



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 03:30 AM UTC

PDB ID : 6LDC / pdb_00006ldc
Title : Structure of Bifidobacterium dentium beta-glucuronidase complexed with C6-nonyl uronic isofagomine
Authors : Lin, H.-Y.; Hsieh, T.-J.; Lin, C.-H.
Deposited on : 2019-11-20
Resolution : 2.18 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

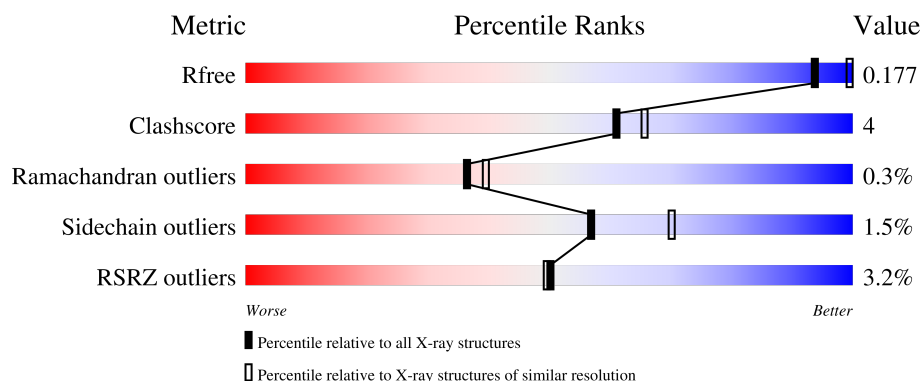
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

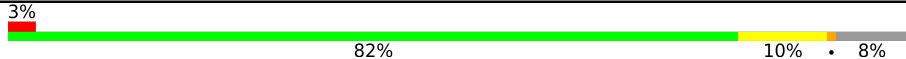
The reported resolution of this entry is 2.18 Å.

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	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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	670	
1	B	670	
1	C	670	
1	D	670	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21662 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.

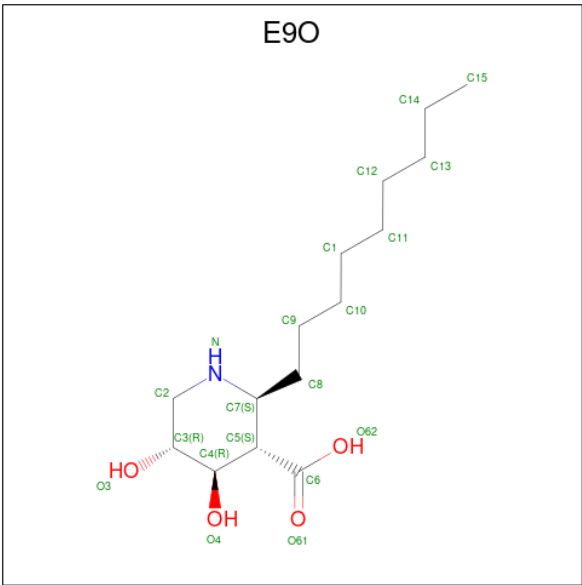
- Molecule 1 is a protein called LacZ1 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	617	Total	C	N	O	S	0	0	0
			4888	3102	852	915	19			
1	B	617	Total	C	N	O	S	0	0	0
			4888	3102	852	915	19			
1	C	617	Total	C	N	O	S	0	0	0
			4888	3102	852	915	19			
1	D	617	Total	C	N	O	S	0	0	0
			4888	3102	852	915	19			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ASN	-	expression tag	UNP D2Q7B1
A	0	GLY	-	expression tag	UNP D2Q7B1
B	-1	ASN	-	expression tag	UNP D2Q7B1
B	0	GLY	-	expression tag	UNP D2Q7B1
C	-1	ASN	-	expression tag	UNP D2Q7B1
C	0	GLY	-	expression tag	UNP D2Q7B1
D	-1	ASN	-	expression tag	UNP D2Q7B1
D	0	GLY	-	expression tag	UNP D2Q7B1

- Molecule 2 is (2 {S},3 {S},4 {R},5 {R})-2-nonyl-4,5-bis(oxidanyl)piperidine-3-carboxylic acid (CCD ID: E9O) (formula: C₁₅H₂₉NO₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			20	15	1	4		
2	B	1	Total	C	N	O	0	0
			20	15	1	4		
2	C	1	Total	C	N	O	0	0
			20	15	1	4		
2	D	1	Total	C	N	O	0	0
			20	15	1	4		

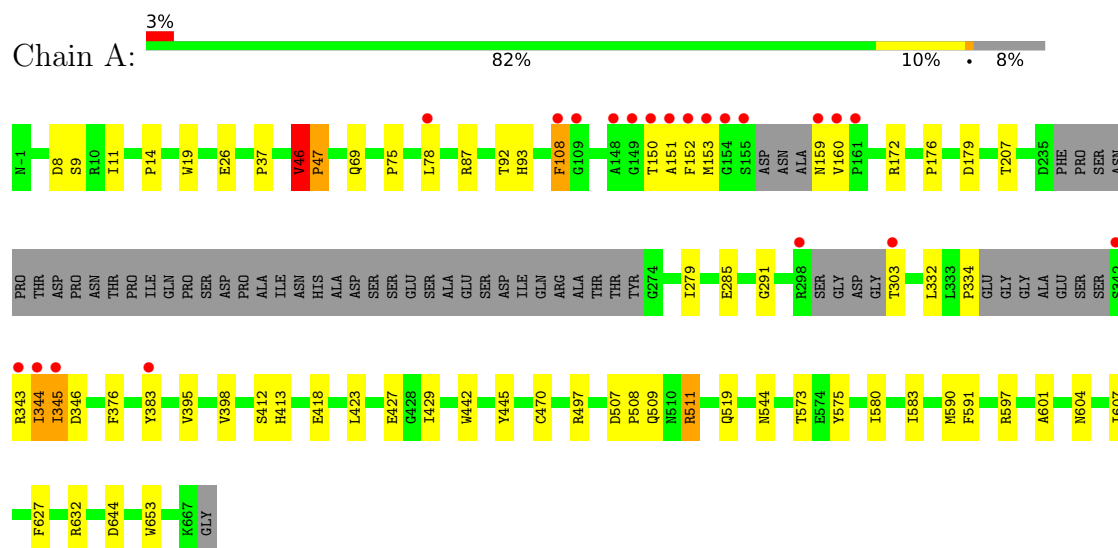
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	553	Total	O	0	0
			553	553		
3	B	462	Total	O	0	0
			462	462		
3	C	574	Total	O	0	0
			574	574		
3	D	441	Total	O	0	0
			441	441		

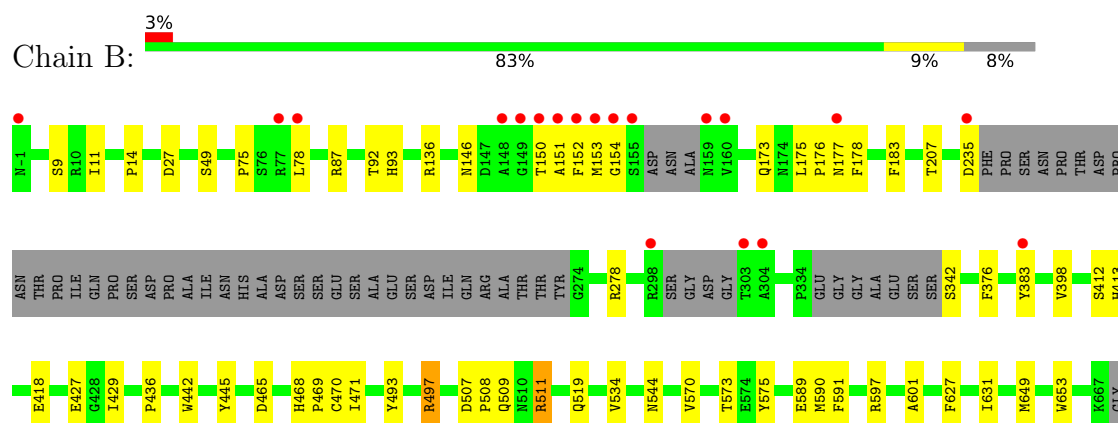
3 Residue-property plots [i](#)

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.

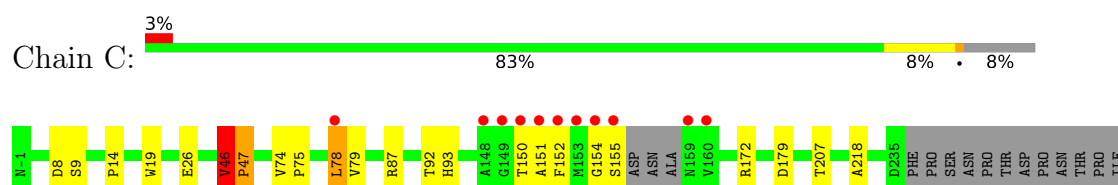
• Molecule 1: LacZ1 Beta-galactosidase

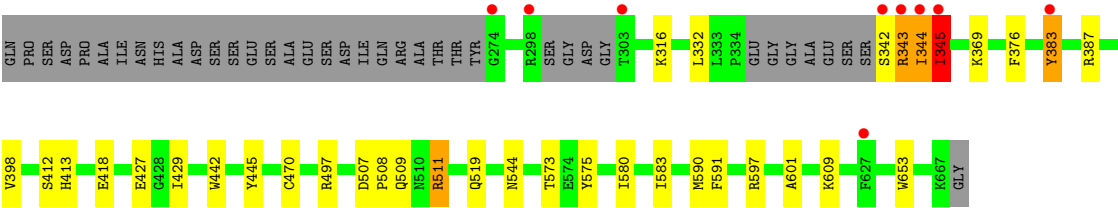


• Molecule 1: LacZ1 Beta-galactosidase

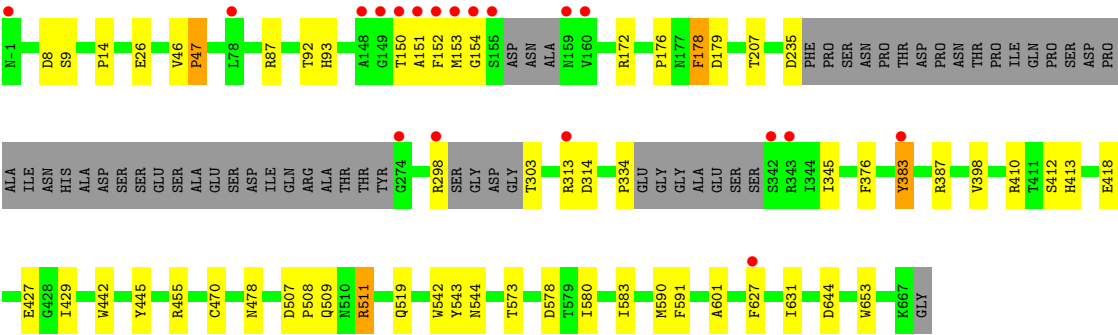
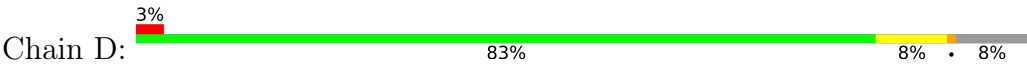


• Molecule 1: LacZ1 Beta-galactosidase





● Molecule 1: LacZ1 Beta-galactosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.66Å 105.14Å 160.97Å 90.00° 91.09° 90.00°	Depositor
Resolution (Å)	28.06 – 2.18 28.06 – 2.18	Depositor EDS
% Data completeness (in resolution range)	98.6 (28.06-2.18) 98.5 (28.06-2.18)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.61 (at 2.18Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.150 , 0.177 0.150 , 0.177	Depositor DCC
R_{free} test set	2001 reflections (1.24%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21662	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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: E9O

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	0.46	0/5018	0.77	5/6825 (0.1%)
1	B	0.45	0/5018	0.78	4/6825 (0.1%)
1	C	0.45	0/5018	0.76	2/6825 (0.0%)
1	D	0.41	0/5018	0.76	4/6825 (0.1%)
All	All	0.44	0/20072	0.77	15/27300 (0.1%)

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	0	2
1	C	0	2
1	D	0	1
All	All	0	5

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	383	TYR	N-CA-C	9.35	122.64	111.33
1	C	383	TYR	N-CA-C	8.90	122.09	111.33
1	B	383	TYR	N-CA-C	8.74	121.91	111.33
1	D	383	TYR	N-CA-C	8.64	121.79	111.33
1	D	47	PRO	N-CA-C	-6.24	99.61	112.47
1	A	47	PRO	N-CA-C	-6.18	99.73	112.47
1	A	108	PHE	N-CA-C	5.71	120.90	113.88
1	A	160	VAL	CA-C-N	5.64	125.32	119.05
1	A	160	VAL	C-N-CA	5.64	125.32	119.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	LEU	CA-C-N	-5.59	114.03	119.90
1	B	175	LEU	C-N-CA	-5.59	114.03	119.90
1	C	47	PRO	N-CA-C	-5.32	101.51	112.47
1	B	178	PHE	N-CA-C	5.27	117.08	109.07
1	D	178	PHE	CA-C-N	-5.18	114.67	122.24
1	D	178	PHE	C-N-CA	-5.18	114.67	122.24

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	46	VAL	Peptide,Mainchain
1	C	46	VAL	Peptide,Mainchain
1	D	46	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	4888	0	4643	49	0
1	B	4888	0	4643	38	0
1	C	4888	0	4643	37	0
1	D	4888	0	4643	42	0
2	A	20	0	0	0	0
2	B	20	0	0	0	0
2	C	20	0	0	0	0
2	D	20	0	0	0	0
3	A	553	0	0	11	0
3	B	462	0	0	6	0
3	C	574	0	0	6	0
3	D	441	0	0	4	0
All	All	21662	0	18572	160	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 4.

All (160) 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:B:11:ILE:HG21	1:B:78:LEU:HD11	1.43	0.99
1:A:344:ILE:HG13	1:A:345:ILE:HG22	1.46	0.97
1:A:11:ILE:HG21	1:A:78:LEU:HD11	1.45	0.95
1:C:344:ILE:HG13	1:C:345:ILE:HG22	1.50	0.93
1:A:627:PHE:O	3:A:801:HOH:O	1.86	0.92
1:D:152:PHE:CE1	1:D:179:ASP:HA	2.14	0.83
1:D:235:ASP:OD2	1:D:298:ARG:NH2	2.15	0.78
1:C:497:ARG:NH1	3:C:801:HOH:O	2.14	0.77
1:A:345:ILE:HG12	1:A:346:ASP:H	1.55	0.71
1:B:342:SER:N	3:B:802:HOH:O	2.22	0.70
1:D:303:THR:N	3:D:804:HOH:O	2.24	0.70
1:B:176:PRO:HG3	1:B:627:PHE:CD2	2.26	0.70
1:C:152:PHE:CE2	1:C:179:ASP:HB2	2.27	0.69
1:A:344:ILE:HG23	1:A:345:ILE:H	1.57	0.69
1:A:632:ARG:HG3	3:A:801:HOH:O	1.97	0.65
1:D:152:PHE:HZ	3:D:819:HOH:O	1.78	0.65
1:A:285:GLU:OE1	3:A:802:HOH:O	2.13	0.65
1:A:176:PRO:HG3	1:A:627:PHE:CE2	2.33	0.64
1:A:159:ASN:O	3:A:803:HOH:O	2.15	0.63
1:A:644:ASP:OD1	3:A:804:HOH:O	2.15	0.62
1:C:597:ARG:NH2	3:C:803:HOH:O	2.26	0.62
1:A:398:VAL:HG11	1:A:427:GLU:HG3	1.82	0.62
1:A:509:GLN:HB3	1:A:511:ARG:HD2	1.82	0.61
1:D:87:ARG:NH2	1:D:418:GLU:OE2	2.33	0.61
1:A:176:PRO:HG2	3:A:1304:HOH:O	2.01	0.60
1:B:398:VAL:HG11	1:B:427:GLU:HG3	1.83	0.60
1:D:152:PHE:CE2	1:D:176:PRO:HG2	2.37	0.59
1:D:398:VAL:HG11	1:D:427:GLU:HG3	1.83	0.58
1:A:87:ARG:NH2	1:A:418:GLU:OE2	2.38	0.57
1:D:153:MET:HB3	1:D:631:ILE:HD11	1.84	0.57
1:D:334:PRO:HB3	1:D:345:ILE:HD11	1.85	0.57
1:D:507:ASP:OD2	1:D:511:ARG:HD3	2.05	0.57
1:B:87:ARG:NH2	1:B:418:GLU:OE2	2.37	0.56
1:A:497:ARG:NH1	3:A:807:HOH:O	2.31	0.56
1:C:507:ASP:OD2	1:C:511:ARG:HD3	2.05	0.56
1:C:609:LYS:NZ	3:C:804:HOH:O	2.27	0.56
1:B:519:GLN:NE2	1:B:544:ASN:HD21	2.03	0.56
1:C:172:ARG:NH2	3:C:809:HOH:O	2.35	0.55
1:D:152:PHE:CE2	1:D:178:PHE:O	2.59	0.55
1:B:589:GLU:OE1	3:B:801:HOH:O	2.17	0.55
1:D:152:PHE:HE2	1:D:176:PRO:HG2	1.72	0.55
1:D:509:GLN:HB3	1:D:511:ARG:HD2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:519:GLN:NE2	1:A:544:ASN:HD21	2.06	0.54
1:C:316:LYS:NZ	3:C:816:HOH:O	2.41	0.53
1:A:152:PHE:HZ	3:A:801:HOH:O	1.89	0.53
1:A:152:PHE:O	1:A:152:PHE:CG	2.62	0.52
1:A:172:ARG:NH2	3:A:817:HOH:O	2.41	0.52
1:A:152:PHE:HD1	1:A:153:MET:HE3	1.75	0.52
1:B:509:GLN:HB3	1:B:511:ARG:HD2	1.92	0.52
1:D:87:ARG:HH22	1:D:418:GLU:CD	2.17	0.52
1:D:207:THR:HG21	1:D:508:PRO:HB2	1.92	0.51
1:D:152:PHE:HE1	1:D:179:ASP:OD1	1.93	0.51
1:C:207:THR:HG21	1:C:508:PRO:HB2	1.93	0.51
1:B:597:ARG:NH2	3:B:814:HOH:O	2.37	0.51
1:C:580:ILE:HB	1:C:583:ILE:HD12	1.92	0.50
1:A:207:THR:HG21	1:A:508:PRO:HB2	1.93	0.50
1:A:26:GLU:OE1	1:A:172:ARG:HD3	2.12	0.50
1:A:507:ASP:OD2	1:A:511:ARG:HD3	2.12	0.49
1:B:27:ASP:OD1	1:B:136:ARG:NH1	2.35	0.49
1:B:173:GLN:NE2	3:B:803:HOH:O	2.25	0.49
1:D:150:THR:HG22	1:D:151:ALA:H	1.78	0.49
1:D:519:GLN:NE2	1:D:544:ASN:HD21	2.09	0.49
1:C:87:ARG:NH2	1:C:418:GLU:OE2	2.45	0.49
1:B:590:MET:O	1:B:591:PHE:HB2	2.13	0.49
1:A:75:PRO:HA	1:C:8:ASP:HA	1.94	0.49
1:A:343:ARG:HG2	1:A:344:ILE:N	2.28	0.49
1:C:509:GLN:HB3	1:C:511:ARG:HD2	1.94	0.49
1:A:344:ILE:HG22	3:A:820:HOH:O	2.14	0.48
1:C:332:LEU:O	1:C:345:ILE:HG21	2.13	0.48
1:D:26:GLU:OE1	1:D:172:ARG:HD3	2.14	0.48
1:D:627:PHE:HB3	3:D:1165:HOH:O	2.13	0.48
1:D:410:ARG:HH12	1:D:478:ASN:ND2	2.11	0.48
1:B:278:ARG:HG3	3:B:845:HOH:O	2.12	0.48
1:C:74:VAL:HB	1:C:78:LEU:HD22	1.95	0.48
1:C:152:PHE:HB3	3:C:1219:HOH:O	2.14	0.48
1:C:155:SER:HB2	1:D:543:TYR:O	2.14	0.47
1:C:342:SER:OG	1:C:343:ARG:N	2.47	0.47
1:D:455:ARG:NH2	3:D:823:HOH:O	2.45	0.47
1:D:150:THR:HG22	1:D:151:ALA:N	2.28	0.47
1:B:150:THR:CG2	1:B:154:GLY:HA3	2.44	0.47
1:B:497:ARG:HA	1:B:497:ARG:HD2	1.59	0.47
1:B:507:ASP:OD2	1:B:511:ARG:HD3	2.15	0.46
1:C:344:ILE:HG23	1:C:345:ILE:H	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:590:MET:O	1:D:591:PHE:HB2	2.14	0.46
1:C:92:THR:HA	1:C:93:HIS:HA	1.75	0.46
1:D:412:SER:HA	1:D:413:HIS:HA	1.69	0.46
1:B:601:ALA:HA	1:B:653:TRP:CH2	2.51	0.46
1:A:92:THR:HA	1:A:93:HIS:HA	1.69	0.46
1:B:412:SER:HA	1:B:413:HIS:HA	1.72	0.46
1:C:398:VAL:HG11	1:C:427:GLU:HG3	1.97	0.46
1:D:429:ILE:O	1:D:470:CYS:HB2	2.16	0.46
1:B:534:VAL:HB	1:B:570:VAL:HG22	1.97	0.45
1:B:150:THR:HG22	1:B:151:ALA:N	2.32	0.45
1:D:152:PHE:CD2	1:D:178:PHE:O	2.69	0.45
1:C:590:MET:O	1:C:591:PHE:HB2	2.16	0.45
1:A:345:ILE:CG1	1:A:346:ASP:H	2.27	0.45
1:B:11:ILE:HG21	1:B:78:LEU:CD1	2.31	0.45
1:C:19:TRP:CD1	1:C:46:VAL:HG13	2.51	0.45
1:D:152:PHE:HE2	1:D:176:PRO:CG	2.28	0.45
1:D:580:ILE:HB	1:D:583:ILE:HD12	1.99	0.45
1:A:597:ARG:NH2	3:A:813:HOH:O	2.35	0.45
1:A:150:THR:HG22	1:A:151:ALA:N	2.32	0.45
1:A:590:MET:O	1:A:591:PHE:HB2	2.17	0.45
1:A:87:ARG:HH22	1:A:418:GLU:CD	2.25	0.45
1:D:150:THR:HG22	1:D:154:GLY:HA3	1.97	0.45
1:D:313:ARG:NH1	1:D:314:ASP:OD2	2.50	0.44
1:B:87:ARG:HH22	1:B:418:GLU:CD	2.26	0.44
1:C:429:ILE:O	1:C:470:CYS:HB2	2.17	0.44
1:A:14:PRO:HG2	1:C:14:PRO:HG2	1.99	0.44
1:B:442:TRP:HA	1:B:445:TYR:CD2	2.53	0.44
1:A:152:PHE:CE1	1:A:179:ASP:OD1	2.71	0.43
1:C:150:THR:HG22	1:C:151:ALA:N	2.33	0.43
1:C:152:PHE:CZ	1:C:179:ASP:HB2	2.53	0.43
1:C:412:SER:HA	1:C:413:HIS:HA	1.72	0.43
1:C:601:ALA:HA	1:C:653:TRP:CH2	2.53	0.43
1:A:601:ALA:HA	1:A:653:TRP:CH2	2.53	0.43
1:D:92:THR:HA	1:D:93:HIS:HA	1.73	0.43
1:B:468:HIS:O	1:B:511:ARG:NH2	2.51	0.43
1:C:442:TRP:HA	1:C:445:TYR:CD2	2.54	0.43
1:A:412:SER:HA	1:A:413:HIS:HA	1.74	0.43
1:B:75:PRO:HA	1:D:8:ASP:HA	1.99	0.43
1:B:413:HIS:O	1:B:436:PRO:HA	2.18	0.43
3:B:1164:HOH:O	1:D:644:ASP:HB3	2.17	0.43
1:A:93:HIS:O	1:A:108:PHE:O	2.36	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:MET:HB3	1:B:631:ILE:HD11	2.02	0.42
1:B:429:ILE:O	1:B:470:CYS:HB2	2.19	0.42
1:D:152:PHE:HD2	1:D:176:PRO:O	2.02	0.42
1:C:218:ALA:HB2	1:C:369:LYS:HE2	2.00	0.42
1:D:442:TRP:HA	1:D:445:TYR:CD2	2.54	0.42
1:A:152:PHE:HE2	1:A:627:PHE:HB2	1.85	0.42
1:B:146:ASN:ND2	1:B:177:ASN:OD1	2.46	0.42
1:B:597:ARG:HG3	1:B:649:MET:HG2	2.01	0.42
1:A:8:ASP:HA	1:C:75:PRO:HA	2.00	0.42
1:A:19:TRP:CD1	1:A:46:VAL:HG13	2.53	0.42
1:A:37:PRO:HA	1:A:69:GLN:OE1	2.20	0.42
1:A:604:ASN:HA	1:A:607:ILE:HG12	2.01	0.42
1:D:26:GLU:OE1	1:D:172:ARG:NH1	2.53	0.42
1:A:395:VAL:HA	1:A:423:LEU:HD21	2.01	0.42
1:C:150:THR:HG22	1:C:154:GLY:HA3	2.02	0.42
1:B:465:ASP:HB3	1:B:471:ILE:CD1	2.49	0.42
1:D:542:TRP:CZ2	1:D:578:ASP:HB2	2.55	0.42
1:B:468:HIS:HA	1:B:469:PRO:HD2	1.95	0.41
1:C:332:LEU:O	1:C:345:ILE:CG2	2.68	0.41
1:A:332:LEU:O	1:A:345:ILE:HG23	2.19	0.41
1:A:442:TRP:HA	1:A:445:TYR:CD2	2.54	0.41
1:B:207:THR:HG21	1:B:508:PRO:HB2	2.02	0.41
1:C:26:GLU:OE1	1:C:172:ARG:HD3	2.20	0.41
1:C:78:LEU:HD23	1:C:79:VAL:HG13	2.03	0.41
1:D:601:ALA:HA	1:D:653:TRP:CH2	2.55	0.41
1:A:279:ILE:O	1:A:291:GLY:HA2	2.21	0.41
1:B:92:THR:HA	1:B:93:HIS:HA	1.71	0.41
1:A:334:PRO:HD3	1:A:345:ILE:HG21	2.02	0.41
1:A:429:ILE:O	1:A:470:CYS:HB2	2.20	0.41
1:B:49:SER:OG	1:B:183:PHE:HB3	2.21	0.41
1:B:176:PRO:HG3	1:B:627:PHE:CE2	2.56	0.41
1:D:410:ARG:HH12	1:D:478:ASN:HD22	1.68	0.41
1:B:14:PRO:HG2	1:D:14:PRO:HG2	2.02	0.41
1:A:580:ILE:HB	1:A:583:ILE:HD12	2.02	0.40
1:C:519:GLN:NE2	1:C:544:ASN:HD21	2.18	0.40
1:B:493:TYR:CZ	1:B:497:ARG:HG2	2.56	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	607/670 (91%)	584 (96%)	20 (3%)	3 (0%)	24	25
1	B	607/670 (91%)	588 (97%)	19 (3%)	0	100	100
1	C	607/670 (91%)	582 (96%)	21 (4%)	4 (1%)	18	17
1	D	607/670 (91%)	585 (96%)	21 (4%)	1 (0%)	43	49
All	All	2428/2680 (91%)	2339 (96%)	81 (3%)	8 (0%)	36	39

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	344	ILE
1	C	343	ARG
1	A	345	ILE
1	C	344	ILE
1	C	345	ILE
1	D	47	PRO
1	A	47	PRO
1	C	47	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	501/543 (92%)	494 (99%)	7 (1%)	59	72
1	B	501/543 (92%)	493 (98%)	8 (2%)	55	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	501/543 (92%)	491 (98%)	10 (2%)	48	61
1	D	501/543 (92%)	495 (99%)	6 (1%)	63	75
All	All	2004/2172 (92%)	1973 (98%)	31 (2%)	57	70

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	46	VAL
1	A	303	THR
1	A	376	PHE
1	A	511	ARG
1	A	573	THR
1	A	575	TYR
1	B	9	SER
1	B	152	PHE
1	B	235	ASP
1	B	376	PHE
1	B	497	ARG
1	B	511	ARG
1	B	573	THR
1	B	575	TYR
1	C	9	SER
1	C	46	VAL
1	C	78	LEU
1	C	345	ILE
1	C	376	PHE
1	C	383	TYR
1	C	387	ARG
1	C	511	ARG
1	C	573	THR
1	C	575	TYR
1	D	9	SER
1	D	376	PHE
1	D	383	TYR
1	D	387	ARG
1	D	511	ARG
1	D	573	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-1	ASN
1	A	350	GLN
1	A	544	ASN
1	B	187	ASN
1	B	350	GLN
1	B	405	HIS
1	B	544	ASN
1	C	-1	ASN
1	C	102	GLN
1	C	544	ASN
1	D	187	ASN
1	D	350	GLN
1	D	478	ASN
1	D	544	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](#)

4 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	E9O	C	701	-	19,20,20	0.58	0	20,25,25	1.49	4 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	E9O	A	701	-	19,20,20	0.58	0	20,25,25	1.58	4 (20%)
2	E9O	B	701	-	19,20,20	0.56	0	20,25,25	1.58	4 (20%)
2	E9O	D	701	-	19,20,20	0.52	0	20,25,25	1.50	4 (20%)

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	E9O	C	701	-	-	3/13/30/30	0/1/1/1
2	E9O	A	701	-	-	3/13/30/30	0/1/1/1
2	E9O	B	701	-	-	5/13/30/30	0/1/1/1
2	E9O	D	701	-	-	3/13/30/30	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	E9O	C2-N-C7	4.38	119.33	109.71
2	A	701	E9O	C2-N-C7	3.89	118.26	109.71
2	C	701	E9O	C2-N-C7	3.88	118.24	109.71
2	D	701	E9O	C2-N-C7	3.88	118.23	109.71
2	A	701	E9O	C9-C8-C7	-3.22	107.82	113.94
2	D	701	E9O	C9-C8-C7	-2.94	108.35	113.94
2	B	701	E9O	C9-C8-C7	-2.90	108.42	113.94
2	D	701	E9O	O62-C6-O61	-2.84	117.63	124.08
2	C	701	E9O	C9-C8-C7	-2.83	108.56	113.94
2	A	701	E9O	O62-C6-O61	-2.75	117.83	124.08
2	B	701	E9O	O62-C6-O61	-2.75	117.84	124.08
2	A	701	E9O	O62-C6-C5	2.63	121.47	113.91
2	D	701	E9O	O62-C6-C5	2.56	121.27	113.91
2	C	701	E9O	O62-C6-C5	2.52	121.17	113.91
2	C	701	E9O	O62-C6-O61	-2.46	118.50	124.08
2	B	701	E9O	O62-C6-C5	2.41	120.84	113.91

There are no chirality outliers.

All (14) torsion outliers are listed below:

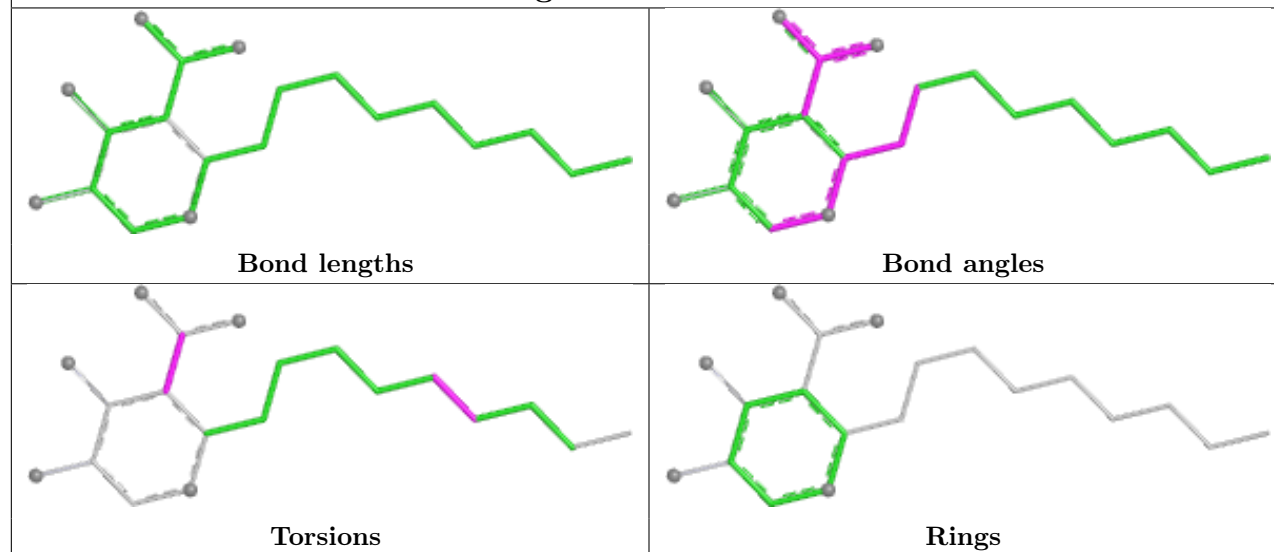
Mol	Chain	Res	Type	Atoms
2	D	701	E9O	C11-C12-C13-C14
2	A	701	E9O	C10-C1-C11-C12
2	B	701	E9O	C1-C10-C9-C8
2	C	701	E9O	C1-C11-C12-C13
2	B	701	E9O	C11-C1-C10-C9
2	B	701	E9O	C7-C8-C9-C10
2	A	701	E9O	C4-C5-C6-O62
2	A	701	E9O	C4-C5-C6-O61
2	B	701	E9O	C4-C5-C6-O61
2	C	701	E9O	C4-C5-C6-O61
2	D	701	E9O	C4-C5-C6-O61
2	B	701	E9O	C4-C5-C6-O62
2	C	701	E9O	C4-C5-C6-O62
2	D	701	E9O	C4-C5-C6-O62

There are no ring outliers.

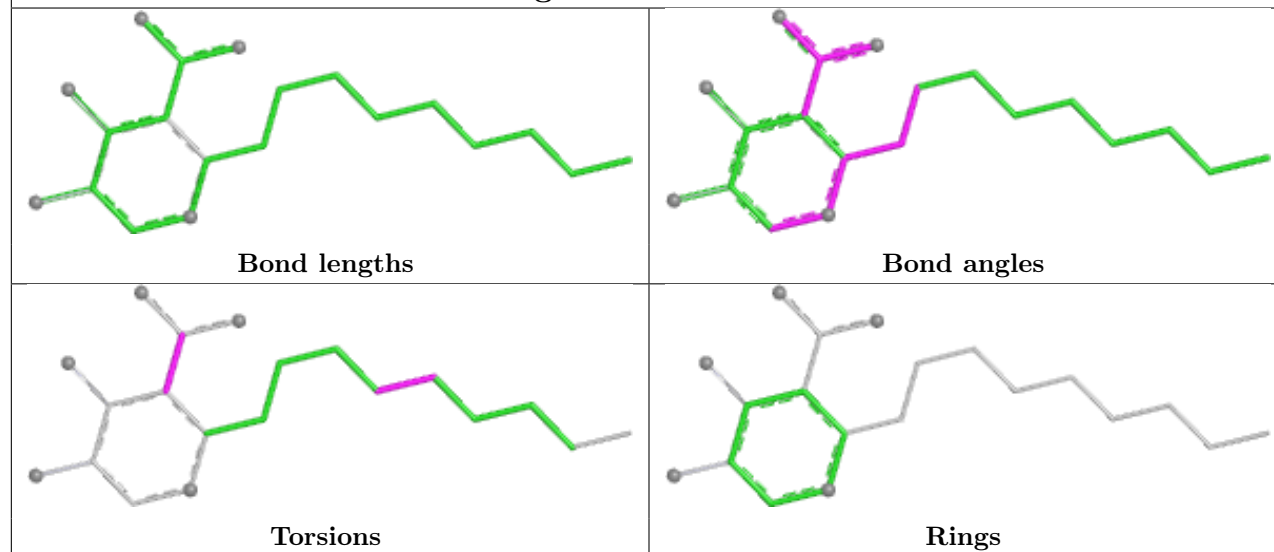
No monomer is involved in short contacts.

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.

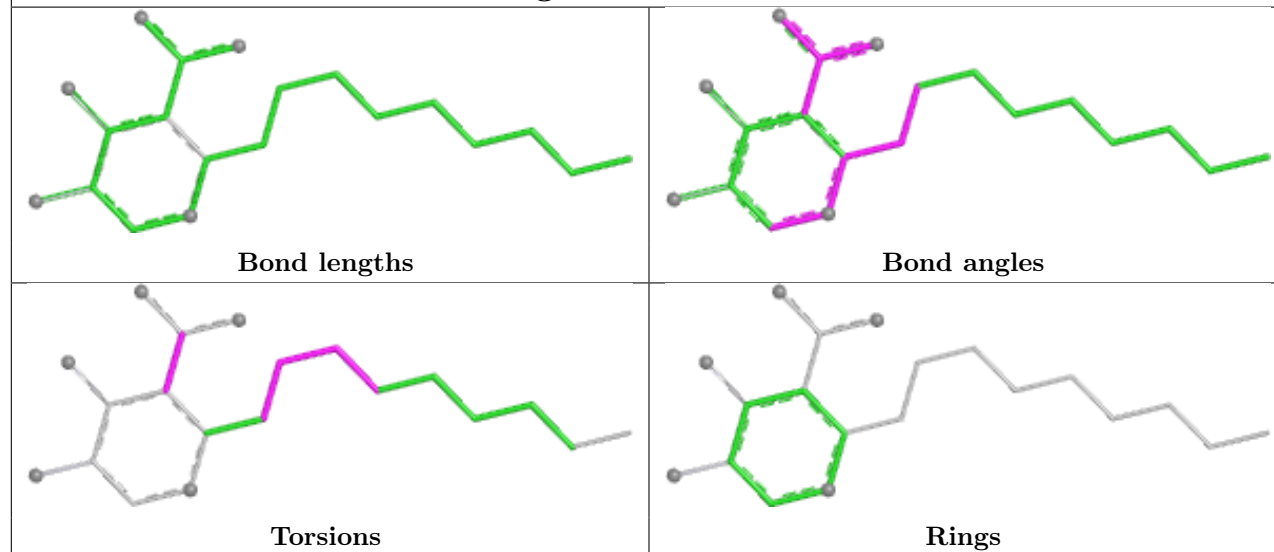
Ligand E9O C 701

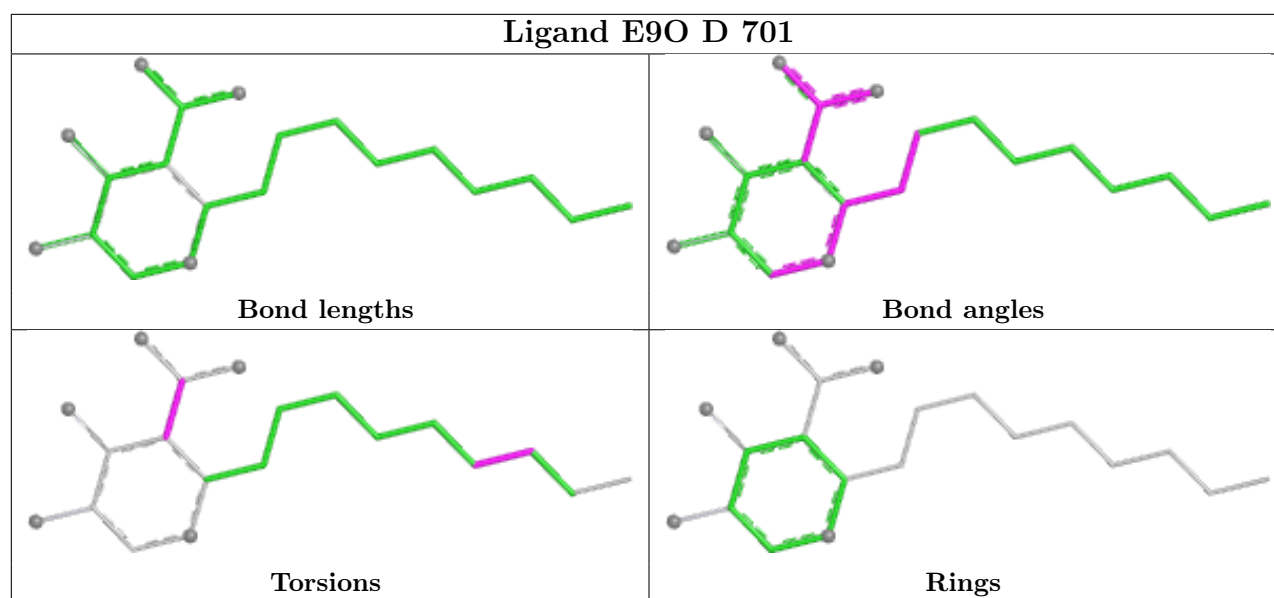


Ligand E9O A 701



Ligand E9O B 701





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	617/670 (92%)	-0.48	21 (3%)	48	47	6, 14, 37, 175	0
1	B	617/670 (92%)	-0.39	19 (3%)	51	51	8, 20, 41, 126	0
1	C	617/670 (92%)	-0.54	20 (3%)	50	49	6, 13, 37, 161	0
1	D	617/670 (92%)	-0.40	19 (3%)	51	51	7, 20, 45, 125	0
All	All	2468/2680 (92%)	-0.45	79 (3%)	50	49	6, 17, 40, 175	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	344	ILE	17.3
1	A	344	ILE	15.7
1	A	345	ILE	15.2
1	B	152	PHE	11.5
1	C	345	ILE	10.6
1	D	152	PHE	9.8
1	A	152	PHE	9.1
1	A	155	SER	8.8
1	C	152	PHE	8.6
1	B	151	ALA	7.8
1	A	151	ALA	7.6
1	B	155	SER	7.3
1	A	150	THR	7.3
1	A	153	MET	7.2
1	D	148	ALA	6.9
1	C	151	ALA	6.8
1	B	153	MET	6.7
1	A	160	VAL	6.7
1	D	155	SER	6.2
1	D	153	MET	6.0
1	B	148	ALA	5.8

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Mol	Chain	Res	Type	RSRZ
1	B	149	GLY	5.7
1	A	148	ALA	5.5
1	D	151	ALA	5.4
1	C	155	SER	5.2
1	C	160	VAL	5.2
1	C	148	ALA	5.2
1	A	149	GLY	5.0
1	A	154	GLY	5.0
1	A	342	SER	4.9
1	C	342	SER	4.9
1	C	150	THR	4.6
1	B	150	THR	4.5
1	D	150	THR	4.4
1	D	149	GLY	4.3
1	A	383	TYR	4.3
1	B	383	TYR	4.3
1	C	159	ASN	4.2
1	B	154	GLY	4.2
1	A	343	ARG	4.1
1	C	149	GLY	4.1
1	D	383	TYR	4.0
1	B	160	VAL	3.8
1	C	343	ARG	3.7
1	A	159	ASN	3.7
1	C	153	MET	3.7
1	D	342	SER	3.7
1	B	159	ASN	3.6
1	D	154	GLY	3.6
1	A	303	THR	3.5
1	B	177	ASN	3.5
1	A	109	GLY	3.4
1	C	383	TYR	3.1
1	D	78	LEU	3.1
1	D	159	ASN	2.9
1	B	235	ASP	2.9
1	A	161	PRO	2.9
1	C	154	GLY	2.9
1	C	627	PHE	2.8
1	D	298	ARG	2.8
1	D	343	ARG	2.8
1	A	108	PHE	2.8
1	C	298	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	160	VAL	2.8
1	D	627	PHE	2.7
1	C	78	LEU	2.7
1	D	-1	ASN	2.7
1	B	298	ARG	2.7
1	B	78	LEU	2.6
1	C	303	THR	2.6
1	B	77	ARG	2.6
1	A	78	LEU	2.4
1	B	304	ALA	2.3
1	C	274	GLY	2.2
1	B	303	THR	2.2
1	A	298	ARG	2.2
1	B	-1	ASN	2.1
1	D	313	ARG	2.1
1	D	274	GLY	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.

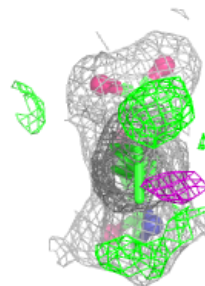
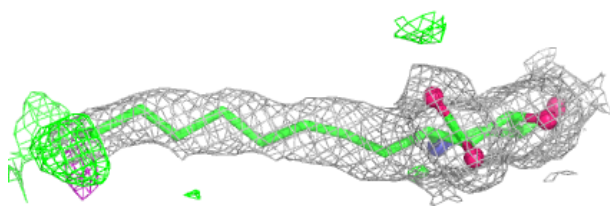
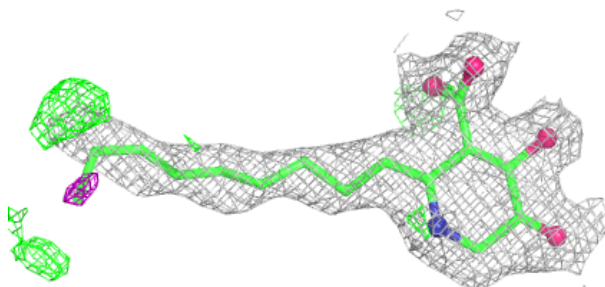
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	E9O	B	701	20/20	0.94	0.10	12,16,42,43	0
2	E9O	A	701	20/20	0.96	0.08	10,16,36,36	0
2	E9O	D	701	20/20	0.96	0.07	14,18,32,32	0
2	E9O	C	701	20/20	0.98	0.06	7,14,36,36	0

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.

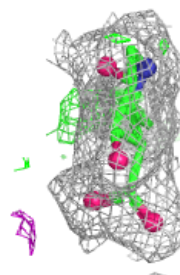
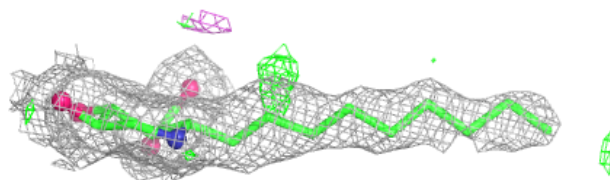
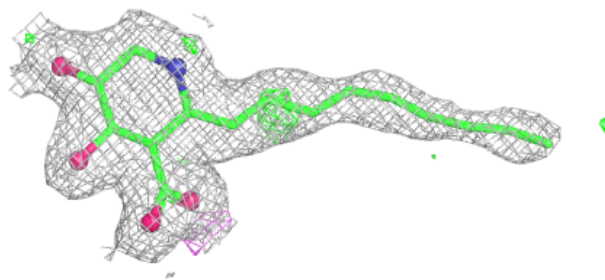
Electron density around E9O B 701:

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



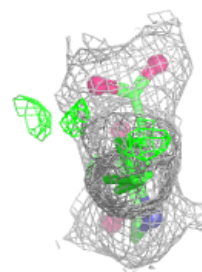
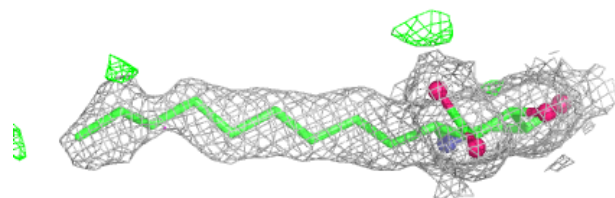
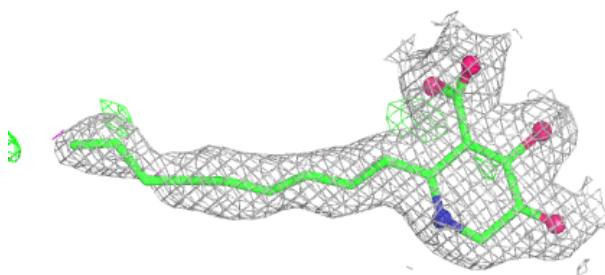
Electron density around E9O A 701:

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

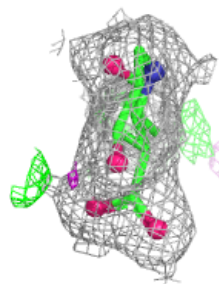
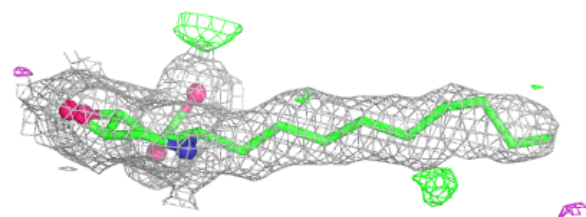
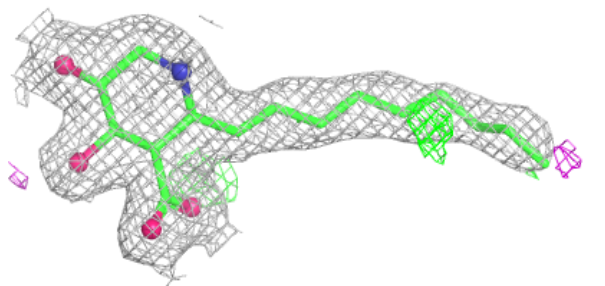


Electron density around E9O D 701:

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

**Electron density around E9O C 701:**

$2mF_o-DF_c$ (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.