



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 04:13 AM BST

PDB ID : 2NZT
Title : Crystal structure of human hexokinase II
Authors : Rabeh, W.M.; Zhu, H.; Nedyalkova, L.; Tempel, W.; Wasney, G.; Landry, R.;
Vedadi, M.; Arrowsmith, C.H.; Edwards, A.M.; Sundstrom, M.; Weigelt, J.;
Bochkarev, A.; Park, H.; Structural Genomics Consortium (SGC)
Deposited on : 2006-11-25
Resolution : 2.45 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

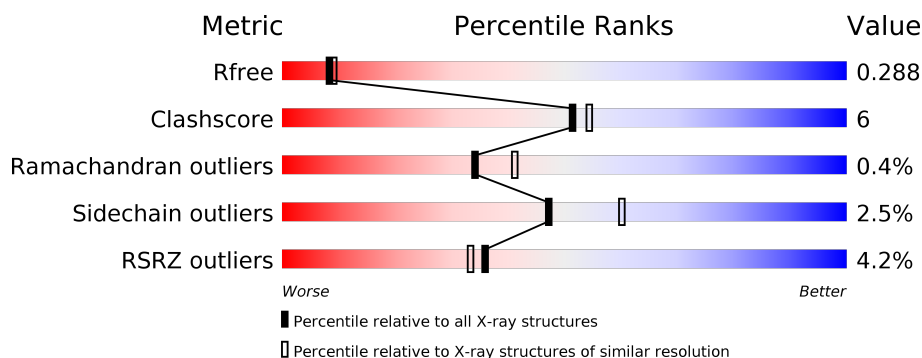
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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}	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 on 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	902	<div> <div>3%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>••</div> </div> </div>
1	B	902	<div> <div>5%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div>••</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	UNX	A	1	-	-	-	X
4	UNX	A	10	-	-	-	X
4	UNX	A	1005	-	-	-	X
4	UNX	A	1006	-	-	-	X
4	UNX	A	1007	-	-	-	X
4	UNX	A	1008	-	-	-	X
4	UNX	A	1009	-	-	-	X
4	UNX	A	11	-	-	-	X
4	UNX	A	4	-	-	-	X
4	UNX	A	6	-	-	-	X
4	UNX	A	8	-	-	-	X
4	UNX	A	9	-	-	-	X
4	UNX	B	1005	-	-	-	X
4	UNX	B	1006	-	-	-	X
4	UNX	B	1007	-	-	-	X
4	UNX	B	1008	-	-	-	X
4	UNX	B	1009	-	-	-	X
4	UNX	B	1010	-	-	-	X
4	UNX	B	1011	-	-	-	X
4	UNX	B	1012	-	-	-	X
4	UNX	B	1013	-	-	-	X
4	UNX	B	12	-	-	-	X
4	UNX	B	13	-	-	-	X
4	UNX	B	14	-	-	-	X
4	UNX	B	2	-	-	-	X
4	UNX	B	3	-	-	-	X
4	UNX	B	5	-	-	-	X
4	UNX	B	7	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 13546 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.

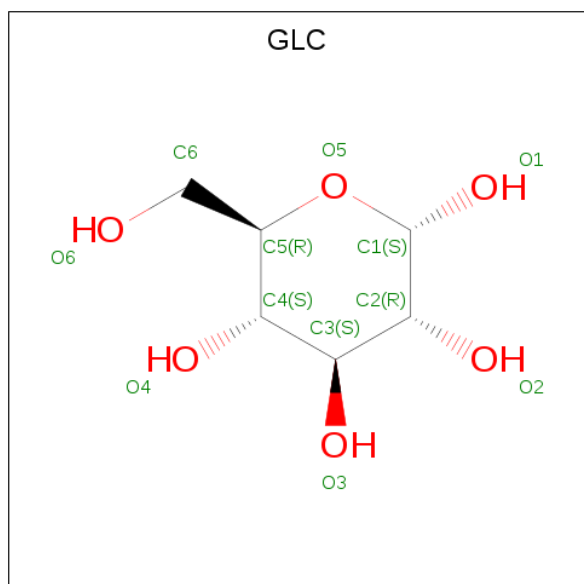
- Molecule 1 is a protein called Hexokinase-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	871	Total	C	N	O	S	0	0	0
			6653	4181	1157	1253	62			
1	B	867	Total	C	N	O	S	0	0	0
			6656	4181	1157	1256	62			

There are 4 discrepancies between the modelled and reference sequences:

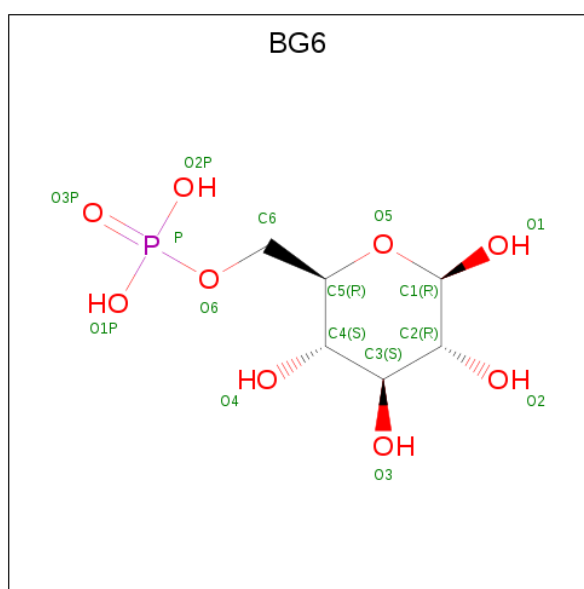
Chain	Residue	Modelled	Actual	Comment	Reference
A	15	GLY	-	cloning artifact	UNP P52789
A	16	SER	-	cloning artifact	UNP P52789
B	15	GLY	-	cloning artifact	UNP P52789
B	16	SER	-	cloning artifact	UNP P52789

- Molecule 2 is alpha-D-glucopyranose (three-letter code: GLC) (formula: C₆H₁₂O₆).



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

- Molecule 3 is 6-O-phosphono-beta-D-glucopyranose (three-letter code: BG6) (formula: $C_6H_{13}O_9P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			16	6	9	1		
3	A	1	Total	C	O	P	0	0
			16	6	9	1		
3	B	1	Total	C	O	P	0	0
			16	6	9	1		
3	B	1	Total	C	O	P	0	0
			16	6	9	1		

- Molecule 4 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	16	Total	X	0	0
			16	16		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	X	0	0
			12	12		

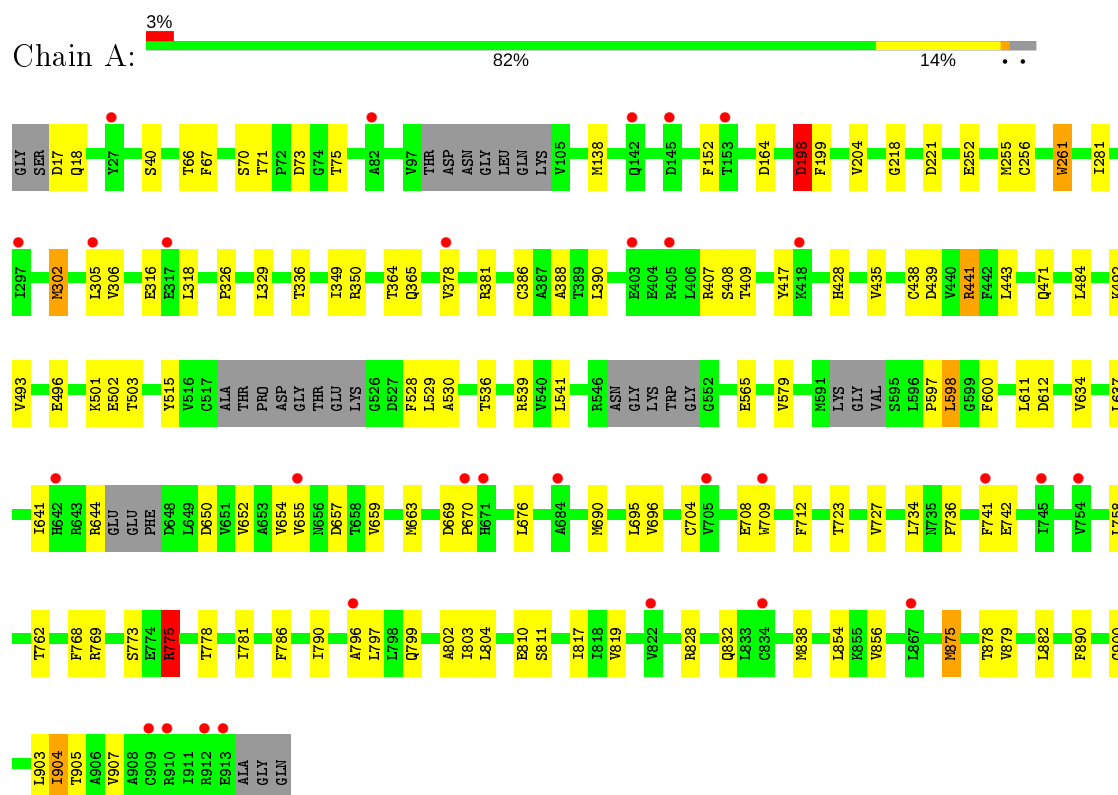
- Molecule 5 is water.

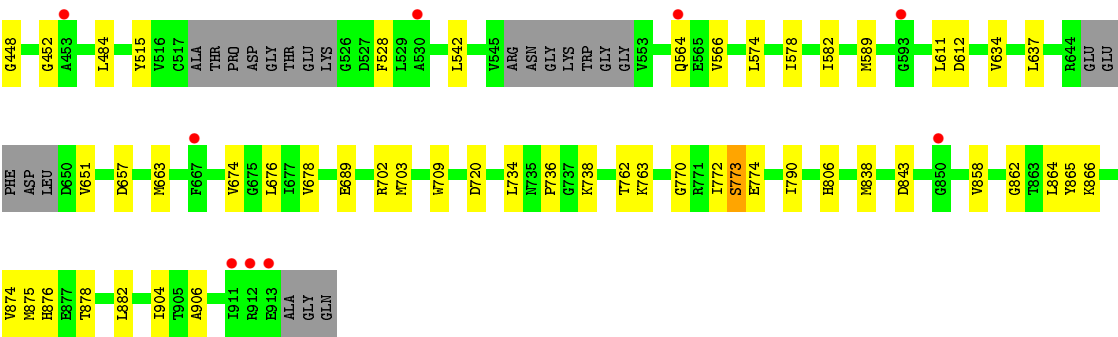
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	49	Total	O	0	0
			49	49		
5	B	48	Total	O	0	0
			48	48		

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: Hexokinase-2





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.89Å 129.34Å 187.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.45 24.74 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.2 (25.00-2.45) 98.2 (24.74-2.45)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.44Å)	Xtriage
Refinement program	REFMAC refmac_5.2.0019, PRODRG	Depositor
R, R_{free}	0.230 , 0.287 0.234 , 0.288	Depositor DCC
R_{free} test set	2083 reflections (2.49%)	wwPDB-VP
Wilson B-factor (Å ²)	51.8	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13546	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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: UNX, GLC, BG6

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	0.79	1/6751 (0.0%)	0.74	2/9091 (0.0%)
1	B	0.82	2/6755 (0.0%)	0.74	2/9095 (0.0%)
All	All	0.80	3/13506 (0.0%)	0.74	4/18186 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	256	CYS	CB-SG	-5.29	1.73	1.81
1	A	810	GLU	CG-CD	5.07	1.59	1.51
1	B	773	SER	CB-OG	5.03	1.48	1.42

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	775	ARG	NE-CZ-NH1	5.95	123.27	120.30
1	B	183	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	B	381	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	A	539	ARG	NE-CZ-NH1	5.04	122.82	120.30

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	6653	0	6557	78	0
1	B	6656	0	6565	73	0
2	A	24	0	23	1	0
2	B	24	0	24	0	0
3	A	32	0	22	0	0
3	B	32	0	22	1	0
4	A	12	0	0	0	0
4	B	16	0	0	0	0
5	A	49	0	0	0	0
5	B	48	0	0	0	0
All	All	13546	0	13213	149	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 6.

All (149) 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:216:THR:HG22	1:B:452:GLY:H	1.13	1.13
1:B:216:THR:HG22	1:B:452:GLY:N	1.83	0.93
1:A:655:VAL:HG11	1:A:904:ILE:HD12	1.52	0.91
1:B:542:LEU:HD23	1:B:589:MET:CE	2.06	0.85
1:B:542:LEU:HD23	1:B:589:MET:HE2	1.57	0.84
1:B:365:GLN:O	1:B:369:VAL:HG23	1.80	0.80
1:B:216:THR:HG21	1:B:448:GLY:O	1.88	0.74
1:A:875:MET:HE1	1:A:890:PHE:CZ	2.31	0.66
1:A:435:VAL:CG1	1:A:438:CYS:SG	2.85	0.65
1:A:435:VAL:HG11	1:A:438:CYS:SG	2.37	0.64
1:A:281:ILE:HG13	1:A:305:LEU:HD13	1.79	0.64
1:A:758:LEU:O	1:A:762:THR:HG23	1.98	0.63
1:A:775:ARG:HH11	1:A:775:ARG:CG	2.12	0.63
1:A:652:VAL:HB	1:A:905:THR:HG23	1.81	0.62
1:A:501:LYS:HB3	1:A:695:LEU:HD23	1.80	0.62
1:A:796:ALA:O	1:A:799:GLN:HB2	1.99	0.62
1:A:612:ASP:O	1:A:634:VAL:HG21	1.99	0.61
1:A:306:VAL:HG22	1:A:378:VAL:HG21	1.82	0.61
1:B:387:ALA:HB2	1:B:427:LEU:HD12	1.83	0.61
1:A:306:VAL:CG2	1:A:378:VAL:HG21	2.30	0.60
1:A:903:LEU:O	1:A:907:VAL:HG23	2.02	0.59
1:B:542:LEU:HD23	1:B:589:MET:HE3	1.84	0.59
1:A:676:LEU:HD13	1:A:838:MET:CE	2.33	0.58
1:B:763:LYS:HB3	1:B:772:ILE:HD11	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:515:TYR:O	1:A:611:LEU:HD13	2.05	0.57
1:B:412:VAL:HG12	1:B:413:ASP:N	2.20	0.57
1:B:862:GLY:O	1:B:866:LYS:HG2	2.04	0.57
1:B:515:TYR:O	1:B:611:LEU:HD13	2.05	0.57
1:A:73:ASP:HB3	1:A:75:THR:HG23	1.86	0.56
1:B:637:LEU:HD23	1:B:651:VAL:HG21	1.88	0.56
1:A:797:LEU:HD11	1:A:817:ILE:HG23	1.87	0.55
1:B:307:ARG:O	1:B:311:VAL:HG23	2.05	0.55
1:B:663:MET:HG3	1:B:904:ILE:HD11	1.87	0.55
1:A:799:GLN:O	1:A:802:ALA:HB3	2.06	0.55
1:B:515:TYR:CD2	1:B:703:MET:HE1	2.42	0.55
1:B:874:VAL:O	1:B:878:THR:OG1	2.21	0.55
1:A:786:PHE:O	1:A:790:ILE:HD12	2.06	0.55
1:B:320:PHE:CE2	1:B:361:LEU:HD13	2.42	0.55
1:A:778:THR:HG22	1:A:781:ILE:HG23	1.89	0.54
1:B:79:GLU:O	1:B:148:LEU:HD22	2.08	0.54
1:A:797:LEU:HD11	1:A:817:ILE:CG2	2.39	0.53
1:B:383:ALA:O	1:B:427:LEU:HD13	2.08	0.53
1:B:542:LEU:CD2	1:B:589:MET:HE3	2.38	0.53
1:B:678:VAL:HG12	1:B:864:LEU:HD23	1.90	0.53
1:B:216:THR:CG2	1:B:452:GLY:H	2.01	0.53
1:A:854:LEU:HD21	1:A:856:VAL:HG21	1.89	0.53
1:B:138:MET:SD	1:B:199:PHE:CD2	3.02	0.53
1:B:353:ARG:NH2	1:B:363:PRO:O	2.42	0.53
1:B:875:MET:HE2	1:B:876:HIS:CD2	2.44	0.53
1:A:407:ARG:HD3	1:A:439:ASP:OD2	2.08	0.52
1:A:854:LEU:HG	1:A:856:VAL:HG23	1.91	0.52
1:A:261:TRP:HH2	1:A:390:LEU:HD13	1.73	0.52
1:A:723:THR:O	1:A:727:VAL:HG23	2.10	0.52
1:B:193:ILE:HG23	1:B:199:PHE:CE1	2.45	0.52
1:A:435:VAL:HG12	1:A:438:CYS:SG	2.50	0.52
1:B:264:PHE:O	1:B:293:PHE:HB2	2.10	0.52
1:A:875:MET:HE1	1:A:879:VAL:HG21	1.92	0.51
1:B:564:GLN:HA	1:B:564:GLN:OE1	2.10	0.51
1:B:295:LYS:HA	1:B:301:TYR:CD2	2.45	0.51
1:B:431:VAL:HG11	1:B:442:PHE:HZ	1.76	0.51
1:A:17:ASP:CG	1:A:18:GLN:H	2.13	0.51
1:B:128:ASP:OD1	1:B:195:ARG:NH1	2.43	0.51
1:B:309:ILE:HG22	1:B:374:ILE:HD13	1.91	0.51
1:A:67:PHE:HA	1:A:255:MET:HE2	1.92	0.51
1:A:875:MET:O	1:A:879:VAL:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:414:GLY:HA2	3:B:1002:BG6:O6	2.11	0.50
1:B:196:ARG:HB3	1:B:198:ASP:HB2	1.92	0.50
1:A:734:LEU:C	1:A:736:PRO:HD3	2.32	0.50
1:B:212:GLY:O	1:B:216:THR:HG23	2.11	0.49
1:B:305:LEU:HD21	1:B:381:ARG:NH2	2.26	0.49
1:B:528:PHE:CZ	1:B:906:ALA:HB2	2.47	0.49
1:A:326:PRO:HA	1:A:329:LEU:HD12	1.94	0.49
1:B:311:VAL:O	1:B:315:LYS:HG3	2.12	0.49
1:A:762:THR:HG21	1:A:768:PHE:HD1	1.78	0.49
1:A:138:MET:SD	1:A:199:PHE:CD2	3.06	0.48
1:A:775:ARG:HH11	1:A:775:ARG:HG3	1.78	0.48
1:A:712:PHE:HB3	1:A:741:PHE:HB2	1.95	0.48
1:B:865:TYR:HE1	1:B:875:MET:HE1	1.78	0.48
1:A:690:MET:CE	1:A:696:VAL:HG12	2.44	0.48
1:B:566:VAL:HG13	1:B:574:LEU:HA	1.96	0.48
1:A:502:GLU:HG2	1:A:503:THR:HG23	1.96	0.47
1:B:689:GLU:OE2	1:B:702:ARG:HG2	2.14	0.47
1:A:349:ILE:HG13	1:A:350:ARG:N	2.28	0.47
1:B:242:MET:HA	1:B:245:ILE:HD12	1.95	0.47
1:A:484:LEU:HB3	1:A:882:LEU:HD11	1.97	0.47
1:B:200:ASP:OD1	1:B:201:ILE:N	2.44	0.47
1:B:578:ILE:O	1:B:582:ILE:HG13	2.15	0.47
1:B:657:ASP:OD1	1:B:657:ASP:N	2.48	0.47
1:B:674:VAL:HB	1:B:858:VAL:HG22	1.96	0.47
1:A:364:THR:O	1:A:365:GLN:C	2.52	0.47
1:B:279:GLN:O	1:B:283:MET:HG2	2.15	0.47
1:A:492:LYS:O	1:A:496:GLU:HG3	2.14	0.46
1:B:209:ASP:N	1:B:209:ASP:OD1	2.45	0.46
1:B:676:LEU:HD13	1:B:838:MET:CE	2.44	0.46
1:B:542:LEU:CD2	1:B:589:MET:CE	2.87	0.46
1:A:409:THR:HG23	1:A:441:ARG:HG2	1.97	0.46
1:B:515:TYR:HA	1:B:703:MET:CE	2.46	0.45
1:A:164:ASP:HA	1:A:204:VAL:O	2.16	0.45
1:B:309:ILE:CG2	1:B:374:ILE:HD13	2.46	0.45
1:A:530:ALA:HB3	1:A:541:LEU:HB2	1.98	0.44
1:A:878:THR:HG22	1:A:882:LEU:HD12	1.99	0.44
1:A:676:LEU:HD13	1:A:838:MET:HE1	1.99	0.44
1:A:690:MET:HE1	1:A:696:VAL:HG12	2.00	0.44
1:A:659:VAL:HG22	1:A:704:CYS:SG	2.58	0.44
1:A:803:ILE:CG1	1:B:58:THR:HG21	2.47	0.44
1:A:828:ARG:O	1:A:832:GLN:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:804:LEU:HD21	1:A:819:VAL:HG11	1.99	0.44
1:B:167:PHE:HA	1:B:183:ARG:O	2.18	0.44
1:B:324:LEU:HD22	1:B:329:LEU:HD21	1.99	0.43
1:B:370:ALA:O	1:B:373:ARG:N	2.50	0.43
1:A:316:GLU:HB3	1:A:318:LEU:HD13	1.99	0.43
1:A:316:GLU:HB3	1:A:318:LEU:CD1	2.48	0.43
1:A:579:VAL:HG12	1:A:644:ARG:HD3	2.00	0.43
1:B:349:ILE:HD13	1:B:369:VAL:HG22	2.00	0.43
1:A:302:MET:CE	1:A:336:THR:HG23	2.49	0.42
1:A:40:SER:CB	1:A:388:ALA:O	2.67	0.42
1:A:803:ILE:HG13	1:B:58:THR:HG21	2.01	0.42
1:B:281:ILE:HD12	1:B:305:LEU:HD13	2.01	0.42
1:A:600:PHE:HB3	1:A:654:VAL:HG22	2.01	0.42
1:A:708:GLU:OE2	2:A:1003:GLC:O3	2.30	0.42
1:B:720:ASP:OD1	1:B:720:ASP:N	2.51	0.42
1:A:218:GLY:HA2	1:A:221:ASP:O	2.19	0.42
1:B:612:ASP:O	1:B:634:VAL:HG21	2.20	0.42
1:B:380:THR:OG1	1:B:426:ARG:CZ	2.67	0.42
1:B:67:PHE:HA	1:B:255:MET:HE2	2.02	0.42
1:A:66:THR:OG1	1:A:256:CYS:HB3	2.20	0.42
1:B:281:ILE:CD1	1:B:305:LEU:HD13	2.50	0.42
1:A:528:PHE:CD2	1:A:597:PRO:HG2	2.55	0.41
1:A:529:LEU:HB3	1:A:598:LEU:HD12	2.01	0.41
1:A:650:ASP:O	1:A:652:VAL:HG23	2.20	0.41
1:A:663:MET:HB2	1:A:900:GLY:HA3	2.02	0.41
1:B:349:ILE:HG23	1:B:368:CYS:HB3	2.02	0.41
1:A:669:ASP:OD1	1:A:669:ASP:C	2.59	0.41
1:A:637:LEU:O	1:A:641:ILE:HG13	2.21	0.41
1:A:612:ASP:HB3	1:A:652:VAL:O	2.21	0.41
1:B:734:LEU:C	1:B:736:PRO:HD3	2.40	0.41
1:A:305:LEU:HD21	1:A:381:ARG:NH2	2.35	0.41
1:A:417:TYR:OH	1:A:428:HIS:NE2	2.33	0.41
1:A:198:ASP:HB3	1:A:199:PHE:HD1	1.86	0.40
1:A:386:CYS:O	1:A:390:LEU:HB2	2.21	0.40
1:A:70:SER:OG	1:A:71:THR:O	2.38	0.40
1:B:302:MET:HG2	1:B:378:VAL:HG12	2.02	0.40
1:B:875:MET:CE	1:B:876:HIS:CD2	3.04	0.40
1:B:261:TRP:CD1	1:B:261:TRP:C	2.94	0.40
1:B:484:LEU:HB3	1:B:882:LEU:HD11	2.03	0.40
1:B:370:ALA:O	1:B:371:THR:C	2.60	0.40
1:A:657:ASP:N	1:A:657:ASP:OD1	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:797:LEU:CD1	1:A:817:ILE:HG22	2.52	0.40
1:B:762:THR:O	1:B:770:GLY:HA2	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	859/902 (95%)	825 (96%)	30 (4%)	4 (0%)	29	34
1	B	853/902 (95%)	815 (96%)	36 (4%)	2 (0%)	47	57
All	All	1712/1804 (95%)	1640 (96%)	66 (4%)	6 (0%)	34	41

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	565	GLU
1	B	198	ASP
1	A	198	ASP
1	A	670	PRO
1	A	769	ARG
1	B	143	ILE

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	702/763 (92%)	683 (97%)	19 (3%)	44	57
1	B	708/763 (93%)	692 (98%)	16 (2%)	50	63
All	All	1410/1526 (92%)	1375 (98%)	35 (2%)	47	60

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	PHE
1	A	198	ASP
1	A	252	GLU
1	A	261	TRP
1	A	302	MET
1	A	408	SER
1	A	441	ARG
1	A	443	LEU
1	A	471	GLN
1	A	493	VAL
1	A	536	THR
1	A	598	LEU
1	A	709	TRP
1	A	742	GLU
1	A	773	SER
1	A	775	ARG
1	A	811	SER
1	A	875	MET
1	A	904	ILE
1	B	20	GLN
1	B	33	ASP
1	B	54	THR
1	B	261	TRP
1	B	279	GLN
1	B	281	ILE
1	B	318	LEU
1	B	324	LEU
1	B	431	VAL
1	B	709	TRP
1	B	738	LYS
1	B	773	SER
1	B	774	GLU
1	B	790	ILE
1	B	806	HIS
1	B	843	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 28 are unknown - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	BG6	A	1004	-	16,16,16	0.85	0	24,24,24	1.47	3 (12%)
3	BG6	B	1002	-	16,16,16	1.12	2 (12%)	24,24,24	1.26	2 (8%)
2	GLC	B	1001	-	12,12,12	0.71	0	17,17,17	0.66	0
2	GLC	A	1001	-	12,12,12	1.23	1 (8%)	17,17,17	1.69	4 (23%)
3	BG6	A	1002	-	16,16,16	0.93	0	24,24,24	0.99	0
2	GLC	B	1003	-	12,12,12	0.80	0	17,17,17	1.15	1 (5%)
3	BG6	B	1004	-	16,16,16	0.58	0	24,24,24	1.48	4 (16%)
2	GLC	A	1003	-	12,12,12	0.89	0	17,17,17	0.91	1 (5%)

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	BG6	A	1004	-	-	2/6/26/26	0/1/1/1
3	BG6	B	1002	-	-	1/6/26/26	0/1/1/1
2	GLC	B	1001	-	-	0/2/22/22	0/1/1/1
2	GLC	A	1001	-	-	0/2/22/22	0/1/1/1
3	BG6	A	1002	-	-	1/6/26/26	0/1/1/1
2	GLC	B	1003	-	-	0/2/22/22	0/1/1/1
3	BG6	B	1004	-	-	1/6/26/26	0/1/1/1
2	GLC	A	1003	-	-	2/2/22/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1002	BG6	P-O3P	2.50	1.58	1.50
3	B	1002	BG6	O3-C3	2.07	1.47	1.43
2	A	1001	GLC	O4-C4	2.03	1.47	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1004	BG6	O5-C1-C2	-4.29	102.63	110.28
2	A	1001	GLC	O5-C1-C2	-4.08	103.01	110.28
3	A	1004	BG6	O1-C1-O5	-3.65	99.44	110.38
2	B	1003	GLC	C6-C5-C4	-3.11	105.72	113.00
3	B	1002	BG6	O1P-P-O6	2.85	114.31	106.73
3	A	1004	BG6	O2P-P-O6	2.82	114.22	106.73
3	B	1004	BG6	O2P-P-O1P	2.68	117.88	107.64
3	B	1002	BG6	O6-P-O3P	-2.60	99.17	106.47
3	B	1004	BG6	C4-C3-C2	-2.57	106.33	110.82
2	A	1001	GLC	O5-C5-C4	2.54	114.31	109.69
2	A	1001	GLC	O3-C3-C4	-2.35	104.91	110.35
2	A	1001	GLC	O3-C3-C2	2.34	115.77	110.35
2	A	1003	GLC	C6-C5-C4	-2.24	107.76	113.00
3	A	1004	BG6	O2P-P-O3P	-2.20	102.08	110.68
3	B	1004	BG6	C1-C2-C3	-2.13	105.89	110.31

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1004	BG6	C6-O6-P-O2P
3	B	1002	BG6	C5-C6-O6-P
2	A	1003	GLC	O5-C5-C6-O6

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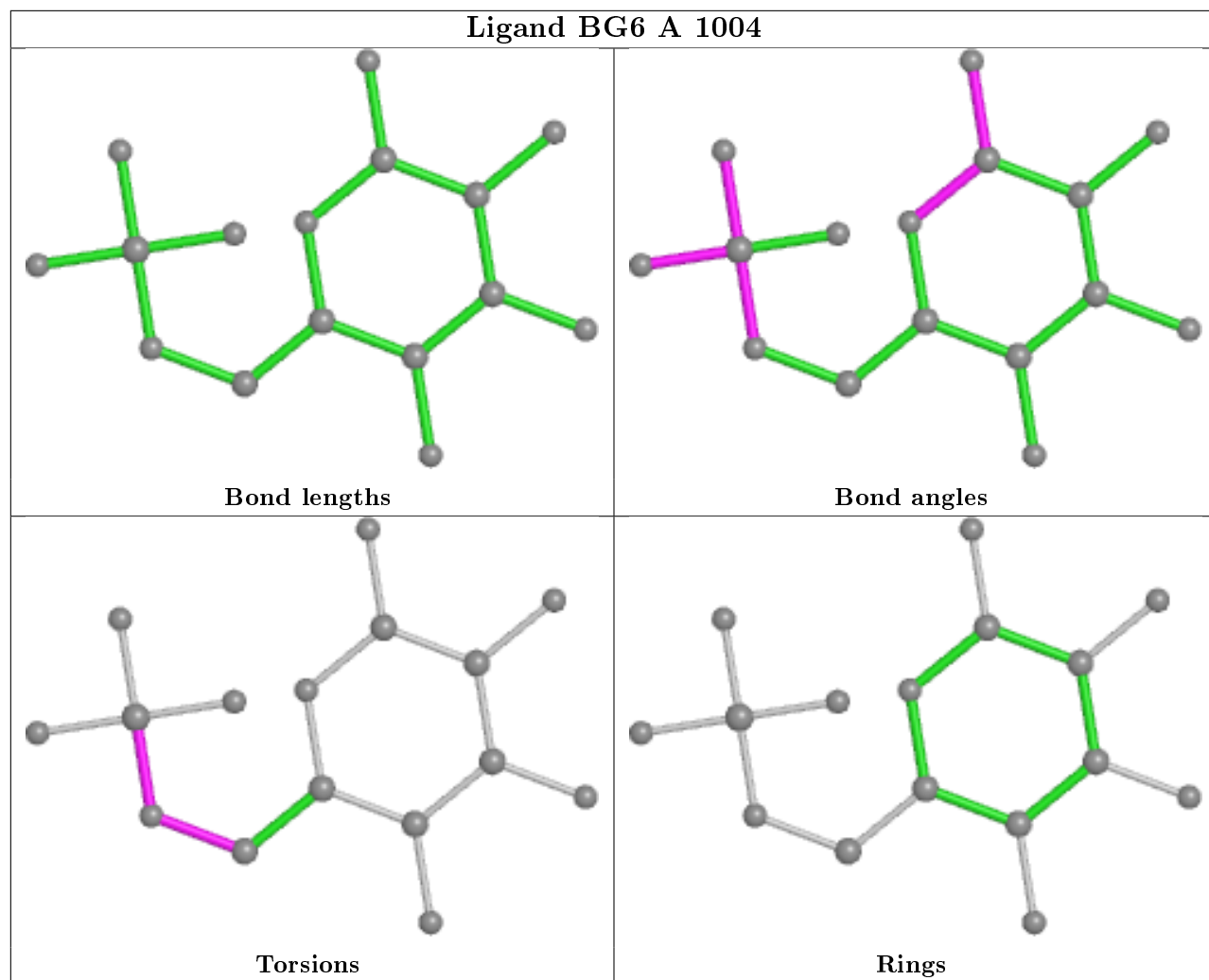
Mol	Chain	Res	Type	Atoms
2	A	1003	GLC	C4-C5-C6-O6
3	A	1002	BG6	C5-C6-O6-P
3	B	1004	BG6	C5-C6-O6-P
3	A	1004	BG6	C5-C6-O6-P

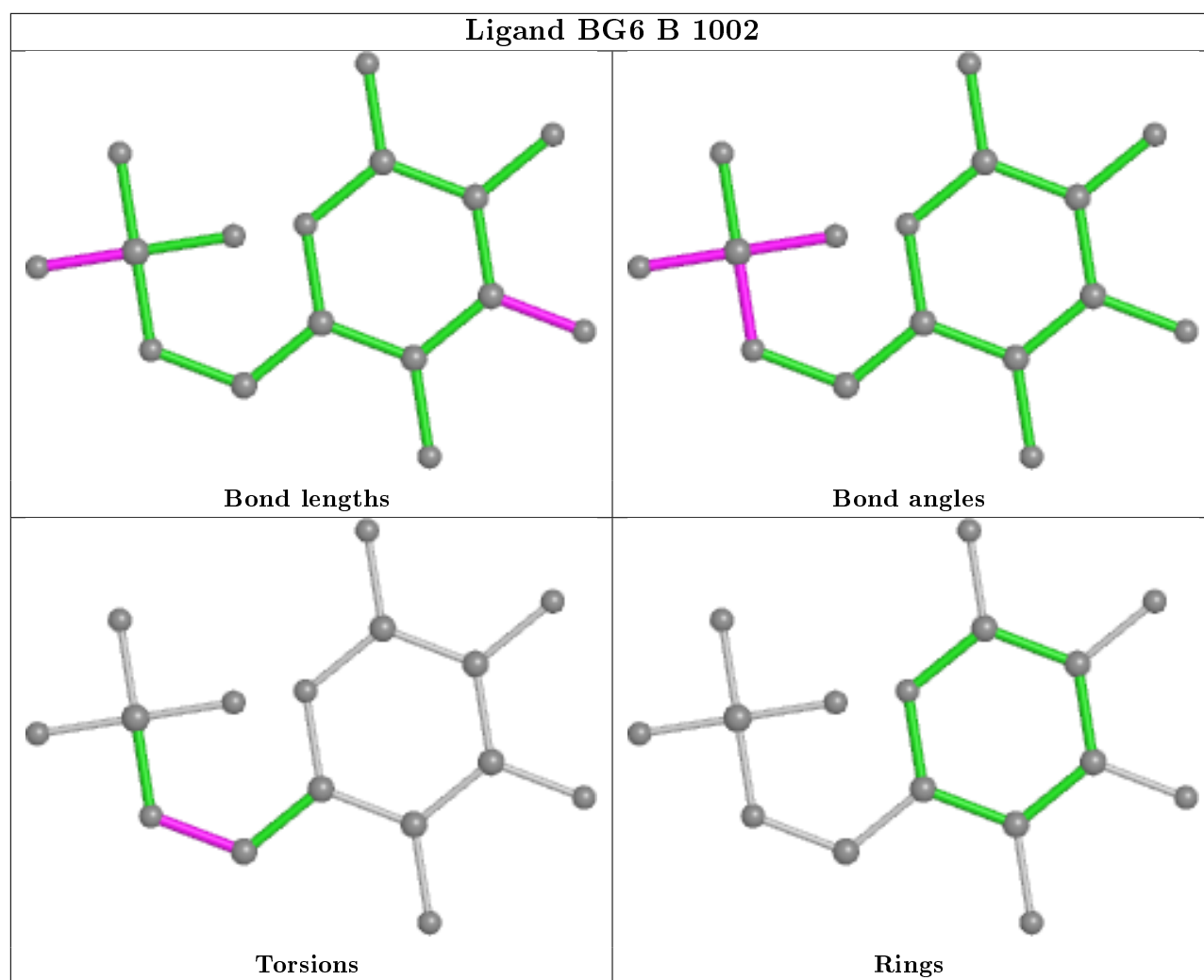
There are no ring outliers.

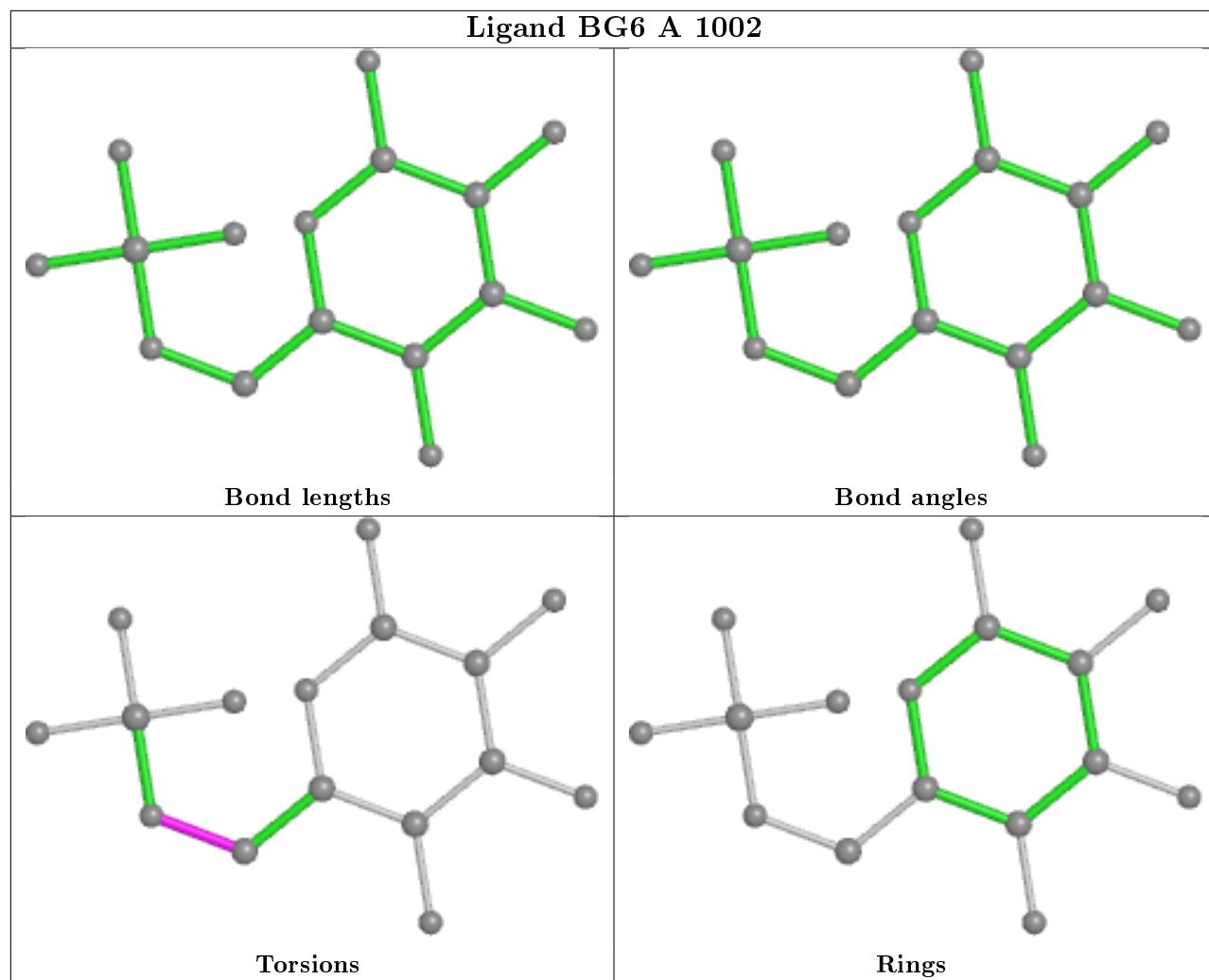
2 monomers are involved in 2 short contacts:

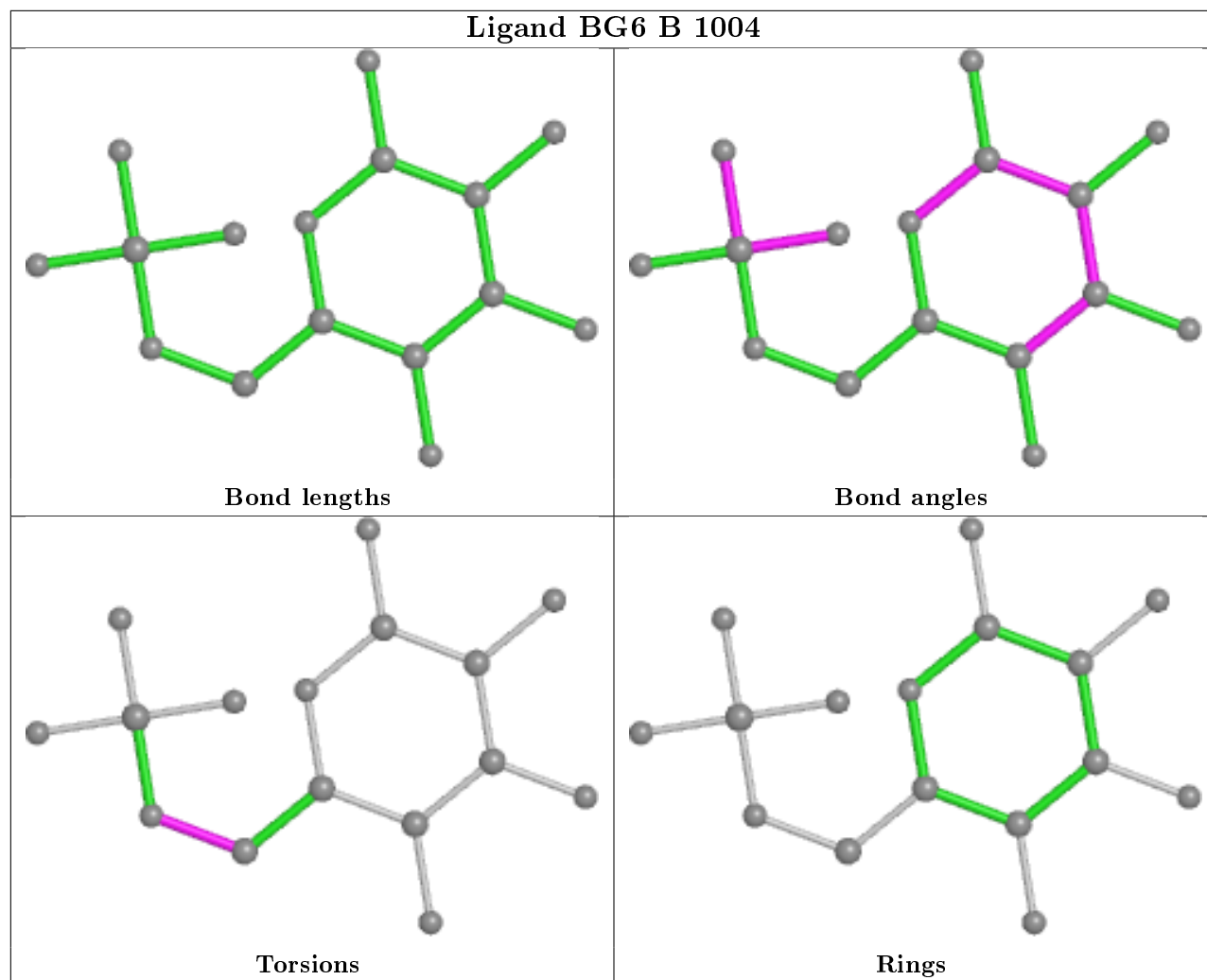
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1002	BG6	1	0
2	A	1003	GLC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [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	871/902 (96%)	0.04	30 (3%) 45 41	31, 49, 68, 85	0
1	B	867/902 (96%)	0.10	43 (4%) 28 26	32, 49, 73, 87	0
All	All	1738/1804 (96%)	0.07	73 (4%) 36 33	31, 49, 71, 87	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	911	ILE	4.9
1	B	27	TYR	4.8
1	B	347	ASP	4.1
1	B	222	HIS	3.9
1	A	378	VAL	3.8
1	B	344	GLY	3.8
1	B	20	GLN	3.5
1	B	437	GLY	3.5
1	B	374	ILE	3.4
1	B	375	CYS	3.3
1	A	317	GLU	3.2
1	B	310	LEU	3.2
1	B	386	CYS	3.1
1	B	19	VAL	3.1
1	A	684	ALA	3.0
1	B	530	ALA	3.0
1	B	145	ASP	2.9
1	B	378	VAL	2.9
1	B	850	GLY	2.8
1	A	912	ARG	2.8
1	B	150	LEU	2.8
1	B	24	GLN	2.8
1	B	453	ALA	2.8
1	A	418	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	82	ALA	2.7
1	A	910	ARG	2.7
1	A	754	VAL	2.7
1	B	667	PHE	2.6
1	A	705	VAL	2.6
1	A	670	PRO	2.6
1	A	867	LEU	2.6
1	B	402	GLY	2.6
1	A	796	ALA	2.6
1	A	655	VAL	2.6
1	B	438	CYS	2.5
1	A	834	CYS	2.5
1	A	142	GLN	2.4
1	A	145	ASP	2.4
1	B	153	THR	2.4
1	A	27	TYR	2.4
1	A	405	ARG	2.4
1	B	390	LEU	2.4
1	A	153	THR	2.4
1	A	305	LEU	2.4
1	A	642	HIS	2.3
1	B	593	GLY	2.3
1	B	237	CYS	2.3
1	B	324	LEU	2.3
1	B	912	ARG	2.3
1	A	745	ILE	2.3
1	A	403	GLU	2.2
1	B	365	GLN	2.2
1	B	74	GLY	2.2
1	B	306	VAL	2.2
1	B	564	GLN	2.2
1	B	81	LEU	2.2
1	A	913	GLU	2.2
1	B	82	ALA	2.2
1	B	317	GLU	2.1
1	B	389	THR	2.1
1	B	83	LEU	2.1
1	B	228	LEU	2.1
1	A	909	CYS	2.1
1	A	709	TRP	2.1
1	A	822	VAL	2.1
1	A	297	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	443	LEU	2.1
1	A	671	HIS	2.1
1	B	28	HIS	2.0
1	B	425	LYS	2.0
1	B	349	ILE	2.0
1	A	741	PHE	2.0
1	B	913	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 monosaccharides 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(\AA^2)	Q<0.9
4	UNX	B	12	1/1	-0.50	3.11	2,2,2,2	1
4	UNX	A	1007	1/1	-0.36	4.71	2,2,2,2	1
4	UNX	A	1008	1/1	-0.33	6.23	2,2,2,2	1
4	UNX	B	1005	1/1	-0.29	2.80	2,2,2,2	1
4	UNX	A	11	1/1	-0.26	2.28	2,2,2,2	1
4	UNX	B	3	1/1	-0.13	2.42	2,2,2,2	1
4	UNX	A	8	1/1	0.00	1.52	2,2,2,2	1
4	UNX	B	1009	1/1	0.03	0.89	2,2,2,2	1
4	UNX	B	13	1/1	0.08	1.28	2,2,2,2	1
4	UNX	B	5	1/1	0.13	0.68	2,2,2,2	1
4	UNX	B	1013	1/1	0.14	1.20	5,5,5,5	1
4	UNX	B	1007	1/1	0.18	1.29	2,2,2,2	1
4	UNX	B	1012	1/1	0.20	2.06	2,2,2,2	1
4	UNX	B	7	1/1	0.20	2.09	2,2,2,2	1
4	UNX	A	1009	1/1	0.26	1.34	2,2,2,2	1

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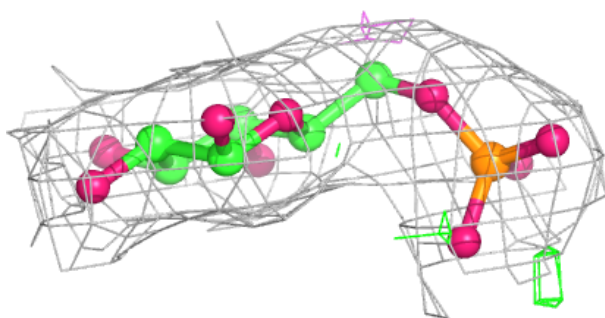
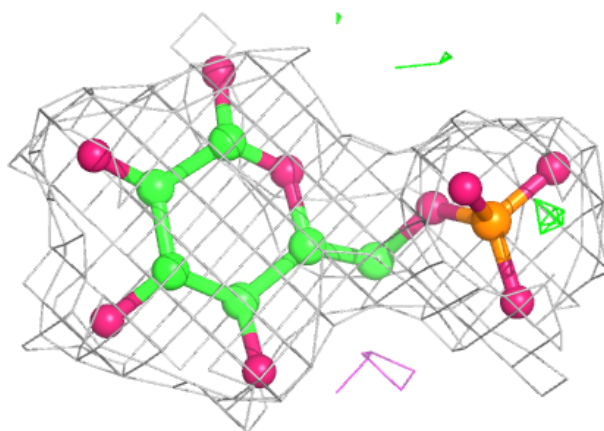
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	UNX	A	4	1/1	0.26	1.18	2,2,2,2	1
4	UNX	B	14	1/1	0.27	2.65	2,2,2,2	1
4	UNX	A	9	1/1	0.28	1.26	2,2,2,2	1
4	UNX	B	1006	1/1	0.31	2.40	2,2,2,2	1
4	UNX	A	1005	1/1	0.32	2.27	2,2,2,2	1
4	UNX	B	1011	1/1	0.34	4.14	2,2,2,2	1
4	UNX	B	1010	1/1	0.40	1.95	2,2,2,2	1
4	UNX	B	2	1/1	0.43	2.60	2,2,2,2	1
4	UNX	B	1008	1/1	0.47	4.46	2,2,2,2	1
4	UNX	A	10	1/1	0.51	3.33	2,2,2,2	1
4	UNX	A	6	1/1	0.58	1.97	2,2,2,2	1
4	UNX	A	1006	1/1	0.59	2.11	2,2,2,2	1
4	UNX	A	1	1/1	0.63	1.42	2,2,2,2	1
2	GLC	A	1001	12/12	0.96	0.19	29,32,34,34	0
3	BG6	A	1004	16/16	0.96	0.10	40,45,48,53	0
3	BG6	B	1002	16/16	0.97	0.08	48,52,54,55	0
2	GLC	A	1003	12/12	0.97	0.17	35,40,41,42	0
3	BG6	B	1004	16/16	0.97	0.11	33,39,43,44	0
3	BG6	A	1002	16/16	0.98	0.10	33,36,38,40	0
2	GLC	B	1003	12/12	0.98	0.20	25,32,34,35	0
2	GLC	B	1001	12/12	0.98	0.13	32,38,39,39	0

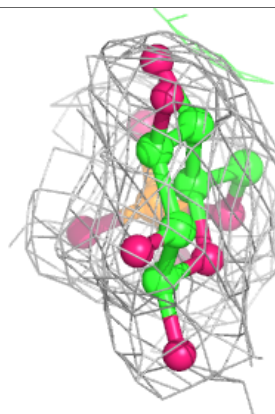
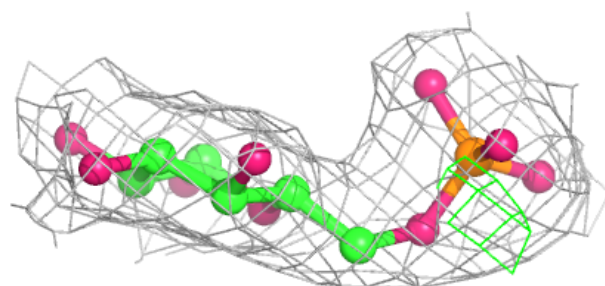
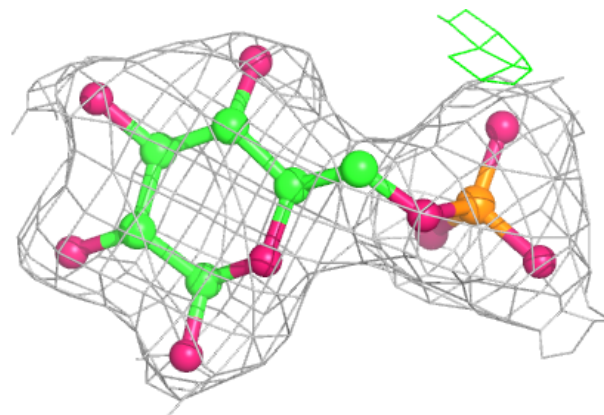
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around BG6 A 1004:

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

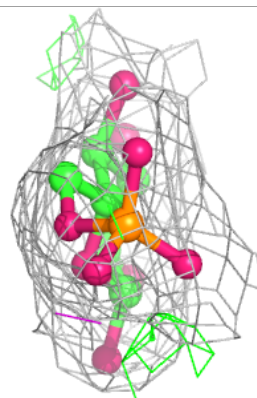
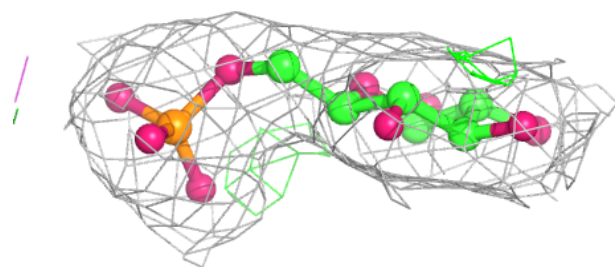
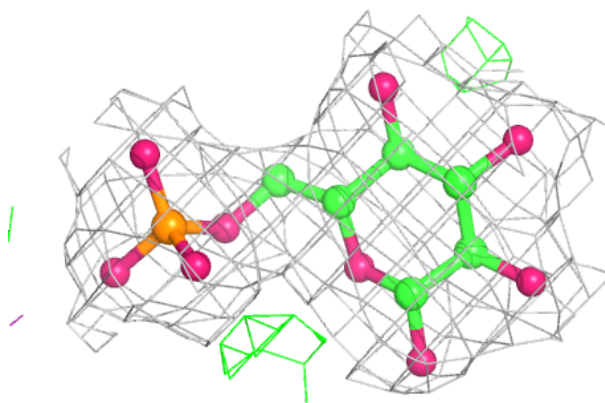
**Electron density around BG6 B 1002:**

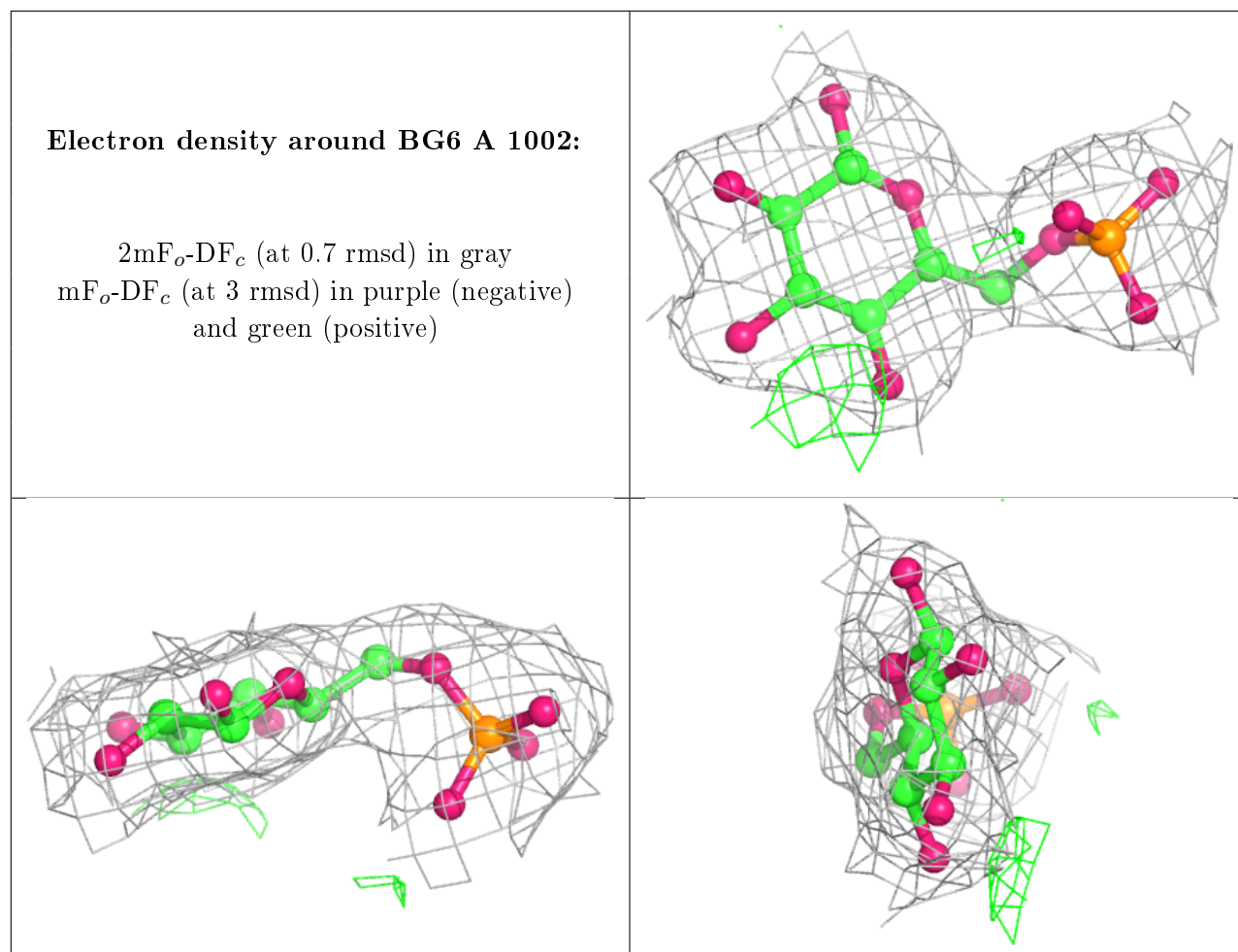
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around BG6 B 1004:

$2mF_o - DF_c$ (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.