



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:38 PM UTC

PDB ID : 3EFK / pdb_00003efk
Title : Structure of c-Met with pyrimidone inhibitor 50
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Deposited on : 2008-09-09
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 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

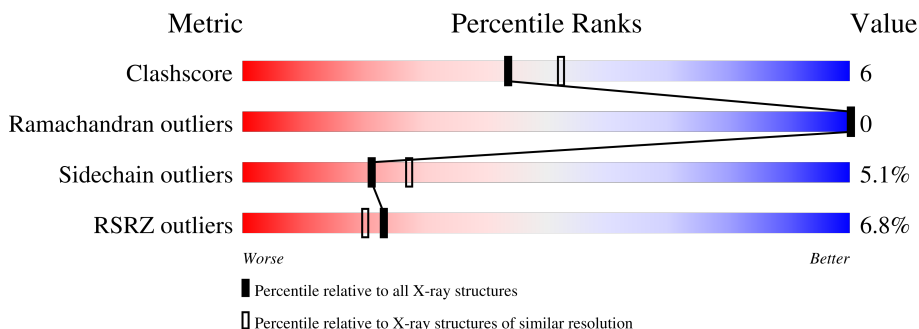
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 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	310	8% 73% 15% • 11%
1	B	310	4% 74% 12% • 12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4486 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.

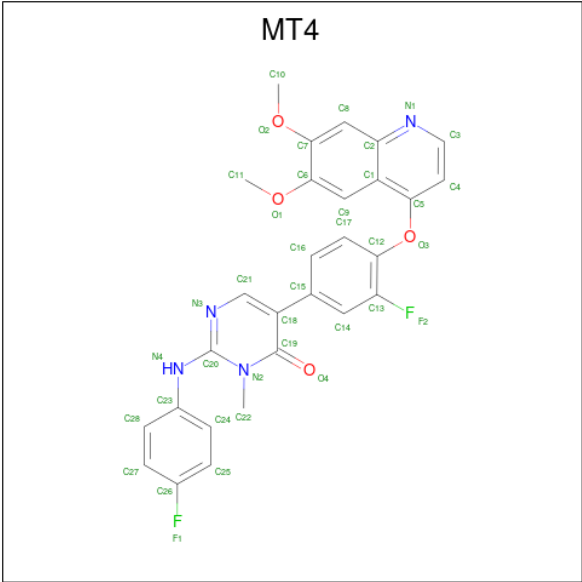
- Molecule 1 is a protein called Hepatocyte growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2166	1402	364	386	14			
1	B	272	Total	C	N	O	S	0	0	0
			2133	1383	359	377	14			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1047	MET	-	initiating methionine	UNP P08581
A	1272	LEU	VAL	engineered mutation	UNP P08581
A	1352	HIS	-	expression tag	UNP P08581
A	1353	HIS	-	expression tag	UNP P08581
A	1354	HIS	-	expression tag	UNP P08581
A	1355	HIS	-	expression tag	UNP P08581
A	1356	HIS	-	expression tag	UNP P08581
B	1047	MET	-	initiating methionine	UNP P08581
B	1272	LEU	VAL	engineered mutation	UNP P08581
B	1352	HIS	-	expression tag	UNP P08581
B	1353	HIS	-	expression tag	UNP P08581
B	1354	HIS	-	expression tag	UNP P08581
B	1355	HIS	-	expression tag	UNP P08581
B	1356	HIS	-	expression tag	UNP P08581

- Molecule 2 is 5-{4-[(6,7-dimethoxyquinolin-4-yl)oxy]-3-fluorophenyl}-2-[(4-fluorophenyl)amino]-3-methylpyrimidin-4(3H)-one (CCD ID: MT4) (formula: C₂₈H₂₂F₂N₄O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 38	C 28	F 2	N 4	O 4	0	0
2	B	1	Total 38	C 28	F 2	N 4	O 4	0	0

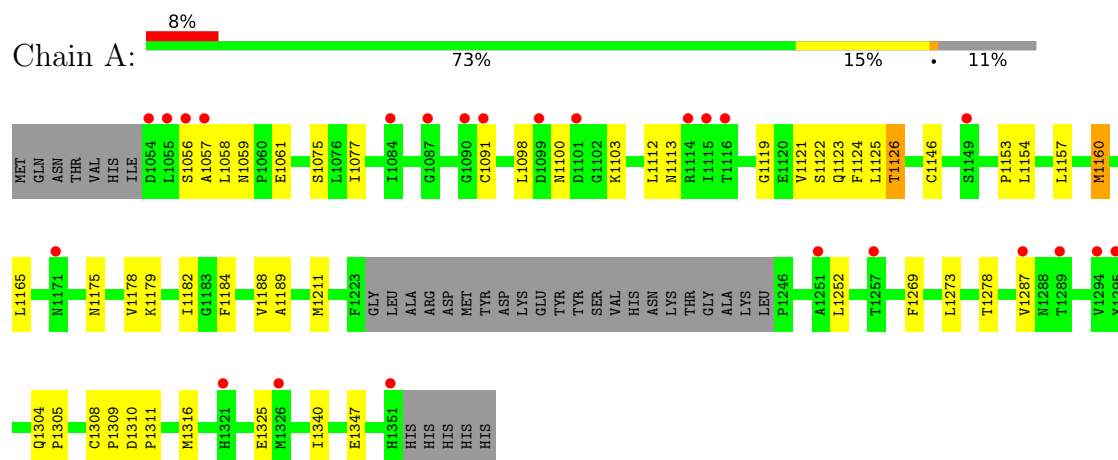
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	62	Total	O	0	0
			62	62		
3	B	49	Total	O	0	0
			49	49		

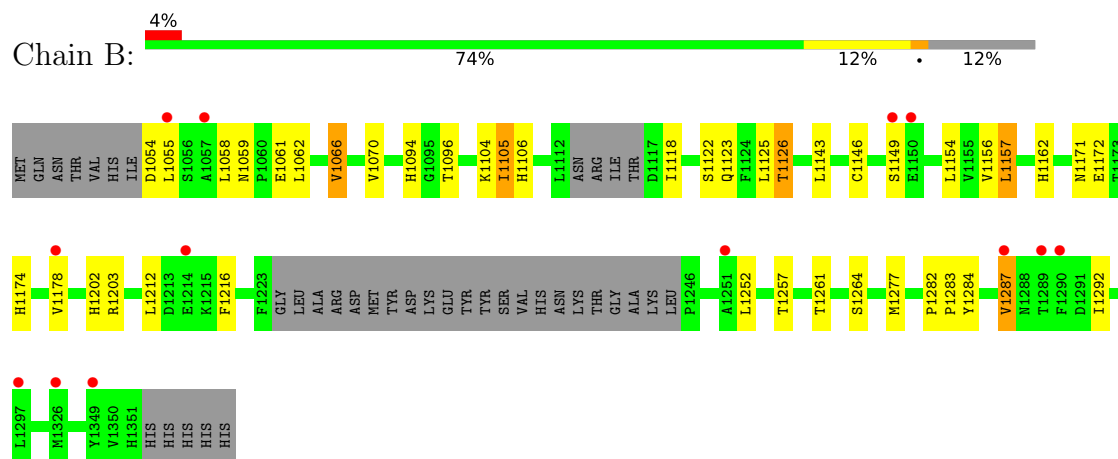
3 Residue-property plots [i](#)

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: Hepatocyte growth factor receptor



• Molecule 1: Hepatocyte growth factor receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.20Å 78.25Å 129.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.64 – 2.20 19.64 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.9 (19.64-2.20) 96.1 (19.64-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.256 , 0.292 0.248 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4486	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.51 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5299e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MT4

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	0/2221	0.85	1/3015 (0.0%)
1	B	0.50	0/2188	0.81	0/2969
All	All	0.49	0/4409	0.83	1/5984 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1057	ALA	N-CA-C	-5.12	107.32	113.21

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	2166	0	2156	29	0
1	B	2133	0	2117	25	0
2	A	38	0	22	3	0
2	B	38	0	22	1	0
3	A	62	0	0	1	0
3	B	49	0	0	0	0
All	All	4486	0	4317	53	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 6.

All (53) 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:1126:THR:HG21	1:B:1126:THR:HG21	1.22	1.15
1:A:1123:GLN:NE2	1:B:1123:GLN:OE1	2.22	0.73
1:A:1126:THR:CG2	1:B:1126:THR:HG21	2.12	0.66
1:A:1119:GLY:O	1:A:1123:GLN:HG2	1.96	0.66
1:B:1062:LEU:O	1:B:1066:VAL:HG23	1.95	0.66
1:A:1122:SER:O	1:A:1126:THR:HG22	1.96	0.64
1:A:1160:MET:HG3	1:A:1211:MET:HB3	1.80	0.64
1:B:1122:SER:O	1:B:1126:THR:HG22	1.97	0.64
1:A:1160:MET:CG	1:A:1211:MET:HB3	2.30	0.62
1:A:1059:ASN:HD21	1:A:1061:GLU:HB2	1.66	0.61
1:B:1212:LEU:HD22	1:B:1216:PHE:HA	1.85	0.58
1:A:1278:THR:HG23	1:A:1305:PRO:HB3	1.86	0.58
1:B:1059:ASN:HD21	1:B:1061:GLU:HB2	1.72	0.55
1:A:1112:LEU:HD12	1:A:1153:PRO:HG2	1.89	0.53
1:A:1146:CYS:HB3	1:A:1154:LEU:HB2	1.90	0.52
1:B:1066:VAL:HG11	1:B:1125:LEU:HD22	1.92	0.52
1:B:1096:THR:CG2	1:B:1104:LYS:HG3	2.40	0.52
1:A:1278:THR:CG2	1:A:1305:PRO:HB3	2.39	0.52
1:B:1094:HIS:HD2	1:B:1106:HIS:NE2	2.10	0.50
1:B:1156:VAL:C	1:B:1157:LEU:HG	2.37	0.49
1:B:1284:TYR:HB3	1:B:1287:VAL:HG12	1.94	0.49
1:A:1112:LEU:HD11	1:A:1124:PHE:CE1	2.47	0.49
1:B:1055:LEU:HD23	1:B:1055:LEU:H	1.78	0.49
1:A:1273:LEU:HB3	1:A:1316:MET:HE2	1.95	0.49
1:A:1184:PHE:O	1:A:1188:VAL:HG23	2.13	0.48
1:A:1056:SER:C	1:A:1058:LEU:H	2.20	0.48
1:B:1162:HIS:HB2	1:B:1212:LEU:HB3	1.96	0.48
1:B:1105:ILE:HD11	1:B:1143:LEU:HD11	1.96	0.47
1:A:1165:LEU:HD21	1:A:1184:PHE:HB3	1.97	0.46
1:A:1121:VAL:O	1:A:1125:LEU:HG	2.16	0.46
1:A:1160:MET:O	2:A:1:MT4:H10A	2.16	0.45
1:B:1059:ASN:HD22	1:B:1062:LEU:H	1.62	0.45
1:A:1160:MET:HB2	2:A:1:MT4:N1	2.31	0.45
1:A:1340:ILE:HG21	3:A:1409:HOH:O	2.16	0.45
1:A:1310:ASP:N	1:A:1311:PRO:HD2	2.32	0.45
1:B:1096:THR:HG23	1:B:1104:LYS:HG3	1.98	0.45
1:B:1171:ASN:HB3	1:B:1174:HIS:ND1	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1160:MET:HG2	1:A:1211:MET:HB3	1.99	0.45
1:A:1304:GLN:HA	1:A:1305:PRO:HD3	1.87	0.44
2:A:1:MT4:H16	2:A:1:MT4:O4	2.17	0.44
1:B:1261:THR:H	1:B:1264:SER:HB3	1.84	0.43
1:B:1202:HIS:O	1:B:1203:ARG:HB2	2.19	0.42
1:A:1179:LYS:HD3	1:A:1347:GLU:HG3	2.00	0.42
1:B:1146:CYS:HB3	1:B:1154:LEU:HB2	2.00	0.42
1:A:1308:CYS:HA	1:A:1309:PRO:HD3	1.91	0.42
1:B:1282:PRO:HA	1:B:1283:PRO:HD3	1.87	0.41
1:B:1058:LEU:HA	1:B:1118:ILE:HD11	2.03	0.41
2:B:2:MT4:O4	2:B:2:MT4:H16	2.20	0.41
1:A:1178:VAL:O	1:A:1182:ILE:HG13	2.21	0.40
1:B:1066:VAL:HG21	1:B:1125:LEU:HD13	2.02	0.40
1:A:1075:SER:HB2	1:A:1098:LEU:O	2.22	0.40
1:A:1189:ALA:HB2	1:A:1269:PHE:CZ	2.57	0.40
1:B:1178:VAL:HG13	1:B:1277:MET:HE1	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	272/310 (88%)	263 (97%)	9 (3%)	0	100	100
1	B	266/310 (86%)	261 (98%)	5 (2%)	0	100	100
All	All	538/620 (87%)	524 (97%)	14 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	239/276 (87%)	227 (95%)	12 (5%)	22	28
1	B	234/276 (85%)	222 (95%)	12 (5%)	21	27
All	All	473/552 (86%)	449 (95%)	24 (5%)	21	27

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1077	ILE
1	A	1091	CYS
1	A	1100	ASN
1	A	1103	LYS
1	A	1113	ASN
1	A	1126	THR
1	A	1157	LEU
1	A	1160	MET
1	A	1175	ASN
1	A	1252	LEU
1	A	1287	VAL
1	A	1325	GLU
1	B	1054	ASP
1	B	1066	VAL
1	B	1070	VAL
1	B	1105	ILE
1	B	1126	THR
1	B	1149	SER
1	B	1157	LEU
1	B	1172	GLU
1	B	1252	LEU
1	B	1257	THR
1	B	1287	VAL
1	B	1292	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1059	ASN
1	A	1162	HIS
1	A	1167	ASN
1	A	1256	GLN
1	A	1258	GLN
1	A	1288	ASN
1	A	1298	GLN
1	B	1059	ASN
1	B	1067	GLN
1	B	1094	HIS
1	B	1162	HIS
1	B	1256	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	MT4	A	1	-	42,42,42	1.55	5 (11%)	58,60,60	2.17	13 (22%)
2	MT4	B	2	-	42,42,42	1.58	6 (14%)	58,60,60	2.08	16 (27%)

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	MT4	A	1	-	-	0/16/16/16	0/5/5/5
2	MT4	B	2	-	-	0/16/16/16	0/5/5/5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	MT4	C15-C18	-4.78	1.41	1.48
2	B	2	MT4	C15-C18	-4.57	1.41	1.48
2	B	2	MT4	C20-N2	4.37	1.42	1.38
2	A	1	MT4	C20-N2	3.90	1.41	1.38
2	A	1	MT4	C18-C19	-3.60	1.40	1.46
2	A	1	MT4	C23-N4	-3.56	1.34	1.41
2	B	2	MT4	C23-N4	-3.42	1.34	1.41
2	B	2	MT4	C18-C19	-3.39	1.41	1.46
2	A	1	MT4	C8-C7	2.15	1.40	1.36
2	B	2	MT4	C21-C18	2.13	1.40	1.37
2	B	2	MT4	C9-C6	2.04	1.40	1.36

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	MT4	N2-C20-N3	-9.75	118.85	124.47
2	B	2	MT4	N2-C20-N3	-8.67	119.47	124.47
2	A	1	MT4	C15-C18-C19	5.46	124.09	119.51
2	B	2	MT4	C15-C18-C19	5.12	123.81	119.51
2	A	1	MT4	N4-C20-N2	4.99	119.16	114.75
2	B	2	MT4	N4-C20-N2	4.78	118.97	114.75
2	A	1	MT4	C20-N2-C19	-3.40	119.50	120.79
2	A	1	MT4	C18-C21-N3	-3.37	122.47	126.19
2	B	2	MT4	C11-O1-C6	-3.26	112.73	117.51
2	B	2	MT4	C18-C21-N3	-3.16	122.70	126.19
2	A	1	MT4	C21-C18-C19	-3.10	117.30	119.70
2	B	2	MT4	C20-N2-C19	-3.07	119.63	120.79
2	A	1	MT4	C3-N1-C2	3.06	121.56	116.93
2	B	2	MT4	O1-C6-C9	-2.91	121.47	125.16
2	A	1	MT4	O4-C19-C18	-2.83	121.67	125.37
2	B	2	MT4	C3-N1-C2	2.82	121.20	116.93
2	A	1	MT4	O1-C6-C9	-2.75	121.67	125.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	MT4	C1-C2-N1	-2.56	120.11	122.82
2	A	1	MT4	C18-C19-N2	2.54	119.54	116.21
2	B	2	MT4	C21-C18-C19	-2.50	117.77	119.70
2	A	1	MT4	C4-C3-N1	-2.48	120.92	124.60
2	B	2	MT4	C4-C3-N1	-2.46	120.95	124.60
2	B	2	MT4	O4-C19-C18	-2.45	122.16	125.37
2	B	2	MT4	O2-C7-C8	-2.30	122.24	125.16
2	B	2	MT4	O1-C6-C7	2.27	118.48	115.40
2	B	2	MT4	C1-C2-N1	-2.21	120.48	122.82
2	A	1	MT4	C11-O1-C6	-2.14	114.37	117.51
2	B	2	MT4	C18-C19-N2	2.11	118.98	116.21
2	B	2	MT4	O2-C7-C6	2.04	118.17	115.40

There are no chirality outliers.

There are no torsion outliers.

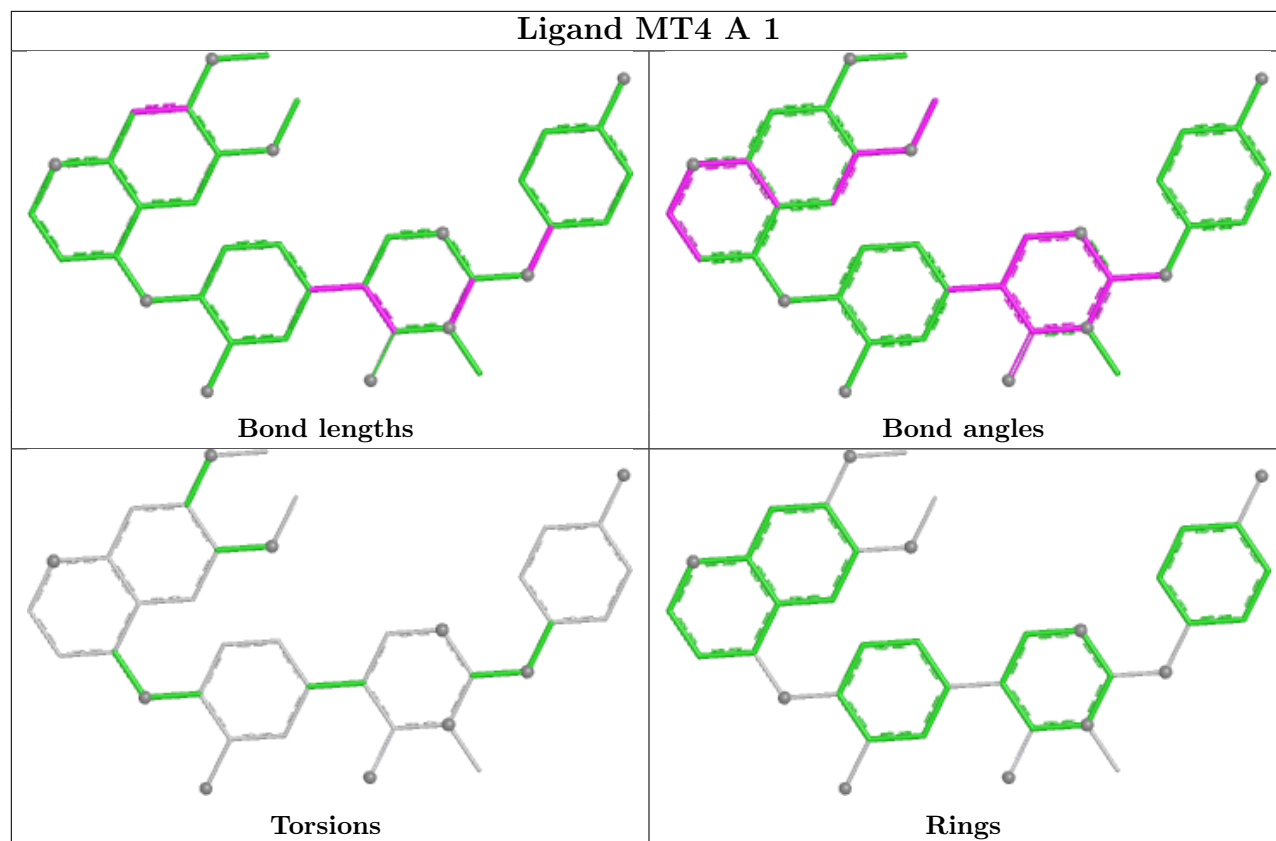
There are no ring outliers.

2 monomers are involved in 4 short contacts:

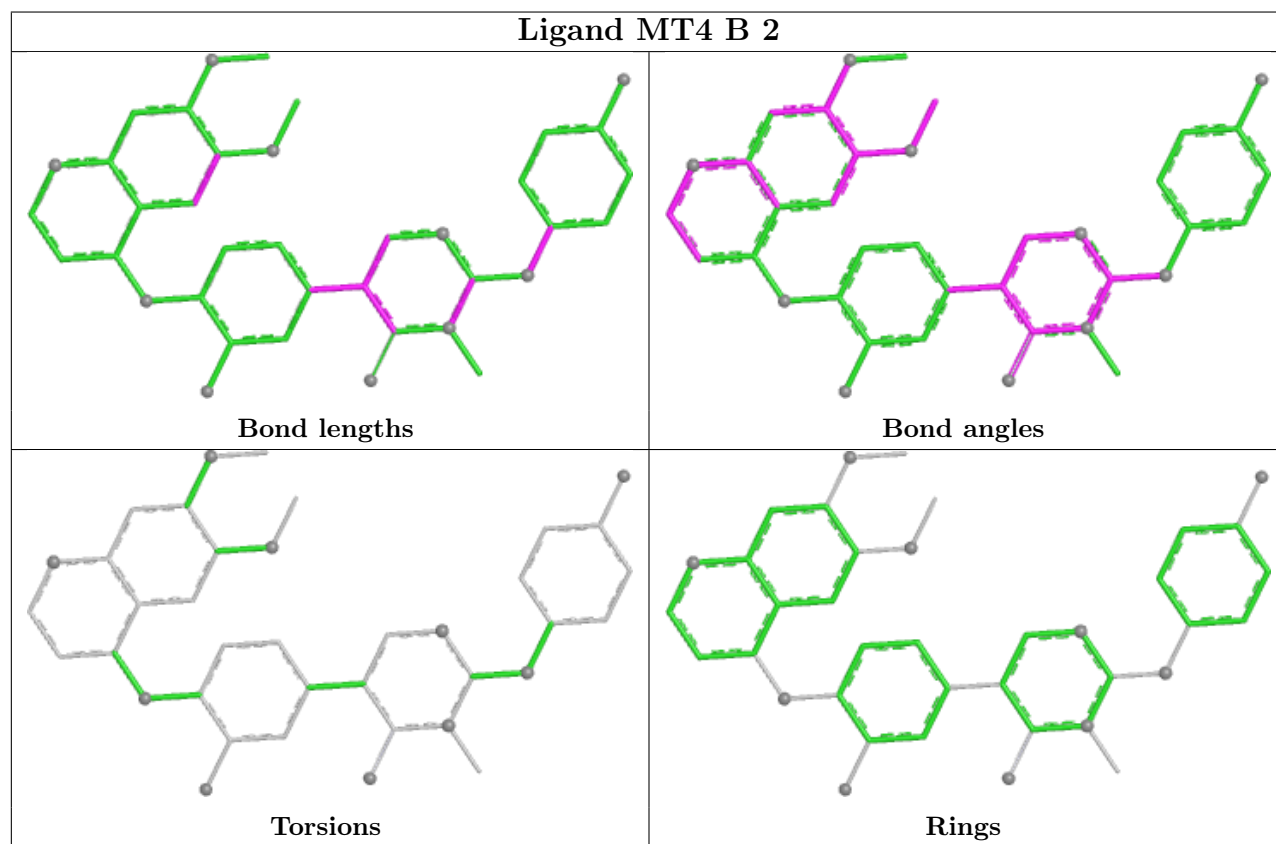
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	MT4	3	0
2	B	2	MT4	1	0

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.

Ligand MT4 A 1



Ligand MT4 B 2



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	276/310 (89%)	0.57	24 (8%) 16 13	23, 35, 53, 61	0
1	B	272/310 (87%)	0.67	13 (4%) 35 32	23, 39, 59, 67	0
All	All	548/620 (88%)	0.62	37 (6%) 23 20	23, 37, 57, 67	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1084	ILE	5.2
1	A	1321	HIS	4.2
1	B	1290	PHE	4.1
1	A	1091	CYS	4.1
1	A	1090	GLY	3.9
1	A	1055	LEU	3.7
1	B	1057	ALA	3.7
1	B	1287	VAL	3.6
1	A	1257	THR	3.4
1	A	1115	ILE	3.1
1	B	1178	VAL	2.9
1	A	1326	MET	2.9
1	A	1087	GLY	2.8
1	A	1057	ALA	2.8
1	A	1056	SER	2.7
1	A	1114	ARG	2.7
1	A	1149	SER	2.7
1	A	1351	HIS	2.5
1	B	1149	SER	2.4
1	B	1349	TYR	2.3
1	A	1287	VAL	2.3
1	A	1054	ASP	2.3
1	A	1251	ALA	2.3
1	B	1326	MET	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	1116	THR	2.3
1	A	1289	THR	2.3
1	B	1289	THR	2.2
1	B	1150	GLU	2.2
1	A	1171	ASN	2.1
1	A	1295	TYR	2.1
1	A	1099	ASP	2.1
1	A	1294	VAL	2.0
1	B	1214	GLU	2.0
1	B	1251	ALA	2.0
1	A	1101	ASP	2.0
1	B	1055	LEU	2.0
1	B	1297	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

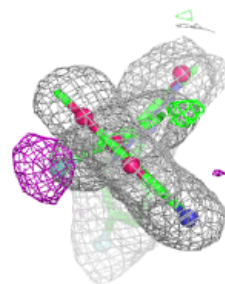
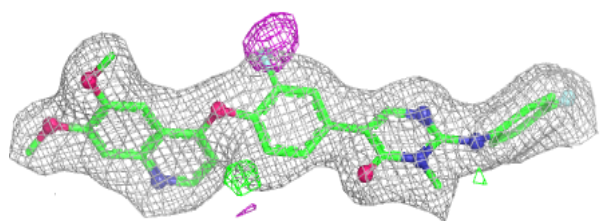
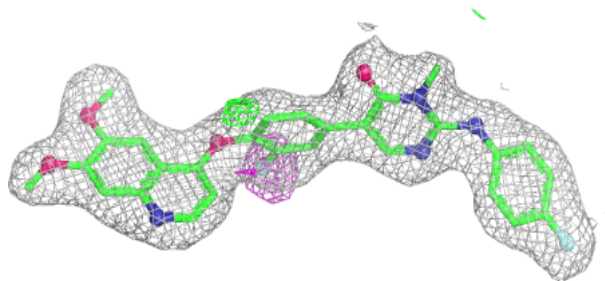
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MT4	A	1	38/38	0.91	0.08	26,28,30,32	0
2	MT4	B	2	38/38	0.91	0.09	26,29,30,31	0

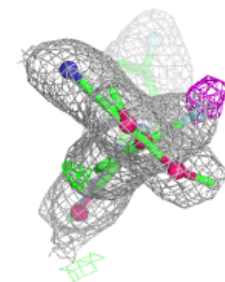
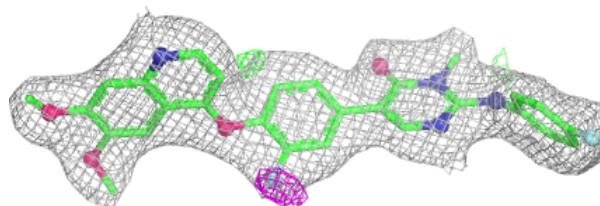
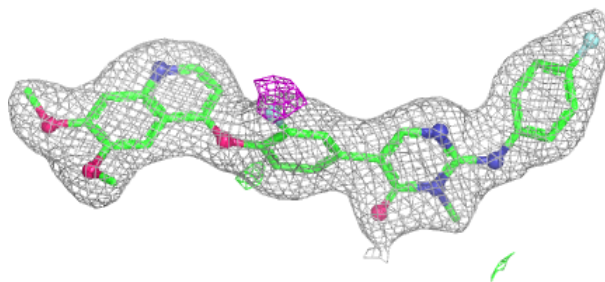
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around MT4 A 1:

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

**Electron density around MT4 B 2:**

$2mF_o-DF_c$ (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.