



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:02 AM UTC

PDB ID : 6FL9 / pdb_00006fl9
Title : The active form of a pentameric ion channel (sTeLIC) gated by alkaline pH - Wild type 2.3 Angstrom resolution
Authors : Hu, H.; Delarue, M.
Deposited on : 2018-01-25
Resolution : 2.30 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

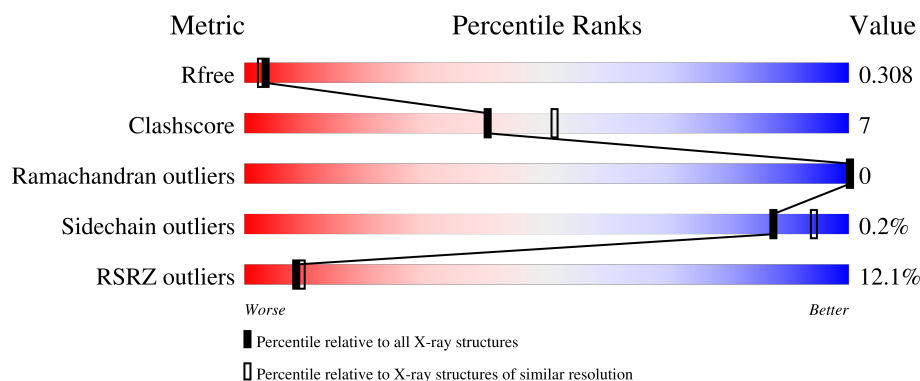
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	320	10% 84% 13% .
1	B	320	8% 83% 13% .
1	C	320	7% 82% 14% ..
1	D	320	12% 85% 12% .
1	E	320	21% 84% 12% ..

2 Entry composition [i](#)

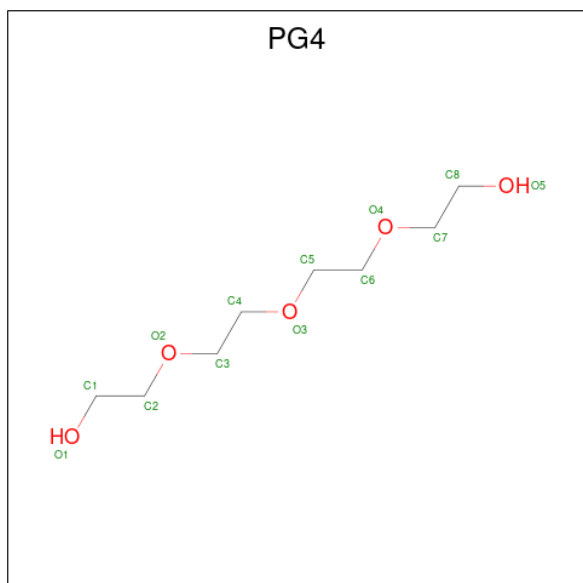
There are 4 unique types of molecules in this entry. The entry contains 13403 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cys-loop ligand-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	0	0	0
			2566	1690	423	448	5			
1	B	310	Total	C	N	O	S	0	0	0
			2566	1690	423	448	5			
1	C	310	Total	C	N	O	S	0	0	0
			2566	1690	423	448	5			
1	D	310	Total	C	N	O	S	0	0	0
			2566	1690	423	448	5			
1	E	310	Total	C	N	O	S	0	0	0
			2566	1690	423	448	5			

- Molecule 2 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



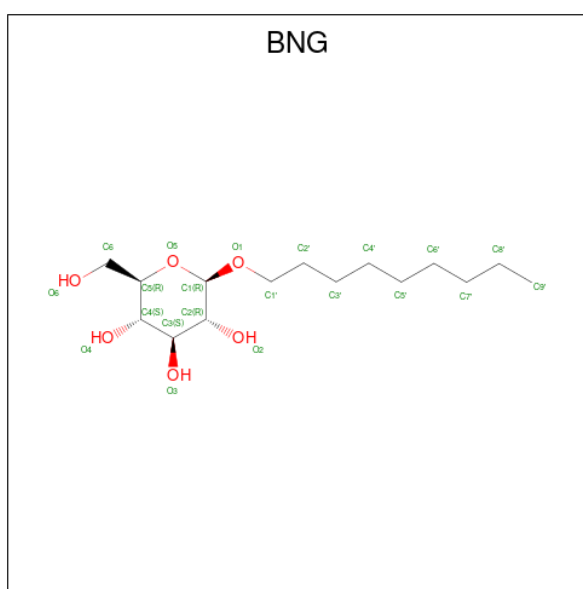
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	8	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	8	5		
2	C	1	Total	C	O	0	0
			13	8	5		
2	C	1	Total	C	O	0	0
			13	8	5		
2	E	1	Total	C	O	0	0
			13	8	5		

- 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	B	1	Total	C	O	0	0
			21	15	6		
3	C	1	Total	C	O	0	0
			21	15	6		
3	D	1	Total	C	O	0	0
			21	15	6		
3	E	1	Total	C	O	0	0
			21	15	6		

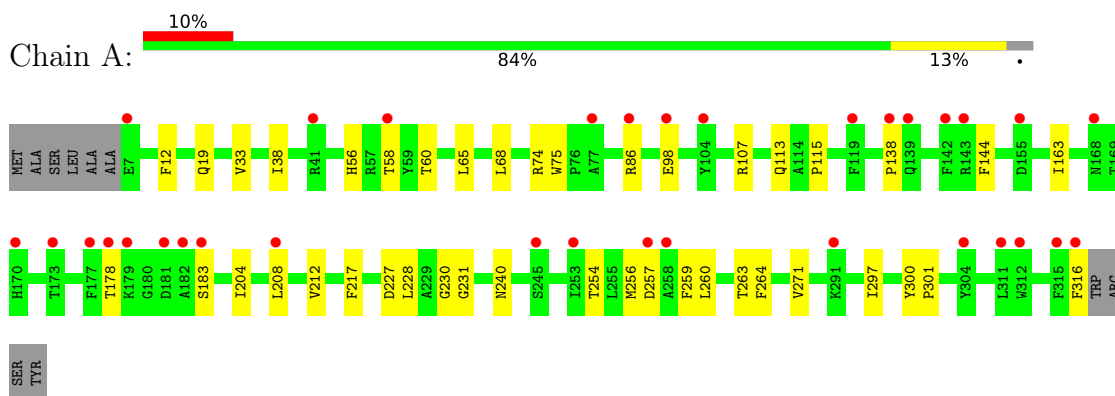
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	74	Total 74	O 74	0	0
4	B	99	Total 99	O 99	0	0
4	C	101	Total 101	O 101	0	0
4	D	88	Total 88	O 88	0	0
4	E	41	Total 41	O 41	0	0

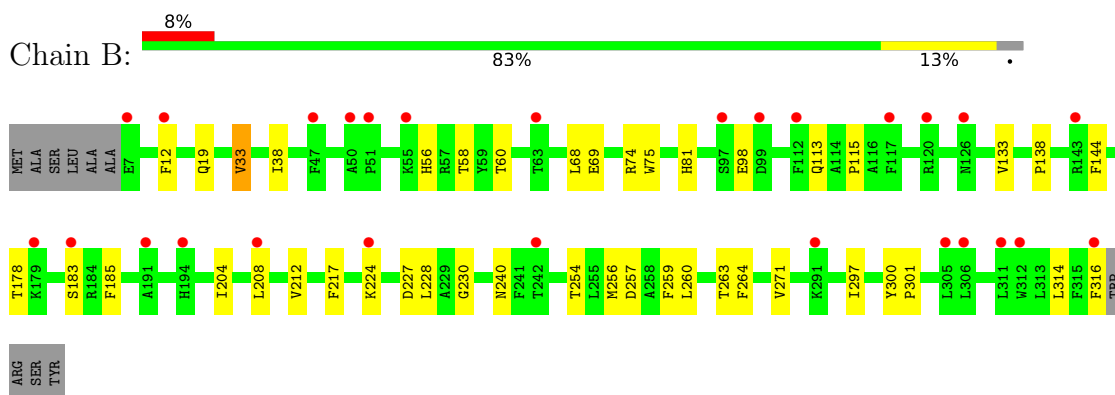
3 Residue-property plots [i](#)

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.

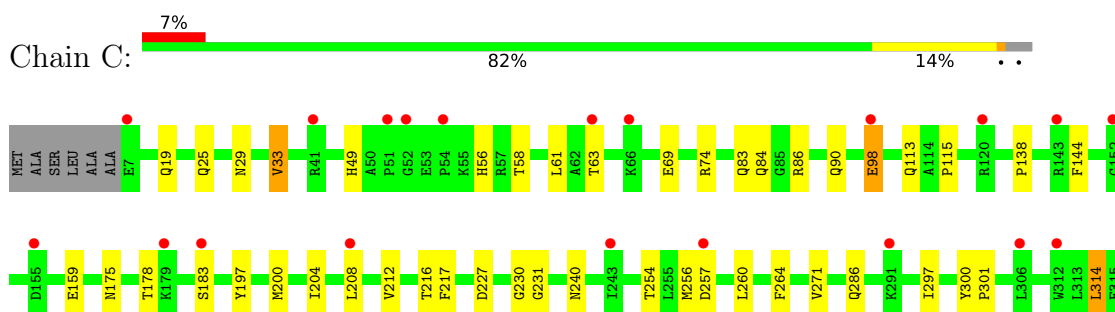
• Molecule 1: Cys-loop ligand-gated ion channel



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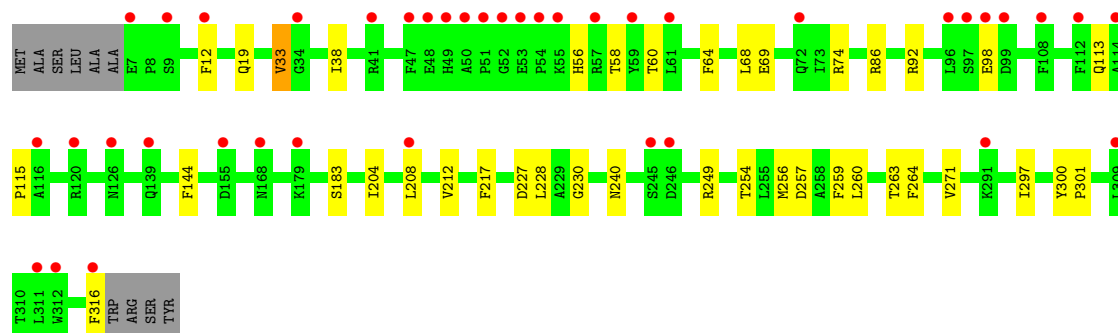
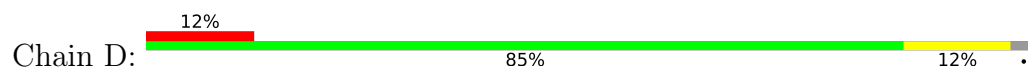


• Molecule 1: Cys-loop ligand-gated ion channel

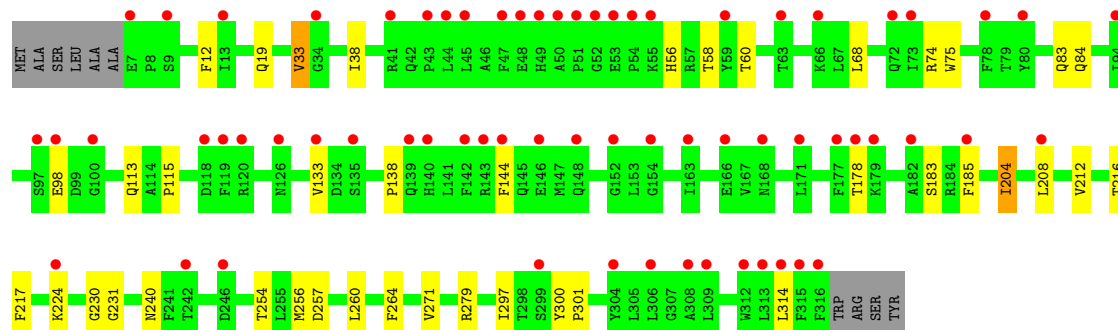
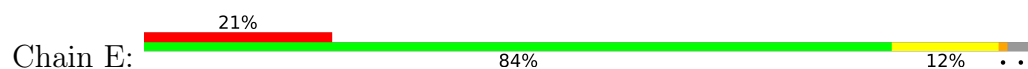




• Molecule 1: Cys-loop ligand-gated ion channel



• Molecule 1: Cys-loop ligand-gated ion channel



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	219.60Å 112.96Å 144.43Å 90.00° 112.05° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.7 (20.00-2.30) 89.5 (20.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.18 (at 2.30Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.209 , 0.236 0.276 , 0.308	Depositor DCC
R_{free} test set	6386 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 59.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13403	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% 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: BNG, PG4

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	0.83	0/2639	1.26	4/3583 (0.1%)
1	B	0.83	0/2639	1.25	7/3583 (0.2%)
1	C	0.85	0/2639	1.26	7/3583 (0.2%)
1	D	0.80	0/2639	1.24	5/3583 (0.1%)
1	E	0.77	0/2639	1.25	5/3583 (0.1%)
All	All	0.82	0/13195	1.25	28/17915 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	98	GLU	CA-C-N	7.70	130.60	120.28
1	B	98	GLU	C-N-CA	7.70	130.60	120.28
1	C	98	GLU	CA-C-N	7.42	130.96	120.28
1	C	98	GLU	C-N-CA	7.42	130.96	120.28
1	E	98	GLU	CA-C-N	7.06	130.45	120.28
1	E	98	GLU	C-N-CA	7.06	130.45	120.28
1	D	98	GLU	CA-C-N	7.00	130.36	120.28
1	D	98	GLU	C-N-CA	7.00	130.36	120.28
1	A	98	GLU	CA-C-N	6.83	130.11	120.28
1	A	98	GLU	C-N-CA	6.83	130.11	120.28
1	C	33	VAL	N-CA-C	-5.95	99.17	107.80
1	E	33	VAL	N-CA-C	-5.71	99.51	107.80
1	B	316	PHE	CA-CB-CG	5.50	119.30	113.80
1	E	60	THR	N-CA-C	-5.48	101.54	110.20
1	A	316	PHE	CA-CB-CG	5.48	119.28	113.80
1	C	175	ASN	CA-C-N	5.31	127.34	120.44
1	C	175	ASN	C-N-CA	5.31	127.34	120.44
1	C	29	ASN	CA-CB-CG	5.29	117.89	112.60
1	D	33	VAL	N-CA-C	-5.28	100.15	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	THR	N-CA-C	-5.26	101.89	110.20
1	B	33	VAL	N-CA-C	-5.20	100.25	107.80
1	E	204	ILE	N-CA-CB	5.19	114.00	110.52
1	D	316	PHE	CA-CB-CG	5.14	118.94	113.80
1	D	60	THR	N-CA-C	-5.07	101.70	109.76
1	B	69	GLU	CA-C-N	5.06	127.02	120.44
1	B	69	GLU	C-N-CA	5.06	127.02	120.44
1	C	316	PHE	CA-CB-CG	5.04	118.84	113.80
1	B	60	THR	N-CA-C	-5.01	102.29	110.20

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	2566	0	2539	40	0
1	B	2566	0	2539	35	0
1	C	2566	0	2539	48	0
1	D	2566	0	2539	38	0
1	E	2566	0	2539	34	0
2	A	26	0	36	4	0
2	C	26	0	36	9	0
2	E	13	0	18	2	0
3	A	21	0	30	7	0
3	B	21	0	30	8	0
3	C	21	0	30	8	0
3	D	21	0	30	3	0
3	E	21	0	30	6	0
4	A	74	0	0	3	0
4	B	99	0	0	3	0
4	C	101	0	0	6	0
4	D	88	0	0	1	0
4	E	41	0	0	2	0
All	All	13403	0	12935	171	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 7.

All (171) 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)
3:C:403:BNG:H2'2	1:D:227:ASP:HA	1.61	0.83
1:C:61:LEU:HD21	2:C:401:PG4:C3	2.11	0.80
1:C:144:PHE:HB2	1:C:183:SER:OG	1.81	0.80
1:E:144:PHE:HB2	1:E:183:SER:OG	1.82	0.79
1:A:144:PHE:HB2	1:A:183:SER:OG	1.82	0.79
1:D:144:PHE:HB2	1:D:183:SER:OG	1.84	0.77
1:B:144:PHE:HB2	1:B:183:SER:OG	1.85	0.76
1:A:230:GLY:HA3	3:E:402:BNG:H1'2	1.67	0.73
1:C:61:LEU:HD21	2:C:401:PG4:H31	1.70	0.72
1:A:228:LEU:HD11	3:A:403:BNG:H6'1	1.72	0.72
1:C:144:PHE:HB2	1:C:183:SER:HG	1.55	0.72
3:C:403:BNG:H1'2	1:D:230:GLY:HA3	1.71	0.72
1:E:231:GLY:HA3	3:E:402:BNG:H61	1.71	0.71
2:C:402:PG4:C2	2:C:402:PG4:H82	2.21	0.70
1:A:58:THR:HG21	1:B:74:ARG:HG3	1.73	0.70
1:B:212:VAL:HG13	1:C:271:VAL:HG11	1.74	0.69
1:C:58:THR:HG21	1:D:74:ARG:HG3	1.74	0.68
1:C:300:TYR:HB3	1:C:301:PRO:HD3	1.75	0.67
1:A:271:VAL:HG11	1:E:212:VAL:HG13	1.78	0.66
1:D:212:VAL:HG13	1:E:271:VAL:HG11	1.77	0.66
1:A:212:VAL:HG13	1:B:271:VAL:HG11	1.78	0.66
1:A:231:GLY:HA3	3:A:403:BNG:O4	1.95	0.65
1:B:300:TYR:HB3	1:B:301:PRO:HD3	1.77	0.65
2:C:402:PG4:H82	2:C:402:PG4:H21	1.79	0.64
1:D:300:TYR:HB3	1:D:301:PRO:HD3	1.79	0.63
3:B:401:BNG:H1'2	1:C:230:GLY:HA3	1.79	0.63
1:A:300:TYR:HB3	1:A:301:PRO:HD3	1.79	0.63
1:B:58:THR:HG21	1:C:74:ARG:HG3	1.81	0.62
1:E:300:TYR:HB3	1:E:301:PRO:HD3	1.80	0.62
1:B:228:LEU:HD11	3:B:401:BNG:H6'1	1.81	0.62
1:A:74:ARG:HG3	1:E:58:THR:HG21	1.81	0.62
1:A:227:ASP:HA	3:E:402:BNG:H2'2	1.80	0.61
2:E:401:PG4:H41	2:E:401:PG4:H81	1.81	0.61
2:A:402:PG4:H11	4:A:544:HOH:O	2.02	0.60
1:C:208:LEU:CD2	1:D:264:PHE:HB3	2.31	0.60
2:C:401:PG4:O4	2:C:401:PG4:H42	2.01	0.60
1:B:212:VAL:CG1	1:C:271:VAL:HG11	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:204:ILE:O	1:D:208:LEU:HD13	2.03	0.59
2:C:402:PG4:H82	2:C:402:PG4:O2	2.03	0.59
1:C:204:ILE:O	1:C:208:LEU:HD13	2.01	0.59
3:C:403:BNG:H3'2	1:D:230:GLY:HA3	1.83	0.58
1:C:61:LEU:HD21	2:C:401:PG4:H32	1.84	0.58
1:B:204:ILE:O	1:B:208:LEU:HD13	2.04	0.57
1:C:208:LEU:HD21	1:D:264:PHE:HB3	1.86	0.57
1:A:208:LEU:CD2	1:B:264:PHE:HB3	2.35	0.57
1:A:204:ILE:O	1:A:208:LEU:HD13	2.04	0.57
1:E:204:ILE:O	1:E:208:LEU:HD13	2.04	0.57
1:A:75:TRP:HB2	4:A:553:HOH:O	2.05	0.56
1:C:159:GLU:HB3	4:C:533:HOH:O	2.06	0.56
3:A:403:BNG:H1'2	1:B:230:GLY:HA3	1.88	0.55
1:A:65:LEU:HD11	2:A:401:PG4:H32	1.88	0.54
1:A:271:VAL:HG11	1:E:212:VAL:CG1	2.37	0.54
1:A:163:ILE:HD12	4:A:526:HOH:O	2.06	0.54
1:A:231:GLY:HA3	3:A:403:BNG:H61	1.89	0.54
1:C:212:VAL:HG13	1:D:271:VAL:HG11	1.90	0.54
3:B:401:BNG:H9'1	4:B:548:HOH:O	2.07	0.54
1:D:208:LEU:CD2	1:E:264:PHE:HB3	2.37	0.54
1:A:230:GLY:HA3	3:E:402:BNG:H3'2	1.90	0.53
3:D:401:BNG:H1'2	1:E:230:GLY:HA3	1.91	0.53
1:A:56:HIS:HB2	1:B:74:ARG:HH22	1.73	0.53
2:E:401:PG4:H41	2:E:401:PG4:C8	2.39	0.53
1:C:231:GLY:HA3	3:C:403:BNG:O4	2.09	0.52
1:A:212:VAL:CG1	1:B:271:VAL:HG11	2.38	0.52
1:A:264:PHE:HB3	1:E:208:LEU:CD2	2.40	0.52
1:A:240:ASN:HA	1:A:260:LEU:HD13	1.90	0.52
1:B:208:LEU:CD2	1:C:264:PHE:HB3	2.40	0.52
1:A:208:LEU:HD21	1:B:264:PHE:HB3	1.91	0.52
1:D:58:THR:HG21	1:E:74:ARG:HG3	1.91	0.51
1:E:240:ASN:HA	1:E:260:LEU:HD13	1.91	0.51
1:B:240:ASN:HA	1:B:260:LEU:HD13	1.92	0.51
1:D:212:VAL:CG1	1:E:271:VAL:HG11	2.39	0.51
1:A:74:ARG:HH22	1:E:56:HIS:HB2	1.75	0.51
1:E:279:ARG:HA	4:E:502:HOH:O	2.09	0.51
1:C:240:ASN:HA	1:C:260:LEU:HD13	1.93	0.50
1:C:144:PHE:CB	1:C:183:SER:HG	2.22	0.50
3:B:401:BNG:H5	4:B:548:HOH:O	2.11	0.50
1:D:240:ASN:HA	1:D:260:LEU:HD13	1.92	0.49
1:B:208:LEU:HD21	1:C:264:PHE:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:VAL:HG12	3:B:401:BNG:H5'1	1.94	0.48
1:B:113:GLN:OE1	1:B:115:PRO:HG3	2.14	0.48
3:D:401:BNG:H5	4:D:520:HOH:O	2.13	0.48
3:C:403:BNG:C1'	1:D:230:GLY:HA3	2.42	0.48
3:B:401:BNG:H2'2	1:C:227:ASP:HA	1.95	0.48
1:B:56:HIS:HB2	1:C:74:ARG:HH22	1.79	0.47
3:B:401:BNG:H9'3	3:B:401:BNG:H1'1	1.96	0.47
1:C:69:GLU:HA	4:C:510:HOH:O	2.13	0.47
3:A:403:BNG:H2'2	1:B:227:ASP:HA	1.96	0.47
1:C:200:MET:HE3	1:D:257:ASP:HB3	1.97	0.47
1:B:212:VAL:HG13	1:C:271:VAL:CG1	2.42	0.46
3:A:403:BNG:H3'2	1:B:230:GLY:HA3	1.97	0.46
1:D:208:LEU:HD21	1:E:264:PHE:HB3	1.95	0.46
1:C:197:TYR:CD2	1:D:249:ARG:HG2	2.50	0.46
1:C:212:VAL:CG1	1:D:271:VAL:HG11	2.45	0.46
1:C:56:HIS:HB2	1:D:74:ARG:HH22	1.80	0.46
1:E:216:THR:HG21	3:E:402:BNG:H4'2	1.98	0.46
1:A:271:VAL:CG1	1:E:212:VAL:HG13	2.46	0.45
1:C:25:GLN:HG2	4:C:513:HOH:O	2.15	0.45
1:A:264:PHE:HB3	1:E:208:LEU:HD21	1.99	0.45
1:C:19:GLN:HB2	1:C:33:VAL:HB	1.99	0.45
1:A:212:VAL:HG13	1:B:271:VAL:CG1	2.45	0.45
1:A:113:GLN:OE1	1:A:115:PRO:HG3	2.17	0.44
1:B:12:PHE:O	1:B:38:ILE:HA	2.17	0.44
1:E:254:THR:HB	1:E:257:ASP:H	1.83	0.44
1:D:228:LEU:HD11	3:D:401:BNG:H8'1	1.99	0.44
1:E:113:GLN:OE1	1:E:115:PRO:HG3	2.18	0.44
1:C:86:ARG:HE	1:C:86:ARG:HB3	1.61	0.44
2:A:401:PG4:H52	2:A:401:PG4:H72	1.79	0.44
1:E:19:GLN:HB2	1:E:33:VAL:HB	2.00	0.44
1:E:144:PHE:CB	1:E:183:SER:OG	2.62	0.44
1:A:254:THR:HB	1:A:257:ASP:H	1.83	0.43
1:C:217:PHE:HB3	1:C:297:ILE:HG12	2.00	0.43
3:C:403:BNG:H1'2	1:D:230:GLY:CA	2.44	0.43
1:A:19:GLN:HB2	1:A:33:VAL:HB	2.00	0.43
1:A:138:PRO:HD3	1:A:178:THR:HG21	2.01	0.43
3:B:401:BNG:H3'2	1:C:230:GLY:HA3	2.00	0.43
1:D:113:GLN:OE1	1:D:115:PRO:HG3	2.17	0.43
1:D:212:VAL:HG13	1:E:271:VAL:CG1	2.46	0.43
1:E:217:PHE:HB3	1:E:297:ILE:HG12	2.00	0.43
1:D:259:PHE:O	1:D:263:THR:HG23	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:138:PRO:HD3	1:C:178:THR:HG21	2.01	0.43
1:C:254:THR:HB	1:C:257:ASP:H	1.84	0.43
2:A:402:PG4:H21	2:A:402:PG4:O5	2.19	0.43
1:C:113:GLN:OE1	1:C:115:PRO:HG3	2.18	0.43
1:D:86:ARG:HE	1:D:86:ARG:HB3	1.60	0.43
1:D:144:PHE:CB	1:D:183:SER:OG	2.63	0.43
1:D:19:GLN:HB2	1:D:33:VAL:HB	2.00	0.42
1:A:86:ARG:HE	1:A:86:ARG:HB3	1.61	0.42
1:B:138:PRO:HD3	1:B:178:THR:HG21	2.00	0.42
1:C:254:THR:HG22	1:C:256:MET:H	1.84	0.42
1:D:217:PHE:HB3	1:D:297:ILE:HG12	2.01	0.42
2:C:402:PG4:H62	2:C:402:PG4:H42	1.97	0.42
1:E:254:THR:HG22	1:E:256:MET:H	1.85	0.42
2:C:402:PG4:H41	1:D:92:ARG:NH2	2.35	0.42
1:B:224:LYS:HD2	4:B:575:HOH:O	2.20	0.42
1:C:286:GLN:HB2	4:C:542:HOH:O	2.20	0.42
1:D:56:HIS:HB2	1:E:74:ARG:HH22	1.85	0.42
1:A:107:ARG:NH2	1:B:81:HIS:HA	2.35	0.41
1:A:259:PHE:O	1:A:263:THR:HG23	2.21	0.41
1:B:68:LEU:HD13	1:B:75:TRP:HB3	2.03	0.41
1:C:231:GLY:HA3	3:C:403:BNG:H61	2.01	0.41
1:D:64:PHE:CE2	1:D:68:LEU:HD11	2.55	0.41
1:E:12:PHE:O	1:E:38:ILE:HA	2.20	0.41
1:E:138:PRO:HD3	1:E:178:THR:HG21	2.02	0.41
1:A:12:PHE:O	1:A:38:ILE:HA	2.21	0.41
1:A:217:PHE:HB3	1:A:297:ILE:HG12	2.02	0.41
1:B:254:THR:HB	1:B:257:ASP:H	1.85	0.41
1:E:68:LEU:HD13	1:E:75:TRP:HB3	2.02	0.41
1:E:83:GLN:HG2	1:E:84:GLN:N	2.36	0.41
1:C:90:GLN:HG3	4:C:502:HOH:O	2.20	0.41
1:C:197:TYR:CD1	1:C:197:TYR:C	2.98	0.41
1:C:216:THR:HG21	3:C:403:BNG:H4'2	2.01	0.41
1:D:254:THR:HG22	1:D:256:MET:H	1.85	0.41
3:E:402:BNG:H9'3	3:E:402:BNG:H1'1	2.03	0.41
1:B:19:GLN:HB2	1:B:33:VAL:HB	2.02	0.41
1:C:58:THR:CG2	1:D:74:ARG:NH1	2.83	0.41
1:B:133:VAL:HB	1:B:185:PHE:HB3	2.03	0.41
1:B:217:PHE:HB3	1:B:297:ILE:HG12	2.02	0.41
1:B:259:PHE:O	1:B:263:THR:HG23	2.21	0.41
1:D:12:PHE:O	1:D:38:ILE:HA	2.20	0.41
1:A:68:LEU:HD13	1:A:75:TRP:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:THR:HG21	1:D:69:GLU:OE1	2.21	0.41
1:D:254:THR:HB	1:D:257:ASP:H	1.86	0.41
1:C:49:HIS:CD2	1:C:98:GLU:CD	2.99	0.41
1:A:144:PHE:CB	1:A:183:SER:OG	2.61	0.40
1:B:254:THR:HG22	1:B:256:MET:H	1.86	0.40
1:E:133:VAL:HB	1:E:185:PHE:HB3	2.03	0.40
1:A:231:GLY:CA	3:A:403:BNG:H61	2.52	0.40
1:C:83:GLN:HG2	1:C:84:GLN:N	2.37	0.40
1:C:314:LEU:HD22	4:C:556:HOH:O	2.20	0.40
1:A:254:THR:HG22	1:A:256:MET:H	1.85	0.40
1:E:224:LYS:HG3	4:E:533:HOH:O	2.21	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	308/320 (96%)	303 (98%)	5 (2%)	0	100	100
1	B	308/320 (96%)	303 (98%)	5 (2%)	0	100	100
1	C	308/320 (96%)	302 (98%)	6 (2%)	0	100	100
1	D	308/320 (96%)	302 (98%)	6 (2%)	0	100	100
1	E	308/320 (96%)	303 (98%)	5 (2%)	0	100	100
All	All	1540/1600 (96%)	1513 (98%)	27 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	275/282 (98%)	275 (100%)	0	100	100
1	B	275/282 (98%)	274 (100%)	1 (0%)	84	92
1	C	275/282 (98%)	274 (100%)	1 (0%)	84	92
1	D	275/282 (98%)	275 (100%)	0	100	100
1	E	275/282 (98%)	274 (100%)	1 (0%)	84	92
All	All	1375/1410 (98%)	1372 (100%)	3 (0%)	87	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	314	LEU
1	C	314	LEU
1	E	314	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	83	GLN
1	A	91	ASN
1	A	139	GLN
1	A	232	ASN
1	B	49	HIS
1	B	72	GLN
1	B	83	GLN
1	B	91	ASN
1	B	139	GLN
1	B	232	ASN
1	C	56	HIS
1	C	72	GLN
1	C	83	GLN
1	C	91	ASN
1	C	139	GLN
1	D	49	HIS
1	D	83	GLN
1	D	91	ASN

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Mol	Chain	Res	Type
1	D	132	HIS
1	D	139	GLN
1	D	220	GLN
1	E	49	HIS
1	E	83	GLN
1	E	91	ASN
1	E	139	GLN
1	E	220	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](#)

10 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	PG4	A	401	-	12,12,12	0.29	0	11,11,11	0.51	0
2	PG4	C	402	-	12,12,12	0.25	0	11,11,11	0.29	0
2	PG4	E	401	-	12,12,12	0.15	0	11,11,11	0.16	0
3	BNG	E	402	-	21,21,21	0.20	0	26,26,26	0.47	0
3	BNG	C	403	-	21,21,21	0.29	0	26,26,26	0.85	1 (3%)
3	BNG	D	401	-	21,21,21	0.29	0	26,26,26	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BNG	A	403	-	21,21,21	0.30	0	26,26,26	0.62	0
3	BNG	B	401	-	21,21,21	0.30	0	26,26,26	0.43	0
2	PG4	C	401	-	12,12,12	0.28	0	11,11,11	0.28	0
2	PG4	A	402	-	12,12,12	0.20	0	11,11,11	0.28	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	PG4	A	401	-	-	7/10/10/10	-
2	PG4	C	402	-	-	7/10/10/10	-
2	PG4	E	401	-	-	5/10/10/10	-
3	BNG	E	402	-	-	6/12/32/32	0/1/1/1
3	BNG	C	403	-	-	8/12/32/32	0/1/1/1
3	BNG	D	401	-	-	8/12/32/32	0/1/1/1
3	BNG	A	403	-	-	7/12/32/32	0/1/1/1
3	BNG	B	401	-	-	8/12/32/32	0/1/1/1
2	PG4	C	401	-	-	5/10/10/10	-
2	PG4	A	402	-	-	6/10/10/10	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	403	BNG	O5-C1-O1	2.52	116.00	110.04

There are no chirality outliers.

All (67) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	402	PG4	C6-C5-O3-C4
2	C	401	PG4	O2-C3-C4-O3
2	C	402	PG4	O2-C3-C4-O3
3	A	403	BNG	O5-C1-O1-C1'
3	B	401	BNG	O5-C1-O1-C1'
3	D	401	BNG	O5-C1-O1-C1'
3	B	401	BNG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	A	403	BNG	C2-C1-O1-C1'
3	B	401	BNG	C2-C1-O1-C1'
3	C	403	BNG	C2-C1-O1-C1'
3	D	401	BNG	C2-C1-O1-C1'
3	C	403	BNG	O5-C1-O1-C1'
2	A	402	PG4	O1-C1-C2-O2
2	C	402	PG4	O1-C1-C2-O2
3	A	403	BNG	C3'-C4'-C5'-C6'
3	C	403	BNG	C3'-C4'-C5'-C6'
3	E	402	BNG	C3'-C4'-C5'-C6'
3	D	401	BNG	C4'-C5'-C6'-C7'
3	D	401	BNG	C3'-C4'-C5'-C6'
3	E	402	BNG	C2'-C3'-C4'-C5'
3	B	401	BNG	C3'-C4'-C5'-C6'
3	A	403	BNG	C4'-C5'-C6'-C7'
3	D	401	BNG	C2'-C3'-C4'-C5'
3	C	403	BNG	C2'-C3'-C4'-C5'
2	A	401	PG4	O1-C1-C2-O2
2	A	401	PG4	C1-C2-O2-C3
2	E	401	PG4	O2-C3-C4-O3
2	E	401	PG4	O4-C7-C8-O5
3	D	401	BNG	O5-C5-C6-O6
3	B	401	BNG	C4-C5-C6-O6
3	B	401	BNG	C2'-C3'-C4'-C5'
3	A	403	BNG	C2'-C3'-C4'-C5'
2	A	401	PG4	C5-C6-O4-C7
3	B	401	BNG	C4'-C5'-C6'-C7'
3	E	402	BNG	C4'-C5'-C6'-C7'
2	C	401	PG4	O4-C7-C8-O5
3	E	402	BNG	C2'-C1'-O1-C1
2	A	402	PG4	O3-C5-C6-O4
3	D	401	BNG	C5'-C6'-C7'-C8'
3	C	403	BNG	C4'-C5'-C6'-C7'
2	A	402	PG4	C3-C4-O3-C5
2	E	401	PG4	C4-C3-O2-C2
3	E	402	BNG	C6'-C7'-C8'-C9'
2	A	401	PG4	C4-C3-O2-C2
2	C	401	PG4	C3-C4-O3-C5
2	C	401	PG4	C6-C5-O3-C4
2	C	402	PG4	C3-C4-O3-C5
2	A	402	PG4	C6-C5-O3-C4
3	C	403	BNG	C6'-C7'-C8'-C9'

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Mol	Chain	Res	Type	Atoms
3	B	401	BNG	C2'-C1'-O1-C1
2	A	401	PG4	C3-C4-O3-C5
2	A	401	PG4	O2-C3-C4-O3
2	C	402	PG4	C1-C2-O2-C3
2	A	402	PG4	O4-C7-C8-O5
2	E	401	PG4	C8-C7-O4-C6
3	A	403	BNG	C6'-C7'-C8'-C9'
3	A	403	BNG	C2'-C1'-O1-C1
3	C	403	BNG	C2'-C1'-O1-C1
3	D	401	BNG	C2'-C1'-O1-C1
2	C	402	PG4	O3-C5-C6-O4
3	C	403	BNG	C5'-C6'-C7'-C8'
2	C	401	PG4	C4-C3-O2-C2
2	C	402	PG4	C4-C3-O2-C2
2	A	402	PG4	O2-C3-C4-O3
2	A	401	PG4	C6-C5-O3-C4
2	E	401	PG4	C5-C6-O4-C7
3	E	402	BNG	C5'-C6'-C7'-C8'

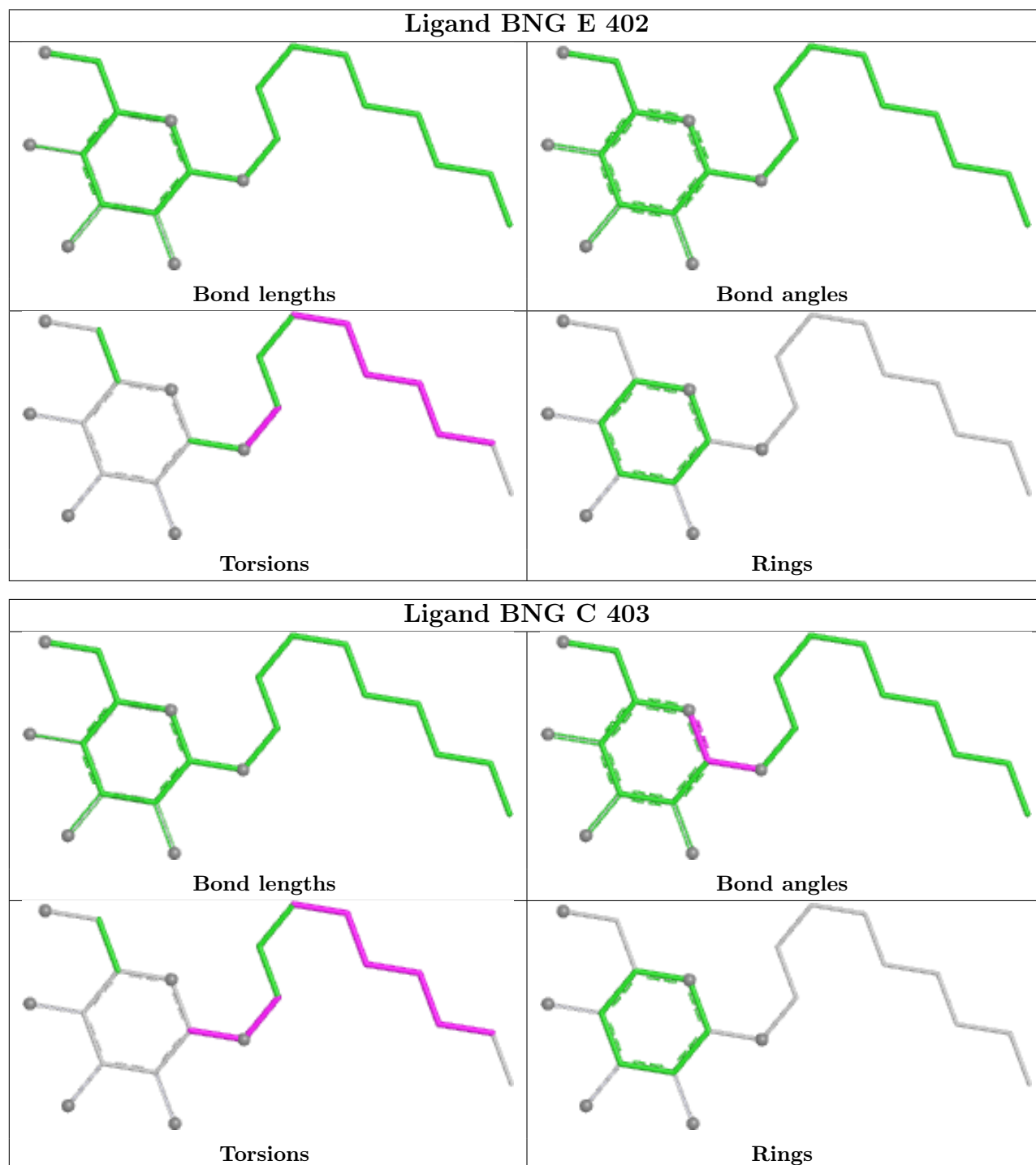
There are no ring outliers.

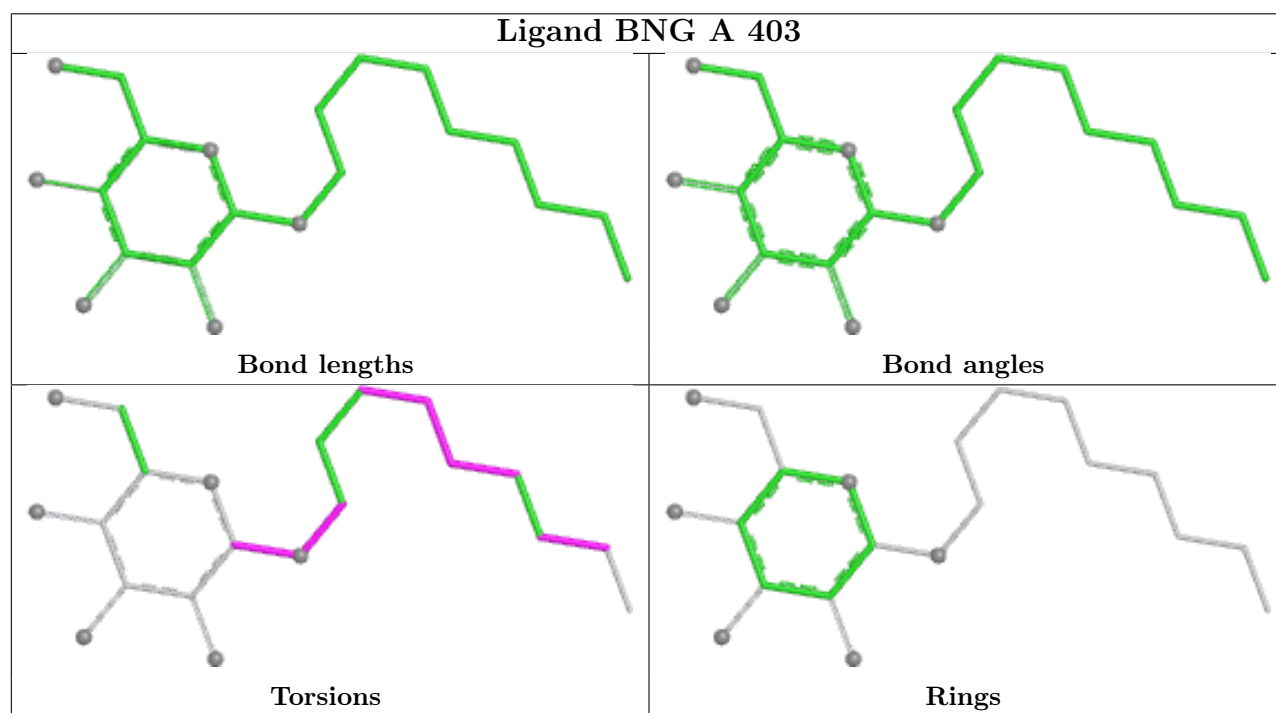
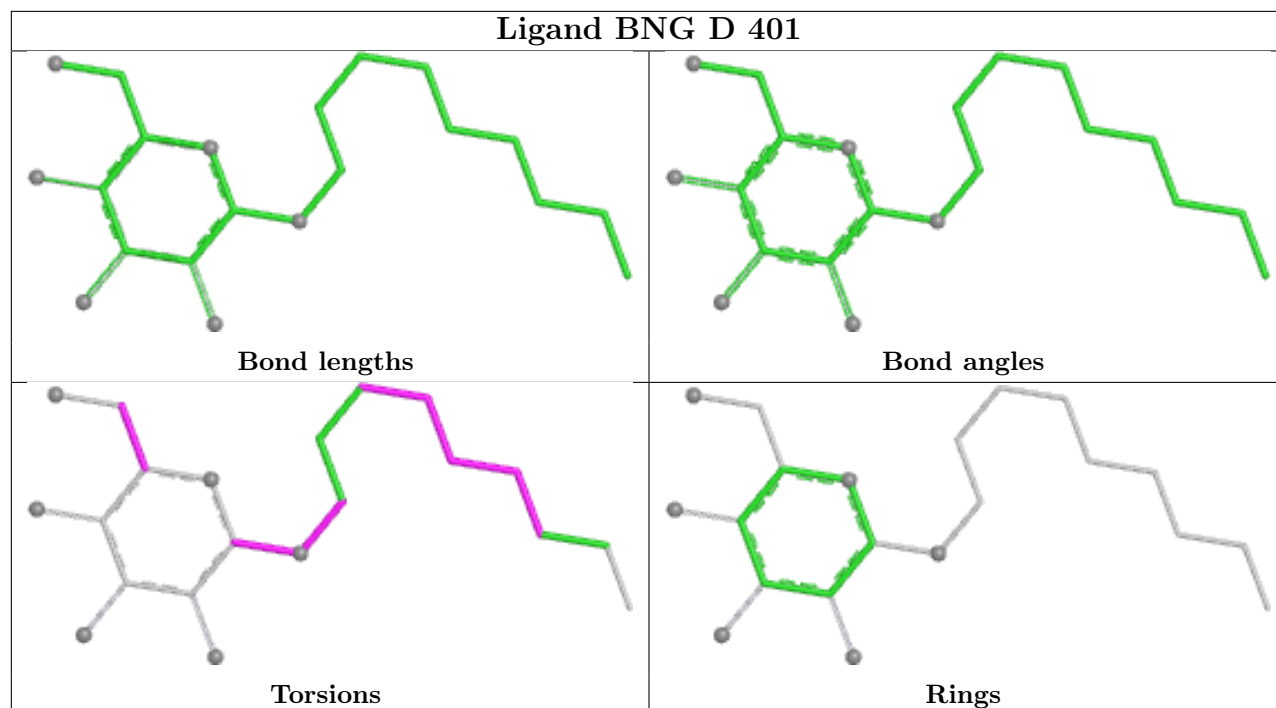
10 monomers are involved in 47 short contacts:

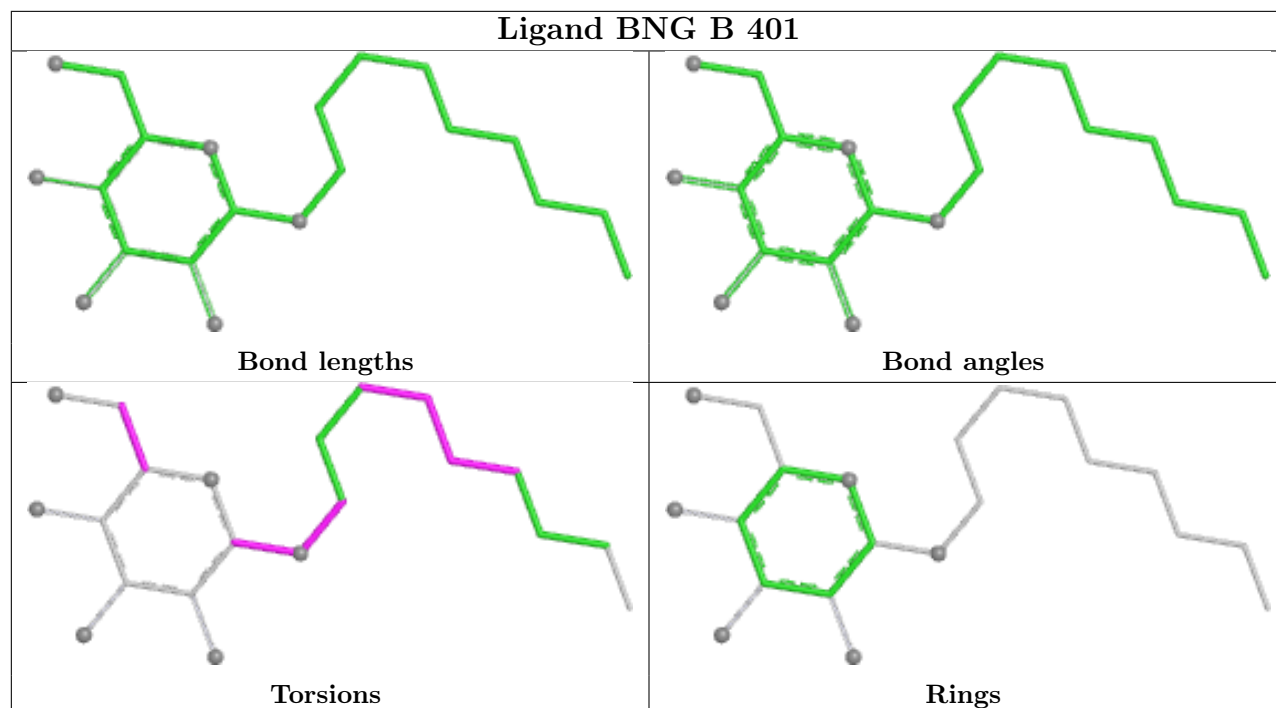
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	PG4	2	0
2	C	402	PG4	5	0
2	E	401	PG4	2	0
3	E	402	BNG	6	0
3	C	403	BNG	8	0
3	D	401	BNG	3	0
3	A	403	BNG	7	0
3	B	401	BNG	8	0
2	C	401	PG4	4	0
2	A	402	PG4	2	0

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	310/320 (96%)	0.94	33 (10%) 11 12	50, 77, 120, 173	0
1	B	310/320 (96%)	0.81	27 (8%) 16 17	45, 72, 115, 165	0
1	C	310/320 (96%)	0.76	21 (6%) 23 25	42, 72, 110, 138	0
1	D	310/320 (96%)	1.09	40 (12%) 7 8	36, 79, 125, 209	0
1	E	310/320 (96%)	1.40	66 (21%) 2 3	51, 88, 138, 186	0
All	All	1550/1600 (96%)	1.00	187 (12%) 8 9	36, 78, 125, 209	0

All (187) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	54	PRO	4.9
1	E	152	GLY	4.8
1	C	312	TRP	4.6
1	E	51	PRO	4.6
1	E	143	ARG	4.5
1	D	291	LYS	4.3
1	C	51	PRO	4.2
1	C	98	GLU	4.2
1	E	55	LYS	4.2
1	A	291	LYS	4.1
1	D	51	PRO	4.1
1	D	48	GLU	4.1
1	A	179	LYS	4.1
1	B	312	TRP	4.0
1	D	7	GLU	4.0
1	C	152	GLY	3.9
1	B	316	PHE	3.8
1	D	53	GLU	3.7
1	E	179	LYS	3.6
1	E	73	ILE	3.6

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Mol	Chain	Res	Type	RSRZ
1	E	50	ALA	3.6
1	D	316	PHE	3.6
1	E	171	LEU	3.5
1	E	246	ASP	3.5
1	B	51	PRO	3.5
1	A	312	TRP	3.5
1	A	181	ASP	3.5
1	E	13	ILE	3.4
1	E	53	GLU	3.4
1	B	50	ALA	3.4
1	E	44	LEU	3.4
1	E	312	TRP	3.3
1	E	120	ARG	3.3
1	E	142	PHE	3.3
1	E	306	LEU	3.3
1	E	133	VAL	3.3
1	C	257	ASP	3.3
1	E	313	LEU	3.2
1	C	179	LYS	3.2
1	D	59	TYR	3.2
1	E	118	ASP	3.1
1	E	78	PHE	3.1
1	E	52	GLY	3.1
1	D	61	LEU	3.1
1	A	177	PHE	3.1
1	D	34	GLY	3.0
1	D	49	HIS	3.0
1	B	242	THR	3.0
1	A	155	ASP	3.0
1	E	316	PHE	3.0
1	D	55	LYS	3.0
1	D	50	ALA	2.9
1	A	182	ALA	2.9
1	D	312	TRP	2.9
1	D	179	LYS	2.9
1	E	208	LEU	2.9
1	B	179	LYS	2.8
1	E	59	TYR	2.8
1	D	96	LEU	2.8
1	D	47	PHE	2.8
1	B	224	LYS	2.8
1	E	168	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	143	ARG	2.8
1	E	135	SER	2.8
1	C	7	GLU	2.8
1	C	120	ARG	2.8
1	D	155	ASP	2.8
1	C	183	SER	2.8
1	E	100	GLY	2.7
1	B	183	SER	2.7
1	C	208	LEU	2.7
1	D	57	ARG	2.7
1	E	308	ALA	2.7
1	E	185	PHE	2.7
1	D	168	ASN	2.7
1	E	154	GLY	2.7
1	E	7	GLU	2.7
1	E	43	PRO	2.7
1	E	54	PRO	2.7
1	E	177	PHE	2.7
1	C	66	LYS	2.7
1	C	155	ASP	2.7
1	E	98	GLU	2.7
1	D	208	LEU	2.7
1	B	12	PHE	2.6
1	B	7	GLU	2.6
1	E	48	GLU	2.6
1	E	34	GLY	2.6
1	A	143	ARG	2.6
1	D	41	ARG	2.6
1	E	314	LEU	2.6
1	A	316	PHE	2.6
1	A	139	GLN	2.6
1	A	311	LEU	2.6
1	D	309	LEU	2.6
1	D	120	ARG	2.6
1	D	114	ALA	2.6
1	C	316	PHE	2.6
1	B	208	LEU	2.5
1	D	311	LEU	2.5
1	D	99	ASP	2.5
1	E	94	ILE	2.5
1	A	98	GLU	2.5
1	D	112	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	257	ASP	2.5
1	B	55	LYS	2.5
1	C	63	THR	2.5
1	E	299	SER	2.5
1	A	41	ARG	2.5
1	E	163	ILE	2.5
1	B	99	ASP	2.5
1	D	12	PHE	2.4
1	B	191	ALA	2.4
1	B	306	LEU	2.4
1	C	306	LEU	2.4
1	A	304	TYR	2.4
1	E	80	TYR	2.4
1	B	126	ASN	2.4
1	A	173	THR	2.4
1	E	242	THR	2.4
1	D	97	SER	2.4
1	D	245	SER	2.4
1	A	208	LEU	2.4
1	A	168	ASN	2.4
1	E	126	ASN	2.4
1	E	97	SER	2.4
1	A	170	HIS	2.4
1	E	166	GLU	2.4
1	C	52	GLY	2.3
1	E	182	ALA	2.3
1	A	58	THR	2.3
1	C	291	LYS	2.3
1	A	258	ALA	2.3
1	C	243	ILE	2.3
1	E	144	PHE	2.3
1	A	178	THR	2.3
1	D	9	SER	2.3
1	E	140	HIS	2.3
1	B	47	PHE	2.3
1	B	291	LYS	2.3
1	A	183	SER	2.3
1	D	52	GLY	2.2
1	A	7	GLU	2.2
1	E	304	TYR	2.2
1	B	311	LEU	2.2
1	E	45	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	315	PHE	2.2
1	B	143	ARG	2.2
1	C	54	PRO	2.2
1	B	63	THR	2.2
1	A	104	TYR	2.2
1	E	119	PHE	2.2
1	E	41	ARG	2.2
1	B	97	SER	2.2
1	E	178	THR	2.2
1	A	138	PRO	2.2
1	A	77	ALA	2.2
1	E	309	LEU	2.1
1	D	98	GLU	2.1
1	E	47	PHE	2.1
1	B	194	HIS	2.1
1	D	72	GLN	2.1
1	D	139	GLN	2.1
1	E	9	SER	2.1
1	D	126	ASN	2.1
1	D	246	ASP	2.1
1	E	146	GLU	2.1
1	E	49	HIS	2.1
1	B	112	PHE	2.1
1	A	86	ARG	2.1
1	D	116	ALA	2.1
1	B	305	LEU	2.1
1	E	72	GLN	2.1
1	A	142	PHE	2.1
1	B	117	PHE	2.1
1	D	108	PHE	2.1
1	E	66	LYS	2.1
1	A	119	PHE	2.1
1	B	120	ARG	2.1
1	E	224	LYS	2.0
1	A	245	SER	2.0
1	E	63	THR	2.0
1	E	139	GLN	2.0
1	A	253	ILE	2.0
1	A	315	PHE	2.0
1	E	148	GLN	2.0
1	C	41	ARG	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](#)

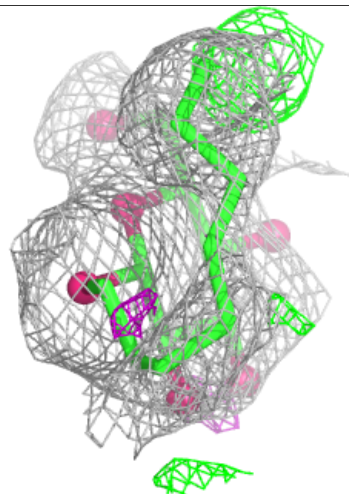
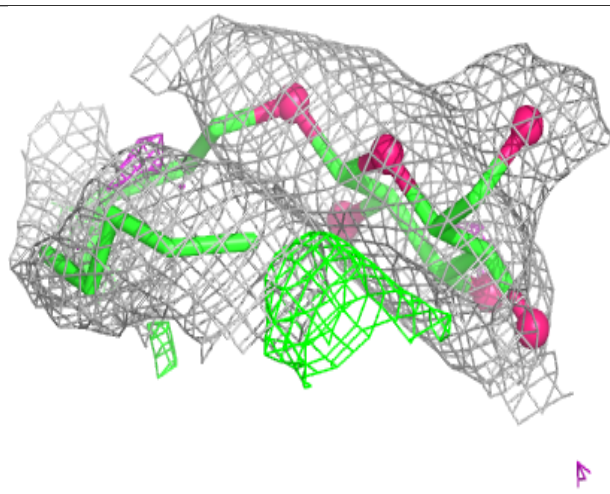
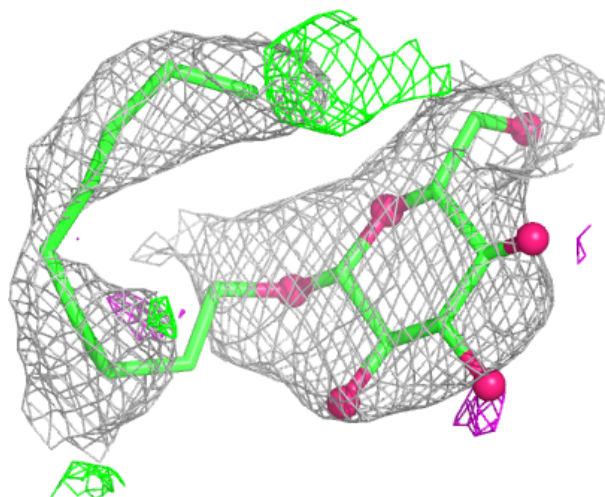
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	PG4	A	401	13/13	0.64	0.27	91,101,106,107	0
2	PG4	C	402	13/13	0.70	0.24	85,97,100,100	0
3	BNG	A	403	21/21	0.73	0.20	48,85,90,92	0
3	BNG	C	403	21/21	0.73	0.22	43,93,102,105	0
2	PG4	C	401	13/13	0.77	0.27	89,103,105,106	0
3	BNG	E	402	21/21	0.77	0.23	55,108,122,123	0
3	BNG	D	401	21/21	0.79	0.19	32,90,101,103	0
2	PG4	A	402	13/13	0.79	0.26	98,101,104,104	0
3	BNG	B	401	21/21	0.80	0.21	31,93,104,107	0
2	PG4	E	401	13/13	0.83	0.18	89,101,109,109	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.

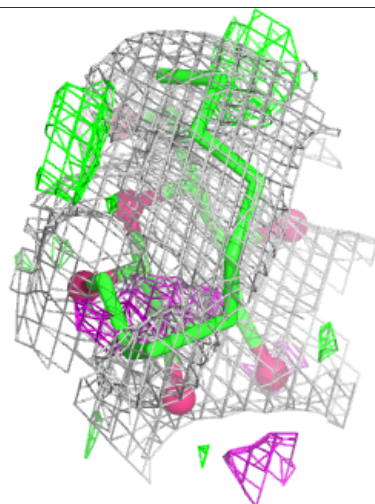
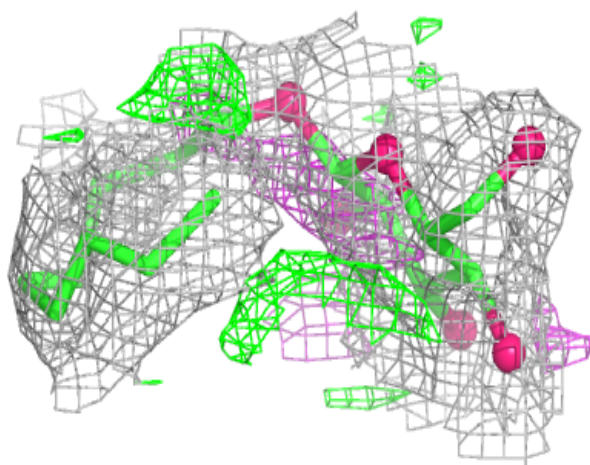
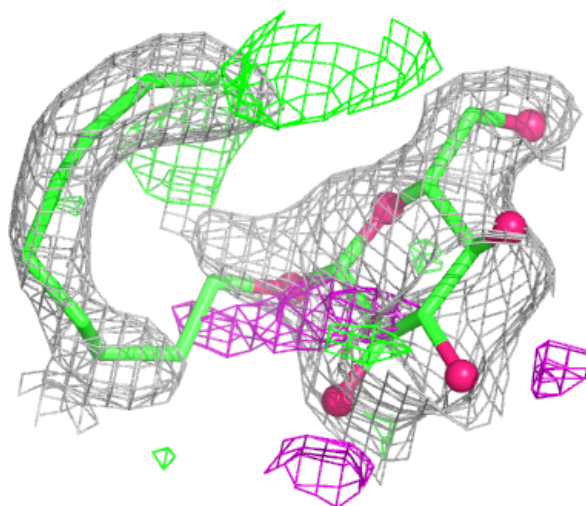
Electron density around BNG A 403:

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



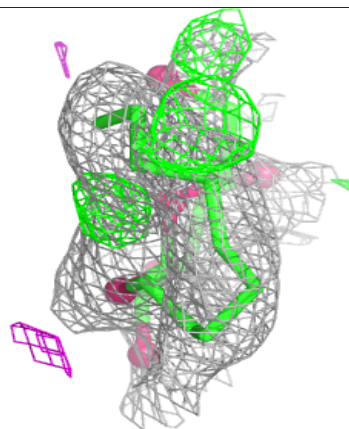
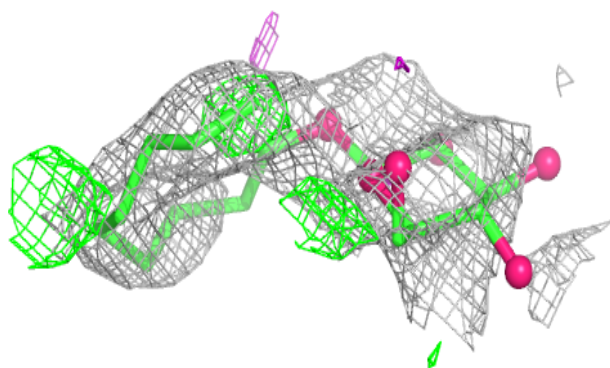
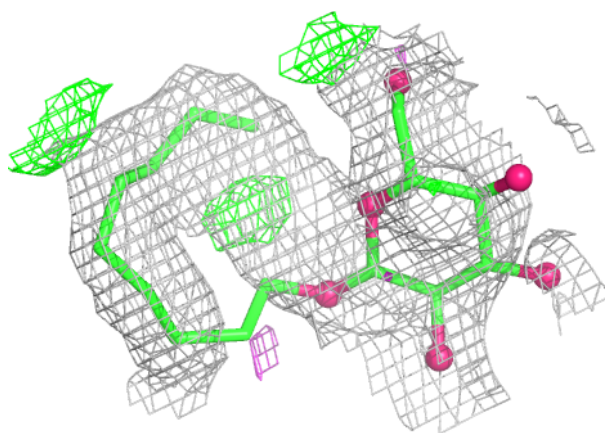
Electron density around BNG C 403:

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

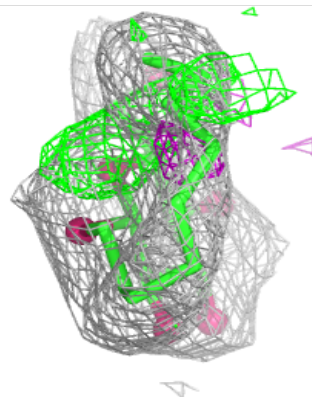
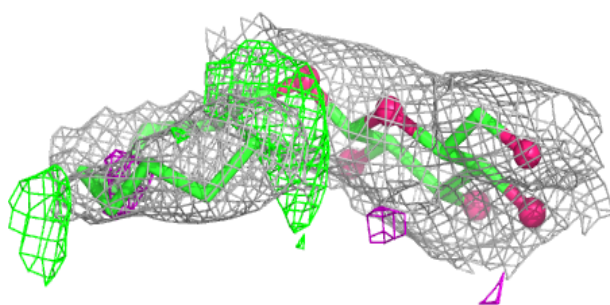
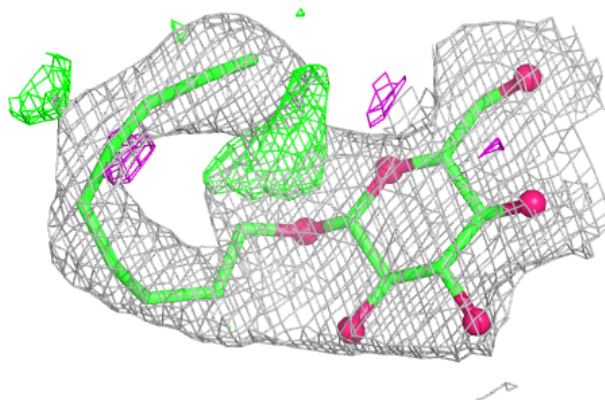


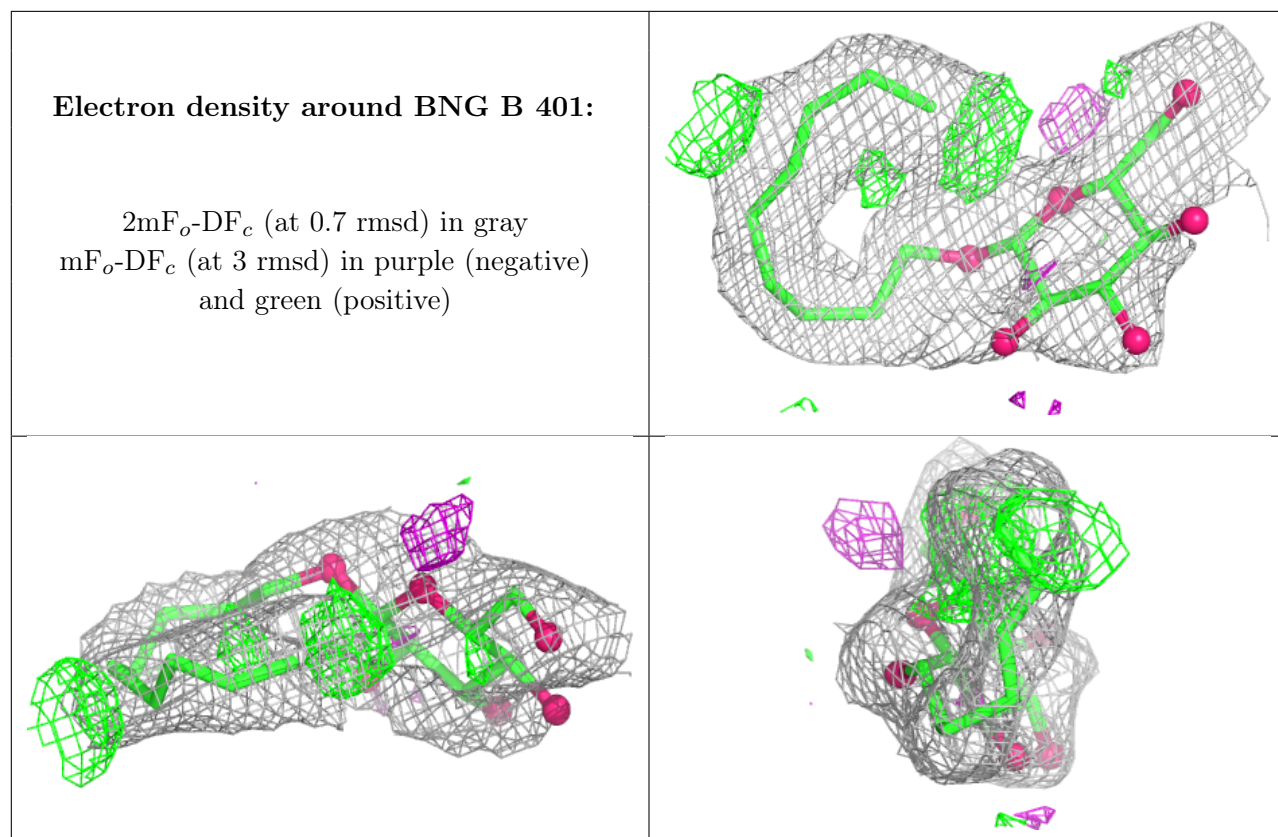
Electron density around BNG E 402:

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

**Electron density around BNG D 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.