



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:03 AM UTC

PDB ID : 4GBZ / pdb_00004gbz
Title : The structure of the MFS (major facilitator superfamily) proton:xylose symporter XylE bound to D-glucose
Authors : Sun, L.F.; Zeng, X.; Yan, C.Y.; Yan, N.
Deposited on : 2012-07-28
Resolution : 2.89 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

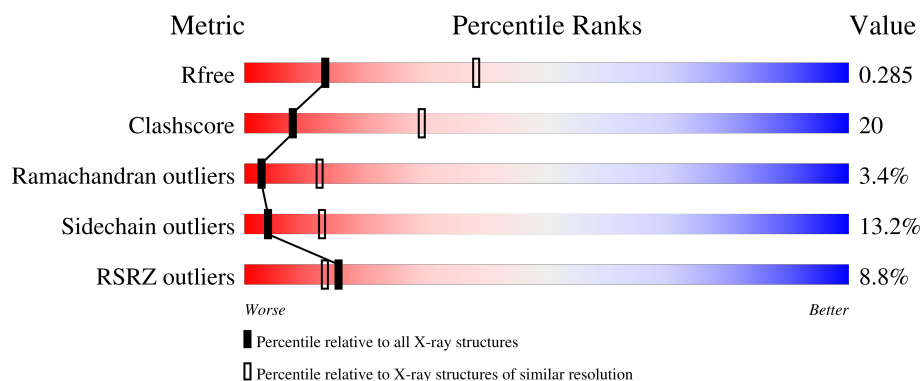
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	491	9% 60% 29% 7% ••

2 Entry composition [i](#)

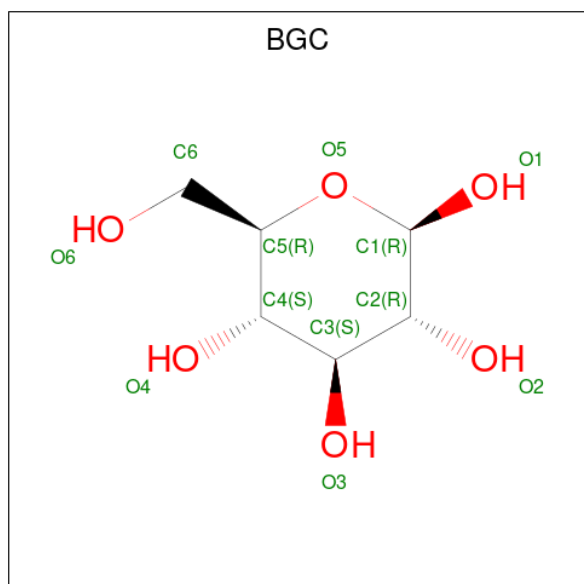
There are 4 unique types of molecules in this entry. The entry contains 3677 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-xylose-proton symporter.

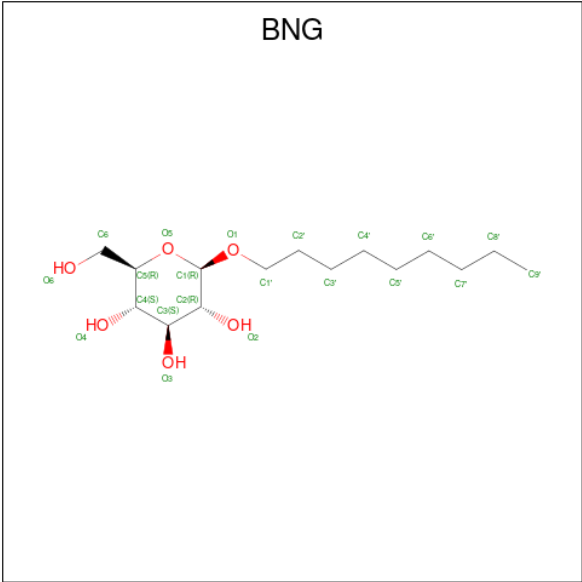
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	475	3579	2369	565	622	23	0	0	0

- Molecule 2 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	12	6	6	0	0

- Molecule 3 is nonyl beta-D-glucopyranoside (CCD ID: BNG) (formula: C₁₅H₃₀O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	15	6		
3	A	1	Total	C	O	0	0
			21	15	6		
3	A	1	Total	C	O	0	0
			21	15	6		
3	A	1	Total	C	O	0	0
			21	15	6		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		

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: D-xylose-proton symporter



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	95.04Å 95.04Å 168.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.52 – 2.89 38.52 – 2.89	Depositor EDS
% Data completeness (in resolution range)	98.1 (38.52-2.89) 98.0 (38.52-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.247 , 0.289 0.246 , 0.285	Depositor DCC
R_{free} test set	899 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	86.6	Xtriage
Anisotropy	0.456	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 67.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3677	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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: BGC, BNG

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	0.57	0/3672	0.97	9/5008 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	336	VAL	CB-CA-C	-6.87	103.18	111.97
1	A	208	ILE	CA-C-N	-6.52	113.08	119.87
1	A	208	ILE	C-N-CA	-6.52	113.08	119.87
1	A	30	VAL	CB-CA-C	-6.13	101.23	111.29
1	A	35	VAL	N-CA-C	5.85	116.50	110.36
1	A	335	THR	N-CA-C	5.80	120.48	113.17
1	A	367	ALA	CA-C-N	5.69	126.95	119.84
1	A	367	ALA	C-N-CA	5.69	126.95	119.84
1	A	29	ALA	N-CA-C	5.17	116.92	111.28

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	3579	0	3586	145	0
2	A	12	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	84	0	120	10	0
4	A	2	0	0	0	0
All	All	3677	0	3718	145	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 20.

All (145) 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:310:SER:CB	1:A:313:ILE:HB	1.72	1.18
1:A:310:SER:HB3	1:A:313:ILE:CB	1.79	1.12
1:A:372:ALA:HB3	3:A:505:BNG:H2'2	1.41	1.01
1:A:372:ALA:CB	3:A:505:BNG:H2'2	1.98	0.93
1:A:470:THR:HG22	1:A:473:GLU:H	1.38	0.88
1:A:182:ASN:OD1	1:A:202:MET:HG2	1.75	0.86
1:A:386:SER:HG	1:A:387:TRP:CD1	1.93	0.85
1:A:197:ASP:OD1	3:A:504:BNG:H61	1.83	0.77
1:A:421:PHE:HA	1:A:425:THR:HG23	1.67	0.75
1:A:310:SER:HB3	1:A:313:ILE:HB	0.84	0.73
1:A:244:MET:SD	1:A:249:ALA:HA	2.32	0.70
1:A:244:MET:HB2	1:A:248:LEU:HB3	1.74	0.70
1:A:310:SER:O	1:A:314:ALA:N	2.22	0.68
1:A:307:LEU:HD23	1:A:307:LEU:N	2.10	0.66
1:A:307:LEU:N	1:A:307:LEU:CD2	2.58	0.66
1:A:307:LEU:HD23	1:A:307:LEU:H	1.62	0.65
1:A:244:MET:HG3	1:A:249:ALA:HB2	1.79	0.64
1:A:234:GLU:CG	1:A:235:GLN:N	2.60	0.64
1:A:79:SER:O	1:A:83:GLY:HA2	1.98	0.64
1:A:307:LEU:HD11	3:A:505:BNG:O2	1.99	0.62
1:A:347:ILE:HG22	1:A:385:MET:HE1	1.81	0.61
1:A:332:ALA:O	1:A:336:VAL:HG22	2.01	0.61
1:A:157:ALA:HB1	1:A:469:LYS:O	2.02	0.60
1:A:234:GLU:HG2	1:A:235:GLN:N	2.16	0.60
1:A:193:TRP:HA	1:A:196:THR:HG22	1.84	0.60
1:A:116:ASN:O	1:A:116:ASN:ND2	2.31	0.59
1:A:470:THR:H	1:A:473:GLU:HB2	1.67	0.58
1:A:160:ARG:HE	1:A:397:GLU:HG2	1.67	0.58
1:A:441:ASN:ND2	3:A:503:BNG:O4	2.36	0.58
1:A:311:THR:HG22	1:A:312:ASP:N	2.16	0.58
1:A:315:LEU:HD22	1:A:319:ILE:HG13	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:VAL:HB	1:A:393:VAL:HG22	1.87	0.57
1:A:341:ARG:NE	1:A:397:GLU:OE1	2.28	0.57
1:A:249:ALA:O	1:A:252:ALA:HB3	2.06	0.56
1:A:390:VAL:O	1:A:394:LEU:HG	2.05	0.56
1:A:31:ILE:HD12	1:A:133:ARG:HD2	1.88	0.56
1:A:194:LEU:HA	1:A:198:GLY:HA3	1.88	0.55
1:A:437:ALA:HA	1:A:441:ASN:HB2	1.88	0.55
1:A:19:LEU:HB3	1:A:213:PHE:CD2	2.42	0.55
1:A:26:TYR:CE2	1:A:206:GLU:HB3	2.43	0.54
1:A:26:TYR:OH	1:A:205:SER:OG	2.25	0.54
1:A:122:LEU:HD11	1:A:200:ARG:NH2	2.23	0.54
1:A:234:GLU:HG2	1:A:235:GLN:H	1.70	0.54
1:A:404:ARG:NH1	1:A:471:LEU:HD13	2.23	0.54
1:A:367:ALA:O	3:A:505:BNG:H62	2.09	0.53
1:A:280:ILE:O	1:A:284:LEU:HB2	2.08	0.53
1:A:149:MET:HE1	1:A:405:GLY:HA2	1.91	0.52
1:A:307:LEU:CD1	3:A:505:BNG:O2	2.57	0.52
1:A:310:SER:O	1:A:311:THR:C	2.52	0.52
1:A:309:ALA:O	1:A:310:SER:CB	2.57	0.52
1:A:470:THR:HG22	1:A:473:GLU:N	2.18	0.52
1:A:241:ARG:C	1:A:243:ILE:H	2.18	0.52
1:A:349:ALA:HB3	1:A:456:ALA:HB2	1.92	0.51
1:A:386:SER:HG	1:A:387:TRP:HD1	1.51	0.51
1:A:386:SER:OG	1:A:387:TRP:CD1	2.62	0.51
1:A:206:GLU:HA	1:A:209:PRO:HD2	1.92	0.51
1:A:224:PRO:HA	1:A:227:LEU:HD12	1.92	0.51
1:A:426:PHE:HB3	1:A:427:PRO:HD3	1.93	0.51
1:A:342:LYS:O	1:A:346:ILE:HG13	2.11	0.51
1:A:345:GLN:HG3	1:A:459:MET:HE2	1.93	0.51
1:A:352:MET:HE1	1:A:383:PHE:HB2	1.93	0.50
1:A:309:ALA:O	1:A:310:SER:OG	2.29	0.50
1:A:241:ARG:C	1:A:243:ILE:N	2.68	0.50
1:A:245:GLY:O	1:A:249:ALA:HB3	2.11	0.50
1:A:330:VAL:HA	1:A:333:ILE:HD12	1.94	0.50
1:A:284:LEU:HD13	1:A:458:PHE:HE1	1.77	0.50
1:A:308:GLY:O	1:A:309:ALA:O	2.30	0.50
1:A:470:THR:HB	1:A:473:GLU:OE1	2.11	0.49
1:A:114:PRO:HG2	1:A:115:ASP:H	1.77	0.49
1:A:332:ALA:HB2	1:A:385:MET:HA	1.94	0.49
1:A:225:ARG:HD3	1:A:401:ASN:ND2	2.27	0.49
1:A:310:SER:CB	1:A:313:ILE:CB	2.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:VAL:O	1:A:144:SER:HB3	2.13	0.49
1:A:228:MET:HE1	1:A:257:LYS:HG3	1.95	0.49
1:A:288:GLN:NE2	2:A:501:BGC:O1	2.46	0.48
1:A:256:ILE:O	1:A:259:SER:HB3	2.14	0.48
1:A:329:THR:HA	1:A:384:ALA:O	2.15	0.47
1:A:372:ALA:HB2	3:A:505:BNG:O6	2.14	0.47
1:A:352:MET:HB2	1:A:379:TYR:CD2	2.49	0.47
1:A:468:GLY:O	1:A:469:LYS:O	2.31	0.47
1:A:315:LEU:HD22	1:A:319:ILE:CD1	2.44	0.47
1:A:357:PHE:O	1:A:361:THR:HG23	2.14	0.47
1:A:244:MET:HB2	1:A:248:LEU:HD13	1.97	0.47
1:A:244:MET:HE2	1:A:248:LEU:HD13	1.97	0.47
1:A:85:ARG:HD3	1:A:89:LYS:HE3	1.97	0.47
1:A:175:GLN:O	1:A:178:VAL:HG23	2.16	0.46
1:A:399:PHE:O	1:A:404:ARG:NH2	2.46	0.46
1:A:404:ARG:O	1:A:408:LEU:HB2	2.14	0.46
1:A:352:MET:HE2	1:A:379:TYR:CE2	2.50	0.46
1:A:149:MET:HE3	1:A:404:ARG:HG2	1.98	0.46
1:A:114:PRO:C	1:A:116:ASN:H	2.23	0.45
1:A:223:SER:HA	1:A:224:PRO:HD2	1.79	0.45
1:A:244:MET:CB	1:A:248:LEU:HD13	2.46	0.45
1:A:44:ALA:C	1:A:46:GLN:H	2.24	0.45
1:A:202:MET:O	1:A:205:SER:HB3	2.17	0.45
1:A:31:ILE:HD13	1:A:61:VAL:HA	1.98	0.45
1:A:154:LEU:HD11	1:A:223:SER:HA	1.98	0.45
1:A:415:GLN:NE2	2:A:501:BGC:O2	2.49	0.45
1:A:469:LYS:HG2	1:A:473:GLU:HB3	1.99	0.45
1:A:240:LEU:O	1:A:243:ILE:HB	2.15	0.45
1:A:399:PHE:O	1:A:404:ARG:NE	2.48	0.45
1:A:26:TYR:HE1	1:A:177:LEU:HD13	1.82	0.45
1:A:367:ALA:HB1	1:A:368:PRO:HD2	1.98	0.45
1:A:248:LEU:HD22	1:A:248:LEU:O	2.17	0.44
1:A:122:LEU:HD11	1:A:200:ARG:HH22	1.82	0.44
1:A:84:ARG:H	1:A:84:ARG:CD	2.31	0.44
1:A:107:LEU:HD22	3:A:502:BNG:H3'2	2.00	0.44
1:A:43:VAL:HG21	1:A:57:LEU:HB2	2.00	0.44
1:A:24:PHE:CD1	1:A:24:PHE:C	2.96	0.44
1:A:308:GLY:O	1:A:309:ALA:C	2.60	0.44
1:A:315:LEU:HD22	1:A:319:ILE:CG1	2.48	0.44
1:A:26:TYR:CZ	1:A:178:VAL:HG13	2.52	0.44
1:A:160:ARG:O	1:A:164:VAL:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:ARG:O	1:A:243:ILE:N	2.51	0.43
1:A:404:ARG:HH12	1:A:471:LEU:HD13	1.82	0.43
1:A:469:LYS:NZ	1:A:477:LEU:HD11	2.34	0.43
1:A:244:MET:HB2	1:A:248:LEU:CD1	2.48	0.43
1:A:118:VAL:HA	1:A:119:PRO:HD3	1.71	0.43
1:A:408:LEU:HA	1:A:408:LEU:HD23	1.75	0.43
1:A:87:SER:HA	1:A:90:ILE:HD12	2.00	0.43
1:A:248:LEU:HD22	1:A:248:LEU:C	2.43	0.43
1:A:404:ARG:HH22	1:A:471:LEU:HB3	1.84	0.43
1:A:470:THR:HG23	1:A:472:GLU:HG2	2.00	0.42
1:A:474:LEU:O	1:A:476:ALA:N	2.53	0.42
1:A:65:LEU:HA	1:A:68:CYS:HB2	2.01	0.42
1:A:81:ARG:HD2	1:A:82:PHE:CZ	2.55	0.42
1:A:197:ASP:OD1	3:A:504:BNB:C6	2.60	0.42
1:A:26:TYR:CE2	1:A:178:VAL:HG13	2.54	0.42
1:A:177:LEU:O	1:A:181:VAL:HG23	2.20	0.42
1:A:243:ILE:O	1:A:243:ILE:HG22	2.19	0.41
1:A:344:LEU:HD23	1:A:344:LEU:HA	1.85	0.41
1:A:41:VAL:HG13	1:A:122:LEU:HB2	2.01	0.41
1:A:250:THR:O	1:A:251:GLN:C	2.63	0.41
1:A:468:GLY:C	1:A:469:LYS:O	2.63	0.41
1:A:310:SER:CB	1:A:313:ILE:CG2	2.99	0.41
1:A:428:MET:HE3	1:A:428:MET:HB3	1.92	0.41
1:A:463:VAL:HA	1:A:464:PRO:HD3	1.82	0.41
1:A:10:ILE:HD11	1:A:154:LEU:O	2.21	0.41
1:A:21:GLY:O	1:A:24:PHE:HB3	2.21	0.41
1:A:149:MET:HE1	1:A:405:GLY:CA	2.51	0.41
1:A:290:PHE:HB3	1:A:451:MET:HG3	2.03	0.41
1:A:173:PHE:O	1:A:177:LEU:HB2	2.20	0.40
1:A:68:CYS:HB3	1:A:416:TRP:CD2	2.56	0.40
1:A:328:PHE:HB2	1:A:384:ALA:HB3	2.02	0.40
1:A:467:LYS:C	1:A:469:LYS:H	2.30	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	473/491 (96%)	410 (87%)	47 (10%)	16 (3%)	3	12

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	107	LEU
1	A	156	PRO
1	A	307	LEU
1	A	309	ALA
1	A	310	SER
1	A	469	LYS
1	A	311	THR
1	A	430	ASP
1	A	475	GLU
1	A	114	PRO
1	A	242	LYS
1	A	368	PRO
1	A	84	ARG
1	A	389	PRO
1	A	397	GLU
1	A	45	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	365/396 (92%)	317 (87%)	48 (13%)	4	13

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ILE

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Mol	Chain	Res	Type
1	A	15	LEU
1	A	19	LEU
1	A	24	PHE
1	A	30	VAL
1	A	40	THR
1	A	46	GLN
1	A	49	SER
1	A	81	ARG
1	A	85	ARG
1	A	107	LEU
1	A	116	ASN
1	A	133	ARG
1	A	144	SER
1	A	145	MET
1	A	177	LEU
1	A	178	VAL
1	A	188	SER
1	A	190	ASP
1	A	196	THR
1	A	197	ASP
1	A	234	GLU
1	A	248	LEU
1	A	266	THR
1	A	278	ILE
1	A	279	VAL
1	A	283	MET
1	A	291	VAL
1	A	307	LEU
1	A	315	LEU
1	A	320	ILE
1	A	331	LEU
1	A	336	VAL
1	A	339	PHE
1	A	347	ILE
1	A	373	LEU
1	A	403	ILE
1	A	408	LEU
1	A	417	LEU
1	A	422	VAL
1	A	425	THR
1	A	432	ASN
1	A	457	LEU

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Mol	Chain	Res	Type
1	A	469	LYS
1	A	470	THR
1	A	472	GLU
1	A	475	GLU
1	A	477	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	GLN
1	A	288	GLN
1	A	289	GLN
1	A	294	ASN
1	A	415	GLN
1	A	441	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](#)

5 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
3	BNG	A	504	-	21,21,21	0.45	0	26,26,26	0.59	0
3	BNG	A	502	-	21,21,21	0.46	0	26,26,26	0.60	0
2	BGC	A	501	-	12,12,12	0.46	0	17,17,17	0.52	0
3	BNG	A	503	-	21,21,21	0.46	0	26,26,26	0.61	0
3	BNG	A	505	-	21,21,21	0.48	0	26,26,26	0.59	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
3	BNG	A	504	-	-	7/12/32/32	0/1/1/1
3	BNG	A	502	-	-	7/12/32/32	0/1/1/1
2	BGC	A	501	-	-	0/2/22/22	0/1/1/1
3	BNG	A	503	-	-	8/12/32/32	0/1/1/1
3	BNG	A	505	-	-	8/12/32/32	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	BNG	O5-C1-O1-C1'
3	A	503	BNG	O5-C1-O1-C1'
3	A	505	BNG	O5-C1-O1-C1'
3	A	505	BNG	O5-C5-C6-O6
3	A	505	BNG	O1-C1'-C2'-C3'
3	A	504	BNG	O5-C1-O1-C1'
3	A	502	BNG	C4-C5-C6-O6
3	A	503	BNG	C4'-C5'-C6'-C7'
3	A	503	BNG	C3'-C4'-C5'-C6'
3	A	503	BNG	C1'-C2'-C3'-C4'
3	A	502	BNG	C1'-C2'-C3'-C4'
3	A	503	BNG	O5-C5-C6-O6
3	A	505	BNG	C2'-C3'-C4'-C5'
3	A	505	BNG	C6'-C7'-C8'-C9'
3	A	504	BNG	C3'-C4'-C5'-C6'
3	A	504	BNG	C1'-C2'-C3'-C4'

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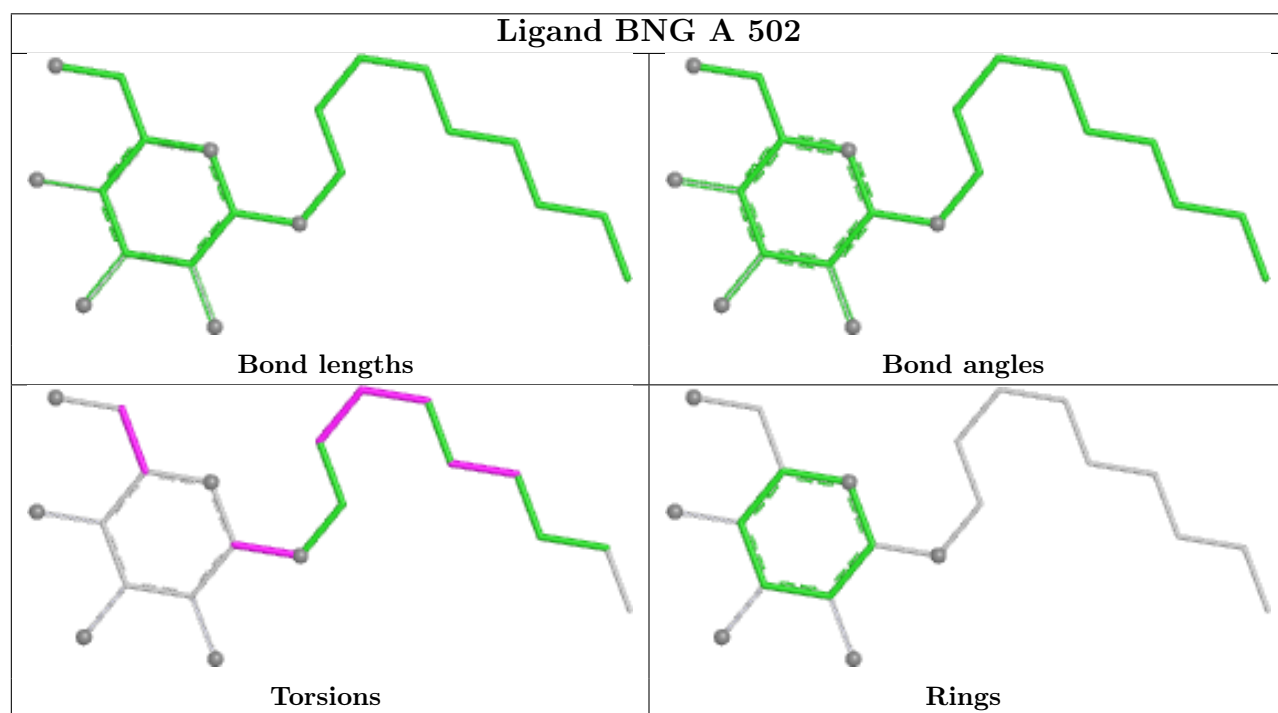
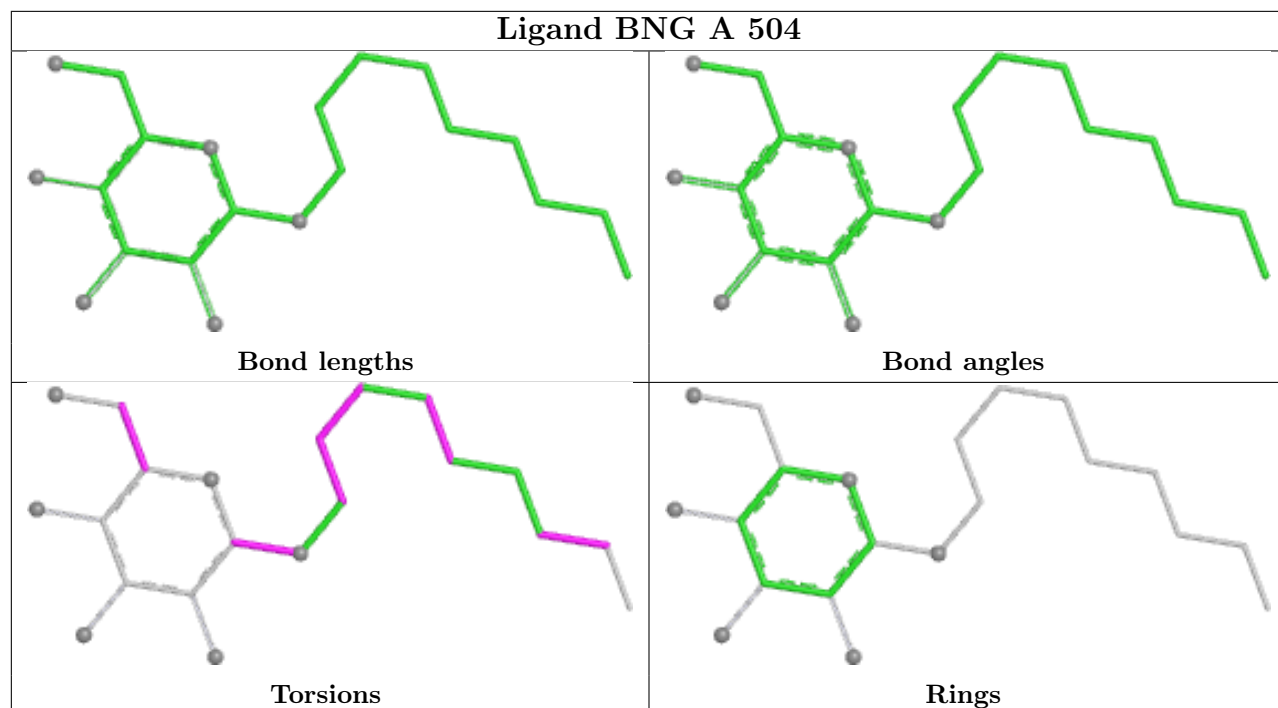
Mol	Chain	Res	Type	Atoms
3	A	503	BNG	C5'-C6'-C7'-C8'
3	A	502	BNG	C2-C1-O1-C1'
3	A	503	BNG	O1-C1'-C2'-C3'
3	A	504	BNG	C6'-C7'-C8'-C9'
3	A	502	BNG	O5-C5-C6-O6
3	A	505	BNG	C4-C5-C6-O6
3	A	502	BNG	C4'-C5'-C6'-C7'
3	A	502	BNG	C2'-C3'-C4'-C5'
3	A	504	BNG	C4-C5-C6-O6
3	A	503	BNG	C2'-C3'-C4'-C5'
3	A	505	BNG	C2'-C1'-O1-C1
3	A	504	BNG	O5-C5-C6-O6
3	A	504	BNG	O1-C1'-C2'-C3'
3	A	505	BNG	C4'-C5'-C6'-C7'

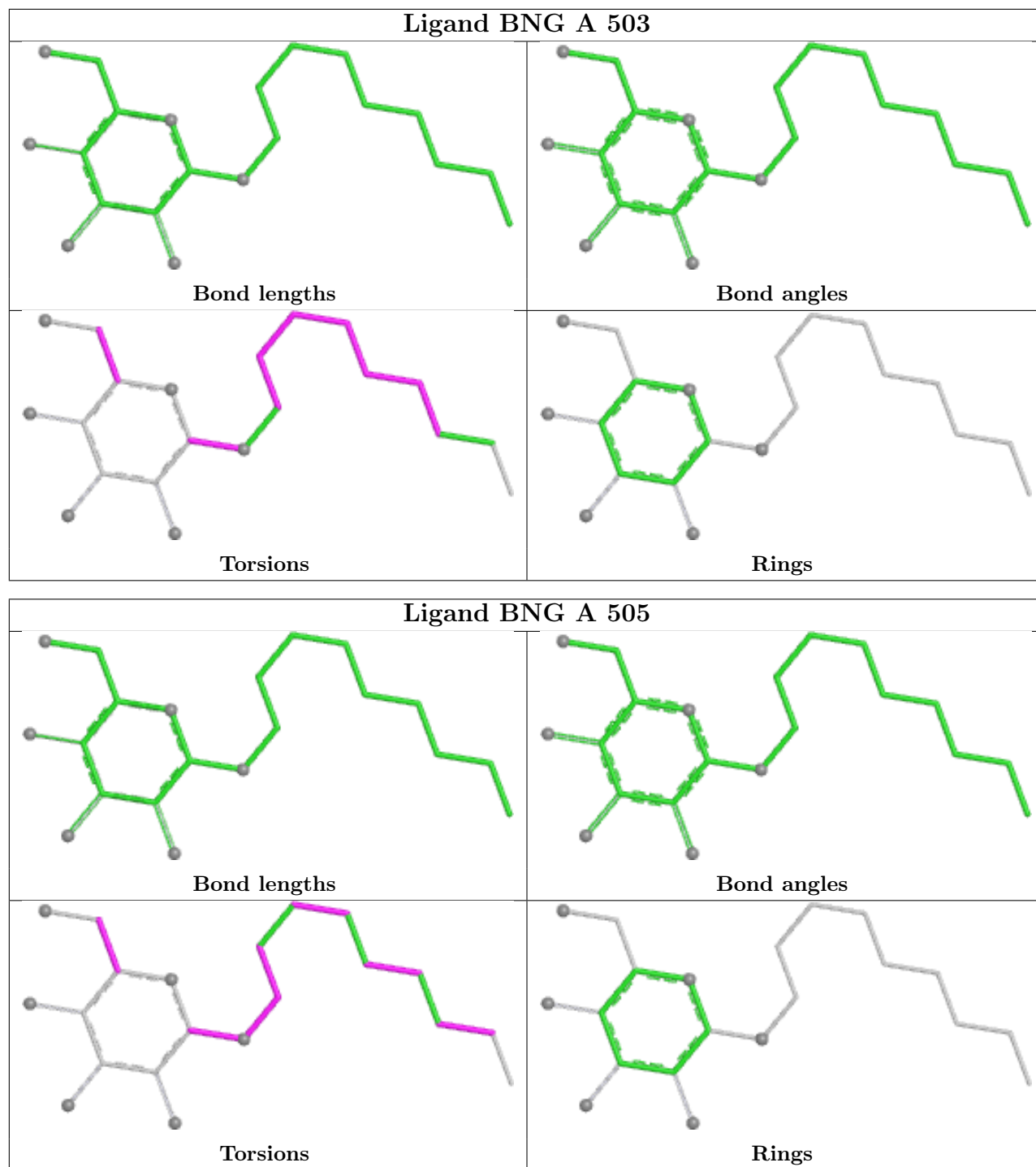
There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	504	BNG	2	0
3	A	502	BNG	1	0
2	A	501	BGC	2	0
3	A	503	BNG	1	0
3	A	505	BNG	6	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	475/491 (96%)	0.65	42 (8%) 15 13	50, 84, 162, 278	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	192	SER	6.6
1	A	271	LEU	4.4
1	A	334	MET	4.3
1	A	108	GLY	4.1
1	A	255	GLU	4.1
1	A	115	ASP	4.0
1	A	468	GLY	3.8
1	A	315	LEU	3.8
1	A	313	ILE	3.8
1	A	244	MET	3.7
1	A	186	ALA	3.5
1	A	229	SER	3.5
1	A	234	GLU	3.3
1	A	312	ASP	3.3
1	A	194	LEU	3.1
1	A	317	GLN	2.9
1	A	270	LEU	2.7
1	A	247	THR	2.7
1	A	471	LEU	2.6
1	A	222	GLU	2.5
1	A	441	ASN	2.5
1	A	107	LEU	2.5
1	A	191	ALA	2.5
1	A	261	ASP	2.4
1	A	343	PRO	2.4
1	A	249	ALA	2.3
1	A	449	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	201	TYR	2.3
1	A	248	LEU	2.3
1	A	319	ILE	2.2
1	A	390	VAL	2.2
1	A	54	ASN	2.2
1	A	117	THR	2.2
1	A	368	PRO	2.1
1	A	246	ASN	2.1
1	A	326	LEU	2.1
1	A	375	SER	2.1
1	A	212	LEU	2.1
1	A	417	LEU	2.1
1	A	100	VAL	2.1
1	A	371	VAL	2.0
1	A	479	GLU	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.

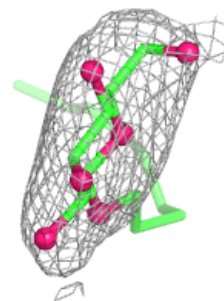
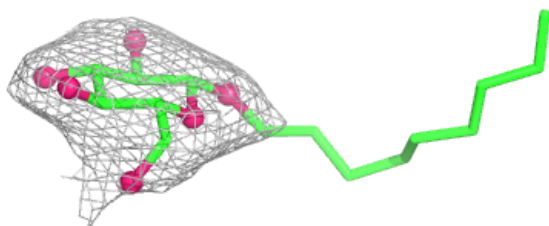
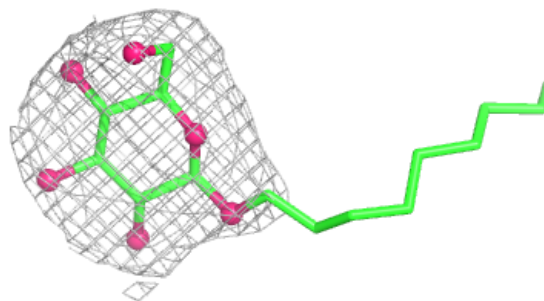
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BNG	A	502	21/21	0.80	0.18	144,151,156,157	0
3	BNG	A	503	21/21	0.82	0.21	134,136,140,142	0
2	BGC	A	501	12/12	0.91	0.16	74,75,78,80	0
3	BNG	A	504	21/21	0.92	0.17	117,128,136,138	0
3	BNG	A	505	21/21	0.92	0.16	108,121,122,123	0

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.

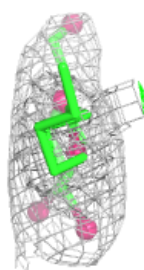
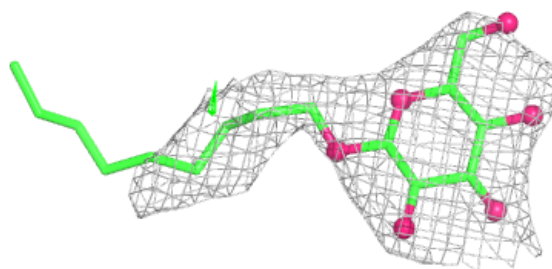
Electron density around BNG A 502:

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

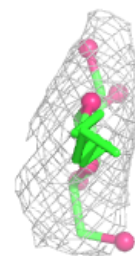
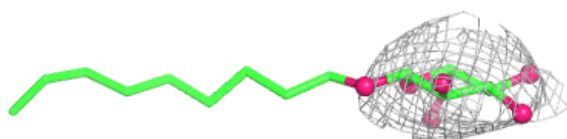
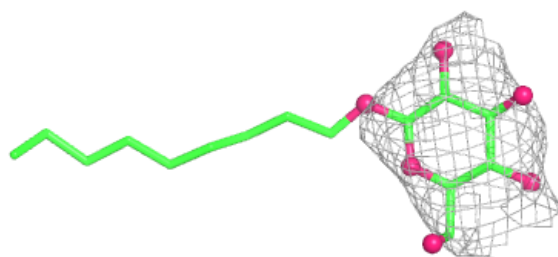


Electron density around BNG A 503:

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

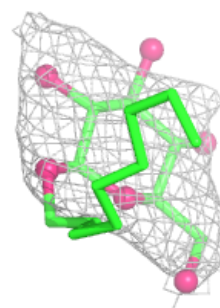
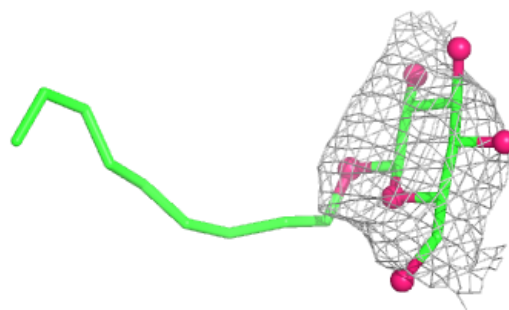
**Electron density around BNG A 504:**

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



Electron density around BNG A 505:

$2mF_o-DF_c$ (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.