



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:48 AM UTC

PDB ID : 3EUV / pdb_00003euv
Title : Crystal structure of FTase(ALPHA-subunit; BETA-subunit DELTA C10, W102T, Y154T) in complex with BiotinGPP
Authors : Guo, Z.; Nguyen, U.T.T.; Delon, C.; Bon, R.S.; Blankenfeldt, W.; Goody, R.S.; Waldmann, H.; Wolters, D.; Alexandrov, K.
Deposited on : 2008-10-11
Resolution : 2.75 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

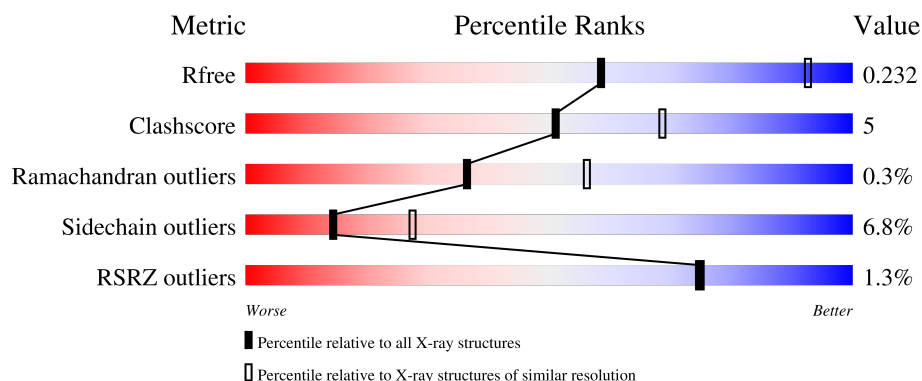
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

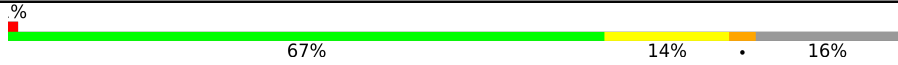
The reported resolution of this entry is 2.75 Å.

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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	 67% 14% • 16%
2	B	427	 79% 13% • 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5966 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	315	Total	C	N	O	S	0	0	0
			2688	1712	473	498	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	402	Total	C	N	O	S	0	2	0
			3159	2014	543	579	23			

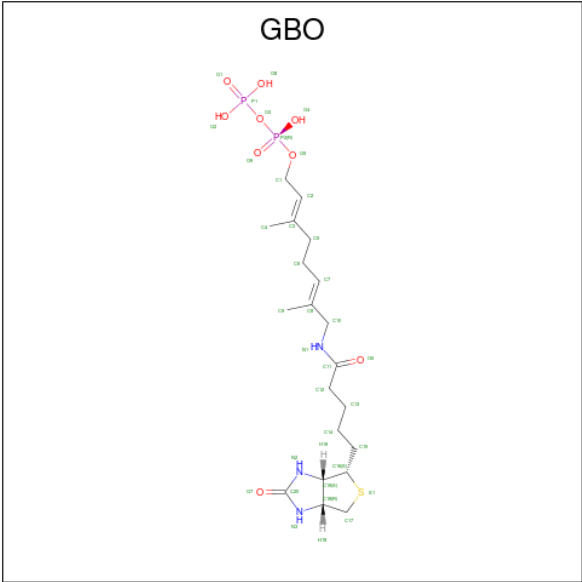
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	102	THR	TRP	engineered mutation	UNP Q02293
B	154	THR	TYR	engineered mutation	UNP Q02293

- 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 (2E,6E)-3,7-dimethyl-8-({5-[(3aS,4S,6aR)-2-oxohexahydro-1H-thieno[3,4-d]imidazol-4-yl]pentanoyl}amino)octa-2,6-dien-1-yl trihydrogen diphosphate (CCD ID: GBO) (formula: C₂₀H₃₅N₃O₉P₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	S	0	0
			35	20	3	9	2	1		

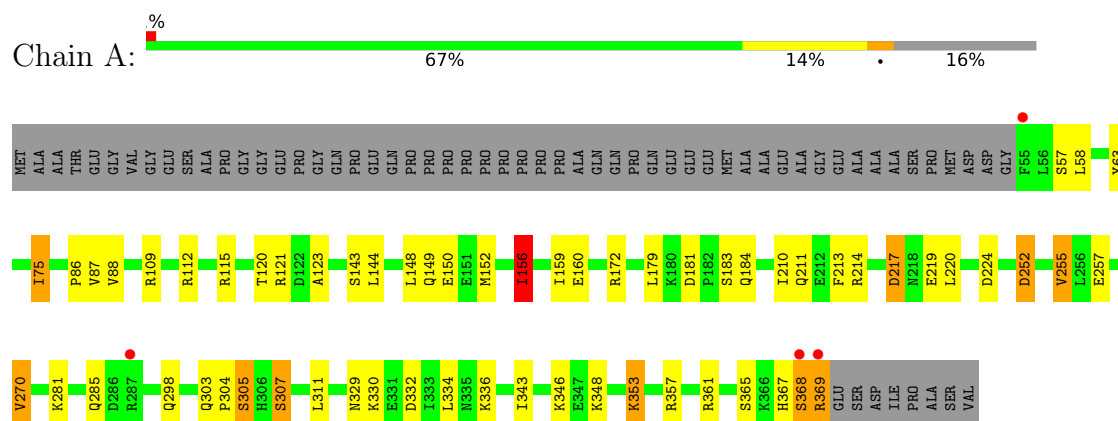
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	33	Total	O	0	0
			33	33		
5	B	50	Total	O	0	0
			50	50		

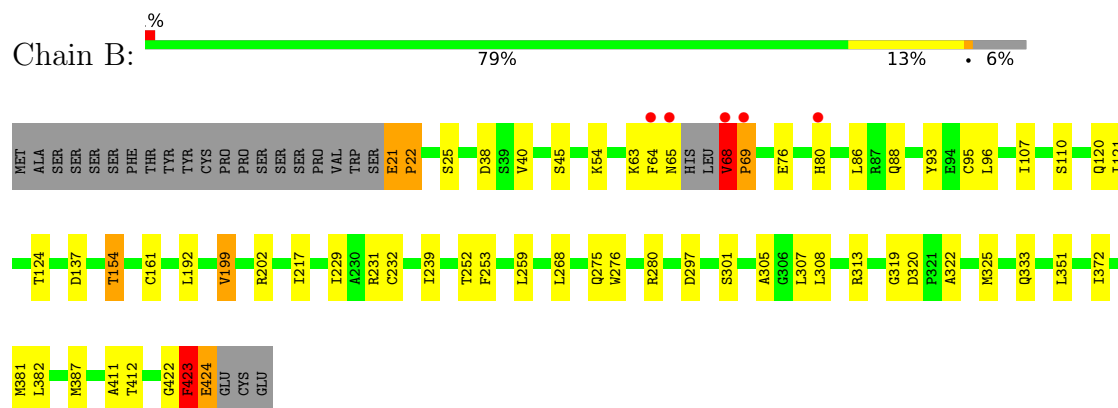
3 Residue-property plots [i](#)

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: Protein farnesyltransferase/geranylgeranyltransferase type-1 subunit alpha



- Molecule 2: Protein farnesyltransferase subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	174.78Å 174.78Å 70.38Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.97 – 2.75 29.97 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.97-2.75) 99.8 (29.97-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.14 (at 2.76Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.169 , 0.231 0.178 , 0.232	Depositor DCC
R_{free} test set	1613 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	53.7	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5966	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% 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: ZN, GBO

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	1.15	1/2754 (0.0%)	1.10	9/3738 (0.2%)
2	B	1.10	1/3247 (0.0%)	1.09	5/4407 (0.1%)
All	All	1.12	2/6001 (0.0%)	1.09	14/8145 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	1
2	B	0	3
All	All	1	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	22	PRO	N-CA	-5.67	1.43	1.47
1	A	123	ALA	CA-CB	-5.21	1.44	1.53

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	285	GLN	N-CA-C	8.64	121.65	111.71
2	B	423	PHE	N-CA-C	-8.51	92.68	110.80
2	B	154	THR	CB-CA-C	-8.20	97.18	110.79
1	A	307	SER	N-CA-C	6.04	118.99	109.39
1	A	270	VAL	CA-C-N	5.86	126.26	119.47
1	A	270	VAL	C-N-CA	5.86	126.26	119.47
1	A	285	GLN	CB-CA-C	5.82	121.40	110.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	THR	N-CA-C	-5.73	105.17	111.82
1	A	156	ILE	CB-CA-C	-5.71	104.58	112.24
2	B	69	PRO	N-CA-C	5.43	123.65	112.47
2	B	161	CYS	N-CA-C	5.11	116.85	111.28
2	B	88	GLN	N-CA-C	5.07	116.94	108.99
1	A	183	SER	N-CA-C	5.00	117.47	111.71
1	A	86	PRO	N-CA-C	5.00	118.91	111.41

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	285	GLN	CA

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	368	SER	Peptide
2	B	21	GLU	Peptide
2	B	423	PHE	Peptide
2	B	68	VAL	Peptide

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	2688	0	2614	30	0
2	B	3159	0	3086	31	0
3	B	1	0	0	0	0
4	B	35	0	32	1	0
5	A	33	0	0	1	0
5	B	50	0	0	4	0
All	All	5966	0	5732	60	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 (60) 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:HD13	1:A:172:ARG:HH12	1.33	0.92
1:A:368:SER:O	1:A:369:ARG:HG2	1.88	0.73
1:A:368:SER:C	1:A:369:ARG:HG2	2.14	0.70
2:B:192:LEU:HD23	2:B:199:VAL:HG23	1.75	0.69
1:A:303:GLN:O	1:A:307:SER:HB2	1.93	0.67
1:A:149:GLN:HE22	1:A:184:GLN:HE22	1.44	0.65
2:B:231:ARG:HD3	5:B:451:HOH:O	2.00	0.61
1:A:156:ILE:CD1	1:A:172:ARG:HH12	2.09	0.60
1:A:152:MET:O	1:A:156:ILE:CG1	2.52	0.57
2:B:76:GLU:OE2	2:B:80[B]:HIS:NE2	2.38	0.56
2:B:93:TYR:HD2	2:B:96:LEU:HD12	1.71	0.55
1:A:210:ILE:HD11	1:A:219:GLU:OE1	2.07	0.54
1:A:152:MET:O	1:A:156:ILE:HG12	2.09	0.53
2:B:325:MET:HB3	2:B:381:MET:HG3	1.91	0.52
2:B:319:GLY:O	2:B:320:ASP:C	2.53	0.52
1:A:353:LYS:O	1:A:357:ARG:HG3	2.10	0.51
1:A:152:MET:O	1:A:156:ILE:HG13	2.11	0.51
2:B:301:SER:O	2:B:305:ALA:HB3	2.11	0.51
5:A:469:HOH:O	2:B:95:CYS:HB3	2.11	0.49
2:B:64:PHE:O	2:B:65:ASN:O	2.30	0.49
1:A:156:ILE:HD13	1:A:172:ARG:NH1	2.16	0.49
1:A:252:ASP:HB3	1:A:255:VAL:HB	1.94	0.49
2:B:423:PHE:CA	2:B:424:GLU:C	2.86	0.48
1:A:361:ARG:NH2	2:B:322:ALA:O	2.45	0.48
2:B:229:ILE:O	2:B:232:CYS:HB2	2.15	0.47
1:A:329:ASN:HB3	1:A:332:ASP:HB3	1.97	0.46
2:B:259:LEU:HD12	2:B:268:LEU:HD11	1.98	0.46
2:B:253:PHE:HA	2:B:307:LEU:HD21	1.97	0.45
2:B:68:VAL:N	2:B:69:PRO:CD	2.78	0.45
2:B:297:ASP:OD1	2:B:297:ASP:C	2.58	0.45
2:B:76:GLU:OE2	2:B:80[B]:HIS:CD2	2.70	0.45
2:B:333:GLN:HG3	2:B:387:MET:SD	2.56	0.45
1:A:181:ASP:OD1	1:A:181:ASP:C	2.59	0.45
2:B:64:PHE:CB	5:B:456:HOH:O	2.66	0.44
2:B:411:ALA:O	2:B:412:THR:C	2.60	0.44
1:A:334:LEU:HD22	1:A:367:HIS:O	2.18	0.44
2:B:423:PHE:HA	2:B:424:GLU:C	2.43	0.44
1:A:329:ASN:O	1:A:330:LYS:C	2.60	0.44
2:B:38:ASP:O	2:B:40:VAL:HG23	2.18	0.44
1:A:75:ILE:HD13	1:A:115:ARG:NH2	2.33	0.44
2:B:239:ILE:HB	2:B:252:THR:HA	1.99	0.44
2:B:68:VAL:N	2:B:69:PRO:HD3	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:LEU:CD2	1:A:367:HIS:O	2.65	0.43
1:A:217:ASP:OD2	1:A:217:ASP:N	2.51	0.43
2:B:276:TRP:O	2:B:280:ARG:HG2	2.19	0.42
1:A:304:PRO:O	1:A:305:SER:CB	2.68	0.42
2:B:64:PHE:HB3	5:B:456:HOH:O	2.18	0.42
2:B:253:PHE:CD2	2:B:253:PHE:C	2.98	0.42
1:A:58:LEU:HD23	1:A:63:TYR:CE1	2.54	0.42
1:A:159:ILE:O	1:A:160:GLU:C	2.61	0.42
1:A:148:LEU:HB2	1:A:179:LEU:HD21	2.02	0.42
1:A:87:VAL:O	1:A:88:VAL:C	2.63	0.41
1:A:121:ARG:O	1:A:121:ARG:HG2	2.12	0.41
1:A:112:ARG:O	1:A:144:LEU:HD21	2.20	0.41
1:A:213:PHE:O	1:A:214:ARG:C	2.63	0.41
2:B:422:GLY:O	2:B:423:PHE:HD1	2.04	0.41
2:B:86:LEU:HB2	2:B:107:ILE:HG21	2.01	0.41
2:B:63:LYS:O	5:B:500:HOH:O	2.22	0.41
2:B:202:ARG:HD2	4:B:429:GBO:H13	2.02	0.41
1:A:311:LEU:C	1:A:311:LEU:HD23	2.45	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	313/377 (83%)	291 (93%)	20 (6%)	2 (1%)	21	38
2	B	400/427 (94%)	379 (95%)	21 (5%)	0	100	100
All	All	713/804 (89%)	670 (94%)	41 (6%)	2 (0%)	36	56

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	305	SER
1	A	252	ASP

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	294/338 (87%)	272 (92%)	22 (8%)	12	24
2	B	340/363 (94%)	319 (94%)	21 (6%)	16	31
All	All	634/701 (90%)	591 (93%)	43 (7%)	14	27

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	SER
1	A	75	ILE
1	A	109	ARG
1	A	143	SER
1	A	150	GLU
1	A	156	ILE
1	A	211	GLN
1	A	217	ASP
1	A	220	LEU
1	A	224	ASP
1	A	255	VAL
1	A	257	GLU
1	A	270	VAL
1	A	281	LYS
1	A	298	GLN
1	A	336	LYS
1	A	343	ILE
1	A	346	LYS
1	A	348	LYS
1	A	353	LYS
1	A	365	SER
1	A	369	ARG

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Mol	Chain	Res	Type
2	B	21	GLU
2	B	22	PRO
2	B	25	SER
2	B	45	SER
2	B	54	LYS
2	B	68	VAL
2	B	110	SER
2	B	120	GLN
2	B	121	ILE
2	B	124	THR
2	B	137	ASP
2	B	154	THR
2	B	199	VAL
2	B	217	ILE
2	B	275	GLN
2	B	308	LEU
2	B	313	ARG
2	B	351	LEU
2	B	372	ILE
2	B	382	LEU
2	B	424	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	HIS
1	A	149	GLN
1	A	221	GLN
1	A	225	GLN
1	A	303	GLN
1	A	325	ASN
1	A	326	GLN
1	A	367	HIS
2	B	56	GLN
2	B	170	ASN
2	B	275	GLN
2	B	332	GLN
2	B	417	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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
4	GB0	B	429	-	34,36,36	2.23	13 (38%)	43,50,50	1.77	10 (23%)

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
4	GB0	B	429	-	-	11/31/52/52	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	429	GB0	O5-C1	-5.99	1.34	1.44
4	B	429	GB0	C1-C2	-4.32	1.37	1.49
4	B	429	GB0	P1-O8	3.74	1.68	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	429	GB0	C12-C11	3.54	1.58	1.51
4	B	429	GB0	P2-O3	-3.40	1.55	1.59
4	B	429	GB0	C6-C7	-3.19	1.40	1.50
4	B	429	GB0	C19-N2	3.02	1.51	1.45
4	B	429	GB0	P1-O2	2.75	1.65	1.54
4	B	429	GB0	C18-N3	2.68	1.50	1.46
4	B	429	GB0	C7-C8	2.66	1.39	1.33
4	B	429	GB0	C2-C3	2.58	1.39	1.33
4	B	429	GB0	C4-C3	2.44	1.56	1.50
4	B	429	GB0	C5-C3	2.40	1.56	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	429	GB0	C6-C7-C8	-3.83	118.85	127.62
4	B	429	GB0	C18-C17-S1	-3.74	100.91	106.06
4	B	429	GB0	C16-C19-N2	3.67	117.22	113.34
4	B	429	GB0	C19-C16-S1	-3.60	100.94	105.03
4	B	429	GB0	C10-N1-C11	-3.44	111.54	121.49
4	B	429	GB0	C17-S1-C16	3.32	96.90	89.98
4	B	429	GB0	C4-C3-C5	2.69	119.90	115.23
4	B	429	GB0	C17-C18-C19	2.56	113.12	109.06
4	B	429	GB0	C1-C2-C3	-2.46	122.16	126.20
4	B	429	GB0	N3-C20-N2	2.05	111.22	108.85

There are no chirality outliers.

All (11) torsion outliers are listed below:

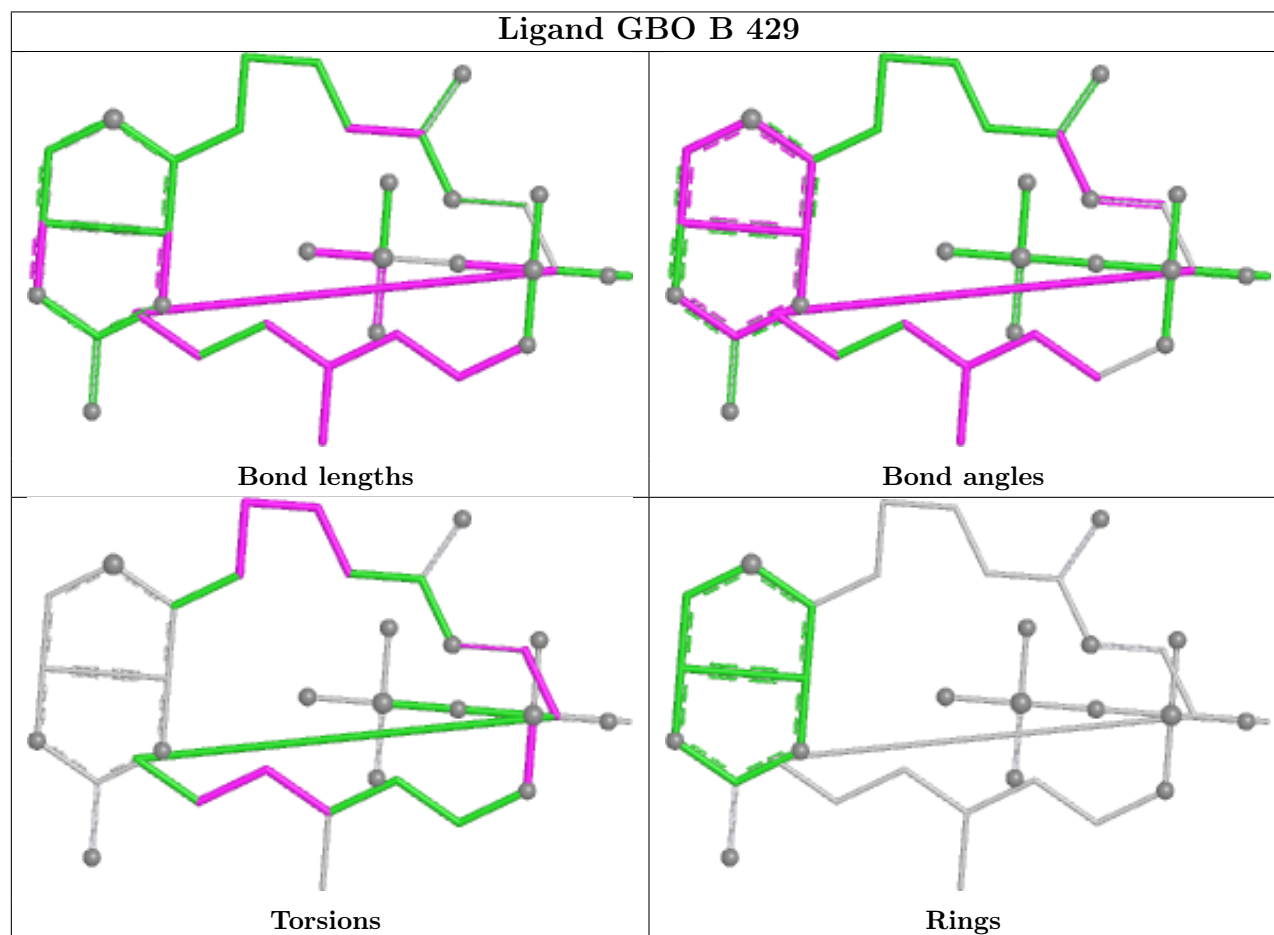
Mol	Chain	Res	Type	Atoms
4	B	429	GB0	C1-O5-P2-O3
4	B	429	GB0	N1-C10-C8-C7
4	B	429	GB0	N1-C10-C8-C9
4	B	429	GB0	C3-C5-C6-C7
4	B	429	GB0	C13-C14-C15-C16
4	B	429	GB0	C11-C12-C13-C14
4	B	429	GB0	C12-C13-C14-C15
4	B	429	GB0	C2-C3-C5-C6
4	B	429	GB0	C4-C3-C5-C6
4	B	429	GB0	C1-O5-P2-O9
4	B	429	GB0	C8-C10-N1-C11

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	429	GBO	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 ⓘ

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	315/377 (83%)	-0.21	4 (1%) 75 75	25, 58, 83, 97	0
2	B	402/427 (94%)	-0.32	5 (1%) 76 76	27, 50, 80, 117	2 (0%)
All	All	717/804 (89%)	-0.27	9 (1%) 75 75	25, 55, 83, 117	2 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	369	ARG	5.0
2	B	68	VAL	4.5
1	A	55	PHE	3.1
2	B	69	PRO	2.8
2	B	65	ASN	2.3
2	B	64	PHE	2.3
1	A	287	ARG	2.3
2	B	80[A]	HIS	2.2
1	A	368	SER	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.

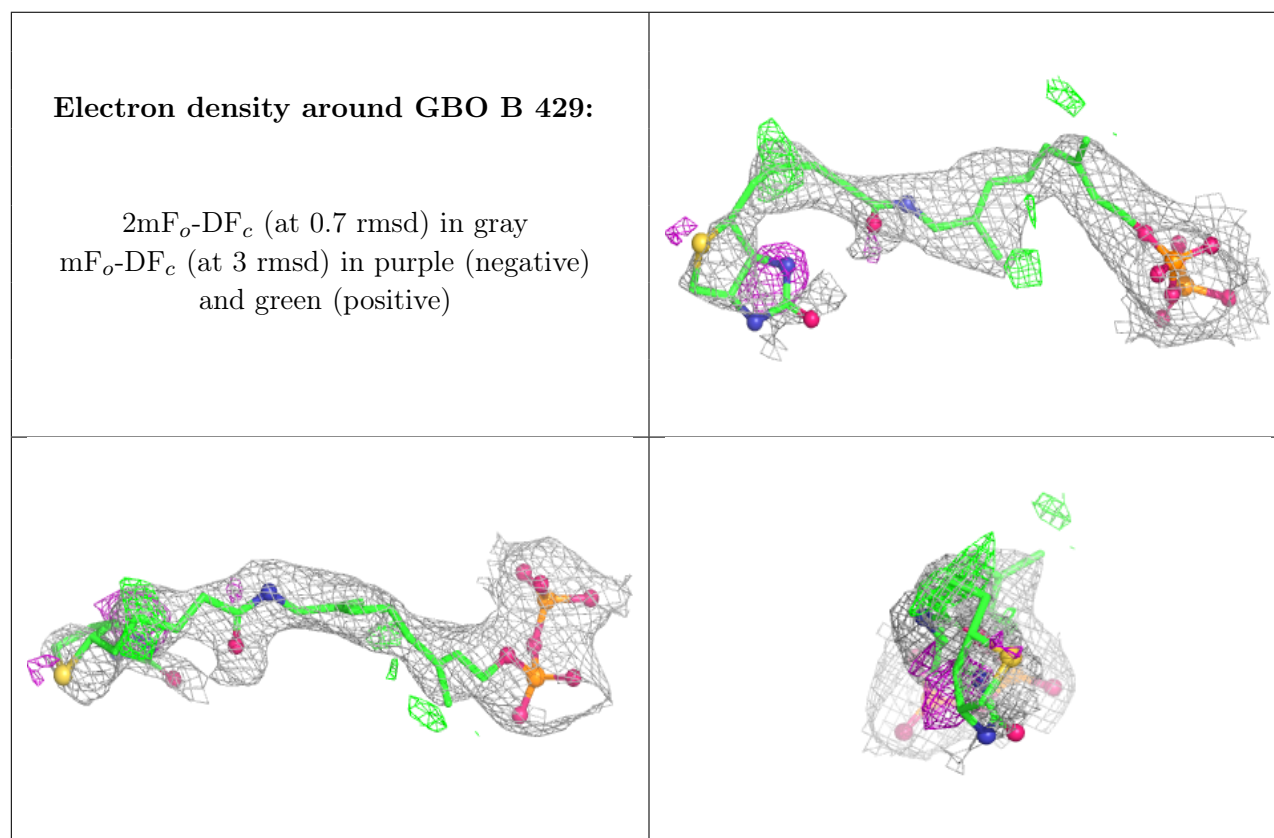
6.4 Ligands [i](#)

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	GBO	B	429	35/35	0.95	0.15	48,66,135,137	0
3	ZN	B	428	1/1	0.98	0.03	41,41,41,41	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.