



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 22, 2025 – 12:10 PM JST

PDB ID : 9V4H / pdb_00009v4h
Title : Prenyltransferase Ord1 Q216A-FSPP
Authors : Oshiro, T.; Uehara, S.; Ito, T.; Tanaka, Y.; Koder, Y.; Matsui, T.
Deposited on : 2025-05-23
Resolution : 2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
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) : 2.46

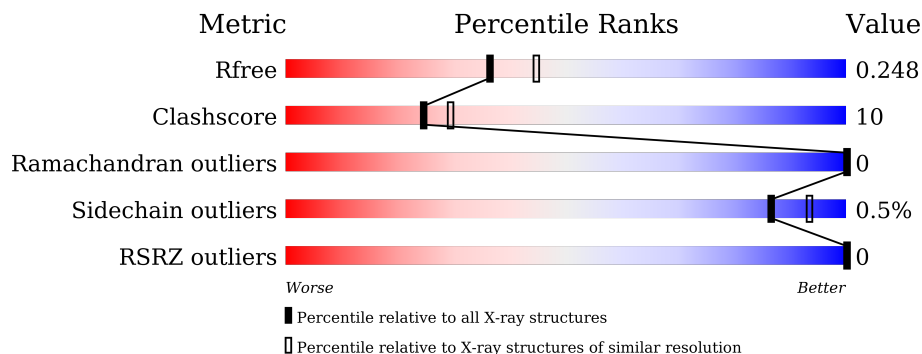
1 Overall quality at a glance

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}	164625	5791 (2.20-2.20)
Clashscore	180529	6634 (2.20-2.20)
Ramachandran outliers	177936	6560 (2.20-2.20)
Sidechain outliers	177891	6561 (2.20-2.20)
RSRZ outliers	164620	5791 (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	337	 79% 12% 9%
1	B	337	 75% 15% 10%
1	C	337	 55% 22% 23%
1	D	337	 55% 20% 25%

2 Entry composition [i](#)

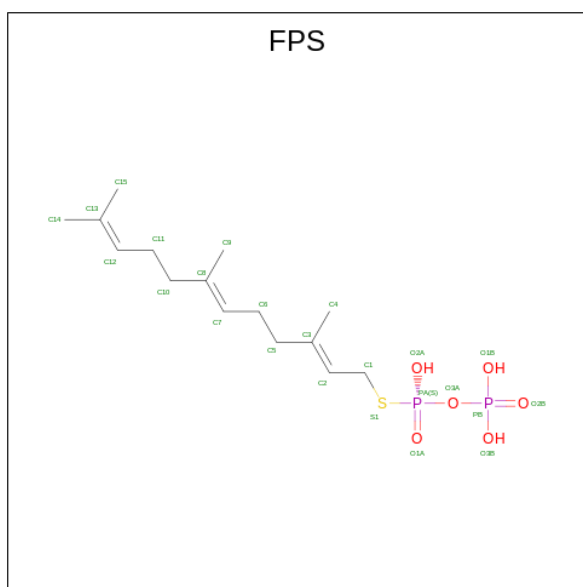
There are 4 unique types of molecules in this entry. The entry contains 8450 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 prenyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	0	0
			2283	1428	413	427	15			
1	B	302	Total	C	N	O	S	0	0	0
			2248	1410	402	422	14			
1	C	259	Total	C	N	O	S	0	0	0
			1926	1207	348	357	14			
1	D	254	Total	C	N	O	S	0	0	0
			1891	1178	345	354	14			

- Molecule 2 is S-[(2E,6E)-3,7,11-TRIMETHYLDODECA-2,6,10-TRIENYL] TRIHYDROGEN THIODIPHOSPHATE (CCD ID: FPS) (formula: C₁₅H₂₈O₆P₂S).



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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		
4	B	18	Total	O	0	0
			18	18		
4	C	10	Total	O	0	0
			10	10		
4	D	13	Total	O	0	0
			13	13		

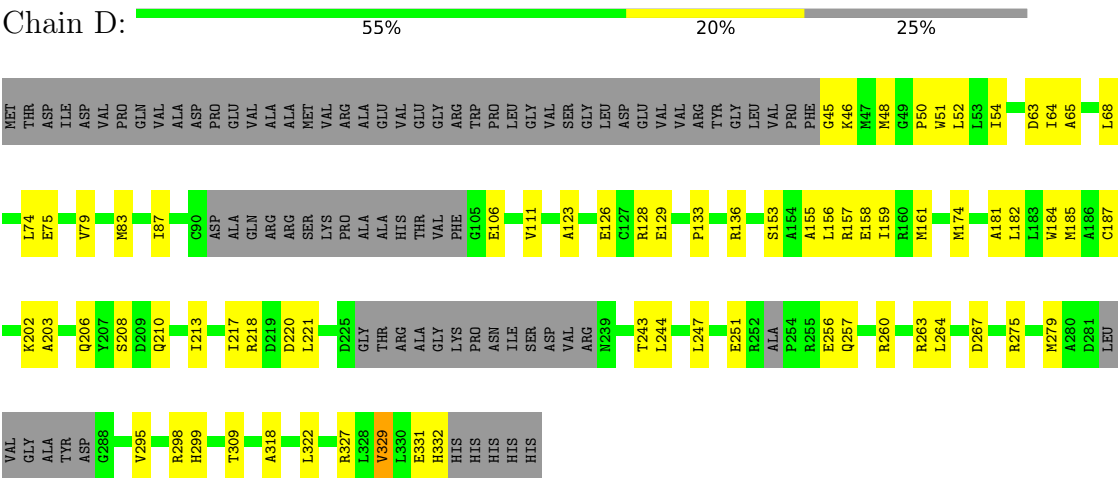
- Molecule 1: prenyltransferase

LYS	PRO	ASP	ASN	ILE	LEU	ASP	VAL	ARG	M239	R255	L259	I261	E262	R263	L264	A270	R275	M279	Q290	E294	V295	R298	H299	V300	P314	T323	V324	P325	GLY	ARG	LEU	VAL	LEU	GLU	HIS	HIS	HIS	HIS	HIS			
THR	THR	ASP	ASP	ILE	ASP	VAL	P77	G78	V9	A10	D11	V14	A15	R19	Y30	M47	P60	W51	L52	R55	A71	V72	A73	C76	K97	E106	L116	G120	Q139	L156	S162	R163	E164	I165	Y171	L175	G225	GLY	THR	ARG	ALA	GLY

[illegible]

V300	Y207	L74	ASP	THR	MET
L307	S208	E75	ASP	THR	
R316	L211		ASP	THR	
L322	L213		ASP	THR	
T323	A214		ASP	THR	
V324	A216		ASP	THR	
P325	L217		ASP	THR	
G326	R218		ASP	THR	
R327	L221		ASP	THR	
L328			ASP	THR	
V329	Y224		ASP	THR	
E331	ASP		ASP	THR	
H333	GLY		ASP	THR	
H334	THR		ASP	THR	
H335	ARG		ASP	THR	
H336	ALA		ASP	THR	
H337	GLY		ASP	THR	
H338	LYS		ASP	THR	
H339	PRO		ASP	THR	
H340	ASN		ASP	THR	
H341	ILE		ASP	THR	
H342	SER		ASP	THR	
H343	ASP		ASP	THR	
H344	VAL		ASP	THR	
H345	ARG		ASP	THR	
H346	N239		ASP	THR	
H347	G240		ASP	THR	
H348	R241		ASP	THR	
H349			ASP	THR	
H350	L244		ASP	THR	
H351	A249		ASP	THR	
H352	R256		ASP	THR	
H353	Q258		ASP	THR	
H354	L259		ASP	THR	
H355	R260		ASP	THR	
H356	L261		ASP	THR	
H357	E262		ASP	THR	
H358	R263		ASP	THR	
H359	L264		ASP	THR	
H360	M279		ASP	THR	
H361			ASP	THR	
H362	V283		ASP	THR	
H363	GLY		ASP	THR	
H364	ALA		ASP	THR	
H365	THR		ASP	THR	
H366	ASP		ASP	THR	
H367	G288		ASP	THR	
H368	A289		ASP	THR	
H369	R293		ASP	THR	
H370	E294		ASP	THR	
H371			ASP	THR	
H372			ASP	THR	
H373			ASP	THR	
H374			ASP	THR	
H375			ASP	THR	
H376			ASP	THR	
H377			ASP	THR	
H378			ASP	THR	
H379			ASP	THR	
H380			ASP	THR	
H381			ASP	THR	
H382			ASP	THR	
H383			ASP	THR	
H384			ASP	THR	
H385			ASP	THR	
H386			ASP	THR	
H387			ASP	THR	
H388			ASP	THR	
H389			ASP	THR	
H390			ASP	THR	
H391			ASP	THR	
H392			ASP	THR	
H393			ASP	THR	
H394			ASP	THR	
H395			ASP	THR	
H396			ASP	THR	
H397			ASP	THR	
H398			ASP	THR	
H399			ASP	THR	
H400			ASP	THR	
H401			ASP	THR	
H402			ASP	THR	
H403			ASP	THR	
H404			ASP	THR	
H405			ASP	THR	
H406			ASP	THR	
H407			ASP	THR	
H408			ASP	THR	
H409			ASP	THR	
H410			ASP	THR	
H411			ASP	THR	
H412			ASP	THR	
H413			ASP	THR	
H414			ASP	THR	
H415			ASP	THR	
H416			ASP	THR	
H417			ASP	THR	
H418			ASP	THR	
H419			ASP	THR	
H420			ASP	THR	
H421			ASP	THR	
H422			ASP	THR	
H423			ASP	THR	
H424			ASP	THR	
H425			ASP	THR	
H426			ASP	THR	
H427			ASP	THR	
H428			ASP	THR	
H429			ASP	THR	
H					

● Molecule 1: prenyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.93Å 71.87Å 84.00Å 86.28° 69.39° 89.99°	Depositor
Resolution (Å)	44.26 – 2.20 44.26 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.0 (44.26-2.20) 97.7 (44.26-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.19.2.4158	Depositor
R, R_{free}	0.236 , 0.253 0.235 , 0.248	Depositor DCC
R_{free} test set	3207 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	40.4	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.477 for h,-k,h-l	Xtriage
Reported twinning fraction	0.500 for h,-k,h-l	Depositor
Outliers	0 of 64136 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8450	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.99% 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, FPS

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	0.12	0/2324	0.31	0/3154
1	B	0.12	0/2288	0.32	0/3107
1	C	0.17	0/1955	0.37	0/2647
1	D	0.13	0/1918	0.36	0/2592
All	All	0.13	0/8485	0.34	0/11500

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 [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	2283	0	2278	26	0
1	B	2248	0	2241	38	0
1	C	1926	0	1943	61	0
1	D	1891	0	1893	48	0
2	C	24	0	25	4	0
2	D	24	0	25	3	0
3	D	1	0	0	0	0
4	A	12	0	0	0	0
4	B	18	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	10	0	0	1	0
4	D	13	0	0	0	0
All	All	8450	0	8405	163	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

The worst 5 of 163 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:178:LYS:HE2	2:C:401:FPS:H2	1.60	0.83
1:C:217:ILE:HD12	1:C:244:LEU:HD22	1.61	0.82
1:C:187:CYS:SG	1:C:208:SER:OG	2.40	0.78
1:B:135:GLU:O	1:B:139:GLN:HG3	1.91	0.70
1:B:158:GLU:HB2	1:B:174:MET:HE2	1.72	0.70

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	302/337 (90%)	300 (99%)	2 (1%)	0	100	100
1	B	296/337 (88%)	291 (98%)	5 (2%)	0	100	100
1	C	251/337 (74%)	247 (98%)	4 (2%)	0	100	100
1	D	244/337 (72%)	241 (99%)	3 (1%)	0	100	100
All	All	1093/1348 (81%)	1079 (99%)	14 (1%)	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	227/254 (89%)	227 (100%)	0	100	100
1	B	224/254 (88%)	224 (100%)	0	100	100
1	C	192/254 (76%)	190 (99%)	2 (1%)	73	84
1	D	188/254 (74%)	186 (99%)	2 (1%)	70	82
All	All	831/1016 (82%)	827 (100%)	4 (0%)	86	93

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	294	GLU
1	C	329	VAL
1	D	329	VAL
1	D	332	HIS

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

Mol	Chain	Res	Type
1	B	297	HIS
1	C	139	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 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	$\# Z > 2$
2	FPS	C	401	-	19,23,23	0.82	1 (5%)	23,31,31	0.64	1 (4%)
2	FPS	D	402	3	19,23,23	0.83	1 (5%)	23,31,31	0.75	1 (4%)

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	FPS	C	401	-	-	2/19/25/25	-
2	FPS	D	402	3	-	0/19/25/25	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	FPS	PA-O2A	-2.76	1.49	1.56
2	D	402	FPS	PA-O2A	-2.75	1.49	1.56

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	402	FPS	O2A-PA-O1A	2.24	115.26	109.82
2	C	401	FPS	O2A-PA-O1A	2.14	115.00	109.82

There are no chirality outliers.

All (2) torsion outliers are listed below:

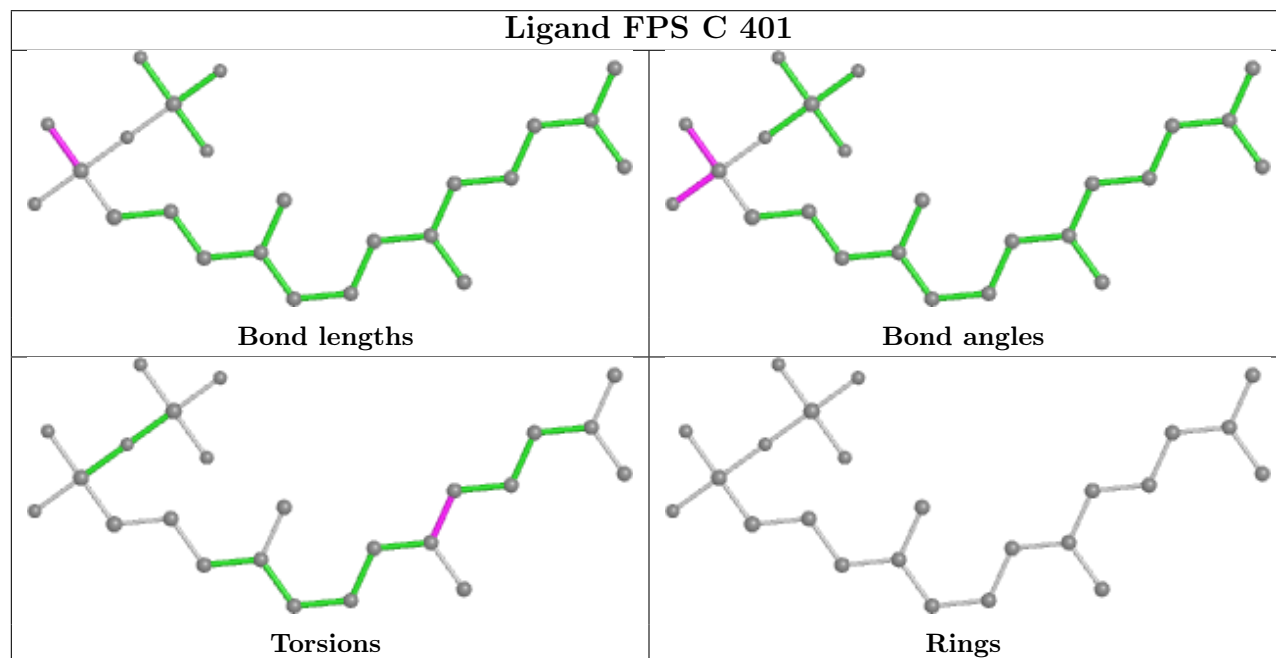
Mol	Chain	Res	Type	Atoms
2	C	401	FPS	C11-C10-C8-C9
2	C	401	FPS	C11-C10-C8-C7

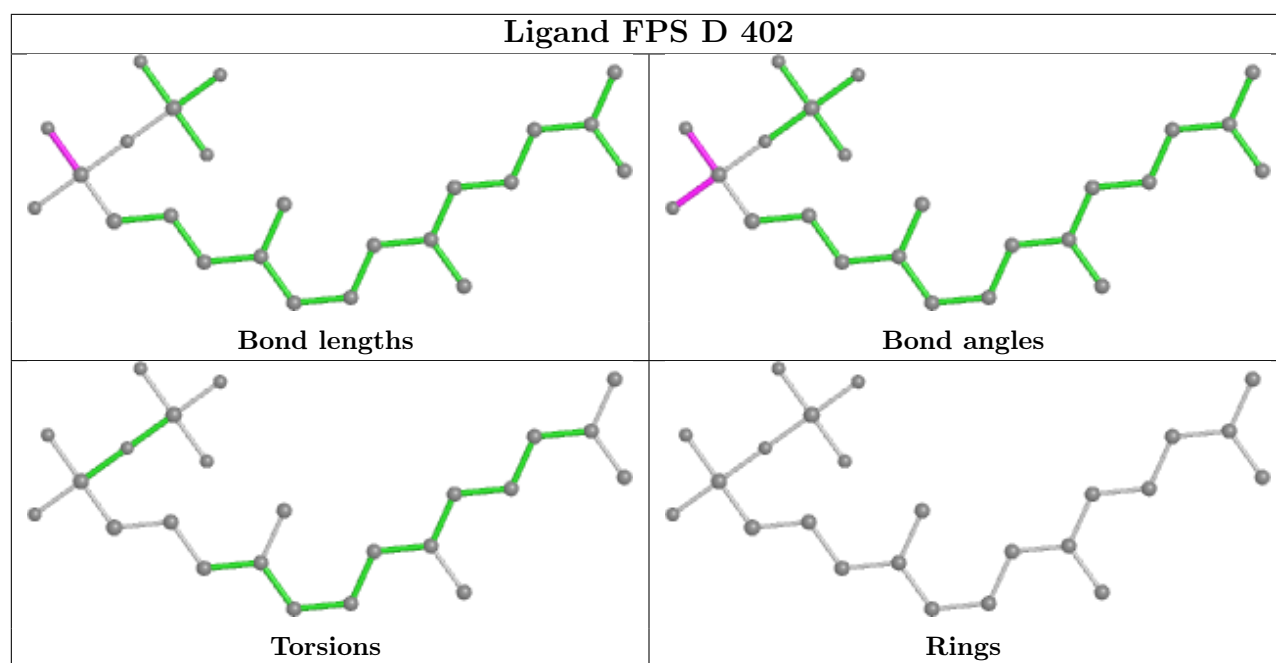
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	FPS	4	0
2	D	402	FPS	3	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 [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	306/337 (90%)	-1.34	0 100 100	32, 41, 87, 123	0
1	B	302/337 (89%)	-1.33	0 100 100	31, 41, 86, 110	0
1	C	259/337 (76%)	-1.23	0 100 100	32, 49, 92, 129	0
1	D	254/337 (75%)	-1.22	0 100 100	31, 52, 89, 113	0
All	All	1121/1348 (83%)	-1.29	0 100 100	31, 46, 88, 129	0

There are no RSRZ outliers to report.

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](#)

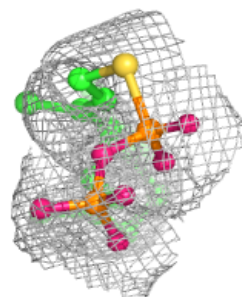
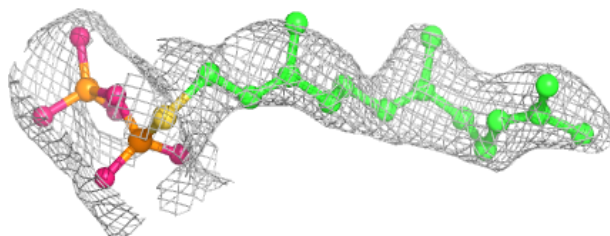
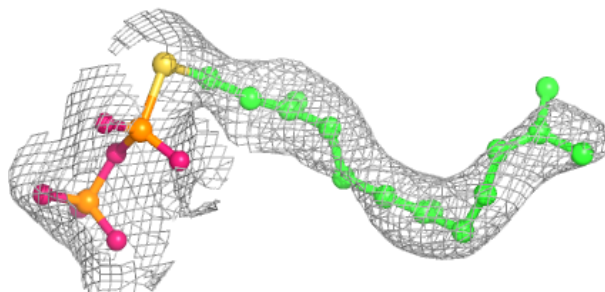
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FPS	C	401	24/24	0.99	0.04	40,51,84,89	0
2	FPS	D	402	24/24	0.99	0.05	43,52,85,103	0
3	MG	D	401	1/1	1.00	0.04	46,46,46,46	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.

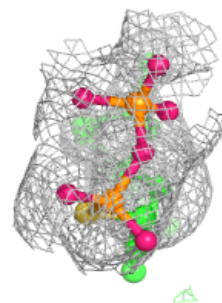
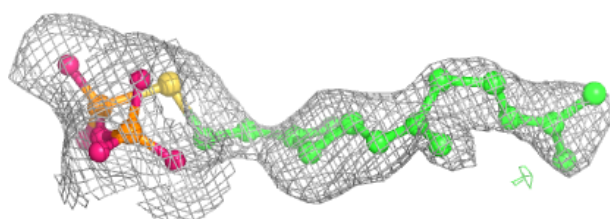
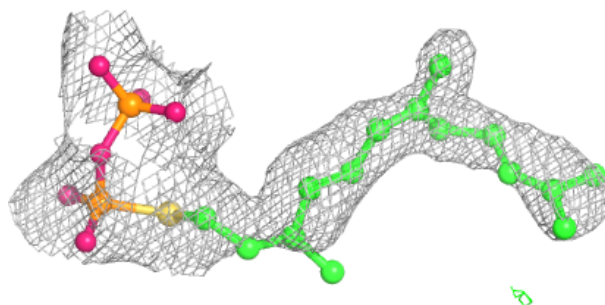
Electron density around FPS C 401:

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



Electron density around FPS D 402:

$2mF_o-DF_c$ (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.