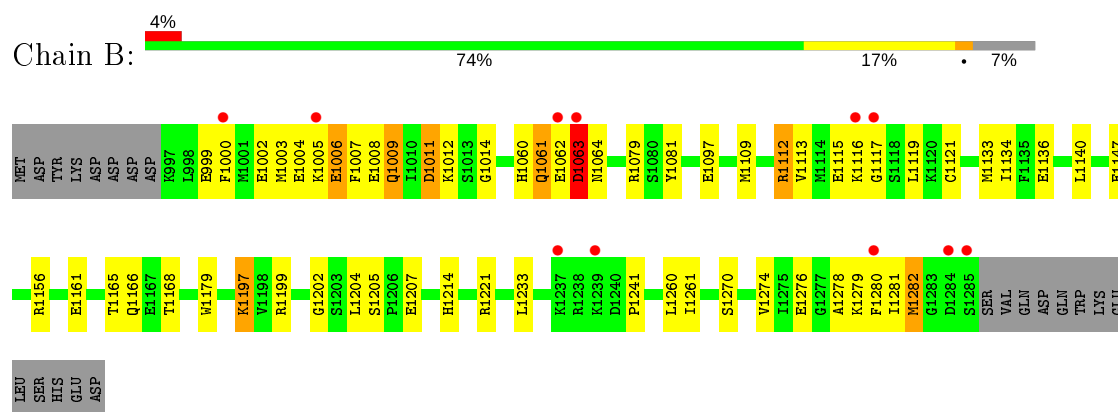
3 Residue-property plots [i](#)

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: Protein-tyrosine phosphatase, non-receptor type 1



- Molecule 1: Protein-tyrosine phosphatase, non-receptor type 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.03Å 87.83Å 138.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.30 28.38 – 1.96	Depositor EDS
% Data completeness (in resolution range)	97.4 (12.00-2.30) 91.2 (28.38-1.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 1.96Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.214 , 0.246 0.215 , 0.246	Depositor DCC
R_{free} test set	3519 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.063 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5018	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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: MG, 600

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	0.49	0/2402	0.63	0/3237
1	B	0.53	1/2399 (0.0%)	0.70	1/3234 (0.0%)
All	All	0.51	1/4801 (0.0%)	0.67	1/6471 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1119	LEU	C-N	-5.74	1.20	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1063	ASP	CA-CB-CG	-10.34	90.64	113.40

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	2349	0	2306	51	0
1	B	2346	0	2298	53	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	58	0	37	2	0
3	B	58	0	37	0	0
4	A	97	0	0	2	0
4	B	108	0	0	2	0
All	All	5018	0	4678	98	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 (98) 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:1161:GLU:HG3	1:B:1168:THR:HG22	1.56	0.88
1:A:500:PHE:HA	1:A:503:MET:HE3	1.61	0.81
1:B:1002:GLU:HG3	1:B:1005:LYS:HE2	1.64	0.78
1:B:1112:ARG:NH1	1:B:1112:ARG:HB2	2.00	0.76
1:B:1011:ASP:O	1:B:1012:LYS:HG3	1.85	0.76
1:A:619:LEU:H	1:A:619:LEU:HD12	1.52	0.74
1:A:727:LEU:HA	1:A:753:MET:HE1	1.73	0.70
1:A:674:PHE:HE2	1:A:697:LYS:HD2	1.57	0.69
1:B:1000:PHE:HE2	1:B:1278:ALA:HB1	1.56	0.69
1:A:513:SER:HB2	1:B:1117:GLY:H	1.58	0.67
1:A:508:GLU:O	1:A:512:LYS:HG3	1.95	0.67
1:A:508:GLU:HB3	1:A:512:LYS:NZ	2.09	0.67
1:A:513:SER:HB2	1:B:1117:GLY:N	2.10	0.67
1:A:626:PRO:HG3	1:A:633:MET:HG3	1.77	0.66
1:B:1064:ASN:C	1:B:1064:ASN:OD1	2.34	0.64
1:A:605:ARG:HG2	1:A:605:ARG:HH11	1.64	0.63
1:B:1270:SER:O	1:B:1274:VAL:HG23	1.99	0.62
1:A:778:ALA:HA	1:A:781:ILE:HG22	1.82	0.62
1:A:502:GLU:HG3	1:A:505:LYS:NZ	2.16	0.61
1:A:613:VAL:HG13	1:A:621:CYS:O	2.00	0.60
1:B:1003:MET:O	1:B:1006:GLU:HG3	2.01	0.60
1:B:1002:GLU:O	1:B:1005:LYS:HG2	2.00	0.60
1:B:1113:VAL:HA	1:B:1121:CYS:HB3	1.84	0.60
1:B:1079:ARG:CZ	1:B:1233:LEU:HD11	2.33	0.59
1:A:502:GLU:HA	1:A:505:LYS:HG2	1.85	0.59
1:B:1112:ARG:HH11	1:B:1112:ARG:HB2	1.66	0.58
1:A:508:GLU:HB3	1:A:512:LYS:HZ3	1.70	0.56
1:A:619:LEU:CD1	1:A:619:LEU:H	2.16	0.56
1:B:1161:GLU:CG	1:B:1168:THR:HG22	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1205:SER:OG	1:B:1207:GLU:HG2	2.06	0.55
1:B:1063:ASP:C	1:B:1063:ASP:OD1	2.44	0.55
1:A:705:SER:OG	1:A:707:GLU:HG2	2.07	0.55
1:B:1062:GLU:OE1	1:B:1062:GLU:HA	2.06	0.54
1:B:1281:ILE:HG13	1:B:1282:MET:SD	2.47	0.54
1:B:1179:TRP:CE2	1:B:1221:ARG:HG2	2.42	0.54
1:A:619:LEU:N	1:A:619:LEU:HD12	2.20	0.54
1:A:661:GLU:HG3	1:A:668:THR:HG22	1.87	0.54
1:B:1136:GLU:HG2	4:B:190:HOH:O	2.09	0.53
1:A:778:ALA:HA	1:A:781:ILE:CG2	2.39	0.53
1:B:1112:ARG:HG2	1:B:1115:GLU:OE2	2.09	0.52
1:A:697:LYS:HD3	1:A:697:LYS:O	2.11	0.51
1:A:739:LYS:HG3	1:A:740:ASP:H	1.75	0.51
1:A:653:TYR:HB2	1:A:675:HIS:O	2.12	0.50
1:A:679:TRP:CE2	1:A:721:ARG:HG2	2.46	0.50
1:A:640:LEU:HD23	1:A:662:ASN:HA	1.93	0.50
1:A:616:LYS:NZ	1:B:1014:GLY:H	2.10	0.49
1:B:1165:THR:O	1:B:1166:GLN:HB2	2.13	0.49
1:B:1281:ILE:HG13	1:B:1282:MET:HE3	1.94	0.49
1:B:1005:LYS:HG3	1:B:1006:GLU:N	2.27	0.49
1:A:513:SER:HB2	1:B:1117:GLY:CA	2.43	0.48
1:B:1109:MET:HG3	1:B:1214:HIS:CE1	2.49	0.48
1:A:548:ASP:O	3:A:801:600:H61	2.15	0.47
1:B:1197:LYS:HE2	1:B:1197:LYS:CA	2.45	0.47
1:A:502:GLU:HG3	1:A:505:LYS:HZ2	1.80	0.46
1:B:1005:LYS:CG	1:B:1006:GLU:N	2.78	0.46
1:A:502:GLU:HG3	1:A:505:LYS:HZ3	1.81	0.46
1:A:572:ILE:HG13	1:A:756:PHE:HB2	1.97	0.46
1:B:1060:HIS:O	1:B:1061:GLN:O	2.34	0.46
1:A:614:MET:HG2	1:A:619:LEU:HG	1.98	0.45
1:B:1133:MET:C	1:B:1134:ILE:HD12	2.37	0.45
1:B:1005:LYS:HA	1:B:1008:GLU:OE2	2.17	0.45
1:B:1063:ASP:OD1	1:B:1063:ASP:O	2.33	0.45
1:A:539:LYS:HE3	4:A:36:HOH:O	2.17	0.45
1:A:515:SER:HA	4:A:29:HOH:O	2.17	0.45
1:B:1202:GLY:O	1:B:1205:SER:HB3	2.17	0.44
1:A:781:ILE:HG23	1:A:782:MET:SD	2.57	0.44
1:B:1004:GLU:O	1:B:1008:GLU:HG3	2.18	0.44
1:B:1112:ARG:CZ	1:B:1112:ARG:HB2	2.48	0.44
1:B:1276:GLU:OE1	1:B:1279:LYS:HE2	2.18	0.44
1:A:739:LYS:HG3	1:A:740:ASP:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:697:LYS:C	1:A:697:LYS:HD3	2.39	0.43
1:A:546:TYR:HB3	3:A:801:600:N47	2.34	0.43
1:A:513:SER:HB2	1:B:1117:GLY:HA3	2.01	0.43
1:B:999:GLU:OE2	1:B:1241:PRO:HB2	2.18	0.43
1:A:727:LEU:CA	1:A:753:MET:HE1	2.46	0.42
1:A:605:ARG:HG2	1:A:605:ARG:NH1	2.31	0.42
1:B:1005:LYS:HA	1:B:1008:GLU:CD	2.40	0.42
1:B:1147:GLU:CB	1:B:1156:ARG:HG2	2.49	0.42
1:B:1278:ALA:O	1:B:1282:MET:HB2	2.19	0.42
1:B:1147:GLU:HG3	1:B:1147:GLU:O	2.19	0.42
1:B:1079:ARG:HG2	1:B:1081:TYR:CZ	2.55	0.42
1:A:647:GLU:HB2	1:A:656:ARG:HG2	2.02	0.42
1:A:777:GLY:O	1:A:781:ILE:HG22	2.20	0.42
1:A:577:ALA:O	1:A:578:GLN:HB2	2.20	0.42
1:A:634:ILE:N	1:A:634:ILE:HD12	2.35	0.42
1:A:602:GLN:O	1:A:709:GLY:HA3	2.19	0.42
1:A:778:ALA:CA	1:A:781:ILE:HG22	2.49	0.41
1:B:1004:GLU:O	1:B:1007:PHE:HB3	2.21	0.41
1:A:618:SER:O	1:A:620:LYS:HG3	2.21	0.41
1:B:1112:ARG:HG2	1:B:1115:GLU:CD	2.41	0.41
1:B:1009:GLN:HB3	1:B:1009:GLN:HE21	1.69	0.41
1:B:1097:GLU:HA	1:B:1140:LEU:HD11	2.03	0.41
1:B:1199:ARG:HG2	1:B:1204:LEU:HD12	2.03	0.40
1:A:674:PHE:CE2	1:A:697:LYS:HD2	2.47	0.40
1:B:1064:ASN:HA	4:B:91:HOH:O	2.21	0.40
1:B:1079:ARG:NE	1:B:1233:LEU:HD11	2.36	0.40
1:A:616:LYS:HE3	1:A:681:ASP:OD1	2.22	0.40
1:A:616:LYS:HZ2	1:B:1014:GLY:H	1.70	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	287/310 (93%)	270 (94%)	14 (5%)	3 (1%)	15	17
1	B	287/310 (93%)	270 (94%)	13 (4%)	4 (1%)	11	11
All	All	574/620 (93%)	540 (94%)	27 (5%)	7 (1%)	13	14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1061	GLN
1	B	1116	LYS
1	A	619	LEU
1	B	1011	ASP
1	A	575	GLU
1	A	761	ILE
1	B	1261	ILE

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	259/283 (92%)	252 (97%)	7 (3%)	44	61
1	B	258/283 (91%)	250 (97%)	8 (3%)	40	55
All	All	517/566 (91%)	502 (97%)	15 (3%)	42	58

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	509	GLN
1	A	511	ASP
1	A	657	GLN
1	A	737	LYS
1	A	760	LEU
1	A	765	ASP
1	A	782	MET
1	B	1006	GLU
1	B	1009	GLN

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Mol	Chain	Res	Type
1	B	1063	ASP
1	B	1112	ARG
1	B	1197	LYS
1	B	1260	LEU
1	B	1280	PHE
1	B	1282	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	1009	GLN
1	B	1139	ASN

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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	600	B	1301	-	57,64,64	2.95	31 (54%)	70,97,97	3.83	18 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	600	A	801	-	57,64,64	2.95	35 (61%)	70,97,97	3.85	14 (20%)

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
3	600	B	1301	-	-	6/46/57/57	0/7/7/7
3	600	A	801	-	-	7/46/57/57	0/7/7/7

All (66) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1301	600	C30-C4	6.48	1.60	1.52
3	B	1301	600	C35-C30	6.19	1.62	1.55
3	A	801	600	C30-C4	6.17	1.60	1.52
3	B	1301	600	C32-C21	6.02	1.61	1.51
3	B	1301	600	C5-C4	5.81	1.48	1.39
3	B	1301	600	C64-C65	5.33	1.49	1.42
3	A	801	600	C67-N66	5.26	1.37	1.32
3	A	801	600	C32-C21	5.21	1.60	1.51
3	A	801	600	C5-C4	5.20	1.47	1.39
3	A	801	600	C35-C30	5.05	1.61	1.55
3	A	801	600	C64-C65	5.03	1.49	1.42
3	A	801	600	C3-C4	5.01	1.47	1.39
3	B	1301	600	P55-O57	-4.53	1.46	1.54
3	A	801	600	O77-C76	-4.45	1.26	1.42
3	B	1301	600	C62-C67	4.43	1.47	1.38
3	A	801	600	C13-C14	4.38	1.46	1.39
3	B	1301	600	C3-C4	4.37	1.46	1.39
3	B	1301	600	O77-C76	-4.19	1.27	1.42
3	A	801	600	C62-C67	3.92	1.46	1.38
3	A	801	600	P55-O57	-3.87	1.47	1.54
3	B	1301	600	C43-C42	3.80	1.48	1.40
3	B	1301	600	C45-C44	3.78	1.45	1.36
3	B	1301	600	P79-O81	-3.77	1.45	1.54
3	A	801	600	P79-O81	-3.51	1.46	1.54
3	A	801	600	C63-C62	3.44	1.43	1.36
3	B	1301	600	C67-N66	3.41	1.35	1.32
3	A	801	600	C12-C11	3.35	1.46	1.38
3	B	1301	600	C20-C21	3.27	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	600	C61-C68	3.26	1.46	1.38
3	A	801	600	C45-C44	3.23	1.44	1.36
3	B	1301	600	C63-C62	3.22	1.43	1.36
3	A	801	600	C43-C42	3.19	1.47	1.40
3	B	1301	600	C35-C11	3.18	1.56	1.51
3	A	801	600	C68-C24	3.17	1.56	1.49
3	A	801	600	C40-C41	3.10	1.43	1.36
3	B	1301	600	C61-C68	3.08	1.46	1.38
3	B	1301	600	C52-C14	3.06	1.53	1.50
3	A	801	600	C10-C11	3.05	1.45	1.38
3	B	1301	600	C40-C41	3.00	1.43	1.36
3	A	801	600	C15-C14	2.94	1.44	1.39
3	B	1301	600	C13-C14	2.93	1.43	1.39
3	B	1301	600	C45-C40	2.88	1.45	1.38
3	A	801	600	P79-O82	2.88	1.55	1.49
3	B	1301	600	C10-C11	2.83	1.44	1.38
3	A	801	600	C13-C12	2.80	1.43	1.38
3	A	801	600	C20-C21	2.77	1.44	1.38
3	B	1301	600	C13-C12	2.76	1.43	1.38
3	A	801	600	O77-C71	-2.65	1.29	1.42
3	A	801	600	C32-C30	2.64	1.58	1.55
3	A	801	600	C25-C24	2.61	1.44	1.39
3	B	1301	600	C12-C11	2.57	1.44	1.38
3	B	1301	600	O77-C71	-2.53	1.30	1.42
3	B	1301	600	C23-C24	2.49	1.44	1.39
3	A	801	600	C22-C21	2.47	1.44	1.38
3	B	1301	600	C6-C5	2.43	1.44	1.38
3	B	1301	600	C22-C21	2.39	1.44	1.38
3	A	801	600	C45-C40	2.38	1.44	1.38
3	B	1301	600	C25-C24	2.36	1.44	1.39
3	A	801	600	C69-C68	2.34	1.43	1.39
3	A	801	600	C6-C5	2.33	1.43	1.38
3	B	1301	600	C2-C1	2.32	1.44	1.38
3	B	1301	600	C25-C20	2.21	1.42	1.38
3	A	801	600	C2-C1	2.21	1.43	1.38
3	A	801	600	C25-C20	2.18	1.42	1.38
3	A	801	600	C23-C24	2.02	1.43	1.39
3	A	801	600	C61-C64	2.02	1.46	1.42

All (32) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	600	C73-C72-C71	25.43	154.66	115.23
3	B	1301	600	C73-C72-C71	24.56	153.31	115.23
3	B	1301	600	C72-C71-C67	-11.40	96.32	111.56
3	A	801	600	C72-C71-C67	-10.41	97.65	111.56
3	B	1301	600	C44-C43-N31	8.13	139.43	131.93
3	A	801	600	C44-C43-N31	7.82	139.15	131.93
3	A	801	600	C67-N66-C65	7.76	125.03	118.07
3	B	1301	600	C67-N66-C65	7.60	124.89	118.07
3	A	801	600	C30-C32-C21	4.51	124.60	115.63
3	B	1301	600	C43-C42-N46	-4.24	103.29	108.58
3	B	1301	600	C30-C32-C21	3.98	123.55	115.63
3	A	801	600	C43-C42-N46	-3.91	103.71	108.58
3	A	801	600	C64-C65-N66	-3.56	117.91	122.64
3	B	1301	600	O77-C71-C72	3.50	122.63	107.48
3	A	801	600	C74-C73-C72	3.49	123.95	111.11
3	B	1301	600	C76-O77-C71	3.41	123.38	113.65
3	A	801	600	O77-C71-C72	3.35	121.97	107.48
3	A	801	600	C76-O77-C71	3.24	122.91	113.65
3	B	1301	600	C64-C65-N66	-3.09	118.53	122.64
3	B	1301	600	C41-C42-N46	2.93	134.96	130.19
3	B	1301	600	C74-C73-C72	2.92	121.84	111.11
3	A	801	600	C62-C67-N66	-2.88	119.92	122.86
3	B	1301	600	C62-C67-N66	-2.85	119.95	122.86
3	B	1301	600	C69-C68-C24	-2.75	116.31	120.86
3	A	801	600	C41-C42-N46	2.72	134.62	130.19
3	A	801	600	C13-C14-C52	2.61	122.21	119.84
3	A	801	600	C30-C35-C11	2.40	120.40	115.63
3	B	1301	600	C32-C21-C20	-2.40	117.87	121.07
3	B	1301	600	C13-C14-C52	2.37	121.99	119.84
3	B	1301	600	C25-C24-C68	-2.22	117.51	121.36
3	B	1301	600	C68-C61-C64	-2.03	118.02	121.53
3	B	1301	600	C30-C35-C11	2.02	119.64	115.63

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1301	600	N66-C67-C71-C72
3	B	1301	600	C67-C71-O77-C76
3	B	1301	600	C72-C71-O77-C76

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Mol	Chain	Res	Type	Atoms
3	B	1301	600	C67-C71-C72-C73
3	A	801	600	C62-C67-C71-O77
3	A	801	600	C67-C71-O77-C76
3	A	801	600	C72-C71-O77-C76
3	A	801	600	C67-C71-C72-C73
3	A	801	600	C20-C21-C32-C30
3	A	801	600	C22-C21-C32-C30
3	B	1301	600	C20-C21-C32-C30
3	A	801	600	C69-C70-P79-O80
3	B	1301	600	C22-C21-C32-C30

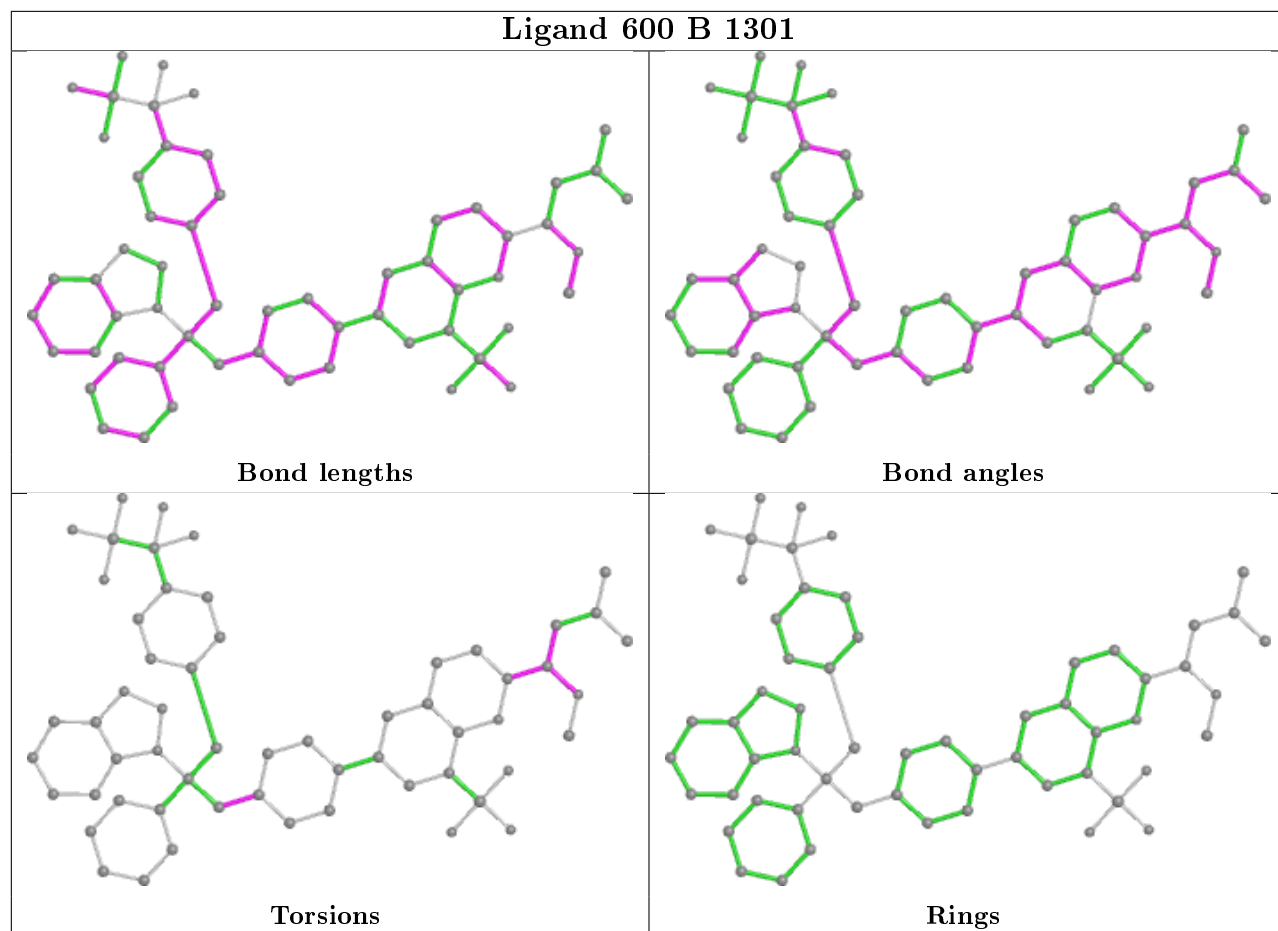
There are no ring outliers.

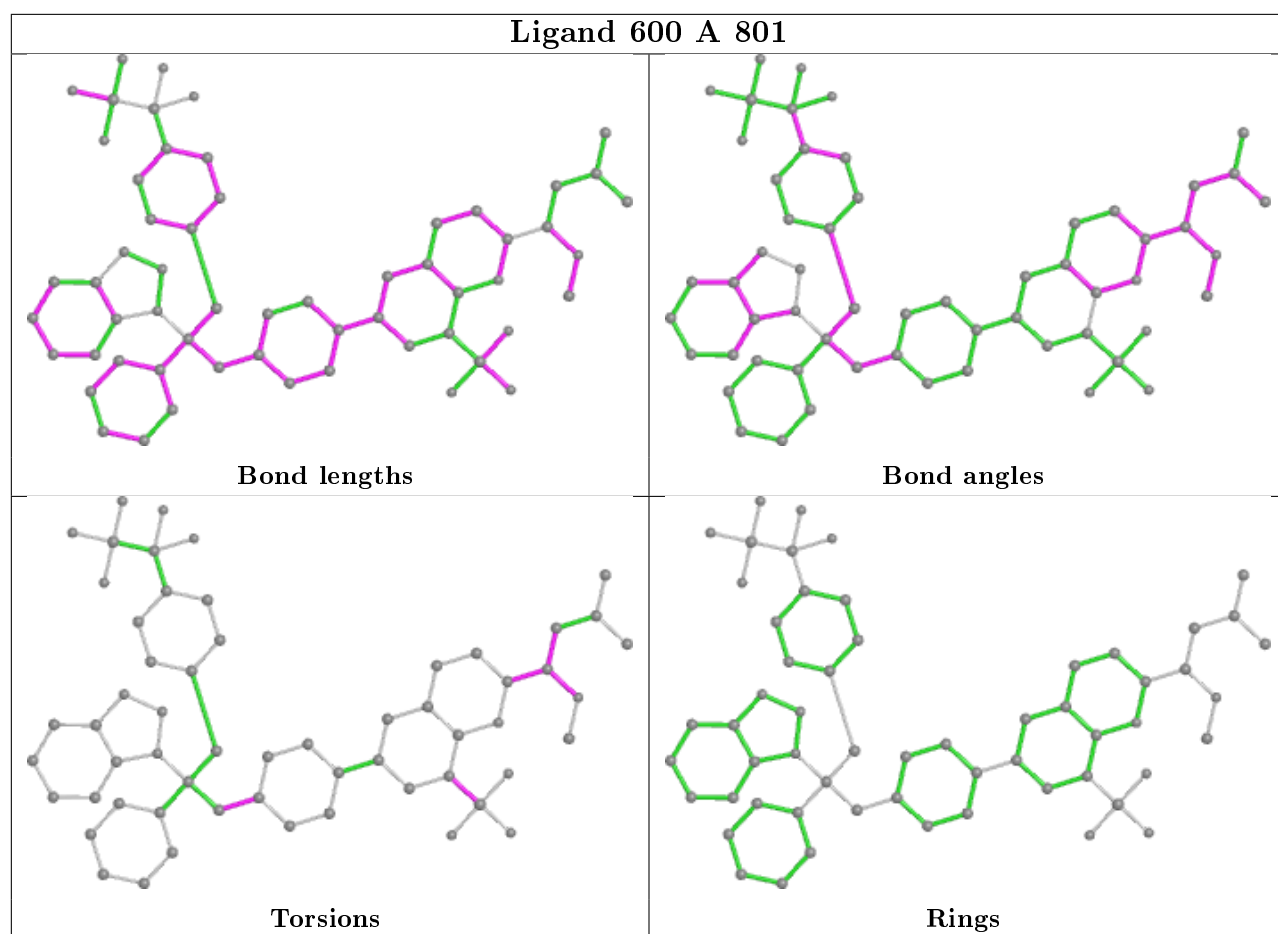
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	600	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand 600 B 1301





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	289/310 (93%)	0.08	14 (4%) 30 37	23, 38, 68, 76	0
1	B	289/310 (93%)	0.13	11 (3%) 40 47	21, 38, 70, 78	0
All	All	578/620 (93%)	0.11	25 (4%) 35 42	21, 39, 70, 78	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1285	SER	11.3
1	A	618	SER	5.9
1	A	619	LEU	5.5
1	B	1000	PHE	4.9
1	B	1117	GLY	4.8
1	B	1063	ASP	3.9
1	A	505	LYS	3.2
1	B	1284	ASP	3.2
1	A	652	TYR	3.1
1	A	616	LYS	3.0
1	A	614	MET	2.9
1	B	1116	LYS	2.9
1	A	617	GLY	2.7
1	B	1062	GLU	2.7
1	A	562	GLU	2.7
1	A	785	SER	2.5
1	A	784	ASP	2.5
1	B	1280	PHE	2.4
1	A	497	LYS	2.4
1	B	1005	LYS	2.4
1	A	500	PHE	2.4
1	A	741	PRO	2.3
1	B	1239	LYS	2.3
1	A	512	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	1237	LYS	2.1

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](#)

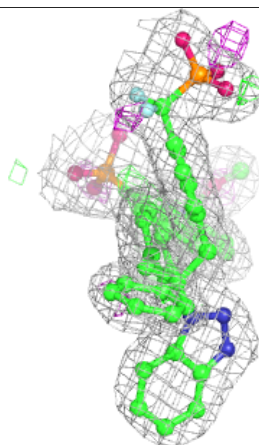
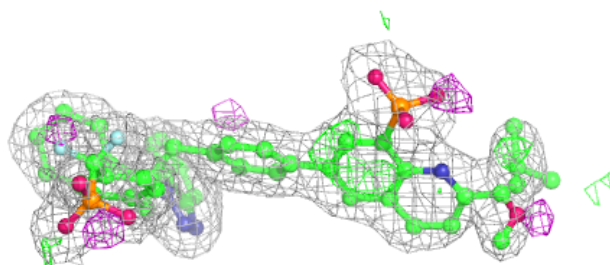
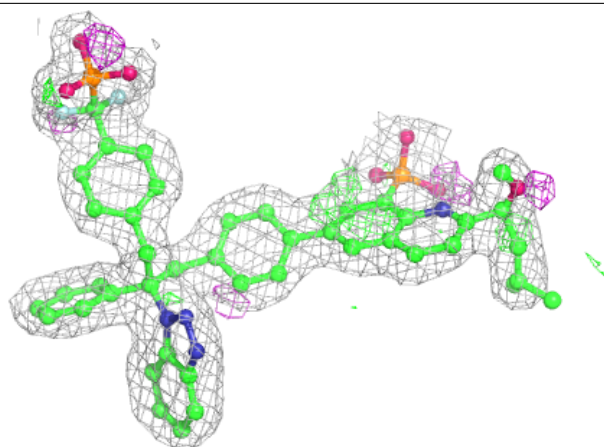
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MG	B	5000	1/1	0.86	0.11	26,26,26,26	1
3	600	B	1301	58/58	0.96	0.14	19,28,44,47	0
3	600	A	801	58/58	0.96	0.16	22,28,55,59	0
2	MG	A	6000	1/1	0.97	0.14	32,32,32,32	1

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.

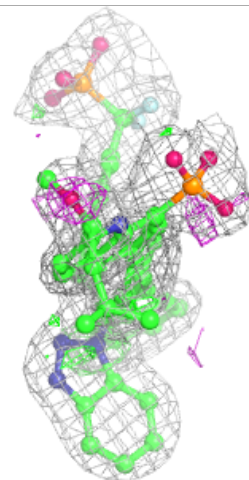
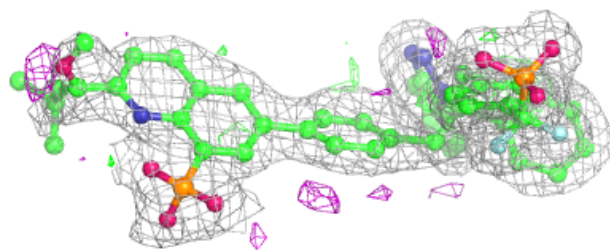
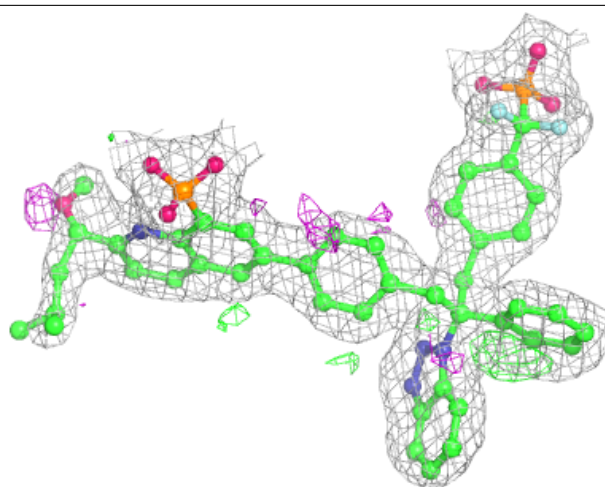
Electron density around 600 B 1301:

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



Electron density around 600 Å 801:

$2mF_o-DF_c$ (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.