



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:56 AM UTC

PDB ID : 5QCU / pdb_00005qcu
Title : Crystal structure of BACE complex with BMC022
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Deposited on : 2017-12-01
Resolution : 1.95 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

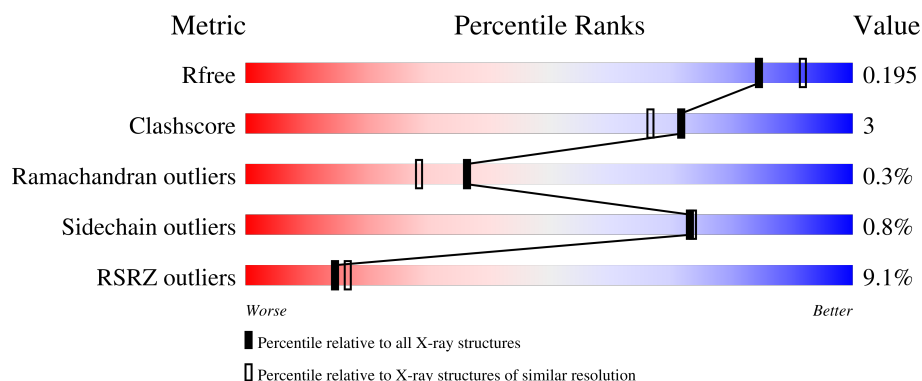
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	402	7% 84% 9% 6%
1	B	402	12% 83% 10% 6%
1	C	402	6% 88% 5% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9594 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.

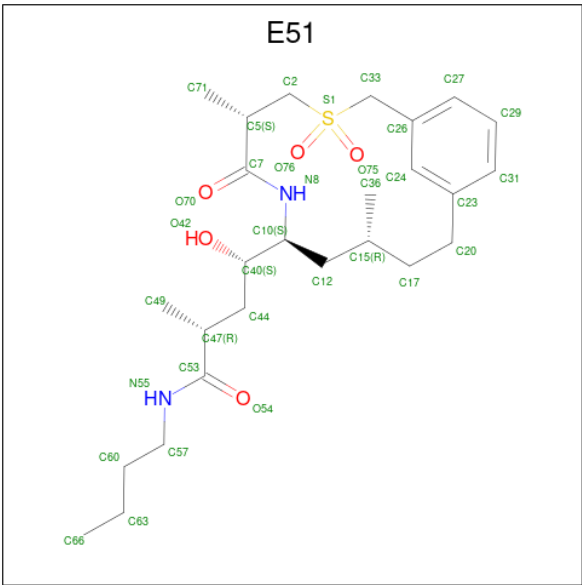
- Molecule 1 is a protein called Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	0
			2960	1895	492	559	14			
1	B	377	Total	C	N	O	S	0	0	0
			2966	1898	493	561	14			
1	C	377	Total	C	N	O	S	0	0	0
			2966	1898	493	561	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	GLY	-	expression tag	UNP P56817
A	-14	PRO	-	expression tag	UNP P56817
B	-15	GLY	-	expression tag	UNP P56817
B	-14	PRO	-	expression tag	UNP P56817
C	-15	GLY	-	expression tag	UNP P56817
C	-14	PRO	-	expression tag	UNP P56817

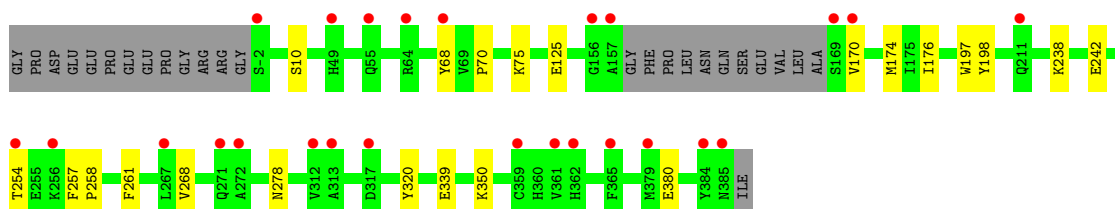
- Molecule 2 is (2R,4S)-N-butyl-4-[(5S,8S,10R)-5,10-dimethyl-3,3,6-trioxo-3lambda 6 -thia-7 -azabicyclo[11.3.1]heptadeca-1(17),13,15-trien-8-yl]-4-hydroxy-2-methylbutanamide (CCD ID: E51) (formula: C₂₆H₄₂N₂O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			34	26	2	5	1		
2	B	1	Total	C	N	O	S	0	0
			34	26	2	5	1		
2	C	1	Total	C	N	O	S	0	0
			34	26	2	5	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	216	Total	O	0	0
			216	216		
3	B	160	Total	O	0	0
			160	160		
3	C	224	Total	O	0	0
			224	224		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.17Å 102.48Å 100.44Å 90.00° 103.67° 90.00°	Depositor
Resolution (Å)	55.29 – 1.95 55.29 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.1 (55.29-1.95) 99.1 (55.29-1.95)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.12_2829, CNX	Depositor
R, R_{free}	0.167 , 0.193 0.169 , 0.195	Depositor DCC
R_{free} test set	11515 reflections (9.92%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.008	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9594	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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: E51

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	0.31	0/3035	0.56	0/4125
1	B	0.29	0/3041	0.55	0/4133
1	C	0.29	0/3041	0.57	0/4133
All	All	0.30	0/9117	0.56	0/12391

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	2960	0	2873	21	0
1	B	2966	0	2878	24	0
1	C	2966	0	2878	15	0
2	A	34	0	0	0	0
2	B	34	0	0	0	0
2	C	34	0	0	0	0
3	A	216	0	0	0	0
3	B	160	0	0	2	0
3	C	224	0	0	1	0
All	All	9594	0	8629	58	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 3.

All (58) 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:267:LEU:HD11	1:A:319:CYS:HB3	1.30	1.06
1:C:320:TYR:OH	3:C:501:HOH:O	1.95	0.77
1:C:258:PRO:HD2	1:C:268:VAL:HG12	1.68	0.76
1:B:153:GLN:OE1	3:B:501:HOH:O	2.07	0.72
1:A:174:MET:HE3	1:A:176:ILE:HD11	1.73	0.70
1:C:68:TYR:HE1	1:C:75:LYS:HD3	1.57	0.68
1:B:215:MET:HE1	1:B:240:VAL:HA	1.79	0.64
1:C:174:MET:HE3	1:C:176:ILE:HD11	1.81	0.62
1:B:110:ILE:HB	1:B:113:SER:HB3	1.82	0.62
1:A:157:ALA:HB2	1:A:170:VAL:HG22	1.83	0.61
1:C:68:TYR:OH	1:C:70:PRO:HB3	2.04	0.57
1:C:257:PHE:HD2	1:C:268:VAL:HG11	1.70	0.55
1:A:364:GLU:HG3	1:A:365:PHE:HD2	1.72	0.54
1:A:278:ASN:HD21	1:C:254:THR:HG23	1.72	0.54
1:A:267:LEU:HD12	1:A:320:TYR:O	2.09	0.53
1:B:252:SER:HA	1:B:279:ILE:HD12	1.91	0.51
1:B:267:LEU:HD22	1:B:309:VAL:HG21	1.92	0.51
1:B:174:MET:HE3	1:B:176:ILE:HD11	1.93	0.51
1:A:254:THR:HG23	1:C:278:ASN:HD21	1.76	0.50
1:B:155:CYS:O	1:B:170:VAL:HG22	2.11	0.50
1:B:209:ASN:HA	3:B:502:HOH:O	2.12	0.50
1:B:197:TRP:CG	1:B:198:TYR:H	2.31	0.49
1:B:258:PRO:HD2	1:B:268:VAL:HG22	1.93	0.49
1:C:197:TRP:CG	1:C:198:TYR:H	2.30	0.49
1:C:261:PHE:CD1	1:C:268:VAL:HG13	2.49	0.48
1:A:197:TRP:CG	1:A:198:TYR:H	2.32	0.47
1:C:238:LYS:O	1:C:242:GLU:HG3	2.15	0.47
1:A:13:GLY:HA3	1:A:30:LEU:HD11	1.97	0.47
1:A:14:TYR:OH	1:A:339:GLU:OE1	2.31	0.47
1:A:314:THR:HG22	1:A:316:GLN:HG3	1.96	0.47
1:B:311:ASP:C	1:B:313:ALA:H	2.23	0.47
1:B:20:VAL:HG12	1:B:85:VAL:HG22	1.97	0.46
1:C:68:TYR:CE1	1:C:75:LYS:HD3	2.44	0.46
1:B:232:THR:O	1:B:336:VAL:HG13	2.16	0.45
1:A:276:PRO:O	1:A:279:ILE:HG12	2.16	0.45
1:B:77:GLU:CG	1:B:104:GLU:HG3	2.47	0.45
1:A:238:LYS:O	1:A:242:GLU:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:SER:HB3	1:C:339:GLU:OE1	2.18	0.44
1:B:13:GLY:HA3	1:B:30:LEU:HD11	1.99	0.43
1:B:310:GLU:HG3	1:B:311:ASP:N	2.34	0.43
1:A:218:LYS:HG3	1:A:381:ASP:O	2.19	0.43
1:B:258:PRO:HB2	1:B:266:GLN:HG2	2.00	0.43
1:A:110:ILE:HB	1:A:113:SER:HB3	2.01	0.43
1:C:258:PRO:CD	1:C:268:VAL:HG12	2.43	0.43
1:C:350:LYS:HE2	1:C:350:LYS:HB3	1.88	0.43
1:A:267:LEU:HD11	1:A:319:CYS:CB	2.22	0.42
1:A:199:TYR:HB3	1:A:352:ILE:HD11	2.02	0.42
1:B:95:VAL:HG11	1:B:140:LEU:HA	2.02	0.42
1:A:222:TYR:HA	1:A:223:ASP:HA	1.72	0.42
1:B:199:TYR:HB3	1:B:352:ILE:HD11	2.02	0.41
1:B:205:ARG:HB3	1:B:286:TYR:HB2	2.03	0.41
1:A:357:SER:O	1:A:360:HIS:HB3	2.20	0.41
1:B:125:GLU:OE2	1:B:195:ARG:NH2	2.26	0.41
1:A:10:SER:HB3	1:A:339:GLU:OE1	2.21	0.40
1:B:8:GLY:C	1:B:170:VAL:HG12	2.46	0.40
1:B:14:TYR:CE2	1:B:170:VAL:HG13	2.56	0.40
1:B:249:LYS:HG3	1:B:262:TRP:CZ2	2.56	0.40
1:A:304:GLN:O	1:A:336:VAL:HB	2.20	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	372/402 (92%)	362 (97%)	9 (2%)	1 (0%)	36	28
1	B	373/402 (93%)	361 (97%)	11 (3%)	1 (0%)	36	28
1	C	373/402 (93%)	365 (98%)	7 (2%)	1 (0%)	36	28
All	All	1118/1206 (93%)	1088 (97%)	27 (2%)	3 (0%)	36	28

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	313	ALA
1	B	312	VAL
1	C	170	VAL

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	321/342 (94%)	320 (100%)	1 (0%)	86	87
1	B	322/342 (94%)	317 (98%)	5 (2%)	55	52
1	C	322/342 (94%)	320 (99%)	2 (1%)	78	79
All	All	965/1026 (94%)	957 (99%)	8 (1%)	73	74

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

Mol	Chain	Res	Type
1	A	267	LEU
1	B	104	GLU
1	B	170	VAL
1	B	242	GLU
1	B	254	THR
1	B	279	ILE
1	C	125	GLU
1	C	380	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	HIS
1	A	233	ASN
1	A	362	HIS
1	A	385	ASN
1	B	89	HIS

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Mol	Chain	Res	Type
1	B	233	ASN
1	B	278	ASN
1	B	385	ASN
1	C	89	HIS
1	C	233	ASN
1	C	385	ASN

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](#)

3 ligands are modelled in this entry.

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	E51	A	401	-	33,35,35	0.53	0	45,48,48	0.42	0
2	E51	C	401	-	33,35,35	0.34	0	45,48,48	0.48	0
2	E51	B	401	-	33,35,35	0.42	0	45,48,48	0.53	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
2	E51	A	401	-	-	3/44/44/44	0/1/2/2
2	E51	C	401	-	-	2/44/44/44	0/1/2/2
2	E51	B	401	-	-	2/44/44/44	0/1/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

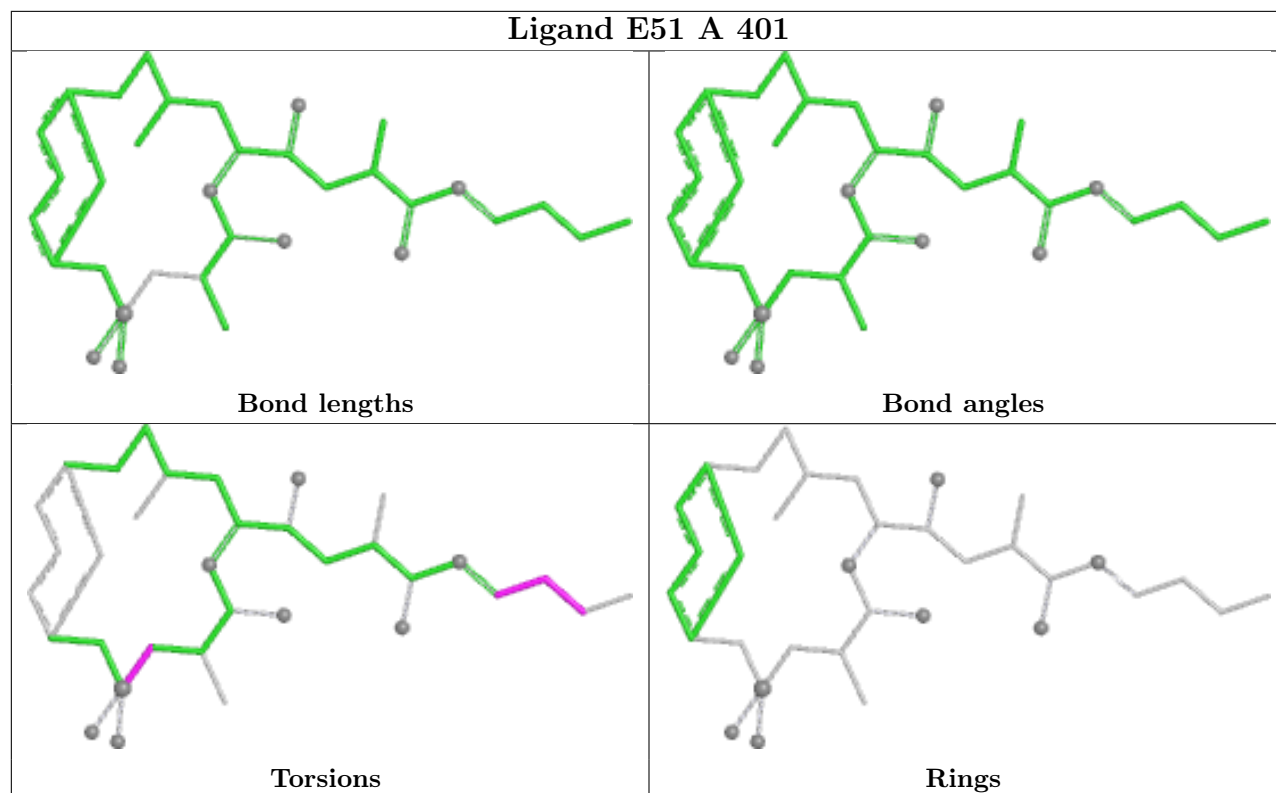
Mol	Chain	Res	Type	Atoms
2	A	401	E51	C57-C60-C63-C66
2	A	401	E51	C5-C2-S1-O75
2	B	401	E51	C5-C2-S1-O75
2	C	401	E51	C5-C2-S1-O75
2	B	401	E51	C5-C2-S1-C33
2	C	401	E51	C71-C5-C7-N8
2	A	401	E51	N55-C57-C60-C63

There are no ring outliers.

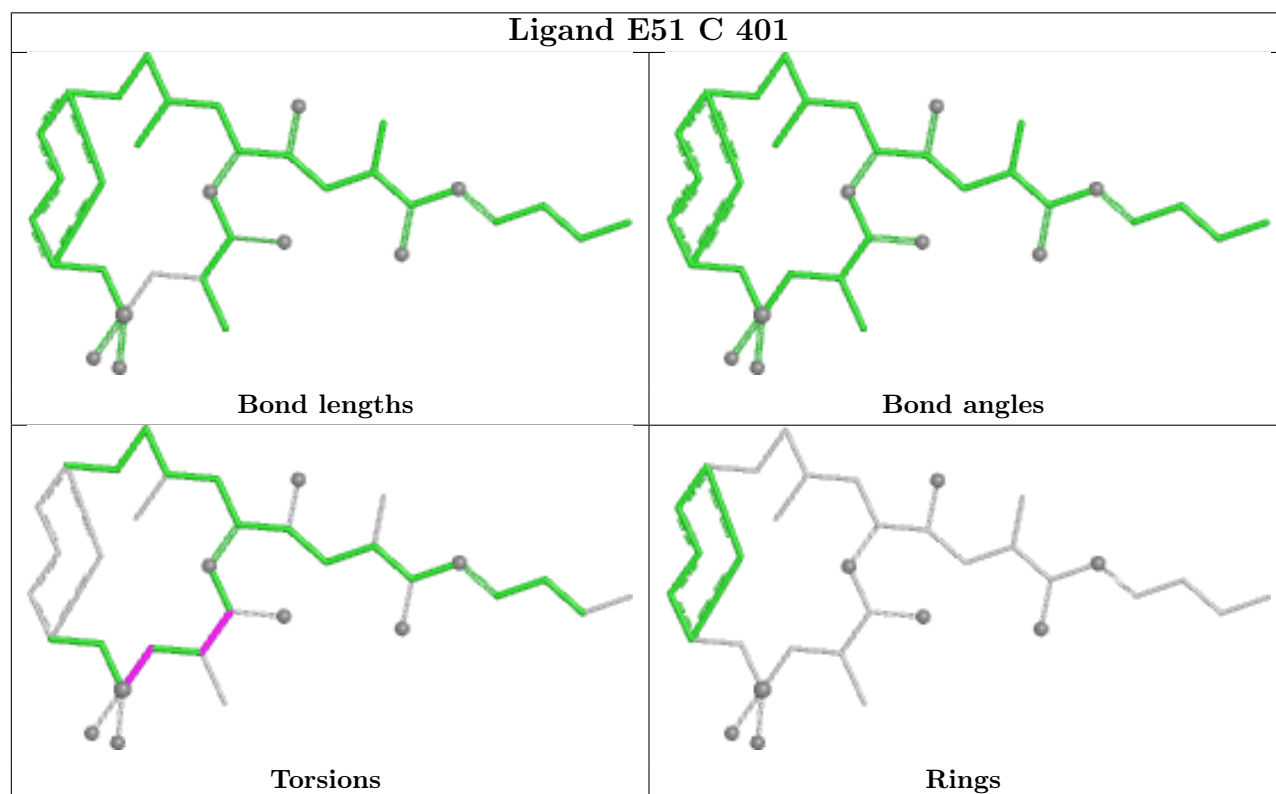
No monomer is involved in short contacts.

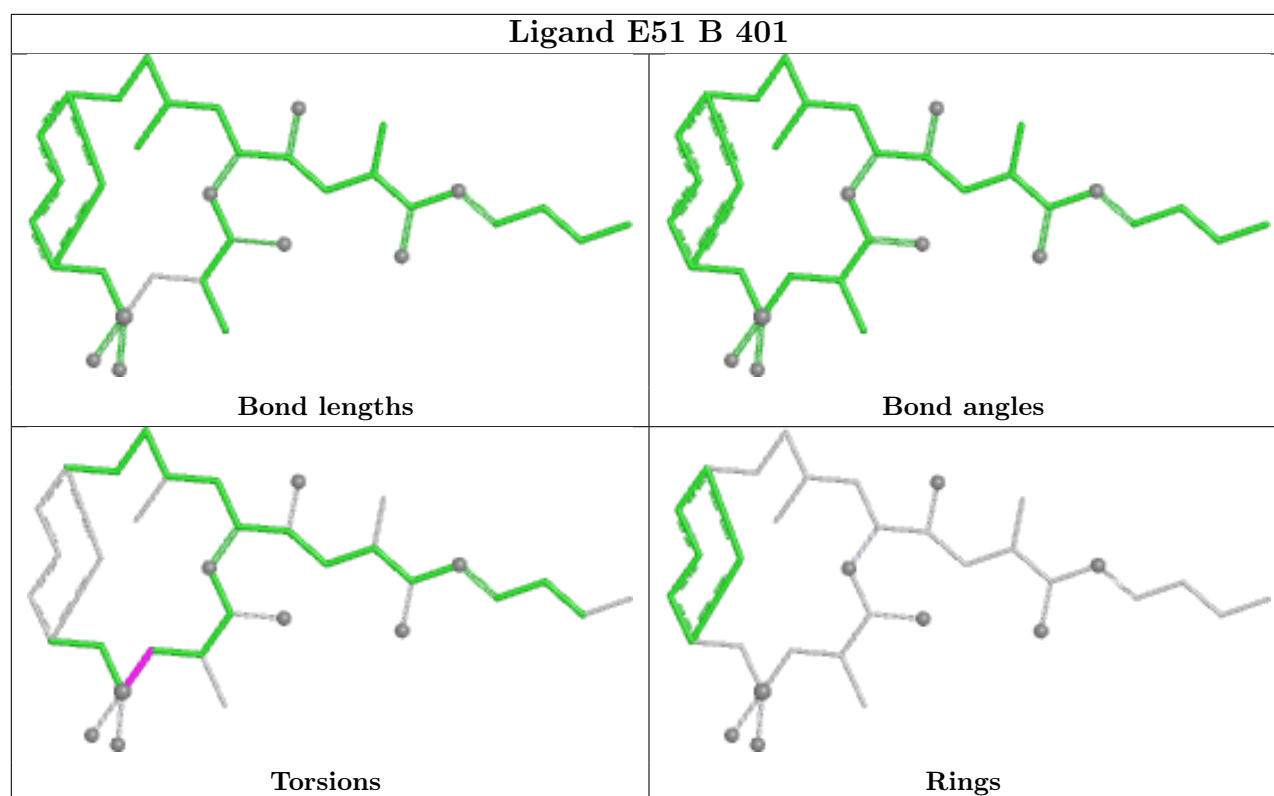
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

Ligand E51 A 401



Ligand E51 C 401





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	376/402 (93%)	0.20	29 (7%) 19 22	23, 36, 76, 148	0
1	B	377/402 (93%)	0.52	49 (12%) 7 8	24, 42, 93, 161	0
1	C	377/402 (93%)	0.18	25 (6%) 24 28	24, 37, 65, 91	0
All	All	1130/1206 (93%)	0.30	103 (9%) 15 17	23, 38, 77, 161	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	170	VAL	9.9
1	B	314	THR	6.2
1	A	312	VAL	6.2
1	B	312	VAL	5.2
1	B	313	ALA	4.9
1	B	260	GLY	4.6
1	C	361	VAL	4.6
1	A	157	ALA	4.6
1	C	68	TYR	4.5
1	A	361	VAL	4.4
1	A	314	THR	4.4
1	A	313	ALA	4.3
1	B	258	PRO	4.3
1	B	257	PHE	4.1
1	B	157	ALA	4.0
1	B	365	PHE	3.9
1	A	362	HIS	3.9
1	B	317	ASP	3.9
1	B	315	SER	3.8
1	A	267	LEU	3.7
1	C	169	SER	3.7
1	C	157	ALA	3.6
1	A	9	LYS	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	266	GLN	3.5
1	C	385	ASN	3.5
1	B	272	ALA	3.5
1	B	362	HIS	3.4
1	A	365	PHE	3.4
1	C	312	VAL	3.4
1	B	-2	SER	3.4
1	A	385	ASN	3.4
1	C	170	VAL	3.3
1	C	362	HIS	3.2
1	B	170	VAL	3.2
1	B	267	LEU	3.2
1	B	279	ILE	3.2
1	A	311	ASP	3.2
1	B	363	ASP	3.2
1	B	367	THR	3.2
1	B	361	VAL	3.1
1	A	360	HIS	2.9
1	C	365	PHE	2.9
1	B	319	CYS	2.9
1	C	272	ALA	2.9
1	B	336	VAL	2.8
1	C	384	TYR	2.8
1	C	271	GLN	2.8
1	B	49	HIS	2.8
1	B	205	ARG	2.7
1	A	156	GLY	2.7
1	C	379	MET	2.7
1	B	9	LYS	2.7
1	B	278	ASN	2.7
1	C	49	HIS	2.7
1	A	49	HIS	2.6
1	B	311	ASP	2.6
1	A	272	ALA	2.6
1	C	211	GLN	2.5
1	B	55	GLN	2.5
1	B	256	LYS	2.5
1	C	55	GLN	2.5
1	C	256	LYS	2.5
1	B	250	ALA	2.5
1	B	259	ASP	2.5
1	B	273	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	254	THR	2.4
1	A	315	SER	2.4
1	B	113	SER	2.4
1	B	251	ALA	2.4
1	B	309	VAL	2.4
1	C	-2	SER	2.4
1	B	263	LEU	2.4
1	C	156	GLY	2.4
1	B	169	SER	2.4
1	A	359	CYS	2.4
1	B	269	CYS	2.4
1	C	313	ALA	2.3
1	B	360	HIS	2.3
1	A	8	GLY	2.3
1	C	359	CYS	2.3
1	B	266	GLN	2.3
1	A	259	ASP	2.3
1	A	55	GLN	2.3
1	B	93	VAL	2.3
1	C	267	LEU	2.2
1	A	316	GLN	2.2
1	B	384	TYR	2.2
1	C	254	THR	2.2
1	A	269	CYS	2.2
1	B	112	GLY	2.2
1	B	253	SER	2.2
1	B	264	GLY	2.2
1	A	310	GLU	2.1
1	A	364	GLU	2.1
1	B	270	TRP	2.1
1	B	268	VAL	2.1
1	A	-2	SER	2.1
1	A	384	TYR	2.1
1	C	317	ASP	2.1
1	C	64	ARG	2.1
1	B	310	GLU	2.0
1	A	271	GLN	2.0
1	B	181	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [i](#)

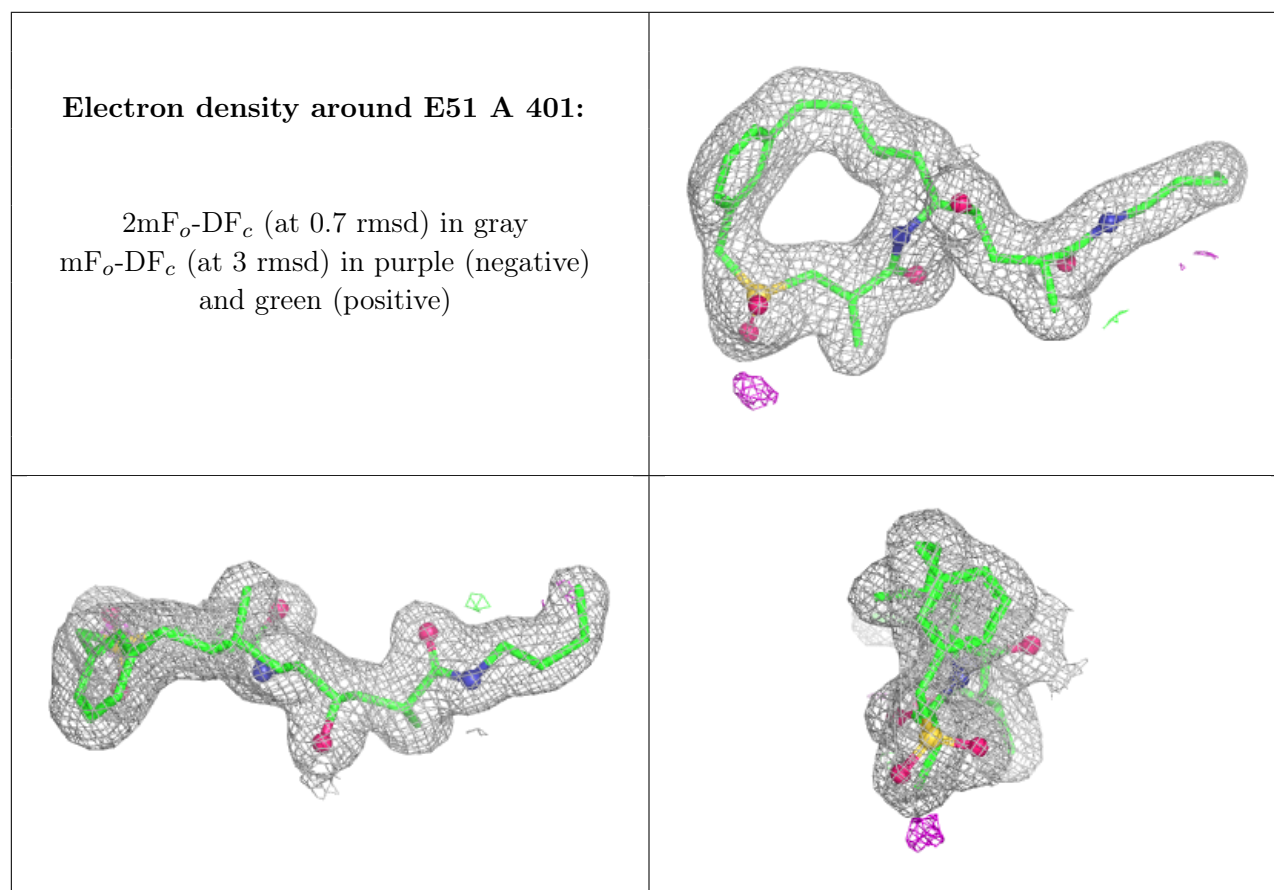
There are no oligosaccharides in this entry.

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.

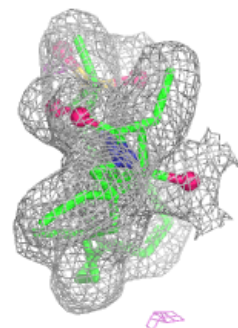
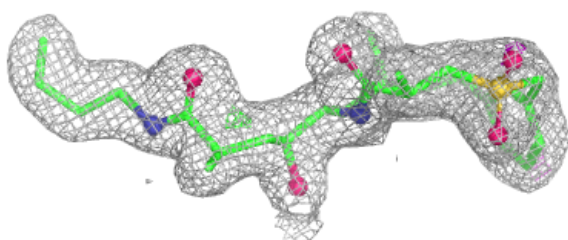
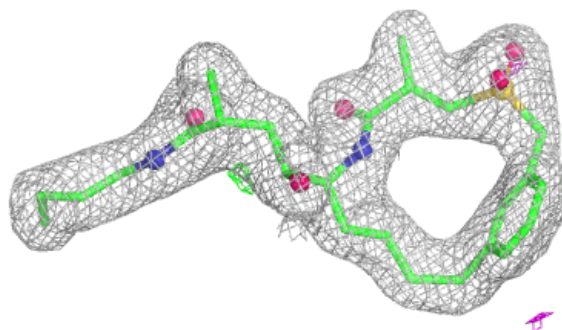
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	E51	A	401	34/34	0.97	0.07	26,31,40,47	0
2	E51	B	401	34/34	0.97	0.08	28,34,44,50	0
2	E51	C	401	34/34	0.98	0.06	25,30,39,41	0

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.

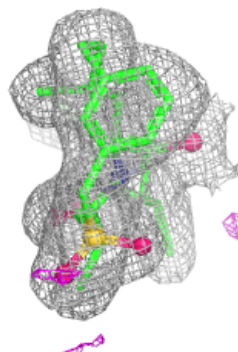
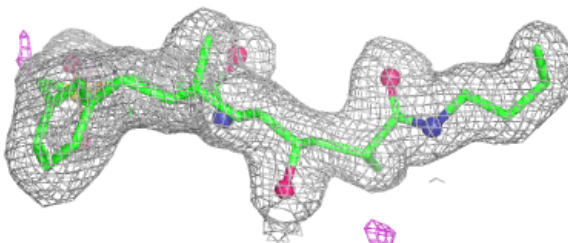
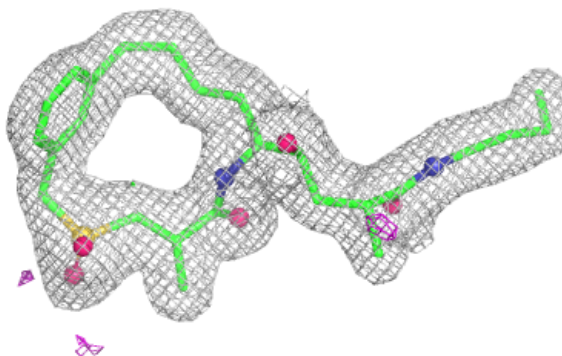


Electron density around E51 B 401:

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

**Electron density around E51 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.