



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2020 – 09:28 AM BST

PDB ID : 1J12
Title : Beta-Amylase from Bacillus cereus var. mycoides in Complex with alpha-EBG
Authors : Oyama, T.; Miyake, H.; Kusunoki, M.; Nitta, Y.
Deposited on : 2002-11-25
Resolution : 2.10 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

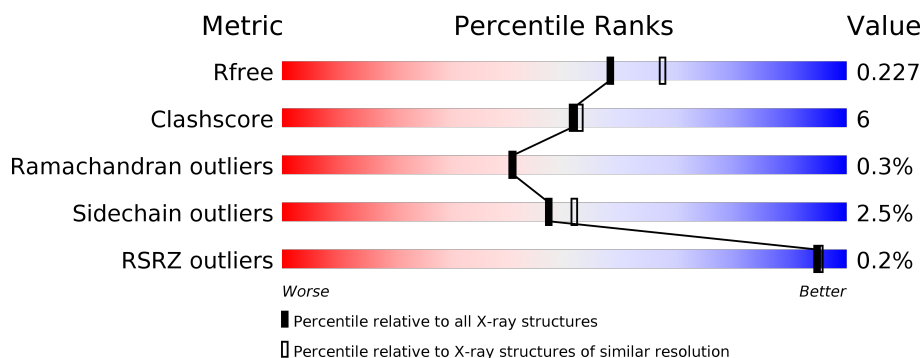
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	516	<div> <div style="width: 86%;"></div> <div style="width: 13%;"></div> <div style="width: 1%;"></div> </div> <div>86% 13% .</div>
1	B	516	<div> <div style="width: 83%;"></div> <div style="width: 16%;"></div> <div style="width: 1%;"></div> </div> <div>83% 16% .</div>
1	C	516	<div> <div style="width: 84%;"></div> <div style="width: 16%;"></div> </div> <div>84% 16%</div>
1	D	516	<div> <div style="width: 83%;"></div> <div style="width: 16%;"></div> </div> <div>83% 16%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EBG	B	601	X	-	-	-
2	EBG	C	601	X	-	-	-
2	EBG	D	601	X	-	-	-

2 Entry composition [i](#)

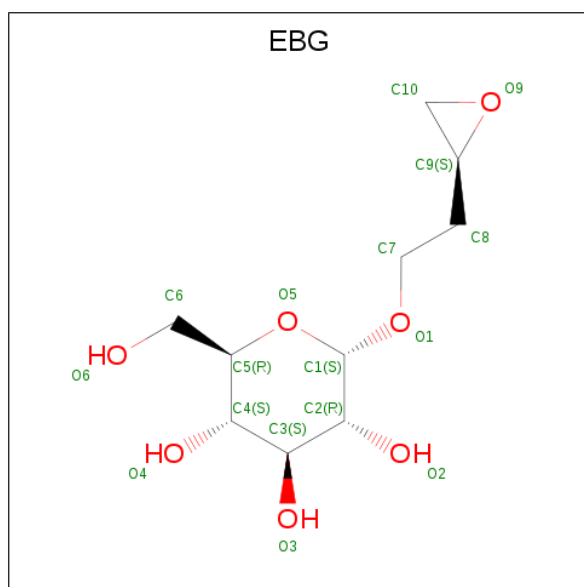
There are 4 unique types of molecules in this entry. The entry contains 17021 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Beta-amylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			
1	B	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			
1	C	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			
1	D	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			

- Molecule 2 is 2-[(2S)-oxiran-2-yl]ethyl alpha-D-glucopyranoside (three-letter code: EBG) (formula: C₁₀H₁₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			17	10	7		
2	B	1	Total	C	O	0	0
			17	10	7		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			17	10	7		
2	D	1	Total	C	O	0	0
			17	10	7		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		
3	A	1	Total	Ca	0	0
			1	1		
3	D	1	Total	Ca	0	0
			1	1		
3	C	1	Total	Ca	0	0
			1	1		

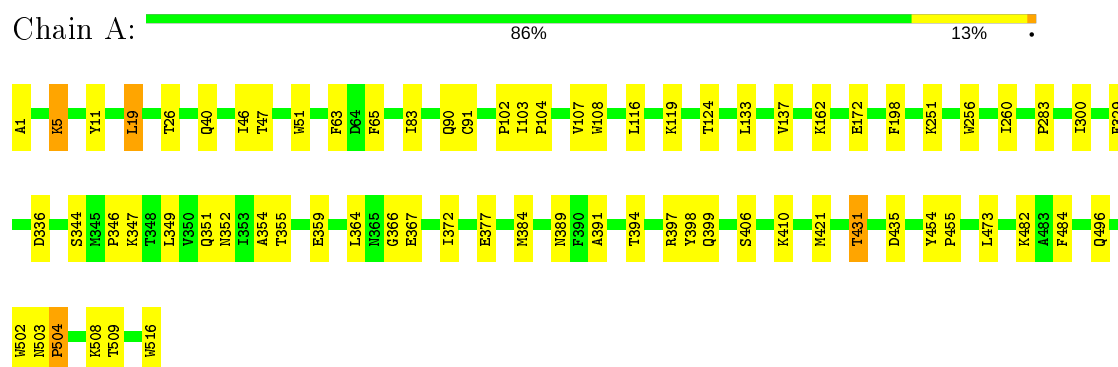
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	128	Total	O	0	0
			128	128		
4	B	101	Total	O	0	0
			101	101		
4	C	131	Total	O	0	0
			131	131		
4	D	113	Total	O	0	0
			113	113		

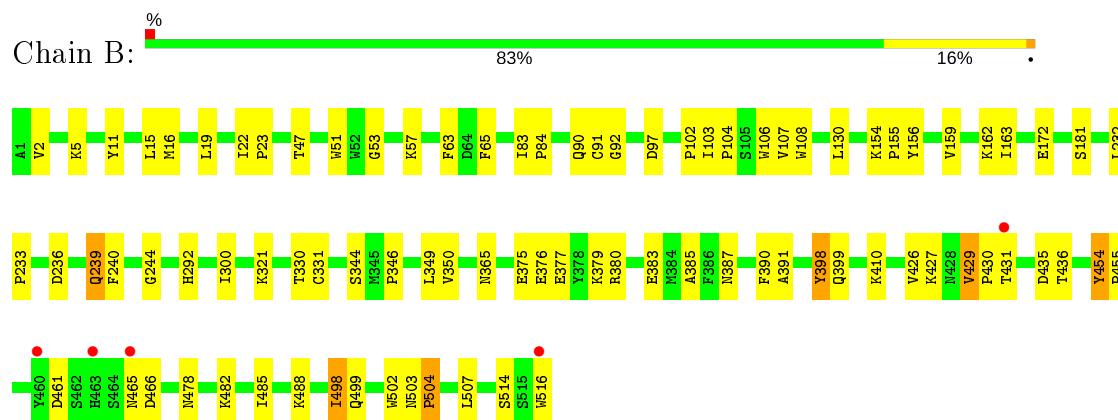
3 Residue-property plots

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.

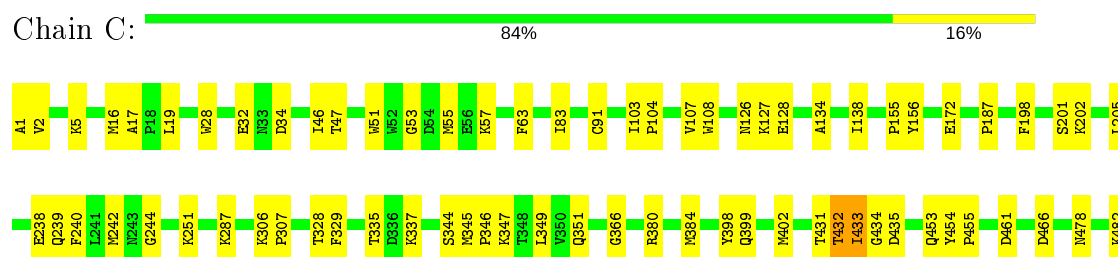
• Molecule 1: Beta-amylase



• Molecule 1: Beta-amylase



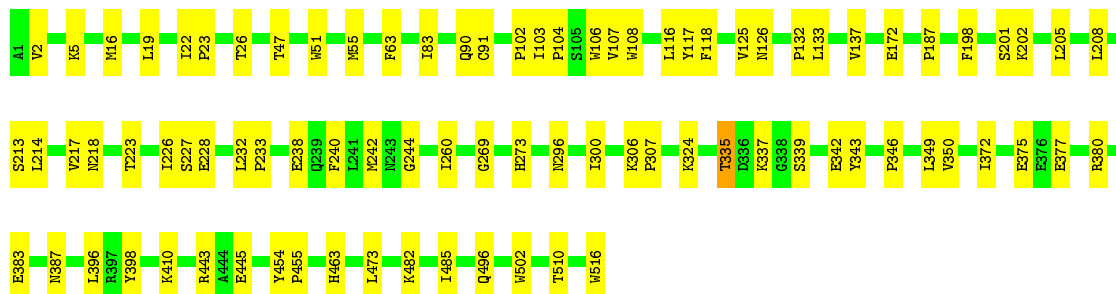
• Molecule 1: Beta-amylase





• Molecule 1: Beta-amylase

Chain D: 83% 16%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.90 Å 112.90 Å 146.20 Å 90.00° 105.80° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10 73.38 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.10) 68.0 (73.38-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.00 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.198 , 0.253 0.181 , 0.227	Depositor DCC
R_{free} test set	6284 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	17.4	Xtriage
Anisotropy	0.479	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 57.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17021	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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: CA, EBG

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	0.42	1/4234 (0.0%)	0.60	0/5751
1	B	0.42	1/4234 (0.0%)	0.59	0/5751
1	C	0.42	1/4234 (0.0%)	0.60	0/5751
1	D	0.42	1/4234 (0.0%)	0.60	0/5751
All	All	0.42	4/16936 (0.0%)	0.60	0/23004

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	172	GLU	CD-OE2	13.54	1.40	1.25
1	D	172	GLU	CD-OE2	13.27	1.40	1.25
1	B	172	GLU	CD-OE2	13.05	1.40	1.25
1	C	172	GLU	CD-OE2	12.49	1.39	1.25

There are no bond angle outliers.

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	4119	0	3984	35	0
1	B	4119	0	3984	50	0
1	C	4119	0	3984	49	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	4119	0	3984	47	0
2	A	17	0	17	0	0
2	B	17	0	17	0	0
2	C	17	0	17	0	0
2	D	17	0	17	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	128	0	0	0	0
4	B	101	0	0	1	0
4	C	131	0	0	1	0
4	D	113	0	0	1	0
All	All	17021	0	16004	181	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 (181) 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:C:432:THR:HG22	1:C:433:ILE:HD13	1.65	0.78
1:C:434:GLY:HA3	1:C:488:LYS:HG2	1.67	0.77
1:D:16:MET:HG2	4:D:712:HOH:O	1.88	0.73
1:C:2:VAL:O	1:C:5:LYS:HG2	1.89	0.72
1:B:431:THR:HG23	1:B:435:ASP:HB2	1.72	0.70
1:A:431:THR:HG23	1:A:435:ASP:HB2	1.75	0.69
1:D:103:ILE:HD11	1:D:108:TRP:CZ2	2.28	0.68
1:D:198:PHE:O	1:D:202:LYS:HG2	1.94	0.68
1:C:482:LYS:HG2	1:C:496:GLN:HB3	1.75	0.67
1:B:485:ILE:HG13	1:B:516:TRP:CH2	2.30	0.66
1:C:1:ALA:HB3	1:C:5:LYS:HG3	1.81	0.62
1:B:482:LYS:HG3	1:B:499:GLN:HA	1.81	0.62
1:D:335:THR:HG23	1:D:380:ARG:HD2	1.81	0.62
1:B:11:TYR:HA	1:B:391:ALA:O	2.00	0.62
1:C:433:ILE:HD13	1:C:433:ILE:H	1.63	0.62
1:D:133:LEU:HB3	1:D:198:PHE:HE2	1.64	0.62
1:B:426:VAL:HA	1:B:514:SER:O	2.00	0.62
1:C:482:LYS:HD3	1:C:496:GLN:O	2.00	0.61
1:A:1:ALA:HB3	1:A:5:LYS:HD2	1.83	0.59
1:B:507:LEU:H	1:B:507:LEU:HD12	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:300:ILE:HG13	1:D:300:ILE:O	2.02	0.59
1:B:498:ILE:H	1:B:498:ILE:HD12	1.68	0.59
1:A:90:GLN:HG3	1:A:102:PRO:HA	1.85	0.58
1:A:482:LYS:HE2	1:A:496:GLN:O	2.04	0.58
1:A:133:LEU:HB3	1:A:198:PHE:HE2	1.69	0.58
1:C:434:GLY:CA	1:C:488:LYS:HG2	2.34	0.57
1:A:347:LYS:NZ	1:A:384:MET:SD	2.78	0.57
1:B:346:PRO:O	1:B:350:VAL:HG23	2.05	0.56
1:D:482:LYS:HG2	1:D:496:GLN:HB3	1.87	0.56
1:D:269:GLY:O	1:D:273:HIS:HD2	1.89	0.56
1:C:510:THR:HG22	1:C:511:SER:H	1.71	0.55
1:C:238:GLU:O	1:C:242:MET:HG3	2.06	0.55
1:D:454:TYR:N	1:D:455:PRO:HD2	2.22	0.55
1:D:201:SER:O	1:D:205:LEU:HG	2.06	0.55
1:D:16:MET:CE	1:D:396:LEU:HB2	2.37	0.55
1:A:454:TYR:N	1:A:455:PRO:HD2	2.22	0.55
1:B:383:GLU:O	1:B:387:ASN:HB2	2.05	0.55
1:C:344:SER:C	1:C:346:PRO:HD3	2.27	0.54
1:C:335:THR:HG23	1:C:380:ARG:HD3	1.88	0.54
1:B:236:ASP:HB3	1:B:239:GLN:HG3	1.90	0.54
1:B:427:LYS:HG2	1:B:466:ASP:HB2	1.89	0.54
1:B:376:GLU:O	1:B:380:ARG:HG3	2.07	0.54
1:D:90:GLN:HG3	1:D:102:PRO:HA	1.90	0.54
1:B:92:GLY:HA2	1:B:97:ASP:HB3	1.89	0.54
1:A:372:ILE:HG23	1:A:377:GLU:HB2	1.90	0.53
1:C:497:THR:OG1	1:C:498:ILE:HD12	2.08	0.53
1:D:117:TYR:HB2	1:D:126:ASN:O	2.09	0.53
1:D:208:LEU:HD11	1:D:228:GLU:HB3	1.89	0.53
1:D:104:PRO:HB2	1:D:107:VAL:HG23	1.90	0.53
1:C:155:PRO:HG2	1:C:156:TYR:CE1	2.43	0.53
1:A:406:SER:O	1:A:410:LYS:HG2	2.09	0.53
1:D:375:GLU:HB3	1:D:410:LYS:HD3	1.90	0.53
1:D:2:VAL:O	1:D:5:LYS:HG2	2.09	0.52
1:C:431:THR:HG22	1:C:432:THR:N	2.25	0.52
1:C:454:TYR:N	1:C:455:PRO:HD2	2.25	0.52
1:B:84:PRO:HG2	1:B:163:ILE:HD13	1.91	0.51
1:B:53:GLY:O	1:B:57:LYS:HB3	2.09	0.51
1:C:329:PHE:O	1:C:366:GLY:HA2	2.09	0.51
1:D:117:TYR:CD1	1:D:125:VAL:HG13	2.45	0.51
1:C:28:TRP:O	1:C:32:GLU:HG3	2.11	0.50
1:D:372:ILE:HG23	1:D:377:GLU:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:SER:O	1:C:205:LEU:HD13	2.11	0.50
1:C:432:THR:H	1:C:435:ASP:HB2	1.76	0.50
1:C:454:TYR:N	1:C:455:PRO:CD	2.75	0.50
1:C:337:LYS:HD3	1:C:344:SER:HB2	1.94	0.50
1:A:355:THR:O	1:A:359:GLU:HG3	2.11	0.50
1:A:367:GLU:HG2	1:A:394:THR:HB	1.93	0.50
1:B:321:LYS:HB3	1:B:321:LYS:NZ	2.27	0.50
1:C:51:TRP:CZ2	1:C:91:CYS:SG	3.05	0.50
1:B:156:TYR:O	1:B:159:VAL:HG22	2.12	0.49
1:D:218:ASN:OD1	1:D:223:THR:HG23	2.12	0.49
1:B:232:LEU:HB3	1:B:233:PRO:HD2	1.95	0.49
1:B:15:LEU:HD11	1:B:398:TYR:HA	1.95	0.49
1:B:461:ASP:O	1:B:465:ASN:N	2.46	0.49
1:A:300:ILE:O	1:A:300:ILE:HG13	2.13	0.49
1:C:351:GLN:HG2	1:C:384:MET:HE1	1.95	0.48
1:C:507:LEU:O	1:C:507:LEU:HD13	2.13	0.48
1:A:116:LEU:HD21	1:A:137:VAL:HG21	1.95	0.48
1:D:133:LEU:HB3	1:D:198:PHE:CE2	2.47	0.48
1:C:47:THR:HG22	1:C:83:ILE:HB	1.95	0.48
1:B:63:PHE:HB3	1:B:65:PHE:CE1	2.48	0.48
1:A:344:SER:C	1:A:346:PRO:HD3	2.34	0.48
1:A:347:LYS:O	1:A:351:GLN:HG3	2.13	0.48
1:B:454:TYR:N	1:B:455:PRO:HD2	2.28	0.48
1:D:337:LYS:HE3	1:D:342:GLU:HB2	1.96	0.47
1:B:300:ILE:HG13	1:B:300:ILE:O	2.14	0.47
1:A:454:TYR:N	1:A:455:PRO:CD	2.78	0.47
1:B:429:VAL:HG12	1:B:516:TRP:HB2	1.97	0.47
1:B:90:GLN:HG3	1:B:102:PRO:HA	1.97	0.47
1:D:51:TRP:CZ2	1:D:91:CYS:SG	3.08	0.46
1:A:47:THR:HG22	1:A:83:ILE:HB	1.97	0.46
1:C:16:MET:HG2	4:C:705:HOH:O	2.14	0.46
1:B:16:MET:SD	1:B:47:THR:OG1	2.68	0.46
1:D:104:PRO:HB3	1:D:106:TRP:NE1	2.31	0.46
1:A:162:LYS:HG3	1:A:283:PRO:O	2.15	0.46
1:C:104:PRO:HB2	1:C:107:VAL:HG23	1.96	0.46
1:D:346:PRO:O	1:D:350:VAL:HG23	2.15	0.46
1:B:454:TYR:N	1:B:455:PRO:CD	2.79	0.46
1:B:162:LYS:HG2	1:B:163:ILE:N	2.31	0.46
1:B:461:ASP:HB3	1:B:466:ASP:OD2	2.16	0.46
1:D:324:LYS:HB3	1:D:324:LYS:NZ	2.31	0.46
1:A:336:ASP:HA	1:A:346:PRO:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:380:ARG:HG2	1:D:380:ARG:HH11	1.80	0.45
1:A:133:LEU:HB3	1:A:198:PHE:CE2	2.49	0.45
1:D:454:TYR:N	1:D:455:PRO:CD	2.78	0.45
1:B:410:LYS:HA	1:B:410:LYS:HE3	1.98	0.45
1:B:240:PHE:HA	1:B:244:GLY:HA3	1.99	0.45
1:D:306:LYS:HB2	1:D:307:PRO:HD3	1.98	0.45
1:B:103:ILE:HD11	1:B:108:TRP:CZ2	2.52	0.45
1:C:126:ASN:OD1	1:C:128:GLU:HG2	2.17	0.45
1:D:118:PHE:HB2	1:D:126:ASN:HB3	1.97	0.45
1:C:55:MET:O	1:C:63:PHE:HA	2.16	0.45
1:A:11:TYR:HA	1:A:391:ALA:O	2.17	0.44
1:B:375:GLU:OE1	1:B:410:LYS:HE2	2.17	0.44
1:C:434:GLY:HA2	1:C:488:LYS:NZ	2.32	0.44
1:D:22:ILE:N	1:D:23:PRO:HD2	2.31	0.44
1:B:22:ILE:HB	1:B:23:PRO:HD3	1.99	0.44
1:A:503:ASN:HA	1:A:504:PRO:HA	1.81	0.44
1:D:485:ILE:HG13	1:D:516:TRP:CH2	2.53	0.44
1:C:482:LYS:HE2	1:C:499:GLN:HA	1.98	0.44
1:D:232:LEU:HB3	1:D:233:PRO:HD2	1.98	0.44
1:B:51:TRP:CZ2	1:B:91:CYS:SG	3.11	0.44
1:D:383:GLU:O	1:D:387:ASN:HB2	2.17	0.44
1:A:329:PHE:O	1:A:366:GLY:HA2	2.18	0.44
1:C:198:PHE:O	1:C:202:LYS:HG2	2.17	0.44
1:B:292:HIS:O	1:B:344:SER:HA	2.18	0.44
1:C:17:ALA:HB2	1:C:46:ILE:HD11	1.99	0.43
1:B:107:VAL:HG11	1:B:130:LEU:HD21	1.99	0.43
1:B:2:VAL:O	1:B:5:LYS:HG2	2.18	0.43
1:A:51:TRP:CZ2	1:A:91:CYS:SG	3.12	0.43
1:D:269:GLY:O	1:D:273:HIS:CD2	2.70	0.43
1:A:256:TRP:O	1:A:260:ILE:HG12	2.18	0.43
1:C:453:GLN:C	1:C:455:PRO:HD2	2.39	0.43
1:B:154:LYS:HB2	1:B:155:PRO:HD3	2.00	0.43
1:C:134:ALA:O	1:C:138:ILE:HG12	2.18	0.43
1:D:55:MET:O	1:D:63:PHE:HA	2.17	0.43
1:B:104:PRO:HB3	1:B:106:TRP:NE1	2.33	0.43
1:A:103:ILE:HD11	1:A:108:TRP:CZ2	2.54	0.43
1:C:306:LYS:HB2	1:C:307:PRO:HD3	2.01	0.43
1:D:132:PRO:HB2	1:D:260:ILE:HG22	2.01	0.43
1:A:421:MET:O	1:A:509:THR:HG23	2.19	0.42
1:C:108:TRP:CZ2	1:C:127:LYS:HB3	2.54	0.42
1:D:296:ASN:OD1	1:D:343:TYR:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:347:LYS:O	1:C:351:GLN:HG3	2.18	0.42
1:D:226:ILE:HG23	1:D:227:SER:N	2.35	0.42
1:B:503:ASN:HA	1:B:504:PRO:HA	1.81	0.42
1:A:104:PRO:HB2	1:A:107:VAL:HG23	2.01	0.42
1:C:345:MET:N	1:C:346:PRO:HD3	2.34	0.42
1:C:461:ASP:HB3	1:C:466:ASP:OD1	2.20	0.42
1:A:119:LYS:HA	1:A:124:THR:O	2.19	0.42
1:A:484:PHE:HA	1:A:516:TRP:HH2	1.85	0.42
1:C:53:GLY:O	1:C:57:LYS:HB2	2.19	0.42
1:C:34:ASP:HB3	1:C:402:MET:HE2	2.02	0.42
1:D:213:SER:O	1:D:217:VAL:HG23	2.20	0.41
1:B:330:THR:O	1:B:331:CYS:HB2	2.20	0.41
1:B:47:THR:HG22	1:B:83:ILE:HB	2.02	0.41
1:D:240:PHE:HA	1:D:244:GLY:HA3	2.03	0.41
1:D:16:MET:HE3	1:D:396:LEU:HB2	2.01	0.41
1:D:443:ARG:HB3	1:D:445:GLU:OE1	2.20	0.41
1:D:47:THR:HG22	1:D:83:ILE:HB	2.02	0.41
1:B:385:ALA:HA	1:B:390:PHE:CD1	2.55	0.41
1:C:287:LYS:HG2	1:C:328:THR:HB	2.03	0.41
1:C:482:LYS:CG	1:C:496:GLN:HB3	2.45	0.41
1:D:238:GLU:O	1:D:242:MET:HG3	2.20	0.41
1:B:365:ASN:HB3	4:B:710:HOH:O	2.20	0.41
1:C:103:ILE:HD11	1:C:108:TRP:CZ2	2.55	0.41
1:C:432:THR:O	1:C:435:ASP:HB2	2.21	0.41
1:A:63:PHE:HB3	1:A:65:PHE:CE1	2.56	0.41
1:B:375:GLU:HG2	1:B:379:LYS:HZ2	1.85	0.41
1:A:19:LEU:HD13	1:A:397:ARG:CZ	2.50	0.41
1:B:429:VAL:HA	1:B:430:PRO:HD3	1.76	0.41
1:B:507:LEU:N	1:B:507:LEU:HD12	2.35	0.41
1:D:337:LYS:HG2	1:D:339:SER:H	1.86	0.41
1:A:5:LYS:HZ1	1:A:389:ASN:HB2	1.85	0.40
1:D:116:LEU:HD21	1:D:137:VAL:HG21	2.03	0.40
1:A:119:LYS:HE3	1:A:119:LYS:HB2	1.93	0.40
1:A:354:ALA:HA	1:A:364:LEU:HD22	2.03	0.40
1:C:240:PHE:HA	1:C:244:GLY:HA3	2.03	0.40
1:C:503:ASN:HA	1:C:504:PRO:HA	1.74	0.40
1:B:427:LYS:HG2	1:B:466:ASP:CB	2.49	0.40
1:B:478:ASN:HB3	1:B:503:ASN:HD22	1.87	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	514/516 (100%)	487 (95%)	25 (5%)	2 (0%)	34	32
1	B	514/516 (100%)	483 (94%)	29 (6%)	2 (0%)	34	32
1	C	514/516 (100%)	487 (95%)	26 (5%)	1 (0%)	47	49
1	D	514/516 (100%)	483 (94%)	29 (6%)	2 (0%)	34	32
All	All	2056/2064 (100%)	1940 (94%)	109 (5%)	7 (0%)	41	41

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	398	TYR
1	A	398	TYR
1	D	398	TYR
1	B	398	TYR
1	A	26	THR
1	D	26	THR
1	B	454	TYR

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	440/440 (100%)	427 (97%)	13 (3%)	41	44
1	B	440/440 (100%)	428 (97%)	12 (3%)	44	48
1	C	440/440 (100%)	430 (98%)	10 (2%)	50	55

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	440/440 (100%)	431 (98%)	9 (2%)	55	60
All	All	1760/1760 (100%)	1716 (98%)	44 (2%)	47	52

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	19	LEU
1	A	40	GLN
1	A	46	ILE
1	A	251	LYS
1	A	349	LEU
1	A	352	ASN
1	A	399	GLN
1	A	431	THR
1	A	473	LEU
1	A	502	TRP
1	A	504	PRO
1	A	508	LYS
1	B	19	LEU
1	B	181	SER
1	B	239	GLN
1	B	349	LEU
1	B	377	GLU
1	B	399	GLN
1	B	429	VAL
1	B	436	THR
1	B	488	LYS
1	B	498	ILE
1	B	502	TRP
1	B	504	PRO
1	C	19	LEU
1	C	187	PRO
1	C	239	GLN
1	C	251	LYS
1	C	349	LEU
1	C	399	GLN
1	C	432	THR
1	C	433	ILE
1	C	478	ASN
1	C	502	TRP
1	D	19	LEU

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Mol	Chain	Res	Type
1	D	187	PRO
1	D	214	LEU
1	D	335	THR
1	D	349	LEU
1	D	463	HIS
1	D	473	LEU
1	D	502	TRP
1	D	510	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	94	ASN
1	A	351	GLN
1	A	453	GLN
1	A	503	ASN
1	B	110	GLN
1	B	239	GLN
1	B	351	GLN
1	B	352	ASN
1	B	465	ASN
1	B	496	GLN
1	B	500	GLN
1	C	40	GLN
1	C	453	GLN
1	D	273	HIS
1	D	302	HIS
1	D	311	ASN
1	D	351	GLN
1	D	352	ASN
1	D	358	ASN
1	D	503	ASN

5.3.3 RNA ⓘ

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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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$
2	EBG	D	601	1	18,18,18	5.70	2 (11%)	23,25,25	4.96	4 (17%)
2	EBG	C	601	1	18,18,18	5.73	2 (11%)	23,25,25	5.04	5 (21%)
2	EBG	B	601	1	18,18,18	5.70	2 (11%)	23,25,25	5.00	5 (21%)
2	EBG	A	601	1	18,18,18	5.86	2 (11%)	23,25,25	5.09	6 (26%)

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	EBG	D	601	1	1/1/6/6	3/8/30/30	0/2/2/2
2	EBG	C	601	1	1/1/6/6	4/8/30/30	0/2/2/2
2	EBG	B	601	1	1/1/6/6	5/8/30/30	0/2/2/2
2	EBG	A	601	1	-	5/8/30/30	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	EBG	O9-C10	24.57	2.44	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	EBG	O9-C10	24.05	2.42	1.43
2	D	601	EBG	O9-C10	23.95	2.42	1.43
2	B	601	EBG	O9-C10	23.93	2.42	1.43
2	A	601	EBG	C10-C9	2.83	1.54	1.44
2	B	601	EBG	C10-C9	2.68	1.53	1.44
2	C	601	EBG	C10-C9	2.67	1.53	1.44
2	D	601	EBG	C10-C9	2.66	1.53	1.44

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	EBG	O9-C9-C10	19.63	110.32	60.04
2	C	601	EBG	O9-C9-C10	19.31	109.50	60.04
2	B	601	EBG	O9-C9-C10	19.23	109.29	60.04
2	D	601	EBG	O9-C9-C10	19.19	109.20	60.04
2	A	601	EBG	O9-C10-C9	-10.60	33.53	59.62
2	C	601	EBG	O9-C10-C9	-10.44	33.92	59.62
2	B	601	EBG	O9-C10-C9	-10.42	33.98	59.62
2	D	601	EBG	O9-C10-C9	-10.38	34.08	59.62
2	A	601	EBG	C10-O9-C9	-8.84	36.15	60.34
2	C	601	EBG	C10-O9-C9	-8.68	36.59	60.34
2	D	601	EBG	C10-O9-C9	-8.63	36.72	60.34
2	B	601	EBG	C10-O9-C9	-8.63	36.73	60.34
2	C	601	EBG	C1-C2-C3	-2.96	103.83	110.00
2	B	601	EBG	C1-C2-C3	-2.95	103.84	110.00
2	C	601	EBG	O1-C1-C2	2.60	112.36	108.30
2	A	601	EBG	O1-C1-C2	2.44	112.12	108.30
2	A	601	EBG	C7-O1-C1	2.18	117.45	113.84
2	D	601	EBG	C1-C2-C3	-2.14	105.53	110.00
2	A	601	EBG	O5-C1-O1	2.11	114.98	109.97
2	B	601	EBG	O1-C1-C2	2.04	111.48	108.30

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	601	EBG	C1
2	D	601	EBG	C1
2	B	601	EBG	C1

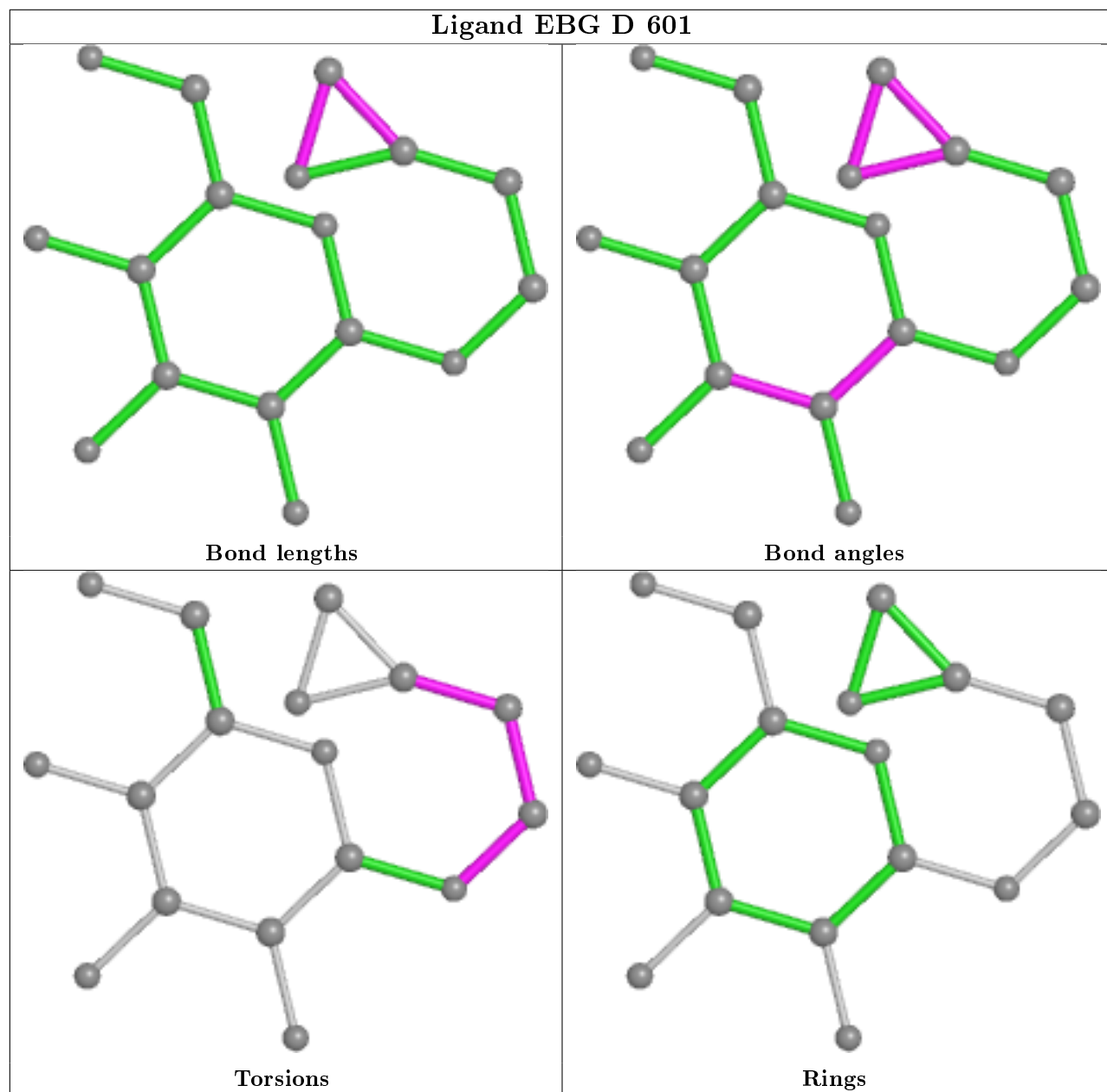
All (17) torsion outliers are listed below:

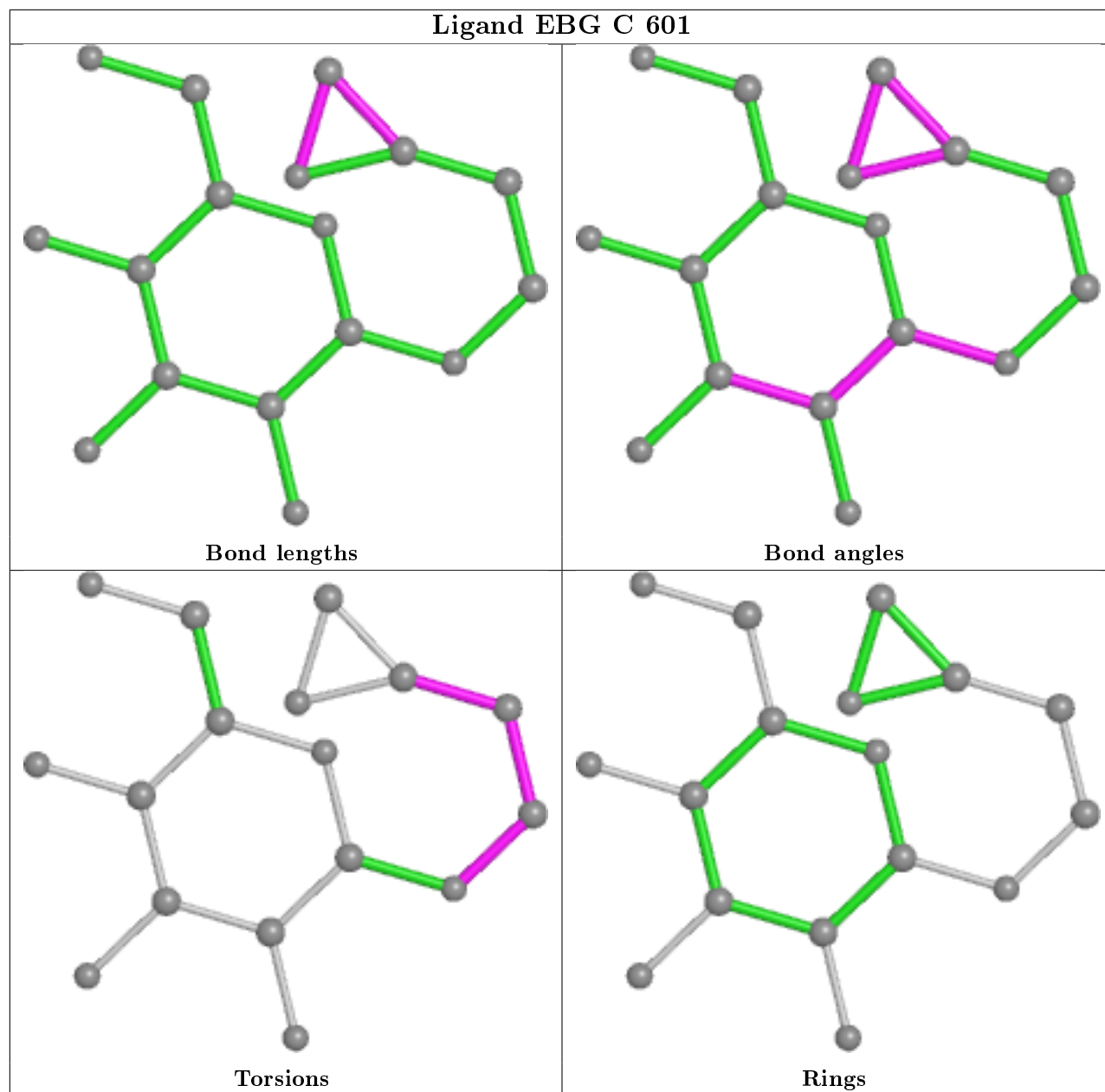
Mol	Chain	Res	Type	Atoms
2	C	601	EBG	O1-C7-C8-C9
2	D	601	EBG	O1-C7-C8-C9
2	B	601	EBG	C2-C1-O1-C7
2	B	601	EBG	C8-C7-O1-C1
2	A	601	EBG	C8-C7-O1-C1
2	A	601	EBG	C7-C8-C9-C10
2	B	601	EBG	O5-C1-O1-C7
2	A	601	EBG	C2-C1-O1-C7
2	D	601	EBG	C7-C8-C9-O9
2	C	601	EBG	C8-C7-O1-C1
2	D	601	EBG	C8-C7-O1-C1
2	B	601	EBG	O1-C7-C8-C9
2	C	601	EBG	C7-C8-C9-O9
2	A	601	EBG	O5-C1-O1-C7
2	A	601	EBG	C7-C8-C9-O9
2	C	601	EBG	C7-C8-C9-C10
2	B	601	EBG	C7-C8-C9-C10

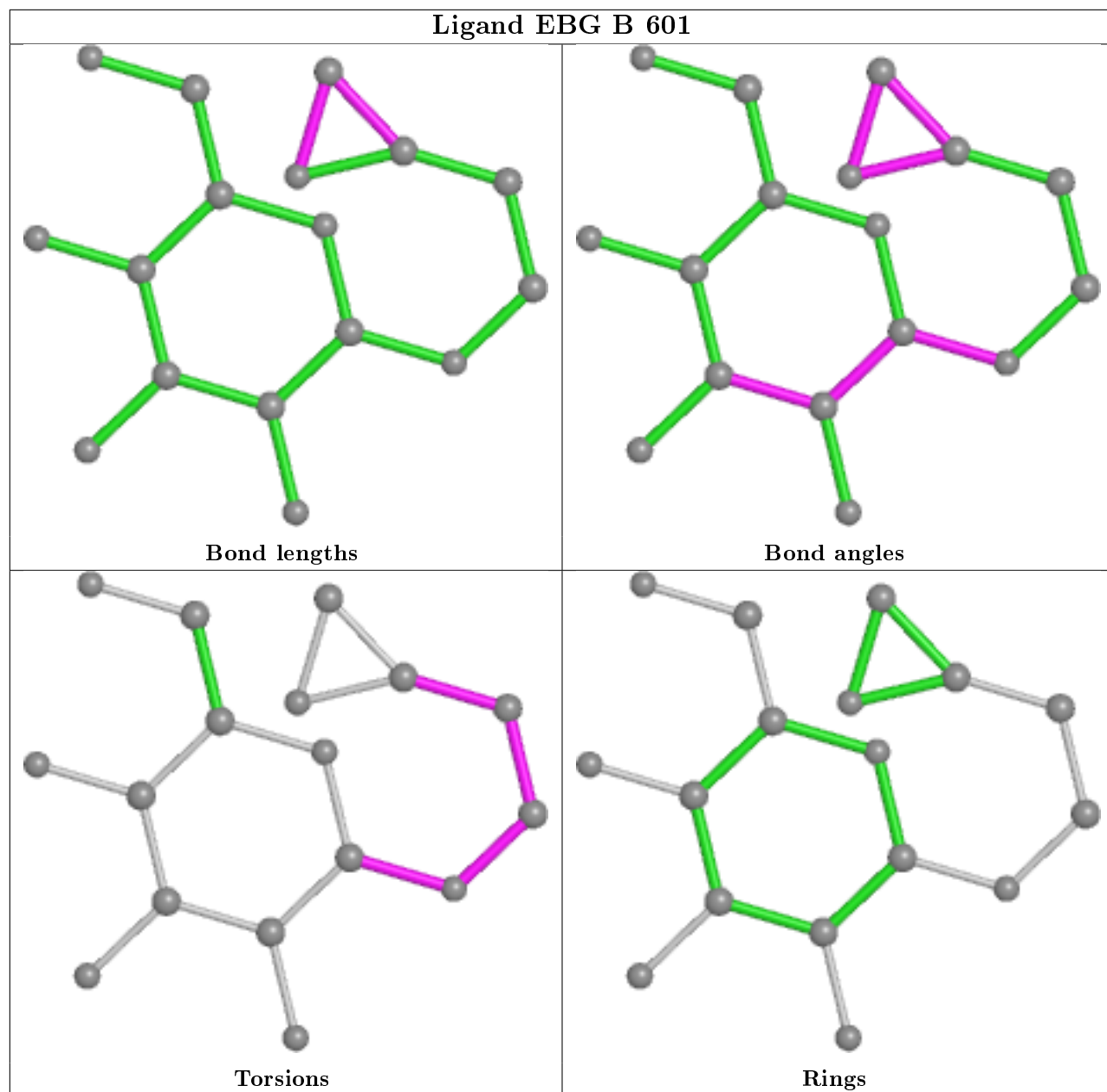
There are no ring outliers.

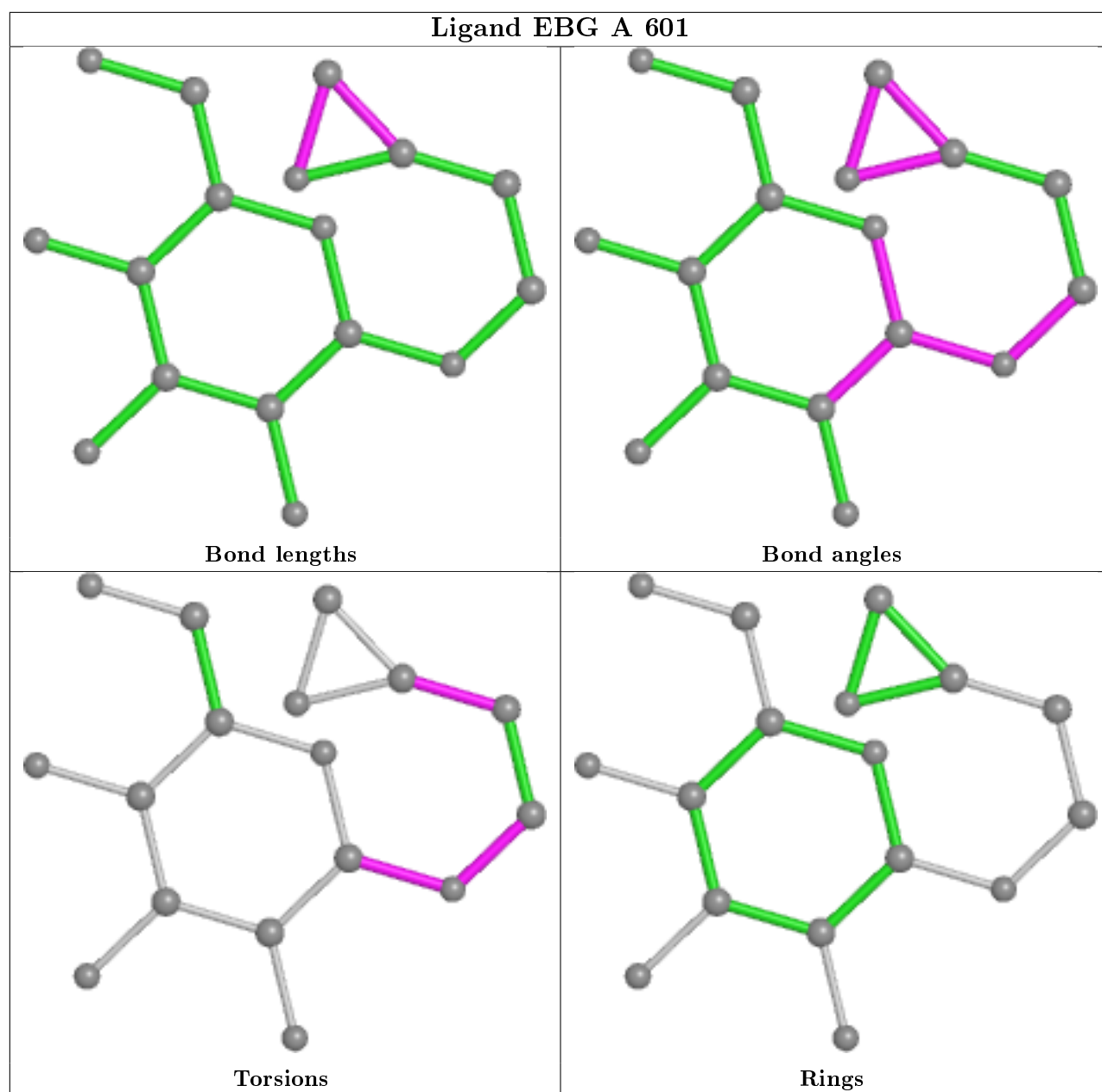
No monomer is involved in short contacts.

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









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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	516/516 (100%)	-0.70	0 100 100	7, 19, 40, 62	0
1	B	516/516 (100%)	-0.54	5 (0%) 82 85	6, 20, 43, 62	0
1	C	516/516 (100%)	-0.58	0 100 100	5, 19, 42, 63	0
1	D	516/516 (100%)	-0.57	0 100 100	6, 20, 41, 61	0
All	All	2064/2064 (100%)	-0.60	5 (0%) 95 95	5, 20, 42, 63	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	465	ASN	4.0
1	B	463	HIS	2.8
1	B	460	TYR	2.7
1	B	516	TRP	2.7
1	B	431	THR	2.3

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.

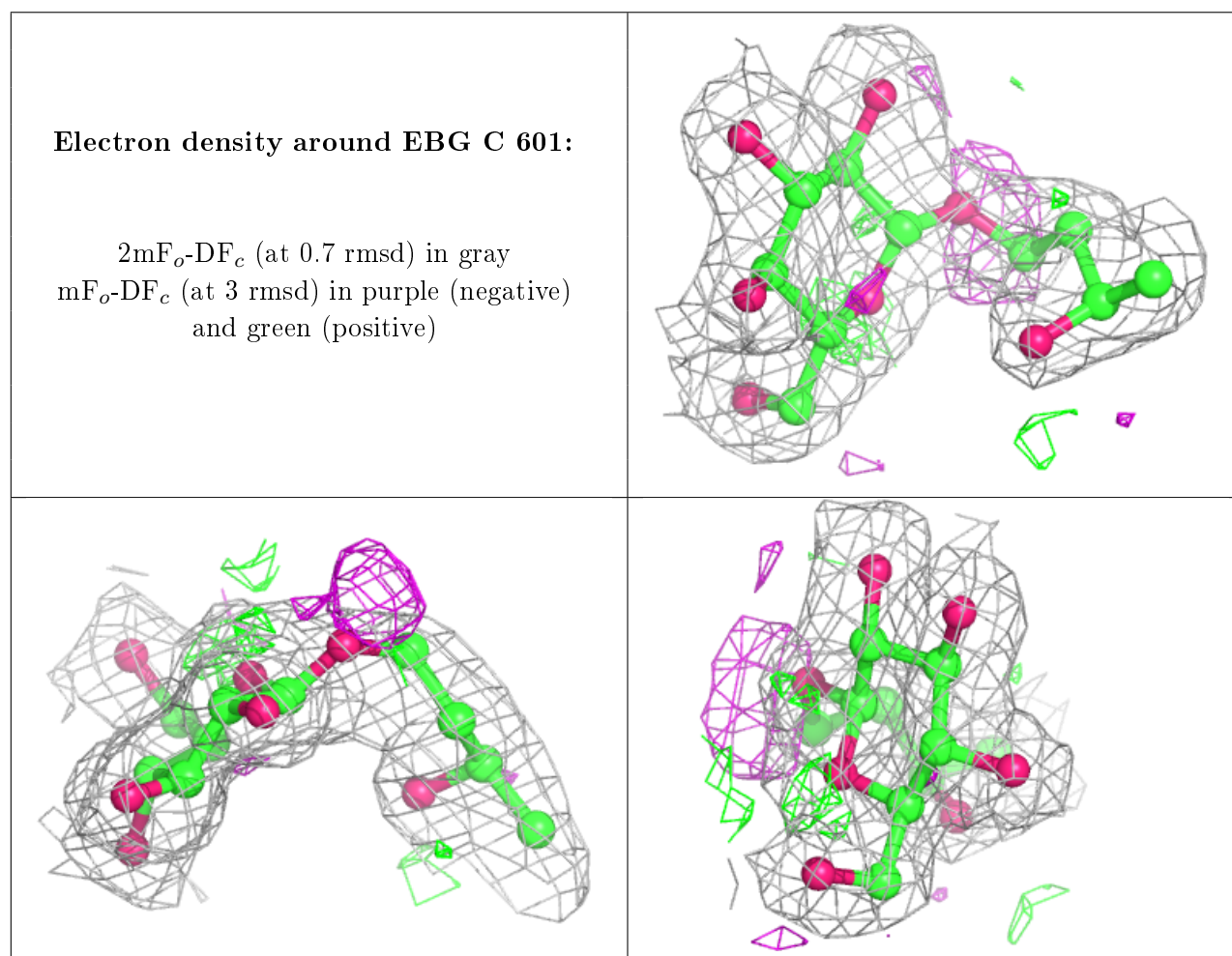
6.4 Ligands [i](#)

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.

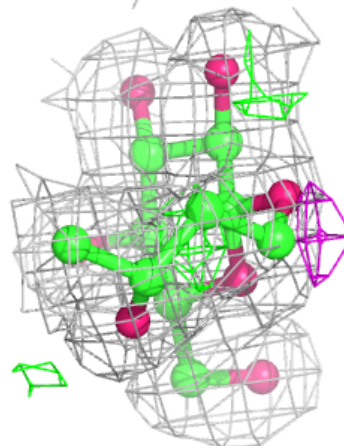
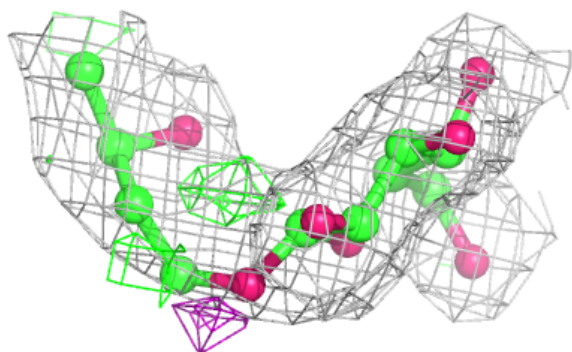
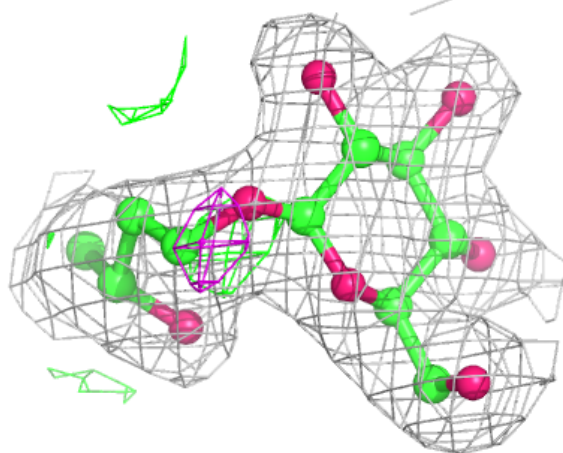
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	EBG	C	601	17/17	0.93	0.11	7,14,29,32	0
2	EBG	D	601	17/17	0.95	0.10	12,20,42,43	0
2	EBG	B	601	17/17	0.95	0.14	5,15,33,52	0
2	EBG	A	601	17/17	0.95	0.09	10,18,37,41	0
3	CA	D	701	1/1	0.99	0.09	14,14,14,14	0
3	CA	B	701	1/1	0.99	0.10	11,11,11,11	0
3	CA	C	701	1/1	1.00	0.15	14,14,14,14	0
3	CA	A	701	1/1	1.00	0.07	10,10,10,10	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.



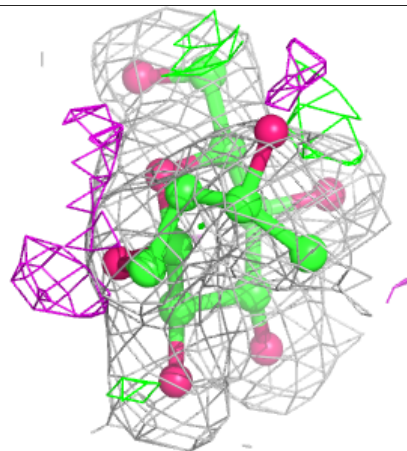
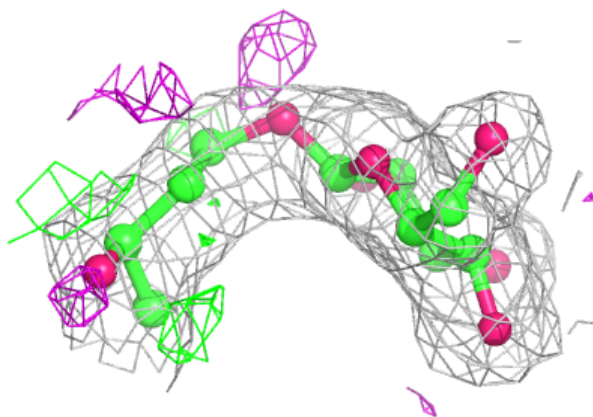
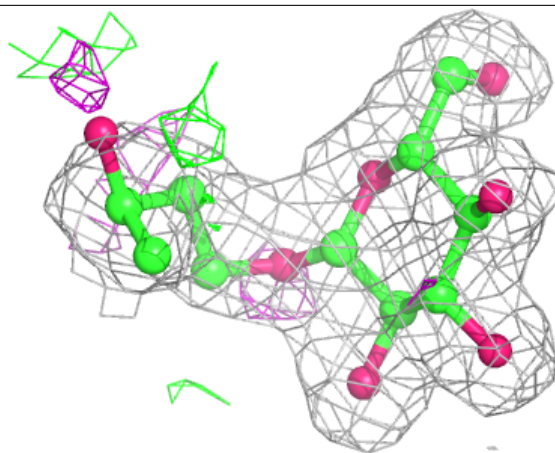
Electron density around EBG D 601:

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



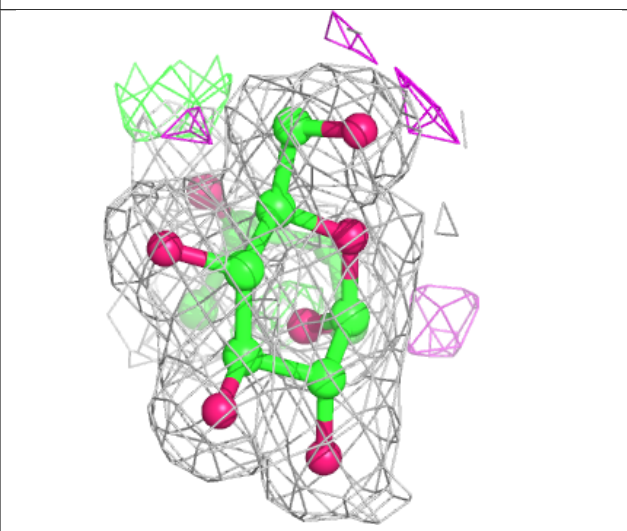
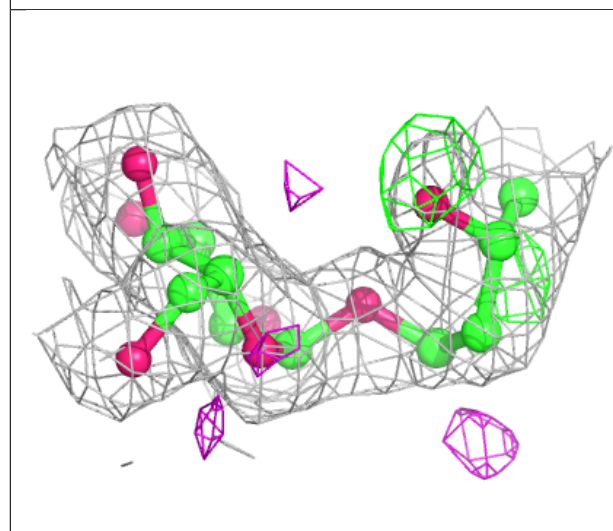
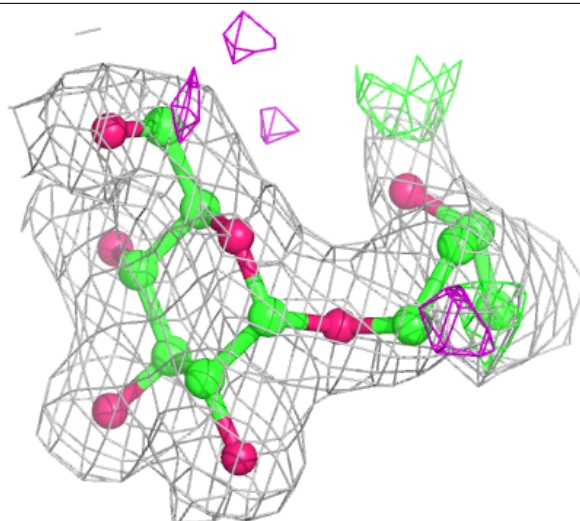
Electron density around EBG B 601:

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



Electron density around EBG A 601:

$2mF_o-DF_c$ (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.