



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:21 PM UTC

PDB ID : 3E32 / pdb_00003e32
Title : Protein farnesyltransferase complexed with FPP and ethylenediamine scaffold inhibitor 2
Authors : Hast, M.A.; Beese, L.S.
Deposited on : 2008-08-06
Resolution : 2.45 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

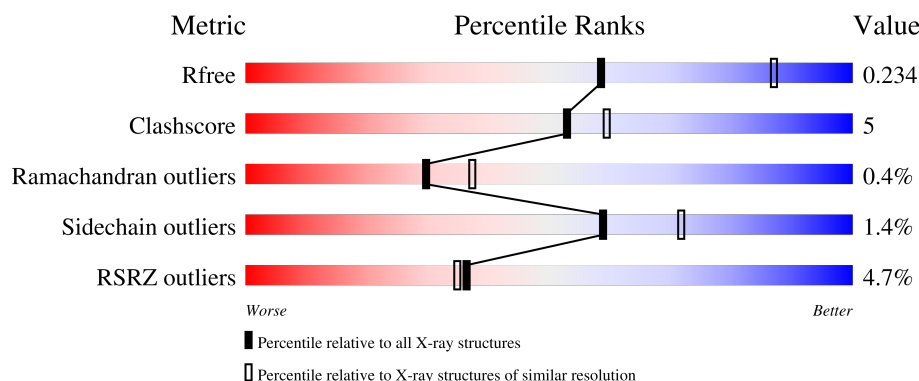
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

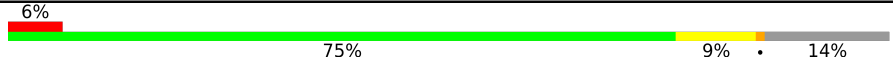
The reported resolution of this entry is 2.45 Å.

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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	437	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6343 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	323	Total	C	N	O	S	0	0	0
			2745	1746	481	513	5			

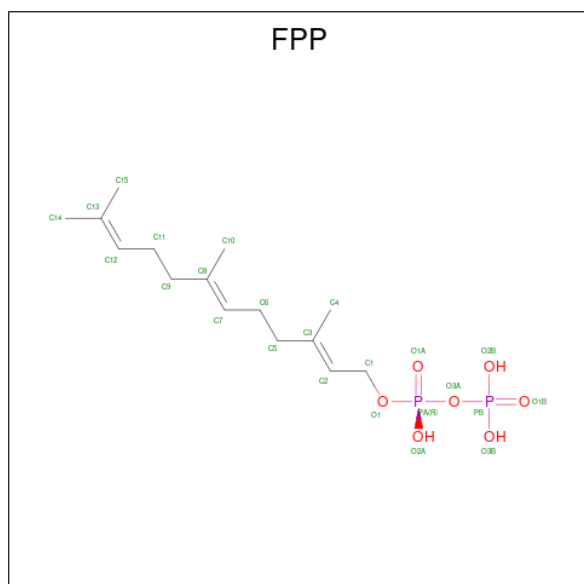
- Molecule 2 is a protein called Protein farnesyltransferase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	407	Total	C	N	O	S	0	0	0
			3203	2050	550	580	23			

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

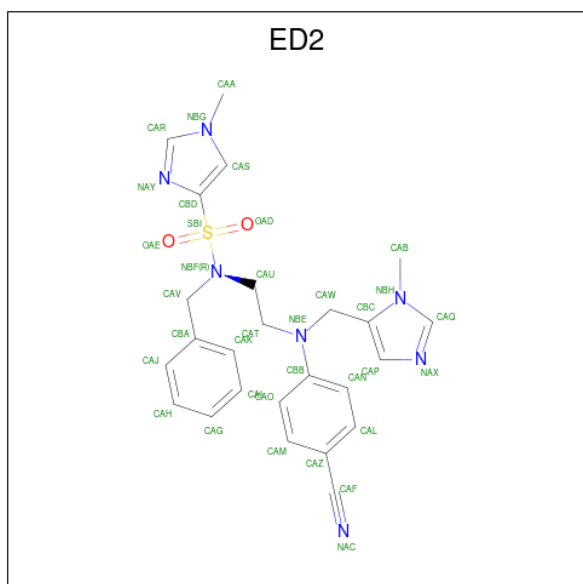
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		

- 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 N-benzyl-N-(2-{(4-cyanophenyl)[(1-methyl-1H-imidazol-5-yl)methyl]amino}ethyl)-1-methyl-1H-imidazole-4-sulfonamide (CCD ID: ED2) (formula: $C_{25}H_{27}N_7O_2S$).

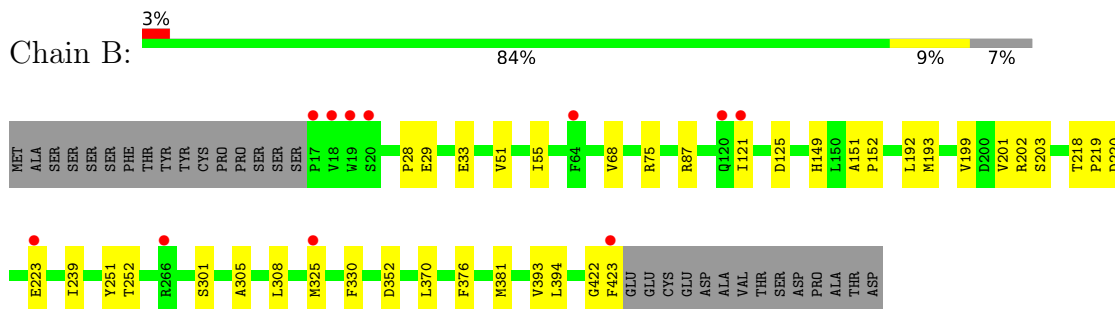


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			35	25	7	2	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	154	Total O 154 154	0	0
6	B	181	Total O 181 181	0	0

- Molecule 1: Protein farnesyltransferase/geranylgeranyltransferase type-1 subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	173.38Å 173.38Å 70.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.50 – 2.45 37.50 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.6 (37.50-2.45) 99.5 (37.50-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.4.0077	Depositor
R, R_{free}	0.208 , 0.237 0.207 , 0.234	Depositor DCC
R_{free} test set	2220 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6343	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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: ED2, FPP, ZN

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	0.39	1/2812 (0.0%)	0.71	4/3817 (0.1%)
2	B	0.37	0/3292	0.69	0/4473
All	All	0.38	1/6104 (0.0%)	0.70	4/8290 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	152	MET	C-O	-5.58	1.17	1.24

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	304	PRO	CA-N-CD	-8.27	100.42	112.00
1	A	85	SER	CA-C-N	6.41	126.38	119.78
1	A	85	SER	C-N-CA	6.41	126.38	119.78
1	A	185	GLU	N-CA-C	5.02	116.56	111.14

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	2745	0	2666	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3203	0	3130	25	0
3	B	1	0	0	0	0
4	B	24	0	25	3	0
5	B	35	0	25	6	0
6	A	154	0	0	0	0
6	B	181	0	0	0	0
All	All	6343	0	5846	59	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 5.

All (59) 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)
5:B:1003:ED2:CAS	5:B:1003:ED2:HAUA	1.45	1.07
5:B:1003:ED2:CAS	5:B:1003:ED2:CAU	2.34	1.05
5:B:1003:ED2:HAUA	5:B:1003:ED2:HAS	1.47	0.95
1:A:152:MET:O	1:A:156:ILE:HG23	1.75	0.86
1:A:152:MET:HE2	1:A:175:LEU:HB3	1.62	0.79
1:A:152:MET:HE2	1:A:175:LEU:CB	2.20	0.71
1:A:152:MET:O	1:A:156:ILE:CG2	2.44	0.64
1:A:262:TYR:O	1:A:265:GLU:HB2	1.97	0.64
1:A:346:LYS:HA	1:A:346:LYS:HE2	1.85	0.59
1:A:152:MET:CE	1:A:175:LEU:HB3	2.33	0.58
1:A:156:ILE:HG22	1:A:172:ARG:HH12	1.68	0.58
1:A:77:PRO:HG3	1:A:102:TYR:CZ	2.40	0.56
2:B:87:ARG:NH2	2:B:125:ASP:OD2	2.37	0.56
1:A:152:MET:CE	1:A:175:LEU:CB	2.83	0.56
2:B:325:MET:SD	2:B:381:MET:HG3	2.47	0.55
2:B:325:MET:HA	2:B:325:MET:HE2	1.92	0.51
1:A:357:ARG:HH11	1:A:357:ARG:HG2	1.76	0.50
2:B:201:VAL:HG11	2:B:251:TYR:HB3	1.94	0.50
2:B:28:PRO:HD2	2:B:29:GLU:OE1	2.12	0.49
2:B:422:GLY:O	2:B:423:PHE:HB2	2.13	0.49
2:B:149:HIS:ND1	2:B:152:PRO:HD2	2.28	0.49
1:A:134:TRP:CD1	1:A:167:GLN:HG2	2.47	0.49
1:A:264:LEU:O	1:A:265:GLU:C	2.55	0.49
2:B:192:LEU:HD23	2:B:199:VAL:HG23	1.94	0.49
2:B:308:LEU:HD13	2:B:330:PHE:HB3	1.94	0.48
2:B:239:ILE:HB	2:B:252:THR:HA	1.95	0.48
2:B:220:ASP:HA	2:B:223:GLU:HG2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:33:GLU:OE2	2:B:33:GLU:HA	2.13	0.47
1:A:263:THR:HG21	1:A:280:LEU:HB2	1.96	0.47
1:A:303:GLN:HB2	1:A:304:PRO:HD2	1.95	0.47
1:A:311:LEU:C	1:A:311:LEU:HD23	2.40	0.47
1:A:152:MET:CE	1:A:175:LEU:HB2	2.45	0.47
2:B:75:ARG:NH2	2:B:393:VAL:O	2.48	0.46
1:A:302:LEU:HB3	1:A:306:HIS:HB2	1.97	0.45
1:A:329:ASN:O	1:A:330:LYS:C	2.58	0.45
1:A:265:GLU:O	1:A:268:LYS:N	2.50	0.44
2:B:370:LEU:HD23	2:B:394:LEU:HD11	2.00	0.44
1:A:87:VAL:O	1:A:88:VAL:C	2.60	0.44
1:A:134:TRP:CD1	1:A:167:GLN:CG	3.01	0.43
2:B:151:ALA:HB3	2:B:152:PRO:CD	2.48	0.43
2:B:51:VAL:O	2:B:55:ILE:HG12	2.19	0.43
4:B:1002:FPP:O1	5:B:1003:ED2:HAA	2.19	0.43
2:B:376:PHE:CD2	2:B:376:PHE:C	2.97	0.43
2:B:193:MET:SD	2:B:203:SER:HB3	2.59	0.43
1:A:96:PHE:HA	1:A:126:LEU:HD13	2.01	0.42
2:B:149:HIS:HB3	2:B:152:PRO:HG2	2.00	0.42
1:A:294:ASN:O	1:A:298:GLN:HG3	2.20	0.42
2:B:29:GLU:H	2:B:29:GLU:CD	2.27	0.42
1:A:156:ILE:HG13	1:A:157:ALA:N	2.34	0.42
1:A:362:SER:O	1:A:366:LYS:HG2	2.20	0.42
2:B:218:THR:HB	2:B:219:PRO:HD2	2.00	0.42
2:B:301:SER:O	2:B:305:ALA:HB3	2.20	0.42
2:B:352:ASP:OD1	2:B:352:ASP:C	2.63	0.41
2:B:149:HIS:HE1	2:B:193:MET:HE2	1.85	0.41
1:A:156:ILE:HG22	1:A:172:ARG:NH1	2.35	0.41
1:A:346:LYS:HE3	1:A:357:ARG:NH2	2.36	0.41
5:B:1003:ED2:HATA	5:B:1003:ED2:HAN	1.40	0.41
2:B:202:ARG:HD2	4:B:1002:FPP:H152	2.02	0.40
4:B:1002:FPP:H61	5:B:1003:ED2:CBD	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	321/377 (85%)	303 (94%)	15 (5%)	3 (1%)	14	17
2	B	405/437 (93%)	394 (97%)	11 (3%)	0	100	100
All	All	726/814 (89%)	697 (96%)	26 (4%)	3 (0%)	30	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	265	GLU
1	A	326	GLN
1	A	304	PRO

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	301/338 (89%)	294 (98%)	7 (2%)	44	59
2	B	344/371 (93%)	342 (99%)	2 (1%)	78	86
All	All	645/709 (91%)	636 (99%)	9 (1%)	59	72

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	ARG
1	A	156	ILE
1	A	167	GLN
1	A	173	ARG
1	A	303	GLN
1	A	324	GLU
1	A	334	LEU
2	B	68	VAL
2	B	121	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	108	GLN
1	A	135	HIS
1	A	145	GLN
1	A	326	GLN
2	B	56	GLN
2	B	327	HIS
2	B	332	GLN
2	B	383	HIS
2	B	410	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	$\# Z > 2$
5	ED2	B	1003	3	36,38,38	4.05	16 (44%)	43,53,53	6.22	19 (44%)
4	FPP	B	1002	-	22,23,23	2.13	8 (36%)	27,31,31	1.25	2 (7%)

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	ED2	B	1003	3	-	13/29/31/31	0/4/4/4
4	FPP	B	1002	-	-	4/25/25/25	-

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1003	ED2	CAV-CBA	-14.92	1.24	1.51
5	B	1003	ED2	CAV-NBF	-9.51	1.33	1.48
5	B	1003	ED2	CAS-NBG	-6.42	1.27	1.37
5	B	1003	ED2	CBC-NBH	-6.09	1.28	1.38
5	B	1003	ED2	OAD-SBI	-5.87	1.37	1.43
5	B	1003	ED2	CAA-NBG	-4.56	1.35	1.46
5	B	1003	ED2	CAP-CBC	-4.16	1.31	1.36
5	B	1003	ED2	CBD-SBI	-4.07	1.66	1.76
5	B	1003	ED2	CAW-CBC	4.01	1.55	1.50
4	B	1002	FPP	C11-C12	-3.58	1.39	1.50
4	B	1002	FPP	PB-O1B	3.48	1.61	1.50
5	B	1003	ED2	CAB-NBH	-3.45	1.40	1.46
4	B	1002	FPP	C6-C7	-3.41	1.40	1.50
5	B	1003	ED2	CBB-NBE	-3.29	1.29	1.38
5	B	1003	ED2	OAE-SBI	-3.23	1.40	1.43
4	B	1002	FPP	C1-C2	-3.14	1.40	1.49
5	B	1003	ED2	CAU-NBF	-3.13	1.43	1.48
5	B	1003	ED2	CBD-NAY	-3.09	1.31	1.38
4	B	1002	FPP	C2-C3	3.08	1.40	1.33
4	B	1002	FPP	PA-O1A	2.98	1.61	1.50
5	B	1003	ED2	CAP-NAX	-2.97	1.32	1.37
4	B	1002	FPP	C7-C8	2.81	1.39	1.33
5	B	1003	ED2	CAQ-NBH	-2.71	1.28	1.35
4	B	1002	FPP	C12-C13	2.47	1.39	1.32

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1003	ED2	NBG-CAR-NAY	-25.62	102.85	112.66
5	B	1003	ED2	CAS-NBG-CAR	18.73	114.11	106.71
5	B	1003	ED2	CBA-CAV-NBF	11.40	134.37	112.14
5	B	1003	ED2	OAE-SBI-OAD	-8.98	106.81	119.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1003	ED2	CAV-NBF-CAU	8.75	132.04	116.98
5	B	1003	ED2	OAD-SBI-CBD	7.41	115.64	108.11
5	B	1003	ED2	CAV-CBA-CAJ	-6.48	108.81	120.75
5	B	1003	ED2	CAU-NBF-SBI	-6.23	105.08	117.88
5	B	1003	ED2	CAR-NAY-CBD	6.07	108.97	103.39
5	B	1003	ED2	OAE-SBI-CBD	5.74	113.94	108.11
5	B	1003	ED2	CAA-NBG-CAS	-5.72	118.52	126.03
5	B	1003	ED2	CAV-CBA-CAK	5.70	131.25	120.75
5	B	1003	ED2	NBH-CAQ-NAX	-5.09	109.91	112.94
5	B	1003	ED2	CAW-NBE-CAT	3.44	123.58	116.11
5	B	1003	ED2	OAE-SBI-NBF	3.29	112.99	107.26
5	B	1003	ED2	CAT-NBE-CBB	-3.05	116.32	121.15
4	B	1002	FPP	C10-C8-C9	2.76	120.02	115.23
5	B	1003	ED2	OAD-SBI-NBF	-2.75	102.48	107.26
4	B	1002	FPP	O3B-PB-O3A	2.54	113.16	104.64
5	B	1003	ED2	CAT-CAU-NBF	2.43	116.57	112.43
5	B	1003	ED2	CAV-NBF-SBI	-2.06	112.60	117.47

There are no chirality outliers.

All (17) torsion outliers are listed below:

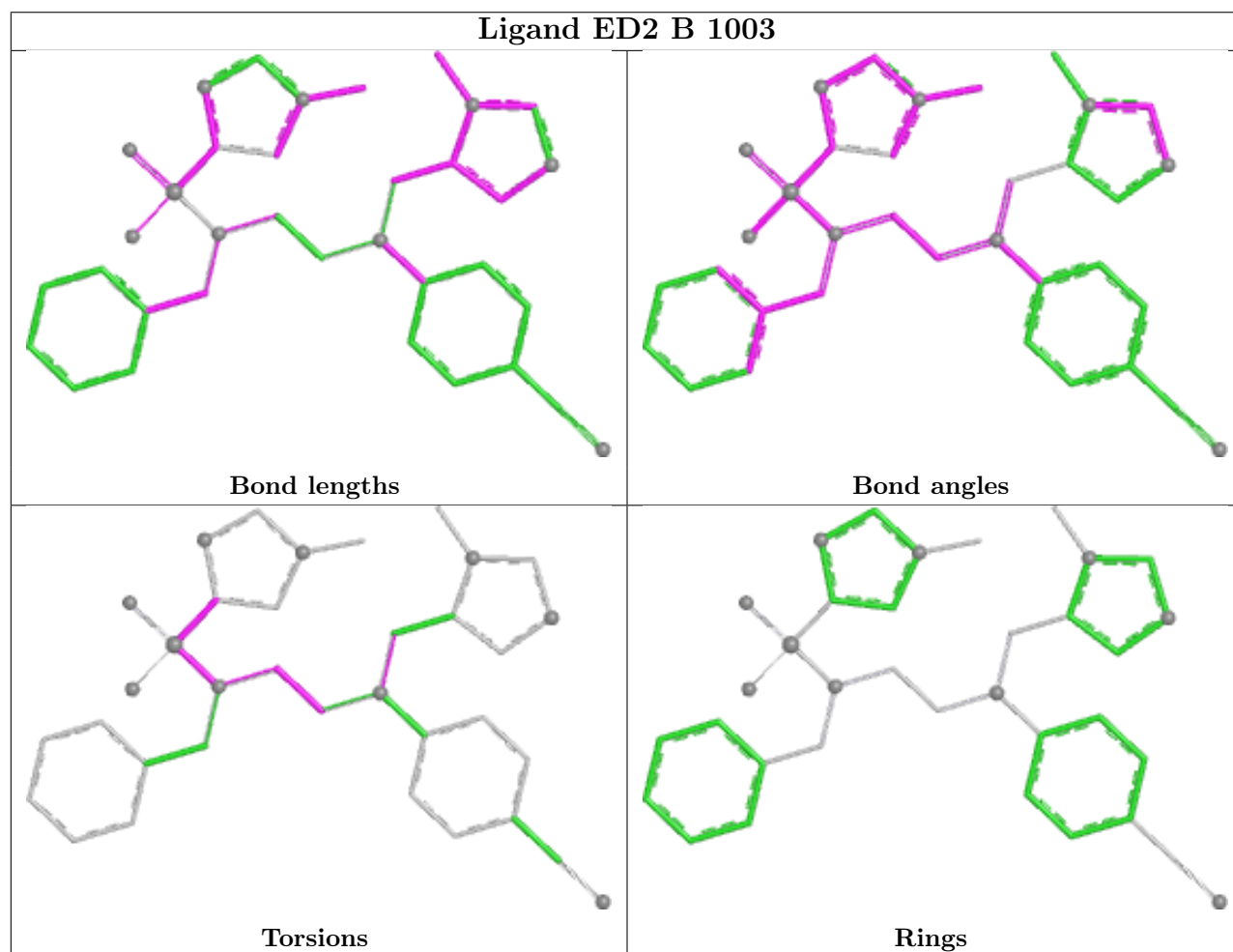
Mol	Chain	Res	Type	Atoms
5	B	1003	ED2	CAT-CAU-NBF-SBI
5	B	1003	ED2	CAT-CAU-NBF-CAV
5	B	1003	ED2	CAU-NBF-SBI-CBD
5	B	1003	ED2	CAS-CBD-SBI-OAD
5	B	1003	ED2	NAY-CBD-SBI-OAD
5	B	1003	ED2	CAS-CBD-SBI-OAE
5	B	1003	ED2	NAY-CBD-SBI-OAE
5	B	1003	ED2	CAV-NBF-SBI-OAD
5	B	1003	ED2	CAV-NBF-SBI-CBD
5	B	1003	ED2	CAU-NBF-SBI-OAD
5	B	1003	ED2	CAU-NBF-SBI-OAE
5	B	1003	ED2	NBE-CAT-CAU-NBF
4	B	1002	FPP	C4-C3-C5-C6
4	B	1002	FPP	C2-C3-C5-C6
5	B	1003	ED2	CBC-CAW-NBE-CAT
4	B	1002	FPP	PB-O3A-PA-O1A
4	B	1002	FPP	C5-C6-C7-C8

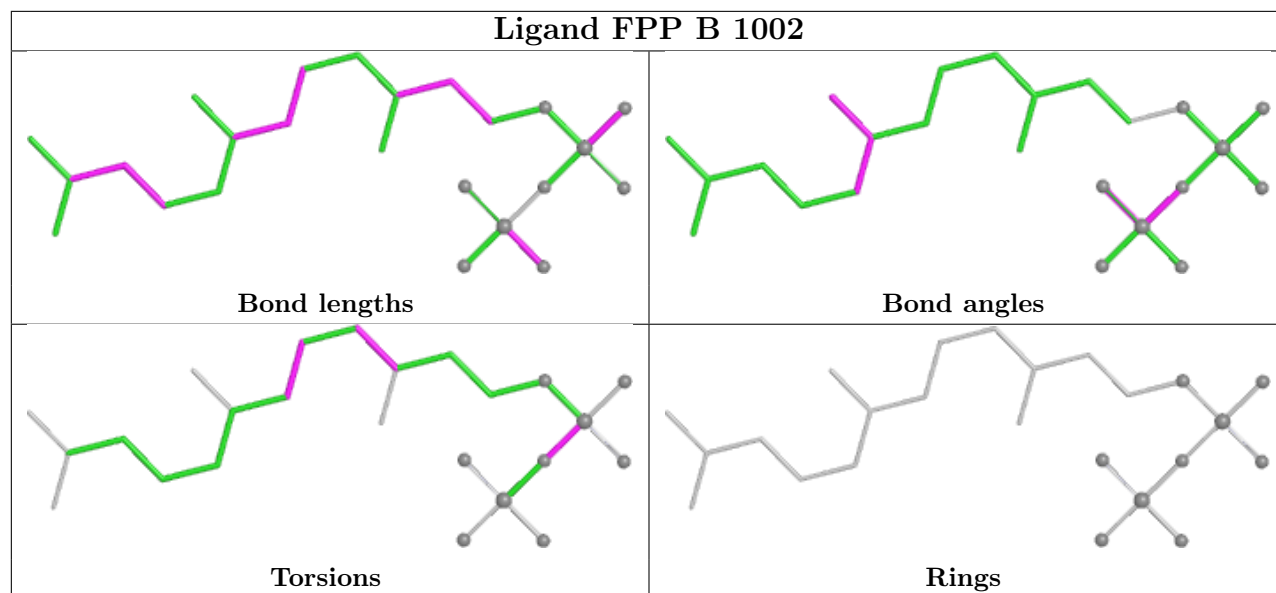
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1003	ED2	6	0
4	B	1002	FPP	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

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	323/377 (85%)	0.38	23 (7%) 22 19	17, 32, 59, 70	0
2	B	407/437 (93%)	0.06	11 (2%) 56 57	18, 29, 44, 56	0
All	All	730/814 (89%)	0.20	34 (4%) 36 35	17, 31, 54, 70	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	375	ALA	7.0
1	A	377	VAL	4.9
2	B	19	TRP	4.7
1	A	374	PRO	4.3
1	A	376	SER	4.1
1	A	55	PHE	3.8
2	B	423	PHE	3.7
1	A	372	ASP	3.5
2	B	20	SER	3.2
2	B	17	PRO	3.2
1	A	373	ILE	3.2
1	A	286	ASP	3.1
2	B	64	PHE	2.9
1	A	370	GLU	2.7
2	B	120	GLN	2.6
2	B	18	VAL	2.6
1	A	305	SER	2.6
1	A	326	GLN	2.6
1	A	364	GLN	2.6
2	B	325	MET	2.6
2	B	266	ARG	2.5
1	A	316	VAL	2.5
1	A	327	CYS	2.4
1	A	330	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	223	GLU	2.4
1	A	328	ASP	2.3
1	A	265	GLU	2.3
1	A	368	SER	2.2
1	A	287	ARG	2.1
1	A	324	GLU	2.1
1	A	371	SER	2.1
1	A	288	GLY	2.1
1	A	334	LEU	2.1
2	B	121	ILE	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](#)

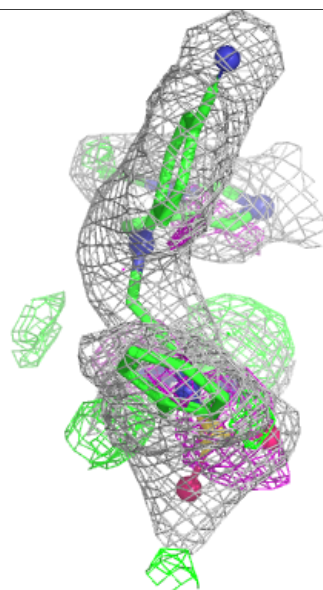
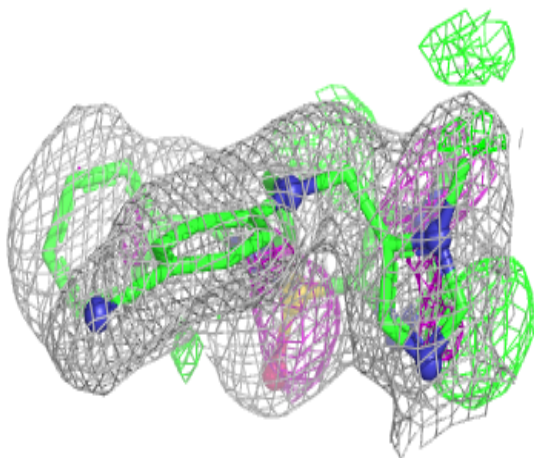
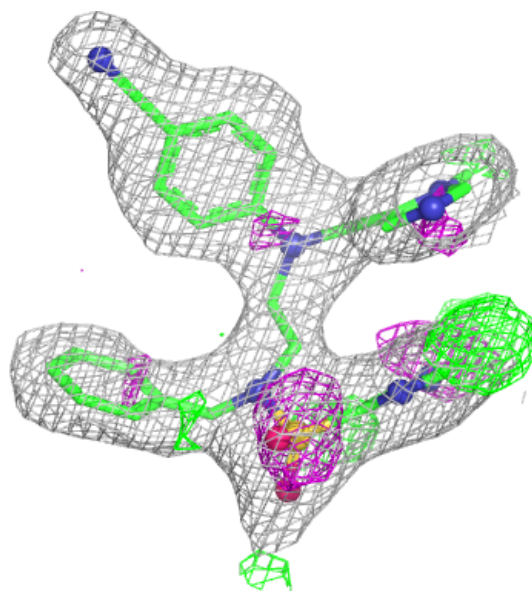
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

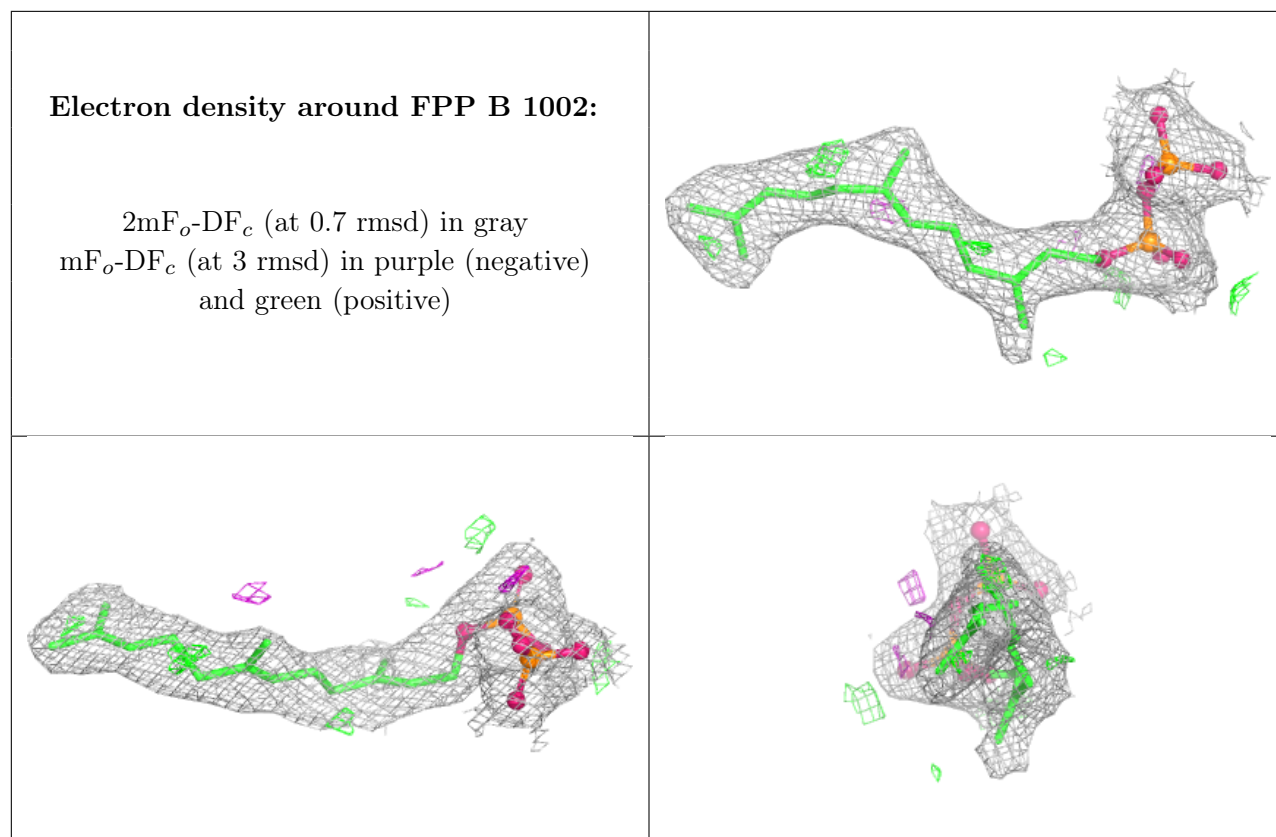
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	ED2	B	1003	35/35	0.88	0.12	18,24,31,32	0
4	FPP	B	1002	24/24	0.98	0.08	18,19,21,22	0
3	ZN	B	1001	1/1	1.00	0.02	19,19,19,19	0

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 ED2 B 1003:

$2mF_o-DF_c$ (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.