



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:15 PM UTC

PDB ID : 4GTM / pdb_00004gtm
Title : FTase in complex with BMS analogue 11
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Deposited on : 2012-08-28
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

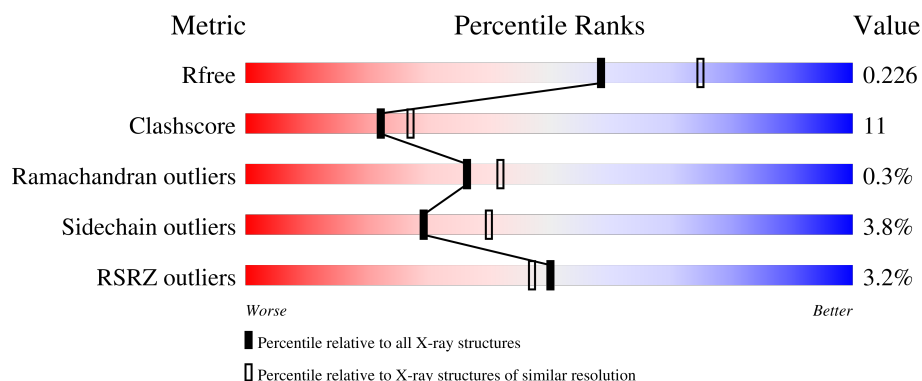
1 Overall quality at a glance

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))
R_{free}	180053	6164 (2.20-2.20)
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	377	
2	B	427	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6589 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 farnesyltransferase/geranylgeranyltransferase type-1 subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	1	0
			2707	1725	476	501	5			

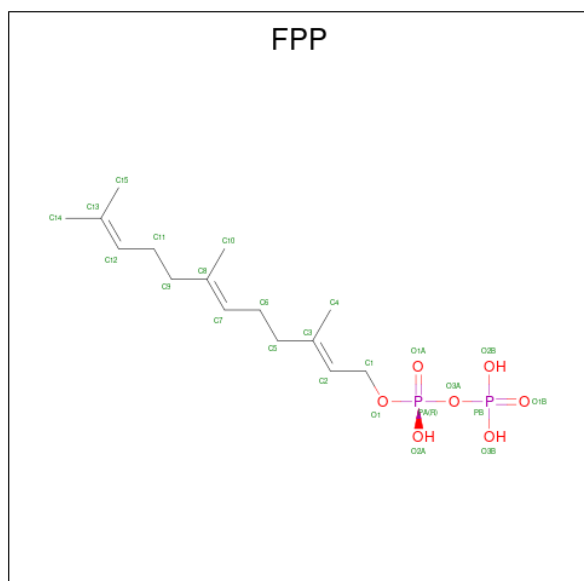
- Molecule 2 is a protein called Protein farnesyltransferase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	405	Total	C	N	O	S	0	3	0
			3200	2046	552	578	24			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

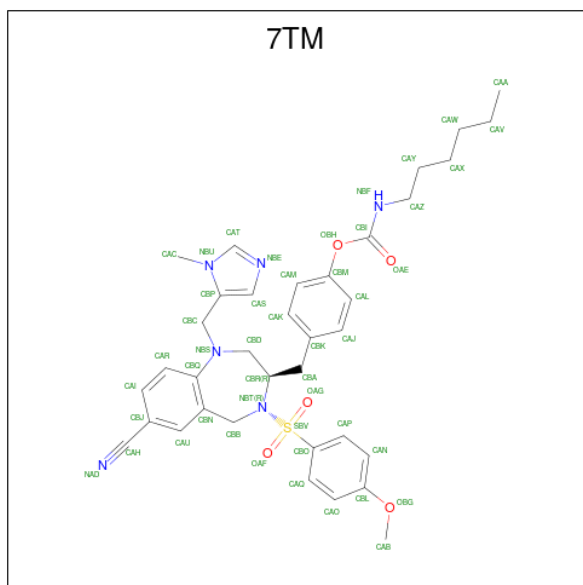
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is FARNESYL DIPHOSPHATE (CCD ID: FPP) (formula: C₁₅H₂₈O₇P₂).



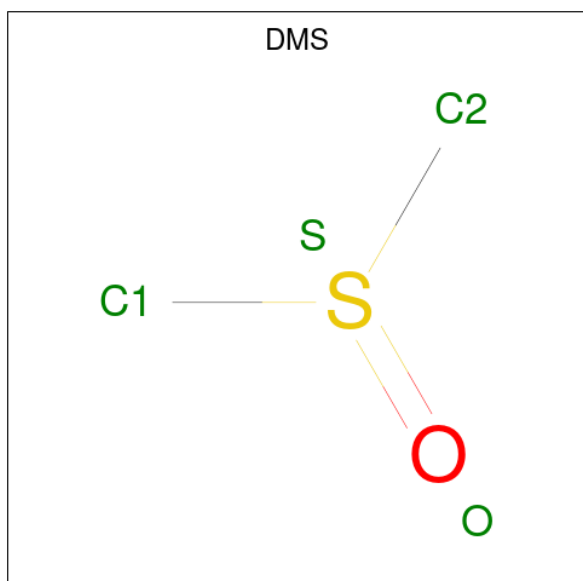
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	O	P	0	0
			24	15	7	2		

- Molecule 5 is 4-({(3R)-7-cyano-4-[(4-methoxyphenyl)sulfonyl]-1-[(1-methyl-1H-imidazol-5-yl)methyl]-2,3,4,5-tetrahydro-1H-1,4-benzodiazepin-3-yl)methyl)phenyl hexylcarbamate (CCD ID: 7TM) (formula: C₃₆H₄₂N₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			48	36	6	5	1		

- Molecule 6 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: $\text{C}_2\text{H}_6\text{OS}$).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	286	Total 286	O 286	0	0
7	B	315	Total 315	O 315	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	173.17Å 173.17Å 70.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.94 – 2.20 19.94 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.94-2.20) 99.8 (19.94-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.87 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.173 , 0.218 0.185 , 0.226	Depositor DCC
R_{free} test set	3048 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6589	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 2.90% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FPP, ZN, 7TM, DMS

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	0.84	0/2776	0.91	1/3768 (0.0%)
2	B	0.81	0/3296	0.92	4/4475 (0.1%)
All	All	0.82	0/6072	0.92	5/8243 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	277	VAL	CB-CA-C	-7.36	102.46	111.88
2	B	409	ILE	CB-CA-C	-6.82	102.98	111.70
2	B	277	VAL	N-CA-CB	6.07	117.25	110.51
2	B	351	LEU	N-CA-C	5.69	117.44	108.96
1	A	185	GLU	N-CA-C	5.26	116.82	111.14

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	2707	0	2632	58	0
2	B	3200	0	3134	73	0
3	B	1	0	0	0	0
4	B	24	0	25	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	48	0	42	20	0
6	B	8	0	12	3	0
7	A	286	0	0	11	0
7	B	315	0	0	11	0
All	All	6589	0	5845	125	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 11.

All (125) 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:156:ILE:HD11	7:A:573:HOH:O	1.18	1.30
1:A:231:VAL:HG21	1:A:266:MET:HE2	1.38	1.05
1:A:328:ASP:O	1:A:329:ASN:HB2	1.63	0.95
1:A:156:ILE:CD1	7:A:573:HOH:O	1.91	0.90
1:A:230:ASP:OD2	2:B:44:THR:HG22	1.71	0.89
1:A:159:ILE:HG12	1:A:168:VAL:HG13	1.56	0.88
2:B:78:HIS:HD1	2:B:349:GLY:H	1.25	0.84
5:B:503:7TM:H40	6:B:505:DMS:H22	1.62	0.81
2:B:99:SER:HA	5:B:503:7TM:H4	1.62	0.81
1:A:296:LEU:HD22	1:A:322:MET:HE1	1.63	0.80
2:B:312:HIS:ND1	7:B:861:HOH:O	2.19	0.76
1:A:221:GLN:NE2	7:A:588:HOH:O	2.19	0.76
2:B:98:ALA:O	5:B:503:7TM:H4	1.86	0.75
1:A:58:LEU:HD22	1:A:95:LYS:HD3	1.69	0.73
2:B:294:LYS:HE3	4:B:502:FPP:O2B	1.88	0.73
2:B:74:GLN:H	2:B:344:GLN:HE22	1.37	0.70
2:B:134:GLN:HE22	2:B:173:ASN:H	1.40	0.69
1:A:159:ILE:CG1	1:A:168:VAL:HG13	2.22	0.69
1:A:256:LEU:O	1:A:259:GLU:N	2.25	0.69
2:B:386:VAL:HG21	2:B:393:VAL:CG1	2.24	0.68
2:B:99:SER:HA	5:B:503:7TM:CAV	2.24	0.67
1:A:266:MET:HE3	1:A:266:MET:HA	1.77	0.66
1:A:114:GLU:OE1	7:A:594:HOH:O	2.12	0.66
2:B:152:PRO:HB3	5:B:503:7TM:H1	1.79	0.65
1:A:90[B]:ILE:HG23	1:A:92:TYR:CZ	2.32	0.65
2:B:152:PRO:HB3	5:B:503:7TM:CAA	2.27	0.64
1:A:312:ILE:HG23	1:A:340:LEU:HD22	1.78	0.64
2:B:149:HIS:HE1	2:B:193[B]:MET:HE3	1.63	0.64
1:A:230:ASP:OD2	2:B:44:THR:CG2	2.46	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:70[A]:ARG:NE	2:B:70[A]:ARG:HA	2.14	0.63
2:B:177:LEU:HD21	2:B:210:VAL:CG1	2.29	0.62
2:B:294:LYS:CE	4:B:502:FPP:O2B	2.47	0.62
2:B:151:ALA:HB1	5:B:503:7TM:OAE	2.00	0.62
2:B:280:ARG:HH11	2:B:289:GLN:HE21	1.47	0.61
1:A:166:TYR:OH	2:B:198:GLU:HG2	2.00	0.61
2:B:280:ARG:HH11	2:B:289:GLN:NE2	1.98	0.61
2:B:34:ARG:HD3	2:B:284:PHE:CE1	2.36	0.61
1:A:207:GLN:HE21	1:A:239:GLN:HE22	1.51	0.59
1:A:90[B]:ILE:HG23	1:A:92:TYR:CE1	2.38	0.58
1:A:90[B]:ILE:HD11	2:B:94:GLU:HG3	1.85	0.58
2:B:423:PHE:C	7:B:833:HOH:O	2.48	0.56
5:B:503:7TM:H39	7:B:818:HOH:O	2.05	0.56
2:B:308:LEU:HD13	2:B:330:PHE:HB3	1.87	0.56
2:B:99:SER:HA	5:B:503:7TM:CAA	2.36	0.55
2:B:386:VAL:HG21	2:B:393:VAL:HG12	1.88	0.55
1:A:207:GLN:HE21	1:A:239:GLN:NE2	2.05	0.55
2:B:149:HIS:CE1	2:B:193[B]:MET:HE3	2.42	0.54
1:A:90[B]:ILE:CG2	1:A:92:TYR:CE1	2.91	0.54
1:A:138:ARG:NH2	7:A:423:HOH:O	2.32	0.54
2:B:152:PRO:HD3	5:B:503:7TM:H5	1.88	0.54
1:A:173:ARG:HD2	2:B:198:GLU:OE2	2.08	0.54
2:B:408:VAL:O	2:B:412:THR:HG23	2.07	0.54
2:B:78:HIS:HD1	2:B:349:GLY:N	2.02	0.53
2:B:312:HIS:CE1	7:B:861:HOH:O	2.59	0.53
2:B:87:ARG:HD3	7:B:710:HOH:O	2.08	0.52
1:A:297:ASN:OD1	1:A:297:ASN:N	2.42	0.52
2:B:102:TRP:CD1	5:B:503:7TM:H2	2.44	0.52
2:B:201:VAL:HG23	2:B:241:GLY:O	2.10	0.52
2:B:99:SER:OG	5:B:503:7TM:H7	2.10	0.52
1:A:345:ALA:O	1:A:350:THR:HG22	2.09	0.51
2:B:118:ILE:HD11	7:B:751:HOH:O	2.09	0.51
2:B:277:VAL:CG1	2:B:307:LEU:HD12	2.40	0.51
1:A:255:VAL:O	1:A:258:ARG:HB3	2.10	0.51
2:B:152:PRO:CG	5:B:503:7TM:H5	2.41	0.51
2:B:275:GLN:NE2	7:B:736:HOH:O	2.42	0.50
2:B:338:TYR:CE2	2:B:343:CYS:SG	3.05	0.50
1:A:166:TYR:OH	2:B:198:GLU:CG	2.59	0.50
2:B:119:PRO:HB2	2:B:122:VAL:HG13	1.93	0.50
1:A:256:LEU:O	1:A:257:GLU:C	2.55	0.49
2:B:70[A]:ARG:HA	2:B:70[A]:ARG:HE	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:ASN:HB3	7:A:493:HOH:O	2.13	0.49
1:A:90[B]:ILE:CD1	2:B:94:GLU:HA	2.42	0.49
2:B:101:PRO:HD2	5:B:503:7TM:H1	1.94	0.48
5:B:503:7TM:CAT	6:B:504:DMS:H12	2.44	0.48
1:A:159:ILE:HG12	1:A:168:VAL:CG1	2.34	0.48
5:B:503:7TM:CAB	6:B:505:DMS:H22	2.40	0.48
1:A:323:LEU:HD22	1:A:367:HIS:HB3	1.96	0.48
2:B:87:ARG:NH2	2:B:125:ASP:OD2	2.42	0.48
2:B:386:VAL:CG2	2:B:393:VAL:HG12	2.43	0.47
1:A:198:LYS:HD3	2:B:294:LYS:HD3	1.95	0.47
2:B:175:GLU:O	2:B:179:GLN:HG3	2.13	0.47
1:A:58:LEU:HD23	1:A:63:TYR:CE1	2.50	0.47
1:A:294:ASN:O	1:A:298:GLN:HG3	2.13	0.47
2:B:102:TRP:CD1	5:B:503:7TM:CAA	2.98	0.47
2:B:202:ARG:HD2	4:B:502:FPP:H152	1.96	0.47
1:A:156:ILE:HG12	1:A:172:ARG:HH12	1.80	0.47
2:B:193[B]:MET:HE2	2:B:202:ARG:HH21	1.81	0.46
2:B:151:ALA:HB3	2:B:152:PRO:CD	2.46	0.46
1:A:121:ARG:HG3	7:A:593:HOH:O	2.16	0.45
2:B:301:SER:O	2:B:305:ALA:HB3	2.15	0.45
1:A:152:MET:O	1:A:156:ILE:HG13	2.17	0.45
1:A:349:ASP:OD2	1:A:352:ARG:HD2	2.17	0.45
1:A:58:LEU:HD22	1:A:95:LYS:CD	2.45	0.44
1:A:334:LEU:O	1:A:334:LEU:HD12	2.17	0.44
2:B:152:PRO:CD	5:B:503:7TM:H5	2.47	0.44
2:B:239:ILE:HB	2:B:252:THR:HA	1.98	0.44
2:B:325:MET:HE2	7:B:706:HOH:O	2.16	0.44
2:B:152:PRO:HD3	5:B:503:7TM:CAV	2.47	0.44
2:B:275:GLN:OE1	7:B:813:HOH:O	2.21	0.44
2:B:177:LEU:HD21	2:B:210:VAL:HG12	1.99	0.43
2:B:202:ARG:CD	4:B:502:FPP:H152	2.49	0.43
2:B:34:ARG:NH1	7:B:869:HOH:O	2.51	0.43
1:A:108:GLN:OE1	7:A:460:HOH:O	2.20	0.43
2:B:56:GLN:NE2	7:B:879:HOH:O	2.47	0.43
2:B:118:ILE:CG2	2:B:122:VAL:HG22	2.48	0.43
2:B:421:PRO:HA	2:B:422:GLY:HA2	1.75	0.43
1:A:240:ARG:HD3	1:A:259:GLU:OE1	2.19	0.43
1:A:328:ASP:O	1:A:329:ASN:CB	2.43	0.42
2:B:277:VAL:HG13	2:B:307:LEU:CD1	2.49	0.42
1:A:312:ILE:HG22	1:A:344:LEU:HD11	2.01	0.42
2:B:375:HIS:HE1	2:B:394:LEU:O	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ARG:NH2	2:B:281:GLN:O	2.52	0.42
1:A:90[B]:ILE:HD13	1:A:92:TYR:CE1	2.55	0.41
1:A:58:LEU:CD2	1:A:95:LYS:HD3	2.44	0.41
1:A:239:GLN:HA	1:A:239:GLN:HE21	1.85	0.41
1:A:280:LEU:HD23	1:A:314:PHE:CE1	2.54	0.41
1:A:173:ARG:CZ	7:A:462:HOH:O	2.68	0.41
2:B:152:PRO:HG3	5:B:503:7TM:H5	2.02	0.41
1:A:97:ARG:NE	7:A:513:HOH:O	2.28	0.41
2:B:308:LEU:HD23	2:B:308:LEU:HA	1.86	0.41
1:A:81:ASN:ND2	7:A:493:HOH:O	2.44	0.41
1:A:250:TYR:C	1:A:252:ASP:H	2.29	0.41
1:A:287:ARG:NE	1:A:287:ARG:N	2.69	0.41
1:A:198:LYS:CD	2:B:294:LYS:HD3	2.51	0.40
1:A:173:ARG:O	1:A:177:GLU:HG2	2.21	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	317/377 (84%)	298 (94%)	17 (5%)	2 (1%)	21	23
2	B	406/427 (95%)	399 (98%)	7 (2%)	0	100	100
All	All	723/804 (90%)	697 (96%)	24 (3%)	2 (0%)	36	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	329	ASN
1	A	256	LEU

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	295/338 (87%)	281 (95%)	14 (5%)	23	31
2	B	344/363 (95%)	334 (97%)	10 (3%)	37	51
All	All	639/701 (91%)	615 (96%)	24 (4%)	29	40

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

Mol	Chain	Res	Type
1	A	71	GLU
1	A	94	GLU
1	A	121	ARG
1	A	145	GLN
1	A	156	ILE
1	A	168	VAL
1	A	251	SER
1	A	252	ASP
1	A	257	GLU
1	A	287	ARG
1	A	297	ASN
1	A	307	SER
1	A	342	GLU
1	A	346	LYS
2	B	20	SER
2	B	34	ARG
2	B	44	THR
2	B	122	VAL
2	B	131	GLU
2	B	263	LYS
2	B	266	ARG
2	B	277	VAL
2	B	280	ARG
2	B	351	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	ASN
1	A	170	HIS
1	A	221	GLN
1	A	239	GLN
1	A	241	HIS
1	A	294	ASN
1	A	329	ASN
1	A	335	ASN
2	B	30	HIS
2	B	56	GLN
2	B	65	ASN
2	B	66	HIS
2	B	134	GLN
2	B	275	GLN
2	B	289	GLN
2	B	332	GLN
2	B	344	GLN
2	B	375	HIS
2	B	414	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is 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	$\# Z > 2$
5	7TM	B	503	3	52,52,52	2.07	8 (15%)	59,72,72	2.58	17 (28%)
4	FPP	B	502	-	22,23,23	2.55	10 (45%)	27,31,31	1.26	2 (7%)
6	DMS	B	505	-	3,3,3	2.67	1 (33%)	3,3,3	0.78	0
6	DMS	B	504	-	3,3,3	2.79	1 (33%)	3,3,3	0.88	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	7TM	B	503	3	-	6/35/51/51	0/4/5/5
4	FPP	B	502	-	-	3/25/25/25	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	503	7TM	OAG-SBV	7.84	1.51	1.43
5	B	503	7TM	OAF-SBV	7.43	1.51	1.43
4	B	502	FPP	PA-O3A	5.64	1.65	1.59
5	B	503	7TM	SBV-NBT	4.93	1.70	1.63
5	B	503	7TM	CBQ-NBS	-4.79	1.34	1.43
6	B	504	DMS	O-S	4.68	1.81	1.50
6	B	505	DMS	O-S	4.45	1.79	1.50
4	B	502	FPP	PB-O1B	4.08	1.63	1.50
5	B	503	7TM	CBC-CBP	3.78	1.55	1.50
4	B	502	FPP	C7-C8	3.59	1.41	1.33
4	B	502	FPP	C11-C12	-3.23	1.40	1.50
4	B	502	FPP	C5-C3	3.23	1.57	1.51
4	B	502	FPP	C2-C3	3.07	1.40	1.33
4	B	502	FPP	C12-C13	3.03	1.41	1.32
5	B	503	7TM	CBP-NBU	-2.99	1.33	1.38
4	B	502	FPP	C1-C2	-2.94	1.40	1.49
4	B	502	FPP	C6-C7	-2.92	1.41	1.50
5	B	503	7TM	CBO-SBV	2.22	1.79	1.76
5	B	503	7TM	CAS-CBP	-2.13	1.34	1.36
4	B	502	FPP	C9-C8	2.05	1.55	1.51

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	503	7TM	NBU-CAT-NBE	-11.05	106.36	112.94
5	B	503	7TM	CBM-OBH-CBI	-7.52	105.62	118.03
5	B	503	7TM	OAG-SBV-OAF	-6.36	109.66	119.59
5	B	503	7TM	CBB-CBN-CAU	-5.67	111.67	119.17
5	B	503	7TM	OBH-CBI-OAE	-4.02	115.89	123.63
5	B	503	7TM	CBB-CBN-CBQ	3.84	128.35	122.45
5	B	503	7TM	CBO-SBV-NBT	3.84	114.06	107.36
5	B	503	7TM	CBD-NBS-CBQ	3.26	124.28	116.96
5	B	503	7TM	CBD-CBR-NBT	2.89	117.91	112.15
5	B	503	7TM	CAB-OBG-CBL	-2.75	111.61	117.50
5	B	503	7TM	CAC-NBU-CBP	-2.74	123.78	126.69
5	B	503	7TM	CAJ-CBK-CAK	2.42	121.83	118.23
5	B	503	7TM	CAS-NBE-CAT	2.40	108.38	105.24
5	B	503	7TM	CAZ-NBF-CBI	2.37	126.00	121.98
5	B	503	7TM	CAO-CAQ-CBO	2.26	121.64	119.44
5	B	503	7TM	CBA-CBR-NBT	-2.21	108.97	112.46
5	B	503	7TM	CBR-NBT-SBV	-2.18	113.92	119.05
4	B	502	FPP	C4-C3-C2	-2.09	118.25	123.63
4	B	502	FPP	C1-C2-C3	2.07	129.59	126.20

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	502	FPP	C1-O1-PA-O2A
4	B	502	FPP	C1-O1-PA-O3A
5	B	503	7TM	NBF-CBI-OBH-CBM
5	B	503	7TM	OAE-CBI-OBH-CBM
5	B	503	7TM	CAN-CBL-OBG-CAB
5	B	503	7TM	CAO-CBL-OBG-CAB
5	B	503	7TM	CAW-CAX-CAY-CAZ
4	B	502	FPP	C1-O1-PA-O1A
5	B	503	7TM	CBB-NBT-SBV-OAG

There are no ring outliers.

4 monomers are involved in 24 short contacts:

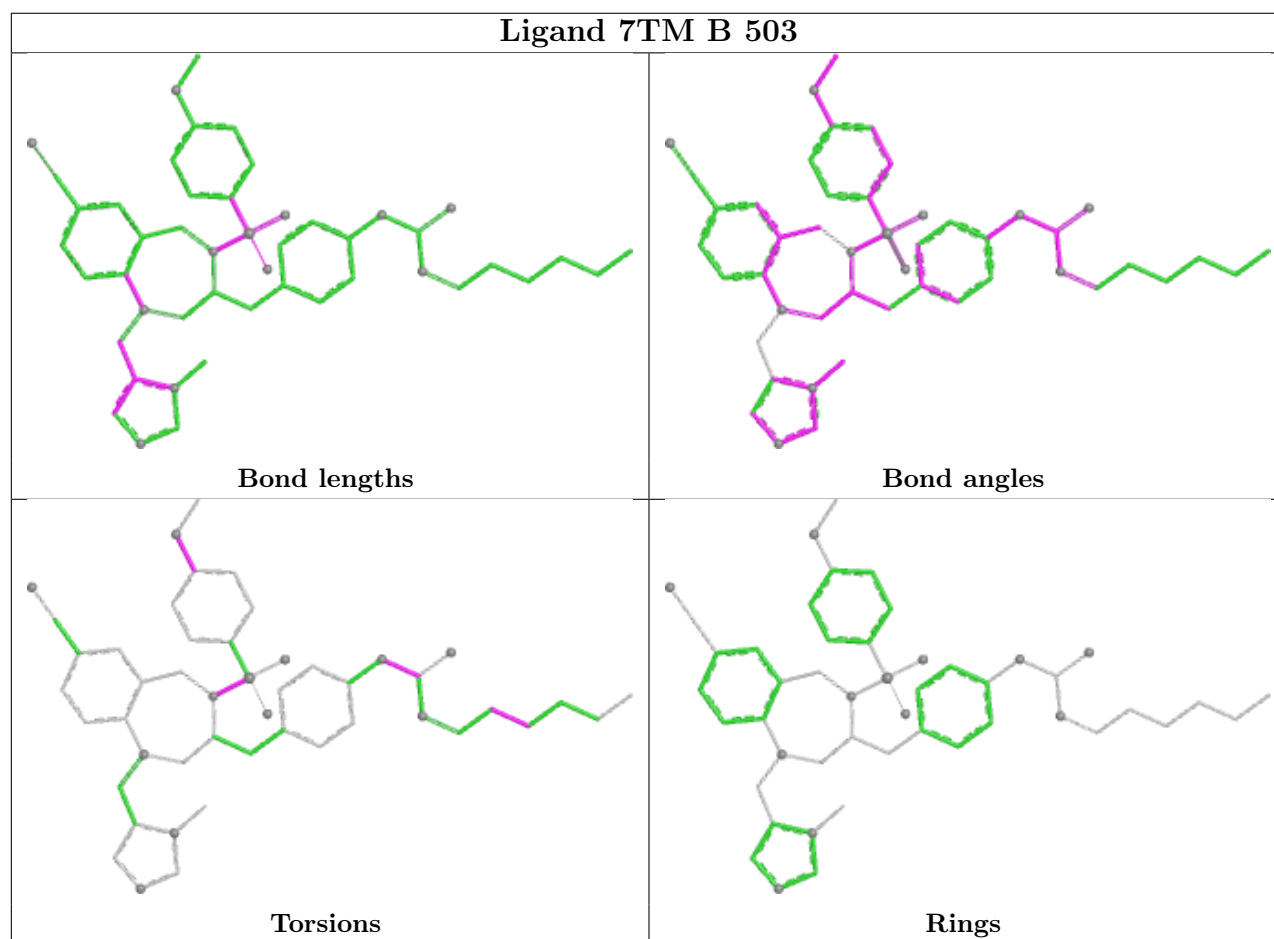
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	503	7TM	20	0
4	B	502	FPP	4	0

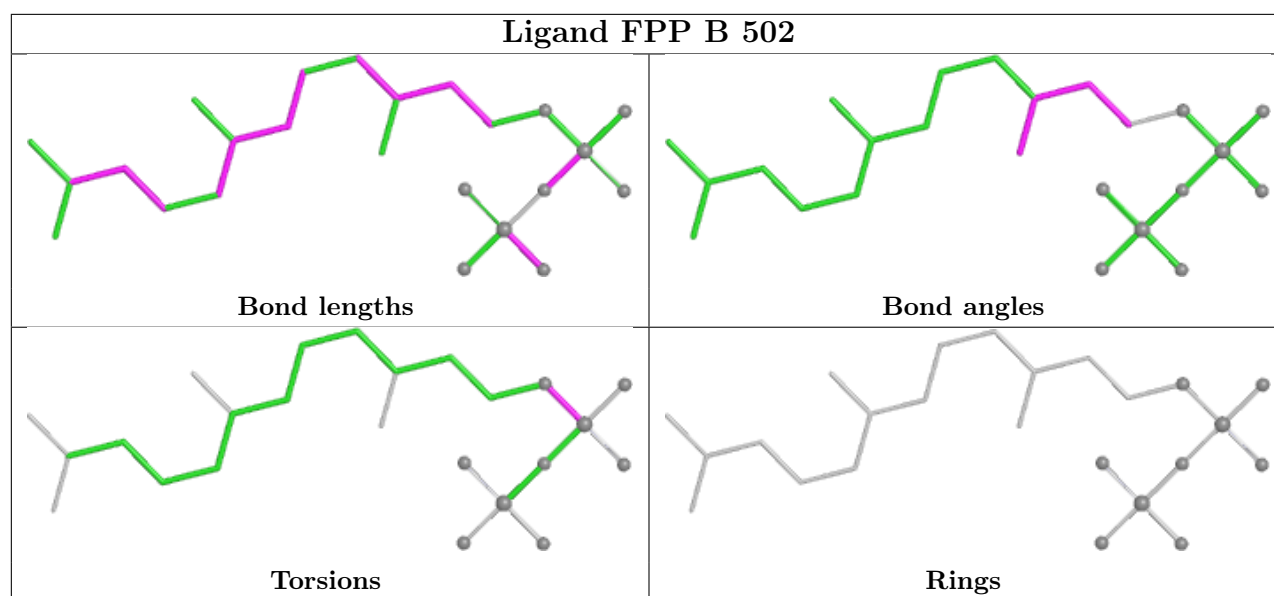
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	505	DMS	2	0
6	B	504	DMS	1	0

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 ⓘ

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	318/377 (84%)	0.06	16 (5%)	34 31	16, 39, 69, 92	1 (0%)
2	B	405/427 (94%)	-0.15	7 (1%)	69 66	19, 35, 59, 83	3 (0%)
All	All	723/804 (89%)	-0.05	23 (3%)	50 47	16, 37, 65, 92	4 (0%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	252	ASP	6.6
1	A	253	ARG	4.4
1	A	371	SER	4.4
2	B	19	TRP	4.0
1	A	254	ALA	3.9
2	B	64	PHE	3.8
1	A	166	TYR	3.1
1	A	256	LEU	3.0
1	A	368	SER	3.0
1	A	55	PHE	2.8
2	B	70[A]	ARG	2.5
1	A	331	GLU	2.5
2	B	121	ILE	2.5
2	B	423	PHE	2.3
2	B	20	SER	2.3
1	A	251	SER	2.2
1	A	54	GLY	2.2
1	A	297	ASN	2.2
2	B	120	GLN	2.2
1	A	287	ARG	2.1
1	A	342	GLU	2.1
1	A	286	ASP	2.1
1	A	369	ARG	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

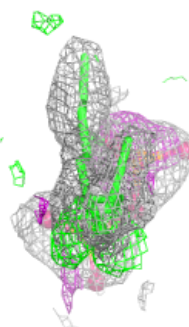
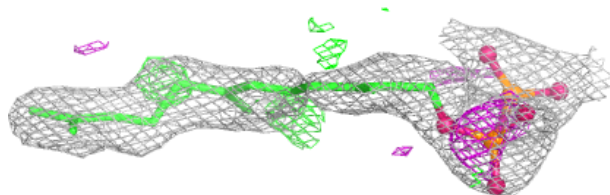
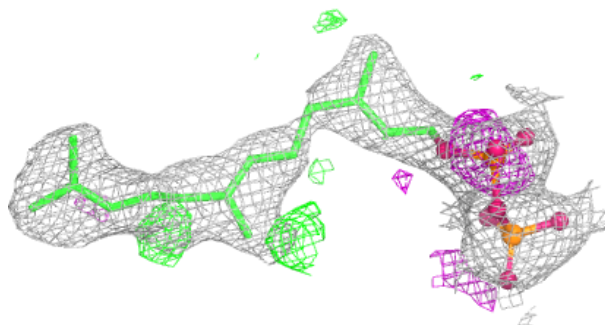
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	FPP	B	502	24/24	0.88	0.15	47,54,60,64	0
5	7TM	B	503	48/48	0.90	0.14	23,51,77,82	0
6	DMS	B	505	4/4	0.91	0.18	65,66,68,71	0
6	DMS	B	504	4/4	0.92	0.20	61,65,65,68	0
3	ZN	B	501	1/1	0.99	0.02	27,27,27,27	0

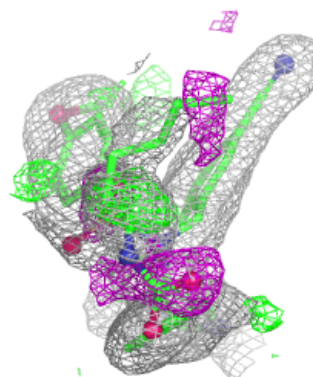
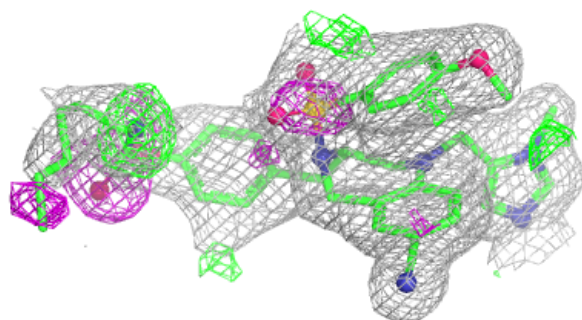
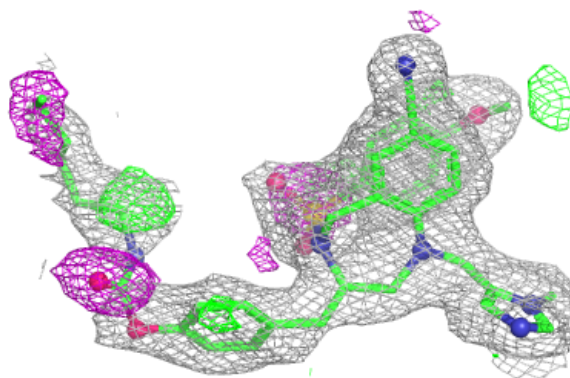
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FPP B 502:

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

**Electron density around 7TM B 503:**

$2mF_o-DF_c$ (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.