



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 07:23 AM EDT

PDB ID : 9R0E / pdb_00009r0e
Title : Structure of the human heterotetrameric cis-prenyltransferase complex harboring NgBR-S249T in complex with magnesium, FsPP and IPP
Authors : Giladi, M.; Haitin, Y.
Deposited on : 2025-04-24
Resolution : 2.82 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

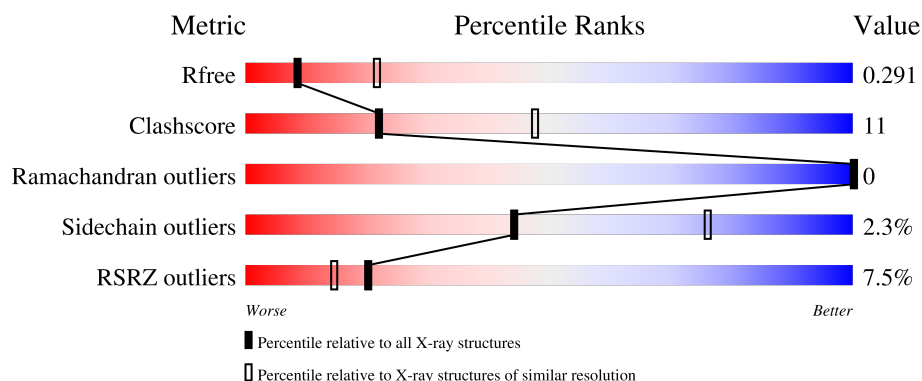
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	340	
2	B	213	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3893 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 Dehydrololichyl diphosphate synthase complex subunit DHDDS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	0
			2475	1568	444	447	16			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	GLY	-	expression tag	UNP Q86SQ9
A	-5	SER	-	expression tag	UNP Q86SQ9
A	-4	GLY	-	expression tag	UNP Q86SQ9
A	-3	SER	-	expression tag	UNP Q86SQ9
A	-2	GLY	-	expression tag	UNP Q86SQ9
A	-1	SER	-	expression tag	UNP Q86SQ9
A	0	GLY	-	expression tag	UNP Q86SQ9

- Molecule 2 is a protein called Dehydrololichyl diphosphate synthase complex subunit NUS1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	192	Total	C	N	O	S	0	0	0
			1379	876	236	259	8			

There are 11 discrepancies between the modelled and reference sequences:

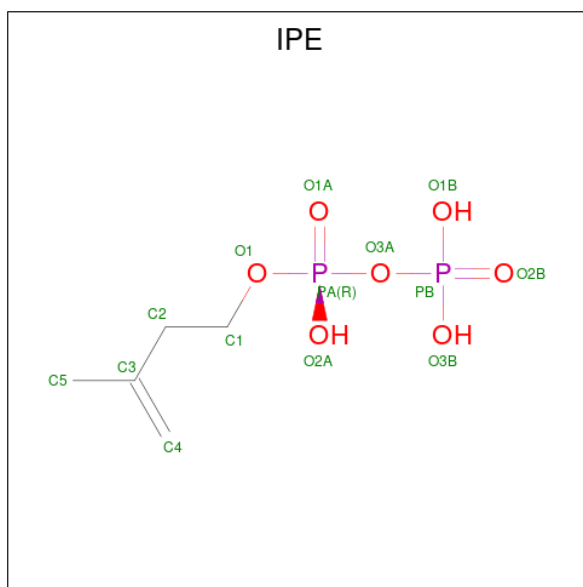
Chain	Residue	Modelled	Actual	Comment	Reference
B	72	GLY	-	expression tag	UNP Q96E22
B	?	-	TYR	deletion	UNP Q96E22
B	?	-	SER	deletion	UNP Q96E22
B	?	-	PRO	deletion	UNP Q96E22
B	?	-	GLU	deletion	UNP Q96E22
B	?	-	PHE	deletion	UNP Q96E22
B	?	-	ALA	deletion	UNP Q96E22
B	?	-	ASN	deletion	UNP Q96E22
B	?	-	SER	deletion	UNP Q96E22

Continued on next page...

Continued from previous page...

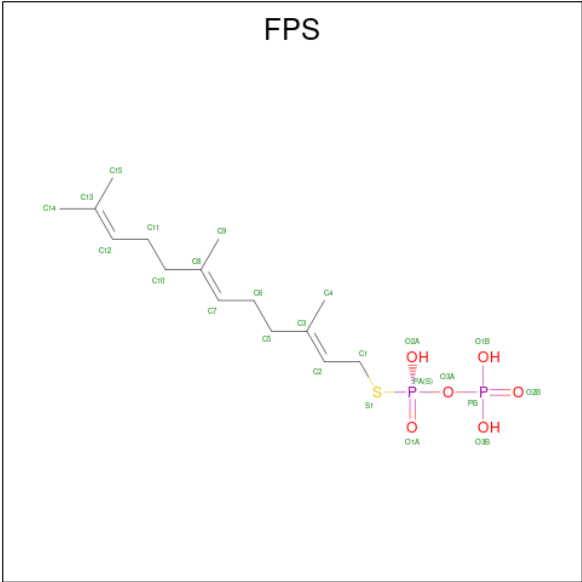
Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	ASN	deletion	UNP Q96E22
B	249	THR	SER	engineered mutation	UNP Q96E22

- Molecule 3 is 3-METHYLBUT-3-ENYL TRIHYDROGEN DIPHOSPHATE (CCD ID: IPE) (formula: $C_5H_{12}O_7P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			14	5	7	2		

- Molecule 4 is S-[(2E,6E)-3,7,11-TRIMETHYLDODECA-2,6,10-TRIENYL] TRIHYDROGEN THIODIPHOSPHATE (CCD ID: FPS) (formula: $C_{15}H_{28}O_6P_2S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	O	P	S	0	0
			24	15	6	2	1		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	184.66Å 184.66Å 112.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.16 – 2.82 46.16 – 2.82	Depositor EDS
% Data completeness (in resolution range)	98.0 (46.16-2.82) 98.0 (46.16-2.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.92 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.240 , 0.290 0.240 , 0.291	Depositor DCC
R_{free} test set	887 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	89.7	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 81.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for -2/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+4/3*l,-1/3*h+1/3*k+1/3*l 0.008 for -h,1/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+1/3*l 0.000 for -1/3*h+1/3*k+4/3*l,-k,2/3*h+1/3*k+1/3*l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3893	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% 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: FPS, MG, IPE

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	0.23	0/2527	0.38	0/3430
2	B	0.23	0/1407	0.38	0/1922
All	All	0.23	0/3934	0.38	0/5352

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2475	0	2264	49	1
2	B	1379	0	1222	33	1
3	A	14	0	9	1	0
4	A	24	0	25	2	0
5	A	1	0	0	0	0
All	All	3893	0	3520	79	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 11.

All (79) 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)
2:B:202:ARG:NH2	2:B:231:SER:HB2	2.09	0.68
1:A:90:VAL:O	1:A:94:MET:HG2	1.95	0.67
2:B:107:GLU:OE1	2:B:269:HIS:ND1	2.25	0.66
2:B:138:GLN:HB2	2:B:140:ILE:HG13	1.78	0.66
1:A:38:ARG:NH2	4:A:402:FPS:O3B	2.31	0.64
2:B:222:ASP:HB3	2:B:225:THR:OG1	2.00	0.62
1:A:162:ILE:HD11	1:A:218:TRP:HB2	1.82	0.61
2:B:222:ASP:OD1	2:B:223:VAL:N	2.36	0.59
1:A:157:THR:OG1	1:A:160:HIS:ND1	2.35	0.58
1:A:260:MET:HE1	2:B:138:GLN:OE1	2.04	0.58
1:A:94:MET:HE2	1:A:128:PRO:HD2	1.87	0.57
1:A:273:ASP:CG	1:A:306:ARG:HH12	2.12	0.56
2:B:202:ARG:HH22	2:B:231:SER:HB2	1.68	0.56
1:A:62:LEU:HD21	1:A:150:LEU:HD13	1.88	0.56
2:B:98:PRO:HG2	2:B:129:ILE:HD12	1.88	0.56
2:B:101:MET:HB2	2:B:129:ILE:HG21	1.89	0.54
2:B:101:MET:CE	2:B:242:LEU:HD22	2.37	0.54
1:A:55:PHE:HB2	4:A:402:FPS:H91	1.90	0.53
1:A:97:ALA:HB3	1:A:131:LEU:HD11	1.88	0.53
2:B:238:PRO:O	2:B:261:THR:OG1	2.26	0.53
2:B:202:ARG:C	2:B:202:ARG:HD2	2.34	0.52
1:A:175:GLN:HB2	1:A:177:LEU:HD23	1.91	0.52
2:B:101:MET:HE1	2:B:242:LEU:HD22	1.91	0.52
1:A:105:MET:HE2	1:A:105:MET:HA	1.92	0.51
1:A:206:THR:O	1:A:207:SER:HB3	2.12	0.50
1:A:94:MET:HE1	1:A:127:LEU:HD22	1.93	0.50
2:B:117:ILE:HG13	2:B:141:PHE:CZ	2.47	0.50
2:B:117:ILE:O	2:B:121:VAL:HG23	2.12	0.49
2:B:130:SER:HB2	2:B:131:TYR:CD2	2.47	0.49
1:A:9:LEU:HD22	1:A:13:GLU:HB3	1.95	0.49
1:A:324:TRP:HA	1:A:324:TRP:CE3	2.48	0.48
1:A:164:ASN:O	1:A:168:GLU:HG2	2.12	0.48
1:A:182:ASP:OD2	1:A:309:ARG:NH2	2.31	0.48
1:A:113:LYS:HE2	1:A:113:LYS:HB2	1.49	0.48
1:A:264:GLU:OE2	1:A:267:ARG:NH1	2.47	0.48
1:A:5:LYS:O	1:A:63:ARG:NE	2.46	0.48
1:A:61:THR:HA	1:A:64:TRP:CE3	2.49	0.47
1:A:79:SER:HB2	3:A:401:IPE:H22	1.97	0.47
2:B:149:MET:HA	2:B:152:ILE:HD12	1.97	0.47
2:B:239:ASP:OD1	2:B:239:ASP:N	2.45	0.47
1:A:31:PHE:HB2	1:A:75:VAL:HG12	1.96	0.47
1:A:185:GLU:HG3	1:A:218:TRP:CZ2	2.50	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:94:LEU:HD11	2:B:278:PHE:CE1	2.51	0.46
2:B:249:THR:HG23	2:B:251:LEU:HG	1.96	0.46
1:A:138:ALA:O	1:A:142:THR:HG23	2.16	0.46
1:A:306:ARG:HA	1:A:309:ARG:NH2	2.31	0.46
2:B:152:ILE:O	2:B:156:GLN:HB2	2.16	0.46
1:A:98:ARG:HD3	1:A:130:ASP:OD2	2.15	0.45
1:A:32:ILE:HB	1:A:205:ARG:HG3	1.98	0.45
1:A:94:MET:HE2	1:A:128:PRO:CD	2.45	0.45
2:B:100:HIS:HB3	2:B:238:PRO:HA	1.98	0.45
2:B:259:ARG:HG3	2:B:260:LEU:N	2.32	0.45
1:A:168:GLU:HB3	1:A:317:LEU:HD11	1.99	0.45
1:A:94:MET:HE1	1:A:127:LEU:CD2	2.47	0.44
1:A:130:ASP:OD1	1:A:130:ASP:N	2.48	0.44
1:A:63:ARG:HA	1:A:63:ARG:HD3	1.85	0.44
2:B:98:PRO:HG3	2:B:240:LEU:HD23	1.99	0.44
2:B:211:VAL:HA	2:B:216:LYS:O	2.17	0.44
2:B:97:LEU:HD22	2:B:129:ILE:HD11	2.00	0.44
2:B:139:GLY:O	2:B:143:ARG:HD3	2.18	0.44
1:A:196:ARG:HH12	1:A:318:GLU:CD	2.27	0.43
1:A:235:TYR:CE1	1:A:239:ASN:HB3	2.52	0.43
2:B:117:ILE:HG13	2:B:141:PHE:HZ	1.82	0.43
1:A:98:ARG:NH1	1:A:130:ASP:OD2	2.52	0.43
1:A:324:TRP:C	1:A:326:ALA:H	2.27	0.42
1:A:213:SER:HA	2:B:260:LEU:HD12	2.02	0.42
1:A:216:LEU:HB3	1:A:219:GLN:HB2	2.01	0.42
2:B:226:LEU:O	2:B:230:LEU:N	2.53	0.42
2:B:142:LYS:HD2	2:B:192:SER:HB3	2.02	0.42
1:A:199:HIS:ND1	1:A:219:GLN:HA	2.35	0.42
2:B:279:SER:HA	2:B:282:ARG:HG3	2.01	0.42
1:A:254:LEU:HD23	1:A:254:LEU:HA	1.84	0.42
1:A:245:LEU:HD23	1:A:245:LEU:HA	1.70	0.41
1:A:116:VAL:O	1:A:142:THR:HB	2.19	0.41
1:A:94:MET:HE3	1:A:131:LEU:HD13	2.03	0.41
1:A:206:THR:HB	1:A:230:VAL:O	2.21	0.41
1:A:220:THR:HA	1:A:223:SER:OG	2.19	0.41
1:A:180:PRO:HG2	2:B:209:GLN:NE2	2.36	0.41
1:A:203:LEU:HG	1:A:220:THR:HG21	2.04	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LYS:NZ	2:B:271:ASN:O[2_665]	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	326/340 (96%)	309 (95%)	17 (5%)	0	100	100
2	B	188/213 (88%)	177 (94%)	11 (6%)	0	100	100
All	All	514/553 (93%)	486 (95%)	28 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	229/295 (78%)	226 (99%)	3 (1%)	61	85
2	B	126/185 (68%)	121 (96%)	5 (4%)	28	61
All	All	355/480 (74%)	347 (98%)	8 (2%)	44	76

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	TRP
1	A	10	SER
1	A	131	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	130	SER
2	B	192	SER
2	B	224	ASP
2	B	235	CYS
2	B	287	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	255	GLN
2	B	205	GLN
2	B	209	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](#)

Of 3 ligands modelled in this entry, 1 is 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	IPE	A	401	5	12,13,13	2.09	2 (16%)	16,19,19	0.98	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FPS	A	402	5	20,23,23	1.06	1 (5%)	24,31,31	1.29	4 (16%)

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	IPE	A	401	5	-	2/13/13/13	-
4	FPS	A	402	5	-	6/19/25/25	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	IPE	PA-O3A	6.23	1.66	1.59
4	A	402	FPS	PA-O1A	2.86	1.52	1.46
3	A	401	IPE	C2-C3	2.15	1.54	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	402	FPS	C6-C7-C8	-2.62	121.62	127.62
4	A	402	FPS	C4-C3-C5	2.58	119.70	115.23
3	A	401	IPE	C5-C3-C2	2.41	121.59	114.96
4	A	402	FPS	C11-C12-C13	-2.22	120.23	127.64
4	A	402	FPS	C15-C13-C14	2.07	119.36	114.59

There are no chirality outliers.

All (8) torsion outliers are listed below:

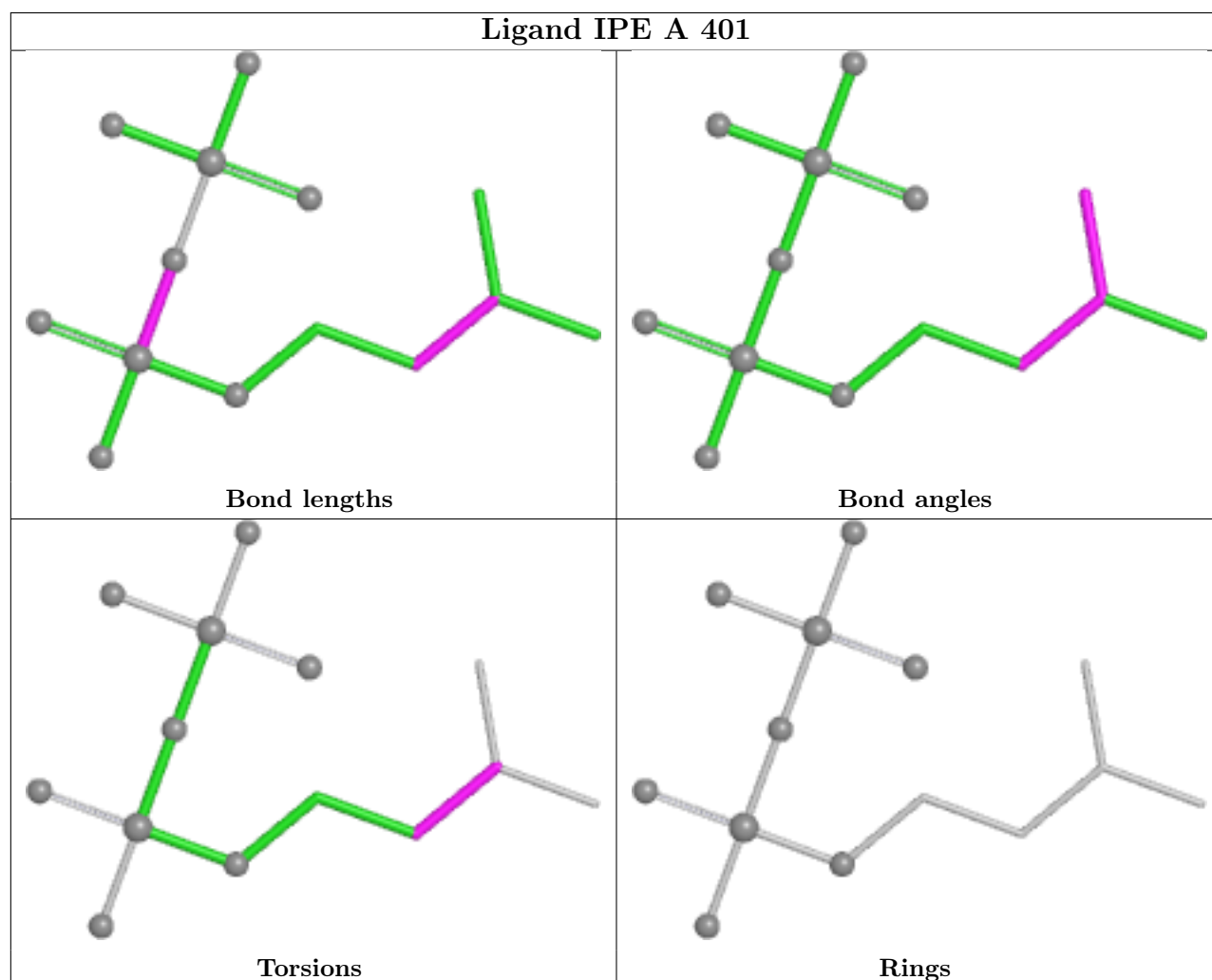
Mol	Chain	Res	Type	Atoms
3	A	401	IPE	C1-C2-C3-C4
3	A	401	IPE	C1-C2-C3-C5
4	A	402	FPS	C3-C5-C6-C7
4	A	402	FPS	C4-C3-C5-C6
4	A	402	FPS	C2-C3-C5-C6
4	A	402	FPS	C8-C10-C11-C12
4	A	402	FPS	PB-O3A-PA-O2A
4	A	402	FPS	PB-O3A-PA-O1A

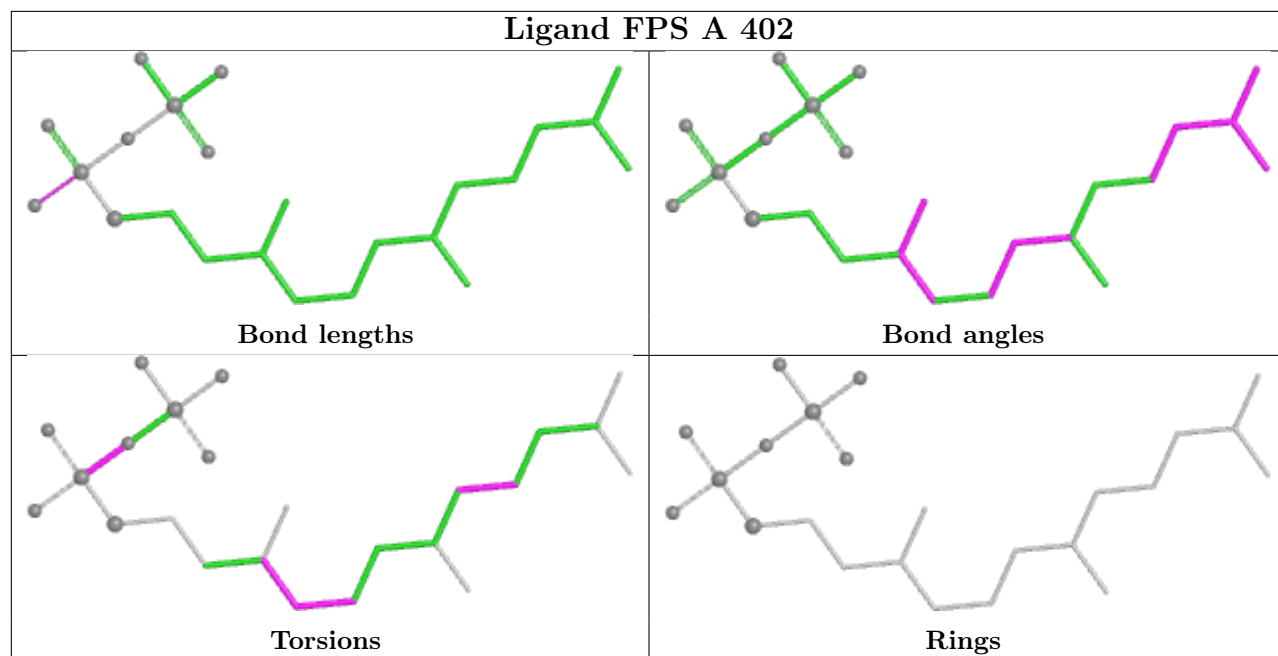
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	IPE	1	0
4	A	402	FPS	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.





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	328/340 (96%)	0.51	19 (5%) 29 21	60, 88, 122, 148	0
2	B	192/213 (90%)	0.79	20 (10%) 11 8	71, 92, 120, 135	0
All	All	520/553 (94%)	0.62	39 (7%) 20 14	60, 90, 122, 148	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	TRP	5.8
1	A	4	ILE	5.2
1	A	292	ALA	4.3
1	A	2	SER	4.2
2	B	181	VAL	4.0
1	A	0	GLY	3.9
1	A	6	GLU	3.6
1	A	326	ALA	3.1
1	A	291	ASP	3.1
1	A	289	SER	3.1
2	B	182	LEU	2.9
2	B	230	LEU	2.9
2	B	219	THR	2.9
1	A	5	LYS	2.8
1	A	324	TRP	2.8
2	B	200	ILE	2.7
2	B	111	GLU	2.7
1	A	327	ARG	2.7
2	B	184	CYS	2.6
2	B	185	HIS	2.6
2	B	231	SER	2.6
2	B	232	SER	2.6
1	A	297	THR	2.5
2	B	220	ASP	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	325	LEU	2.4
2	B	109	GLU	2.3
2	B	82	HIS	2.3
2	B	108	VAL	2.3
1	A	288	ALA	2.2
2	B	157	GLN	2.2
1	A	56	ASN	2.2
1	A	111	LEU	2.2
2	B	183	ASN	2.2
1	A	1	MET	2.2
2	B	233	ASN	2.2
2	B	141	PHE	2.2
2	B	221	LEU	2.1
1	A	294	LEU	2.1
2	B	84	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](#)

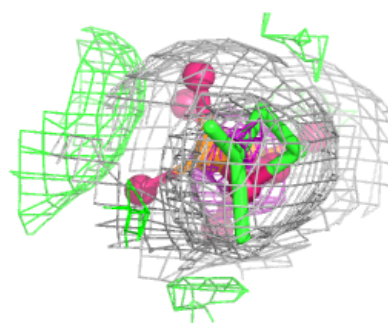
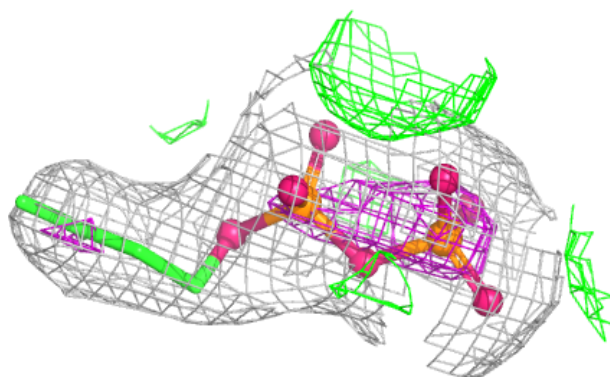
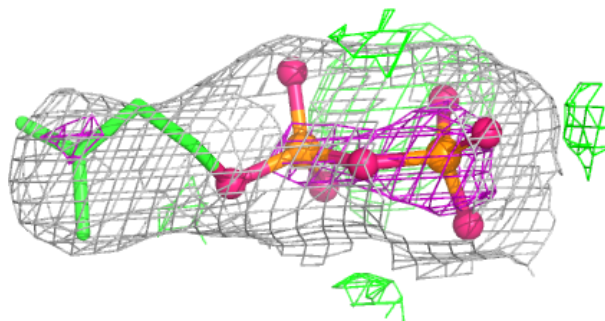
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	IPE	A	401	14/14	0.95	0.10	38,44,46,48	0
4	FPS	A	402	24/24	0.95	0.10	42,48,51,63	0
5	MG	A	403	1/1	0.95	0.10	42,42,42,42	0

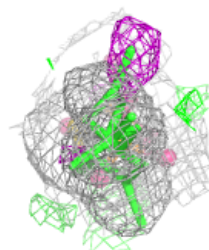
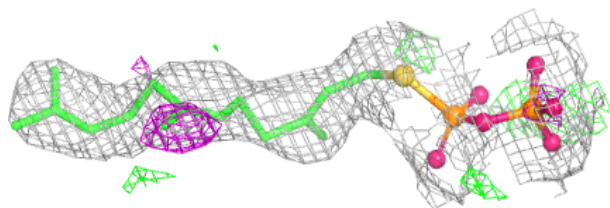
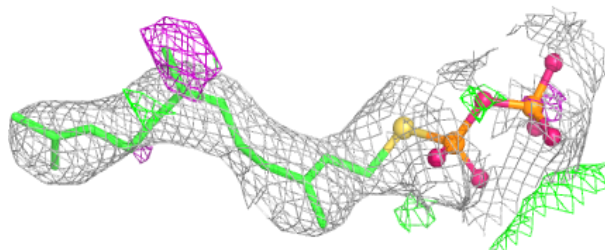
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around IPE A 401:

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

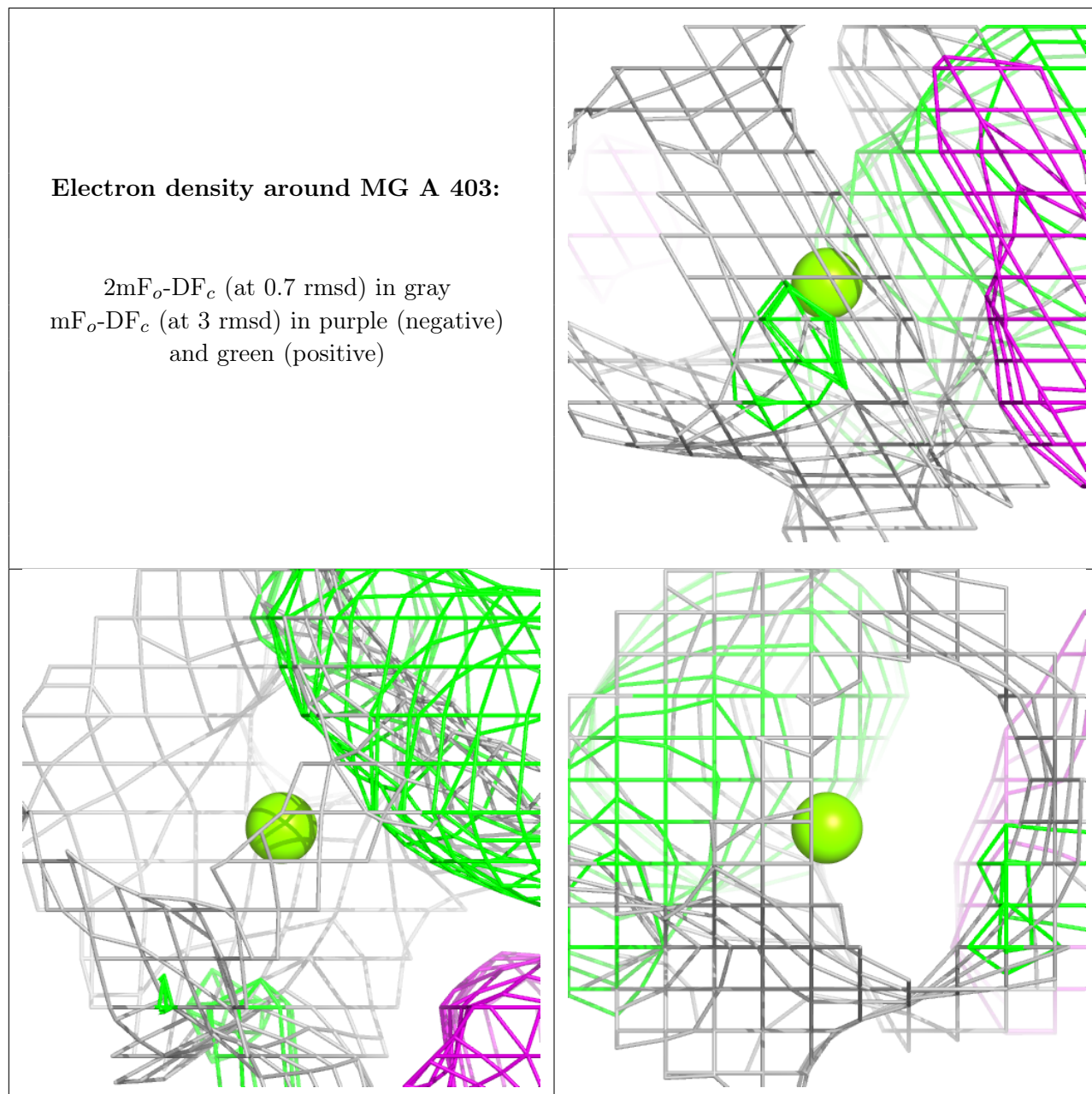
**Electron density around FPS A 402:**

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



Electron density around MG A 403:

$2mF_o-DF_c$ (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.