



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:21 PM UTC

PDB ID : 4RXD / pdb_00004rx
Title : T. Brucei Farnesyl Diphosphate Synthase Complexed with Risedronate
Authors : Cao, R.; Liu, Y.-L.; Oldfield, E.
Deposited on : 2014-12-10
Resolution : 2.00 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

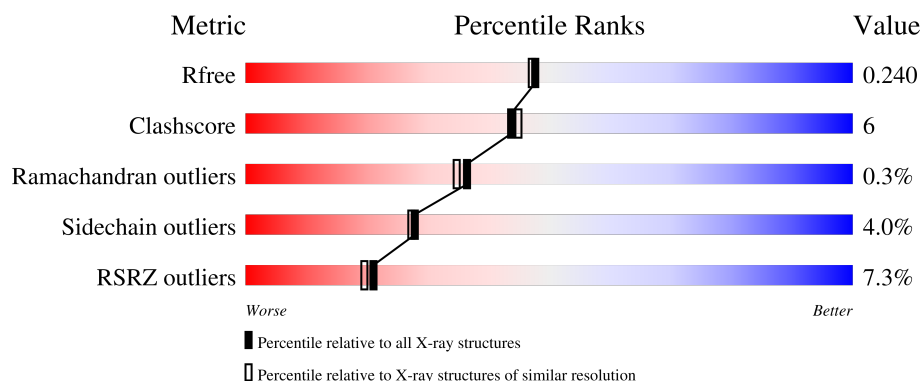
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	390	7% 76% 13% • 9%
1	B	390	5% 81% 9% • 8%
1	C	390	8% 80% 10% • 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9091 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 Farnesyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	S	0	0	0
			2839	1809	467	536	27			
1	B	357	Total	C	N	O	S	0	0	0
			2854	1819	469	538	28			
1	C	357	Total	C	N	O	S	0	0	0
			2854	1819	469	538	28			

There are 69 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP Q86C09
A	-21	GLY	-	expression tag	UNP Q86C09
A	-20	SER	-	expression tag	UNP Q86C09
A	-19	SER	-	expression tag	UNP Q86C09
A	-18	HIS	-	expression tag	UNP Q86C09
A	-17	HIS	-	expression tag	UNP Q86C09
A	-16	HIS	-	expression tag	UNP Q86C09
A	-15	HIS	-	expression tag	UNP Q86C09
A	-14	HIS	-	expression tag	UNP Q86C09
A	-13	HIS	-	expression tag	UNP Q86C09
A	-12	SER	-	expression tag	UNP Q86C09
A	-11	SER	-	expression tag	UNP Q86C09
A	-10	GLY	-	expression tag	UNP Q86C09
A	-9	LEU	-	expression tag	UNP Q86C09
A	-8	VAL	-	expression tag	UNP Q86C09
A	-7	PRO	-	expression tag	UNP Q86C09
A	-6	ARG	-	expression tag	UNP Q86C09
A	-5	GLY	-	expression tag	UNP Q86C09
A	-4	SER	-	expression tag	UNP Q86C09
A	-3	HIS	-	expression tag	UNP Q86C09
A	-2	MET	-	expression tag	UNP Q86C09
A	-1	ALA	-	expression tag	UNP Q86C09
A	0	SER	-	expression tag	UNP Q86C09

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	MET	-	expression tag	UNP Q86C09
B	-21	GLY	-	expression tag	UNP Q86C09
B	-20	SER	-	expression tag	UNP Q86C09
B	-19	SER	-	expression tag	UNP Q86C09
B	-18	HIS	-	expression tag	UNP Q86C09
B	-17	HIS	-	expression tag	UNP Q86C09
B	-16	HIS	-	expression tag	UNP Q86C09
B	-15	HIS	-	expression tag	UNP Q86C09
B	-14	HIS	-	expression tag	UNP Q86C09
B	-13	HIS	-	expression tag	UNP Q86C09
B	-12	SER	-	expression tag	UNP Q86C09
B	-11	SER	-	expression tag	UNP Q86C09
B	-10	GLY	-	expression tag	UNP Q86C09
B	-9	LEU	-	expression tag	UNP Q86C09
B	-8	VAL	-	expression tag	UNP Q86C09
B	-7	PRO	-	expression tag	UNP Q86C09
B	-6	ARG	-	expression tag	UNP Q86C09
B	-5	GLY	-	expression tag	UNP Q86C09
B	-4	SER	-	expression tag	UNP Q86C09
B	-3	HIS	-	expression tag	UNP Q86C09
B	-2	MET	-	expression tag	UNP Q86C09
B	-1	ALA	-	expression tag	UNP Q86C09
B	0	SER	-	expression tag	UNP Q86C09
C	-22	MET	-	expression tag	UNP Q86C09
C	-21	GLY	-	expression tag	UNP Q86C09
C	-20	SER	-	expression tag	UNP Q86C09
C	-19	SER	-	expression tag	UNP Q86C09
C	-18	HIS	-	expression tag	UNP Q86C09
C	-17	HIS	-	expression tag	UNP Q86C09
C	-16	HIS	-	expression tag	UNP Q86C09
C	-15	HIS	-	expression tag	UNP Q86C09
C	-14	HIS	-	expression tag	UNP Q86C09
C	-13	HIS	-	expression tag	UNP Q86C09
C	-12	SER	-	expression tag	UNP Q86C09
C	-11	SER	-	expression tag	UNP Q86C09
C	-10	GLY	-	expression tag	UNP Q86C09
C	-9	LEU	-	expression tag	UNP Q86C09
C	-8	VAL	-	expression tag	UNP Q86C09
C	-7	PRO	-	expression tag	UNP Q86C09
C	-6	ARG	-	expression tag	UNP Q86C09
C	-5	GLY	-	expression tag	UNP Q86C09
C	-4	SER	-	expression tag	UNP Q86C09

Continued on next page...

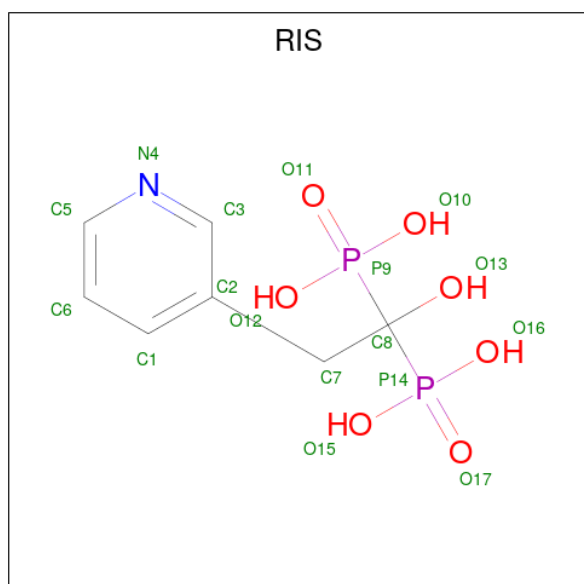
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	HIS	-	expression tag	UNP Q86C09
C	-2	MET	-	expression tag	UNP Q86C09
C	-1	ALA	-	expression tag	UNP Q86C09
C	0	SER	-	expression tag	UNP Q86C09

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 3 is 1-HYDROXY-2-(3-PYRIDINYL)ETHYLIDENE BIS-PHOSPHONIC ACID (CCD ID: RIS) (formula: C₇H₁₁NO₇P₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O P 17 7 1 7 2	0	0
3	B	1	Total C N O P 17 7 1 7 2	0	0
3	C	1	Total C N O P 17 7 1 7 2	0	0

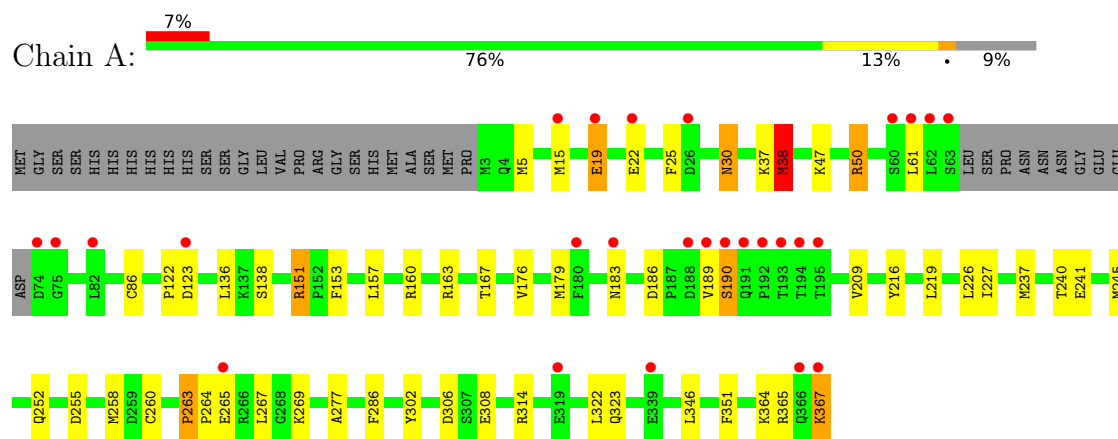
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	166	Total 166	O 166	0	0
4	B	159	Total 159	O 159	0	0
4	C	159	Total 159	O 159	0	0

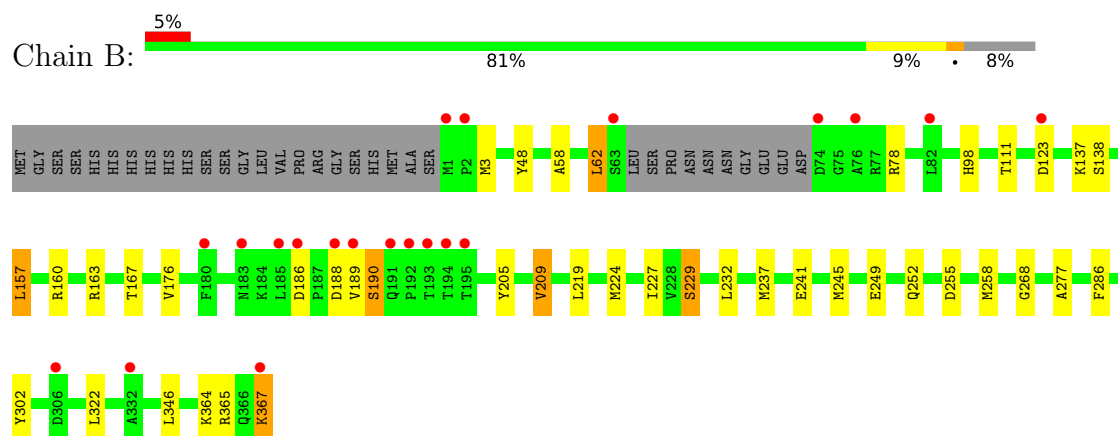
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

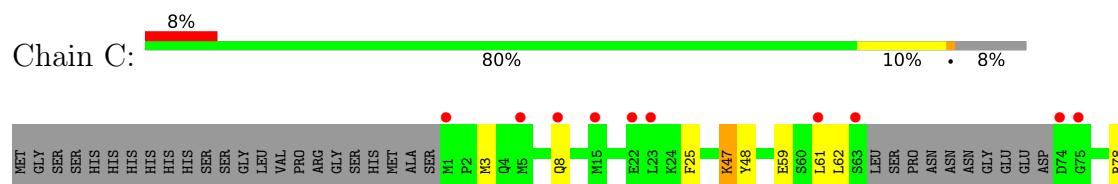
• Molecule 1: Farnesyl pyrophosphate synthase

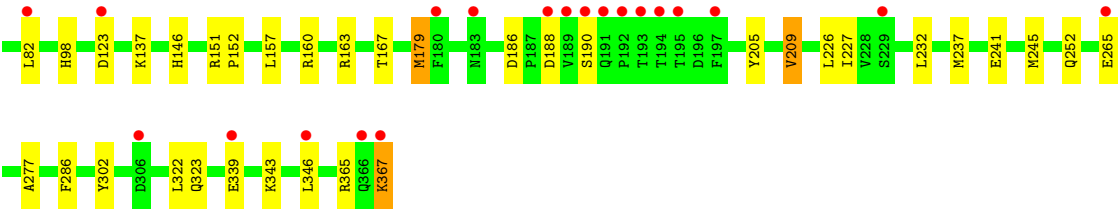


• Molecule 1: Farnesyl pyrophosphate synthase



• Molecule 1: Farnesyl pyrophosphate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	156.34Å 90.31Å 88.71Å 90.00° 89.91° 90.00°	Depositor
Resolution (Å)	29.55 – 2.00 29.55 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.55-2.00) 99.6 (29.55-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.71 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.211 , 0.251 (Not available) , 0.240	Depositor DCC
R_{free} test set	4142 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.137	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.016 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.017 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.486 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.486 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.015 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9091	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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, RIS

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	0.60	1/2897 (0.0%)	0.84	2/3917 (0.1%)
1	B	0.58	0/2913	0.83	0/3939
1	C	0.60	0/2913	0.83	0/3939
All	All	0.60	1/8723 (0.0%)	0.84	2/11795 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	38	MET	SD-CE	-6.54	1.63	1.79

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	PRO	CA-C-N	5.21	125.26	119.32
1	A	263	PRO	C-N-CA	5.21	125.26	119.32

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	2839	0	2789	51	0
1	B	2854	0	2808	29	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2854	0	2808	31	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
3	A	17	0	7	0	0
3	B	17	0	7	0	0
3	C	17	0	7	0	0
4	A	166	0	0	4	0
4	B	159	0	0	2	0
4	C	159	0	0	1	0
All	All	9091	0	8426	104	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 6.

All (104) 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:C:365:ARG:HH11	1:C:367:LYS:HG3	1.04	1.14
1:B:365:ARG:HH11	1:B:367:LYS:HG3	1.13	1.06
1:A:365:ARG:HH11	1:A:367:LYS:HG3	1.13	1.05
1:C:365:ARG:NH1	1:C:367:LYS:HG3	1.79	0.96
1:B:365:ARG:NH1	1:B:367:LYS:HG3	1.87	0.90
1:A:365:ARG:NH1	1:A:367:LYS:HG3	1.86	0.90
1:A:186:ASP:O	1:C:123:ASP:CB	2.22	0.87
1:A:237:MET:HE3	1:A:240:THR:HB	1.57	0.84
1:B:209:VAL:HG22	1:B:252:GLN:HG2	1.61	0.83
1:A:38:MET:HE2	1:A:38:MET:C	2.04	0.82
1:C:209:VAL:HG22	1:C:252:GLN:HG2	1.62	0.80
1:A:123:ASP:CB	1:C:186:ASP:O	2.31	0.78
1:A:50:ARG:HG2	4:A:1527:HOH:O	1.83	0.77
1:A:237:MET:HE2	1:A:241:GLU:HG3	1.68	0.75
1:B:365:ARG:HH11	1:B:367:LYS:CG	1.98	0.75
1:A:61:LEU:CD1	1:A:226:LEU:HD23	2.16	0.74
1:B:163:ARG:O	1:B:167:THR:HG23	1.90	0.72
1:A:167:THR:HG22	1:C:25:PHE:HE1	1.55	0.70
1:B:232:LEU:HD11	1:B:237:MET:CE	2.23	0.69
1:A:38:MET:HE2	1:A:38:MET:O	1.93	0.68
1:A:209:VAL:HG22	1:A:252:GLN:HG2	1.76	0.68
1:A:163:ARG:O	1:A:167:THR:HG23	1.94	0.67
1:A:167:THR:HG22	1:C:25:PHE:CE1	2.31	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:PHE:HE1	1:C:167:THR:HG22	1.61	0.65
1:A:151:ARG:HG2	1:A:153:PHE:CZ	2.31	0.65
1:C:209:VAL:HG22	1:C:252:GLN:CG	2.26	0.64
1:B:157:LEU:C	1:B:157:LEU:HD23	2.23	0.64
1:C:163:ARG:O	1:C:167:THR:HG23	1.98	0.63
1:C:232:LEU:HD11	1:C:237:MET:CE	2.29	0.62
1:B:157:LEU:HD21	1:B:224:MET:HE3	1.82	0.62
1:A:25:PHE:CE1	1:C:167:THR:HG22	2.35	0.61
1:A:61:LEU:HD13	1:A:226:LEU:HD23	1.83	0.61
1:A:237:MET:HE2	1:A:241:GLU:CG	2.30	0.61
1:C:160:ARG:HD2	1:C:227:ILE:HD11	1.83	0.60
1:A:47:LYS:HD3	1:A:367:LYS:HE2	1.84	0.58
1:A:237:MET:CE	1:A:240:THR:HB	2.33	0.58
1:C:47:LYS:HG2	1:C:367:LYS:HE2	1.85	0.58
1:B:232:LEU:HD11	1:B:237:MET:HE1	1.88	0.56
1:A:38:MET:HE1	1:A:136:LEU:HD21	1.88	0.56
1:A:190:SER:HB3	4:A:1541:HOH:O	2.06	0.55
1:B:157:LEU:HD23	1:B:157:LEU:O	2.06	0.55
1:B:160:ARG:HD2	1:B:227:ILE:HD11	1.88	0.55
1:B:3:MET:HE1	1:B:48:TYR:CE2	2.43	0.53
1:A:260:CYS:O	1:A:314:ARG:NH1	2.42	0.53
1:A:160:ARG:HD2	1:A:227:ILE:HD11	1.91	0.52
1:B:209:VAL:HG22	1:B:252:GLN:CG	2.37	0.52
1:A:30:ASN:HD22	1:A:30:ASN:H	1.57	0.52
1:C:3:MET:HE1	1:C:48:TYR:CE2	2.45	0.52
1:C:47:LYS:HE3	1:C:47:LYS:HA	1.91	0.52
1:B:157:LEU:C	1:B:157:LEU:CD2	2.83	0.52
1:A:364:LYS:HA	4:A:1552:HOH:O	2.09	0.51
1:C:61:LEU:CD1	1:C:226:LEU:HD23	2.41	0.51
1:A:306:ASP:OD1	1:A:308:GLU:HG2	2.11	0.50
1:C:205:TYR:CZ	1:C:209:VAL:HG21	2.46	0.50
1:A:209:VAL:CG1	1:A:252:GLN:HB3	2.42	0.50
1:C:146:HIS:HE1	4:C:1794:HOH:O	1.95	0.49
1:C:98:HIS:CD2	1:C:137:LYS:HE3	2.48	0.49
1:A:50:ARG:NH1	1:A:216:TYR:CE1	2.81	0.48
1:C:232:LEU:HD11	1:C:237:MET:HE1	1.95	0.48
1:C:286:PHE:HB2	1:C:322:LEU:HD11	1.94	0.48
1:B:205:TYR:CZ	1:B:209:VAL:HG21	2.48	0.48
1:A:30:ASN:H	1:A:30:ASN:ND2	2.12	0.47
1:A:47:LYS:CD	1:A:367:LYS:HE2	2.45	0.47
1:B:209:VAL:CG1	1:B:249:GLU:HA	2.44	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:LEU:HD21	1:B:78:ARG:HA	1.96	0.47
1:A:15:MET:O	1:A:19:GLU:HB3	2.15	0.47
1:B:189:VAL:O	1:B:190:SER:CB	2.64	0.46
1:A:209:VAL:HG13	1:A:252:GLN:CB	2.46	0.46
1:A:61:LEU:HD11	1:A:226:LEU:HD23	1.93	0.45
1:C:59:GLU:HG3	1:C:82:LEU:HD21	1.98	0.45
1:B:364:LYS:HA	4:B:1676:HOH:O	2.16	0.45
1:B:111:THR:HG22	1:B:268:GLY:O	2.16	0.45
1:A:151:ARG:HG2	1:A:153:PHE:CE2	2.50	0.45
1:C:78:ARG:O	1:C:82:LEU:HD23	2.17	0.45
1:A:61:LEU:HD21	1:A:351:PHE:HB2	2.00	0.44
1:B:157:LEU:HD21	1:B:224:MET:CE	2.48	0.44
1:C:277:ALA:HA	1:C:302:TYR:CE2	2.53	0.44
1:C:241:GLU:O	1:C:245:MET:HG2	2.18	0.43
1:A:50:ARG:HG2	1:A:50:ARG:H	1.67	0.43
1:B:286:PHE:HB2	1:B:322:LEU:HD11	2.00	0.43
1:A:209:VAL:HG11	1:A:252:GLN:HB3	1.99	0.43
1:B:232:LEU:HD11	1:B:237:MET:HE3	1.99	0.43
1:B:241:GLU:O	1:B:245:MET:HG2	2.18	0.43
1:A:277:ALA:HA	1:A:302:TYR:CE2	2.54	0.43
1:A:122:PRO:HG2	1:C:188:ASP:C	2.44	0.42
1:C:232:LEU:HD11	1:C:237:MET:HE3	1.97	0.42
1:A:267:LEU:HG	1:A:269:LYS:HG2	2.01	0.42
1:B:189:VAL:O	4:B:1743:HOH:O	2.21	0.42
1:A:189:VAL:O	1:A:190:SER:CB	2.68	0.42
1:B:98:HIS:CD2	1:B:137:LYS:HE3	2.55	0.42
1:B:58:ALA:O	1:B:229:SER:OG	2.38	0.41
1:A:269:LYS:HE3	4:A:1647:HOH:O	2.20	0.41
1:A:255:ASP:HA	1:A:258:MET:HE2	2.03	0.41
1:A:286:PHE:HB2	1:A:322:LEU:HD11	2.03	0.41
1:A:237:MET:HE3	1:A:237:MET:O	2.20	0.41
1:B:255:ASP:HA	1:B:258:MET:HE2	2.03	0.41
1:C:151:ARG:HA	1:C:152:PRO:HD3	1.98	0.41
1:A:365:ARG:HH11	1:A:367:LYS:CG	2.05	0.40
1:A:5:MET:HG3	1:A:86:CYS:SG	2.62	0.40
1:A:241:GLU:O	1:A:245:MET:HG2	2.22	0.40
1:C:179:MET:SD	1:C:179:MET:N	2.91	0.40
1:B:277:ALA:HA	1:B:302:TYR:CE2	2.57	0.40
1:C:339:GLU:O	1:C:343:LYS:HG3	2.21	0.40
1:A:263:PRO:HA	1:A:264:PRO:HD3	1.94	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:ASP:CB	1:B:186:ASP:O[2_657]	2.19	0.01

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	351/390 (90%)	345 (98%)	5 (1%)	1 (0%)	36	35
1	B	353/390 (90%)	346 (98%)	6 (2%)	1 (0%)	36	35
1	C	353/390 (90%)	347 (98%)	5 (1%)	1 (0%)	36	35
All	All	1057/1170 (90%)	1038 (98%)	16 (2%)	3 (0%)	36	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	SER
1	B	190	SER
1	C	190	SER

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	308/340 (91%)	291 (94%)	17 (6%)	19	17
1	B	310/340 (91%)	300 (97%)	10 (3%)	34	35

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	310/340 (91%)	300 (97%)	10 (3%)	34	35
All	All	928/1020 (91%)	891 (96%)	37 (4%)	28	27

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	22	GLU
1	A	30	ASN
1	A	37	LYS
1	A	38	MET
1	A	50	ARG
1	A	138	SER
1	A	151	ARG
1	A	157	LEU
1	A	176	VAL
1	A	179	MET
1	A	183	ASN
1	A	219	LEU
1	A	265	GLU
1	A	323	GLN
1	A	346	LEU
1	A	367	LYS
1	B	62	LEU
1	B	138	SER
1	B	157	LEU
1	B	176	VAL
1	B	188	ASP
1	B	209	VAL
1	B	219	LEU
1	B	229	SER
1	B	346	LEU
1	B	367	LYS
1	C	8	GLN
1	C	47	LYS
1	C	62	LEU
1	C	157	LEU
1	C	179	MET
1	C	209	VAL
1	C	265	GLU
1	C	323	GLN
1	C	346	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	367	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	49	ASN
1	A	121	HIS
1	A	131	ASN
1	A	146	HIS
1	A	162	ASN
1	A	183	ASN
1	A	191	GLN
1	A	336	GLN
1	B	8	GLN
1	B	121	HIS
1	B	131	ASN
1	B	146	HIS
1	B	162	ASN
1	B	172	GLN
1	B	336	GLN
1	C	8	GLN
1	C	121	HIS
1	C	131	ASN
1	C	146	HIS
1	C	162	ASN
1	C	323	GLN
1	C	336	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 9 are monoatomic - leaving 3 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	RIS	A	1404	2	17,17,17	1.35	1 (5%)	27,27,27	1.37	4 (14%)
3	RIS	C	1600	2	17,17,17	1.25	1 (5%)	27,27,27	1.34	4 (14%)
3	RIS	B	1504	2	17,17,17	1.37	1 (5%)	27,27,27	1.51	5 (18%)

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	RIS	A	1404	2	-	3/23/23/23	0/1/1/1
3	RIS	C	1600	2	-	3/23/23/23	0/1/1/1
3	RIS	B	1504	2	-	2/23/23/23	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1504	RIS	P9-C8	-3.04	1.83	1.85
3	A	1404	RIS	P9-C8	-2.99	1.83	1.85
3	C	1600	RIS	P9-C8	-2.57	1.83	1.85

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1404	RIS	C5-N4-C3	3.16	122.38	116.85
3	B	1504	RIS	C5-N4-C3	3.11	122.30	116.85
3	C	1600	RIS	O15-P14-O16	2.87	116.34	108.24
3	C	1600	RIS	C5-N4-C3	2.70	121.58	116.85
3	B	1504	RIS	C7-C2-C1	-2.69	117.45	121.06

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1504	RIS	C8-C7-C2	2.56	120.41	116.21
3	B	1504	RIS	O10-P9-O12	2.51	115.31	108.24
3	A	1404	RIS	C7-C2-C1	-2.51	117.70	121.06
3	B	1504	RIS	O15-P14-O16	2.41	115.05	108.24
3	C	1600	RIS	C7-C2-C1	-2.38	117.87	121.06
3	A	1404	RIS	C8-C7-C2	2.32	120.03	116.21
3	A	1404	RIS	O15-P14-O16	2.19	114.43	108.24
3	C	1600	RIS	C8-C7-C2	2.14	119.72	116.21

There are no chirality outliers.

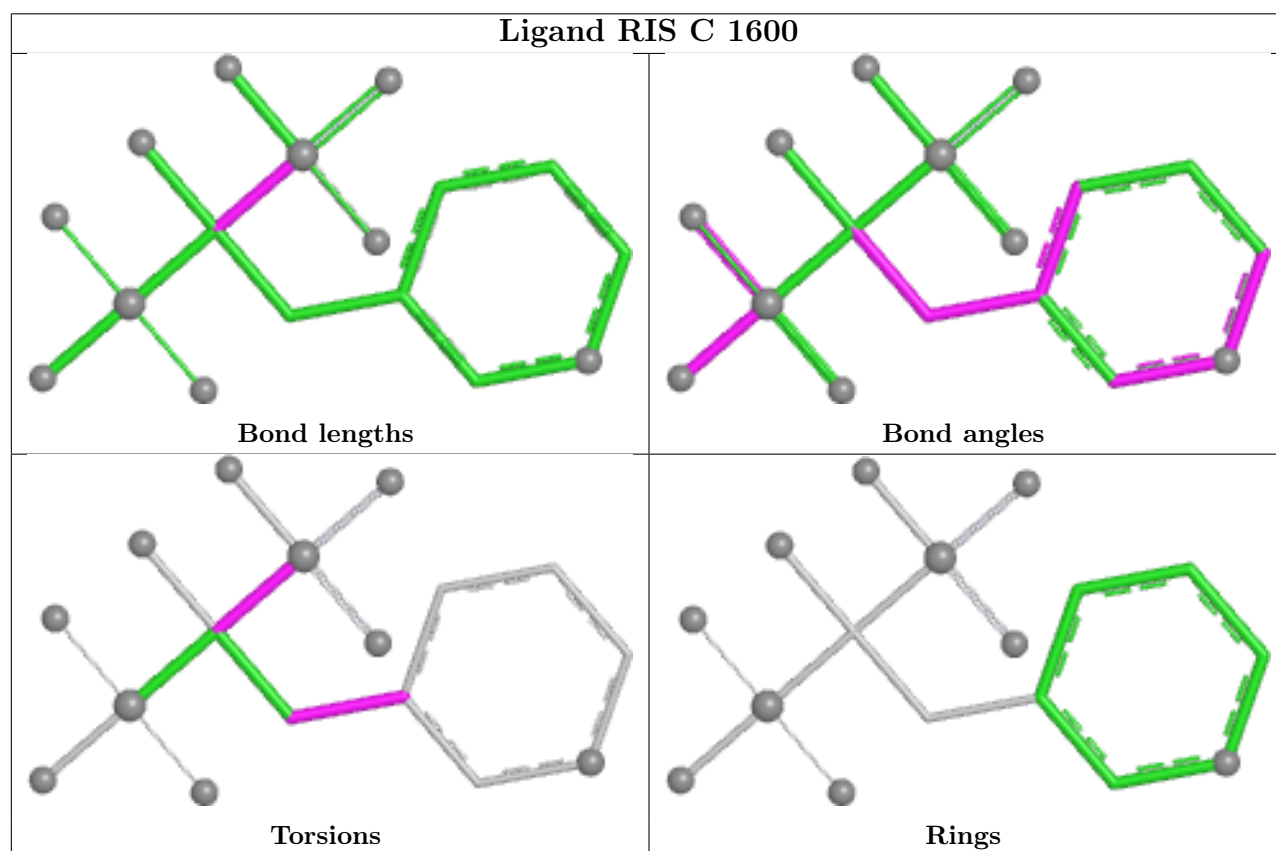
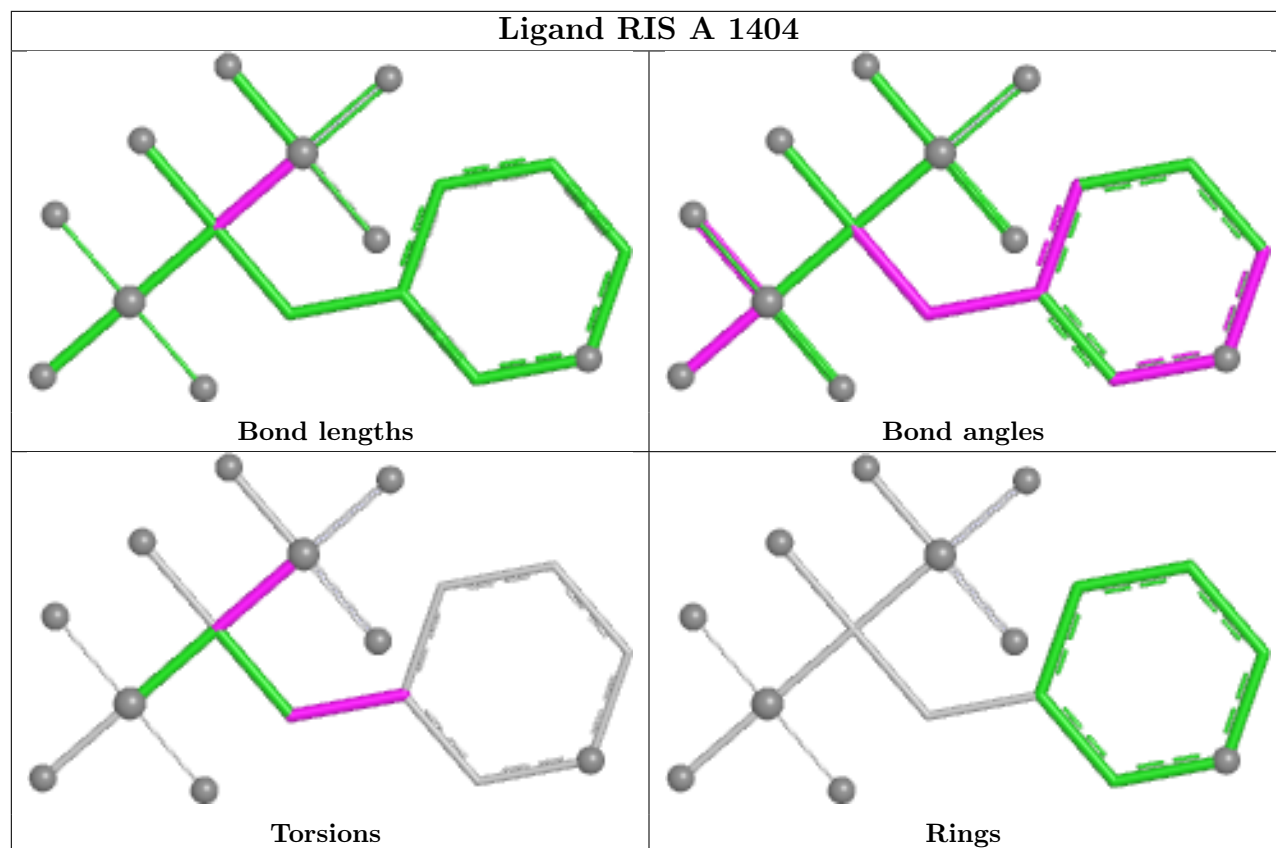
All (8) torsion outliers are listed below:

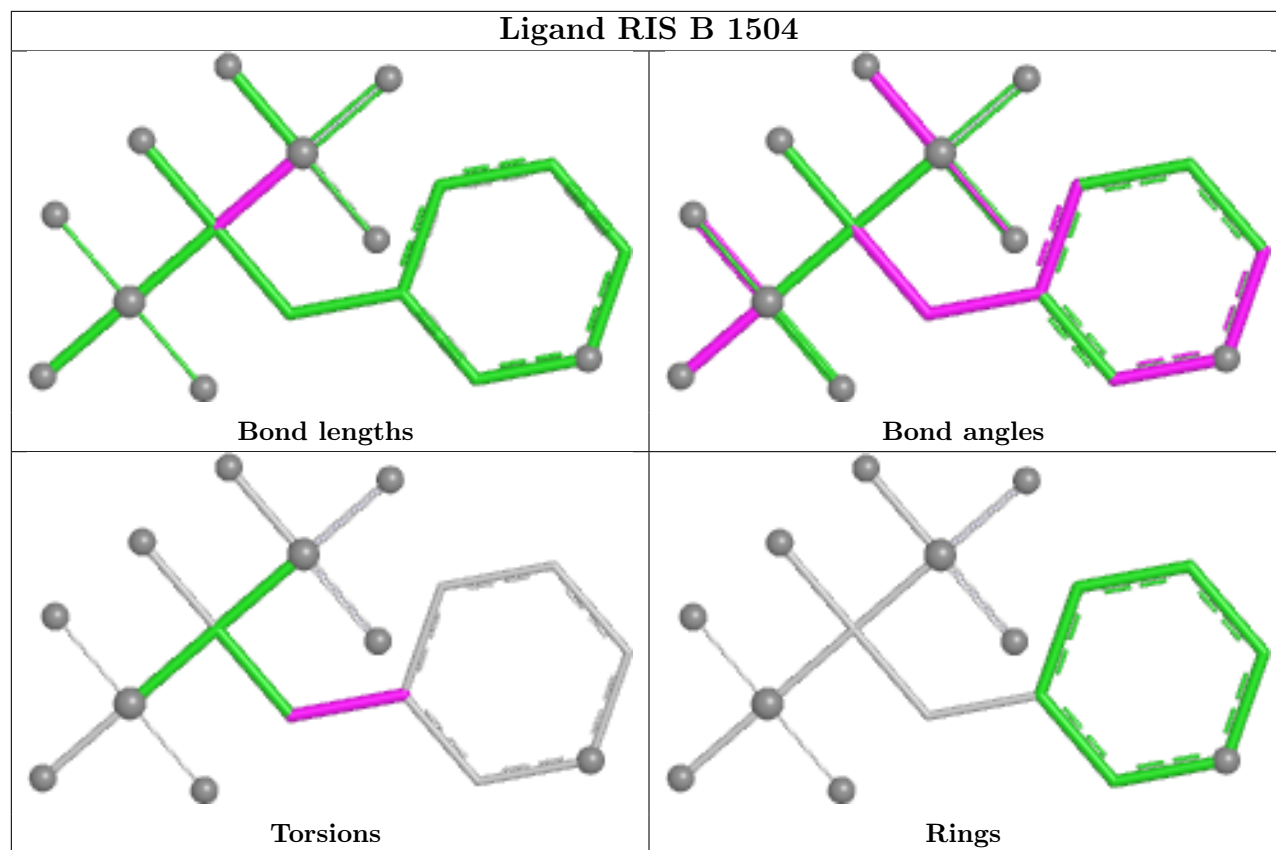
Mol	Chain	Res	Type	Atoms
3	A	1404	RIS	C1-C2-C7-C8
3	A	1404	RIS	C3-C2-C7-C8
3	B	1504	RIS	C1-C2-C7-C8
3	B	1504	RIS	C3-C2-C7-C8
3	C	1600	RIS	C1-C2-C7-C8
3	C	1600	RIS	C3-C2-C7-C8
3	A	1404	RIS	C7-C8-P9-O11
3	C	1600	RIS	C7-C8-P9-O11

There are no ring outliers.

No monomer is involved in short contacts.

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	355/390 (91%)	0.34	27 (7%)	20 18	15, 27, 44, 52	0
1	B	357/390 (91%)	0.28	21 (5%)	28 27	15, 27, 44, 51	0
1	C	357/390 (91%)	0.30	30 (8%)	17 15	15, 27, 44, 55	0
All	All	1069/1170 (91%)	0.31	78 (7%)	21 19	15, 27, 44, 55	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	367	LYS	4.8
1	A	63	SER	4.6
1	A	19	GLU	4.4
1	C	189	VAL	4.3
1	C	193	THR	4.2
1	C	367	LYS	4.1
1	B	193	THR	4.0
1	B	192	PRO	3.9
1	A	191	GLN	3.9
1	B	189	VAL	3.8
1	A	192	PRO	3.7
1	B	1	MET	3.6
1	C	1	MET	3.5
1	A	193	THR	3.5
1	C	192	PRO	3.3
1	A	61	LEU	3.3
1	B	195	THR	3.3
1	C	74	ASP	3.3
1	A	62	LEU	3.2
1	C	190	SER	3.1
1	B	367	LYS	3.1
1	A	189	VAL	3.0
1	B	2	PRO	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	74	ASP	3.0
1	C	180	PHE	3.0
1	C	191	GLN	2.9
1	C	82	LEU	2.9
1	C	265	GLU	2.9
1	A	195	THR	2.8
1	B	194	THR	2.8
1	A	190	SER	2.8
1	A	180	PHE	2.8
1	B	123	ASP	2.8
1	A	194	THR	2.8
1	B	185	LEU	2.8
1	B	191	GLN	2.7
1	C	63	SER	2.7
1	B	183	ASN	2.7
1	C	306	ASP	2.7
1	A	265	GLU	2.6
1	A	183	ASN	2.6
1	A	26	ASP	2.6
1	A	123	ASP	2.6
1	B	180	PHE	2.6
1	C	188	ASP	2.6
1	B	306	ASP	2.5
1	A	15	MET	2.5
1	C	15	MET	2.4
1	C	8	GLN	2.4
1	C	366	GLN	2.4
1	B	188	ASP	2.4
1	A	366	GLN	2.4
1	B	82	LEU	2.3
1	C	61	LEU	2.3
1	C	194	THR	2.3
1	C	195	THR	2.3
1	B	186	ASP	2.3
1	C	75	GLY	2.3
1	C	5	MET	2.3
1	C	22	GLU	2.3
1	A	82	LEU	2.2
1	C	23	LEU	2.2
1	A	319	GLU	2.2
1	C	183	ASN	2.2
1	B	76	ALA	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	346	LEU	2.2
1	A	188	ASP	2.2
1	C	123	ASP	2.2
1	A	22	GLU	2.2
1	A	339	GLU	2.2
1	A	75	GLY	2.1
1	C	339	GLU	2.1
1	B	63	SER	2.1
1	A	60	SER	2.1
1	A	74	ASP	2.1
1	C	197	PHE	2.1
1	C	229	SER	2.0
1	B	332	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 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.

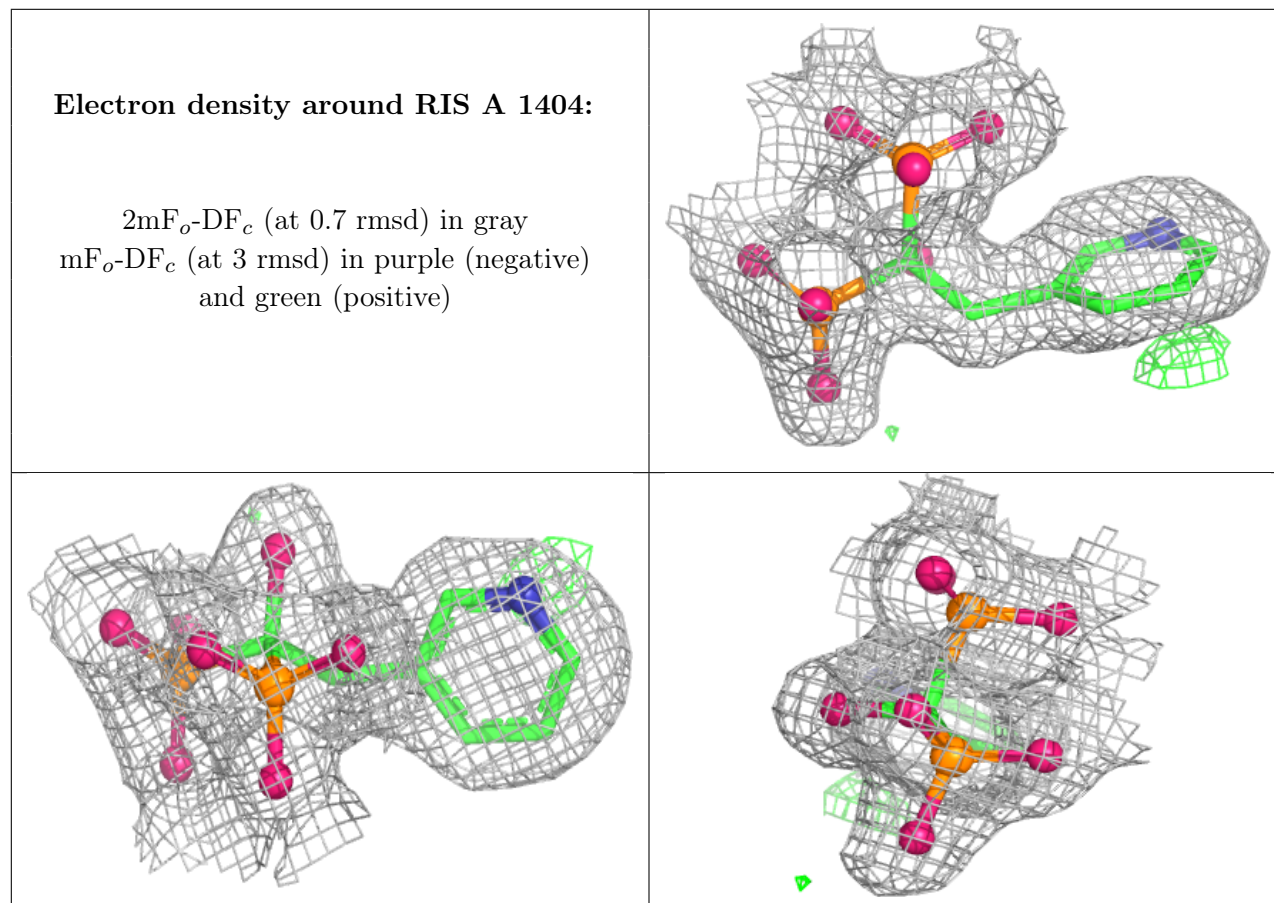
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	B	1503	1/1	0.94	0.05	17,17,17,17	0
2	MG	C	1603	1/1	0.95	0.04	16,16,16,16	0
2	MG	A	1403	1/1	0.96	0.03	16,16,16,16	0
2	MG	C	1601	1/1	0.97	0.03	18,18,18,18	0
2	MG	A	1402	1/1	0.98	0.02	22,22,22,22	0
2	MG	B	1501	1/1	0.98	0.04	17,17,17,17	0
2	MG	C	1602	1/1	0.98	0.03	21,21,21,21	0
2	MG	B	1502	1/1	0.98	0.02	22,22,22,22	0
3	RIS	A	1404	17/17	0.98	0.05	16,18,19,19	0
3	RIS	B	1504	17/17	0.98	0.05	15,17,18,19	0

Continued on next page...

Continued from previous page...

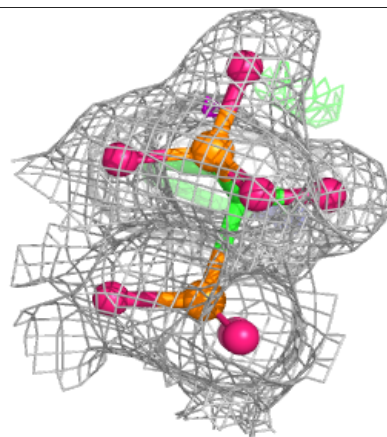
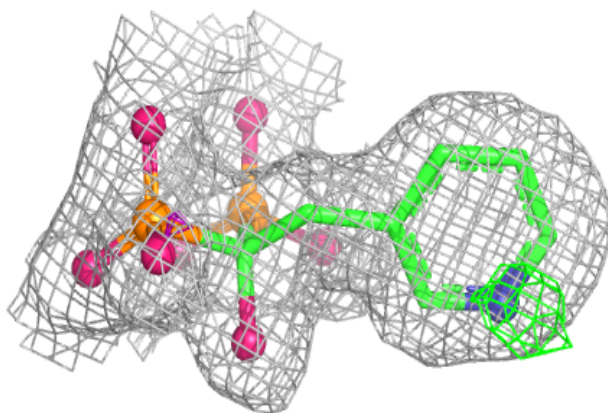
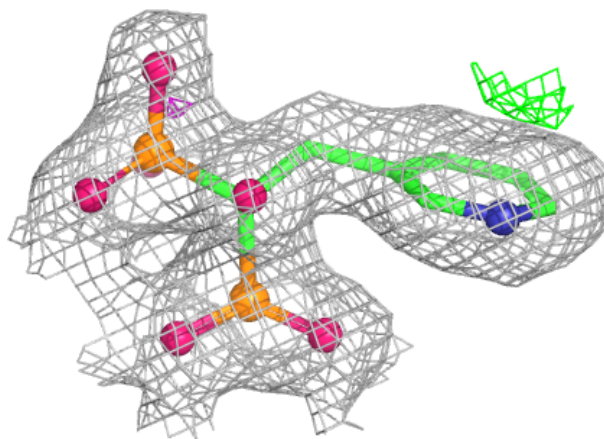
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	RIS	C	1600	17/17	0.98	0.06	17,18,20,20	0
2	MG	A	1401	1/1	0.99	0.02	17,17,17,17	0

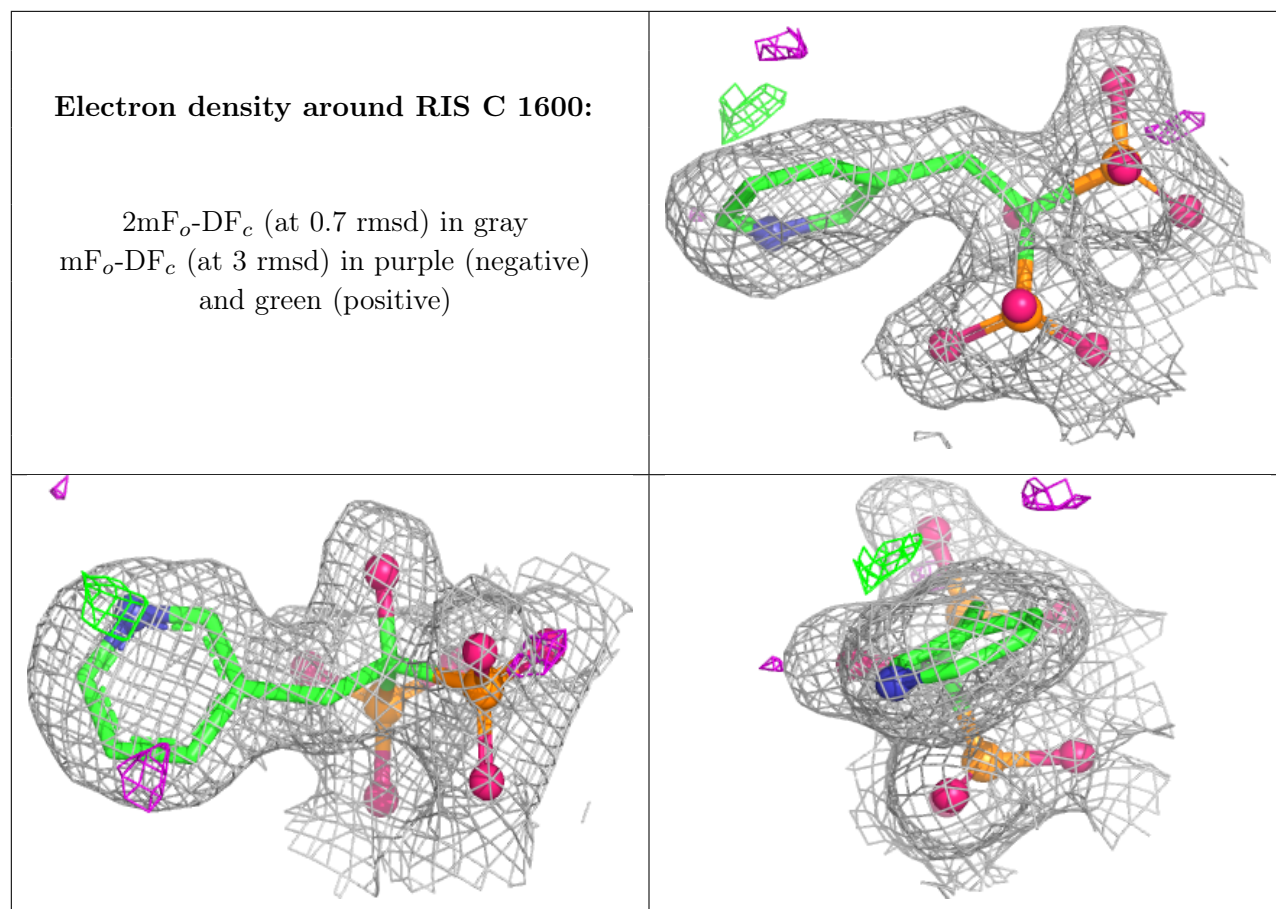
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



Electron density around RIS B 1504:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.