Full wwPDB X-ray Structure Validation Report (i)

Sep 27, 2023 – 01:36 AM EDT

PDB ID	:	6D50
Title	:	Bacteroides uniforms beta-glucuronidase 2 bound to D-glucaro-1,5-lactone
Authors	:	Walton, W.G.; Pellock, S.J.; Redinbo, M.R.
Deposited on	:	2018-04-19
Resolution	:	2.50 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 (i) 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 (i)) were used in the production of this report:

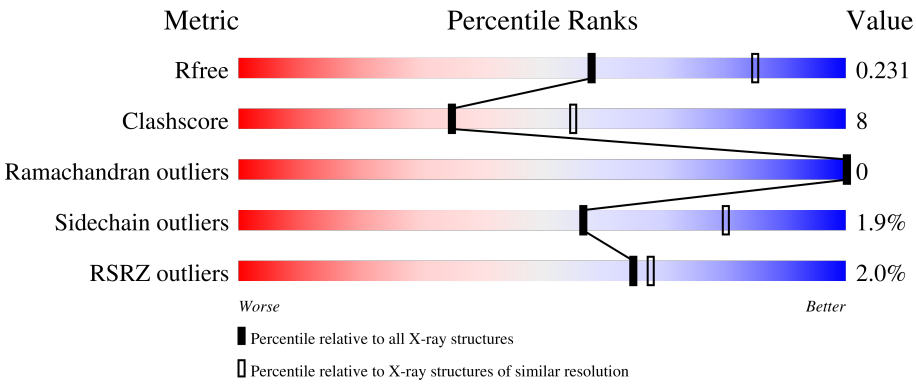
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 <=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	886	3%, 74%, 20%, • 6%
1	B	886	%, 78%, 16%, • 5%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 14207 atoms, of which 14 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 Glycosyl hydrolases family 2, sugar binding domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	837	6802	4329	1168	1284	21	0	4	0
1	B	840	6830	4345	1176	1288	21	0	4	0

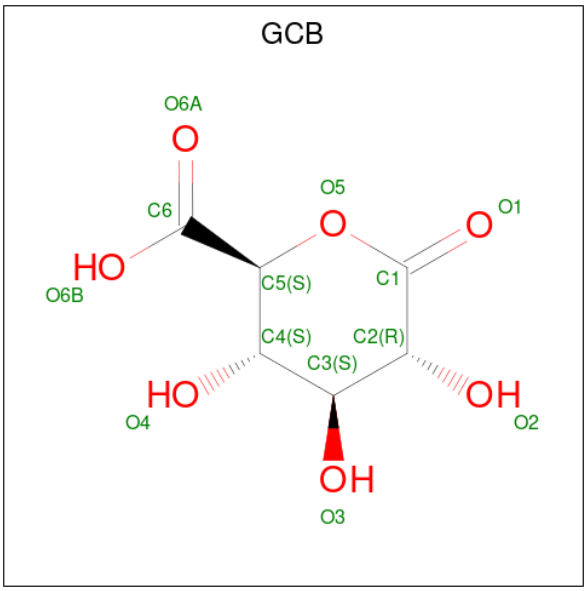
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
2	A	1	1	1	0	0
2	B	1	1	1	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
3	A	1	1	1	0	0
3	B	1	1	1	0	0

- Molecule 4 is (2S,3S,4S,5R)-3,4,5-trihydroxy-6-oxo-oxane-2-carboxylic acid (three-letter code: GCB) (formula: $C_6H_8O_7$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
4	A	1	20	6	7	7	0	0
			Total	C	H	O		
4	B	1	20	6	7	7	0	0

- Molecule 5 is water.

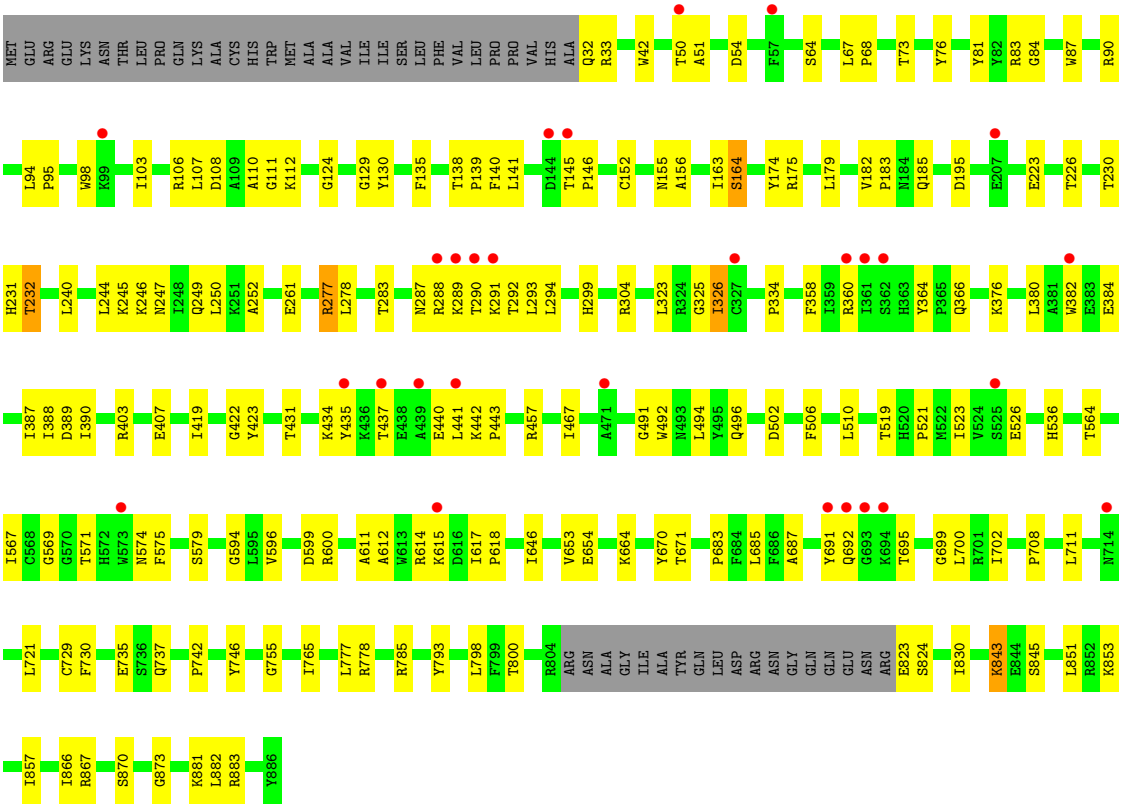
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	225	225	225	0	0
			Total	O		
5	B	306	306	306	0	0

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.

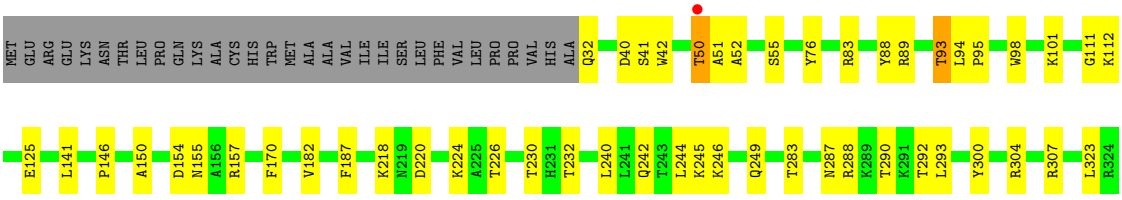
• Molecule 1: Glycosyl hydrolases family 2, sugar binding domain protein

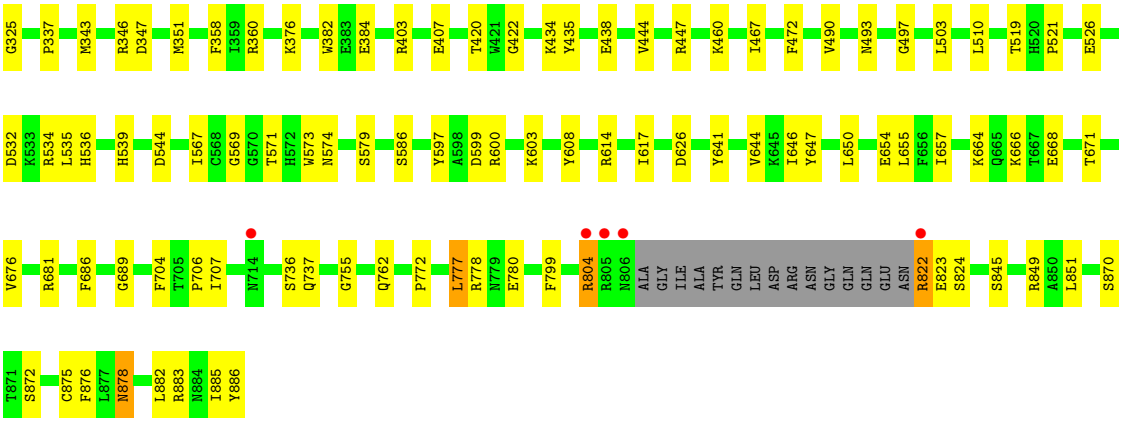
Chain A:



• Molecule 1: Glycosyl hydrolases family 2, sugar binding domain protein

Chain B:




G325 P337 M343 R346 D347 M351 F358 I359 R360 K376 W382 E383 E384 R403 E407 T420 W421 G422 K434 Y435 E438 V444 R447 K460 I467 F472 V490 N493 G497 L503 L510 T519 H520 P521 E526
D532 K533 R534 L535 H536 H539 D544 I567 C568 G569 G570 T571 H572 W573 N574 S579 S586 Y597 A598 D599 R600 K603 Y608 R614 I617 D626 Y641 V644 K645 I646 Y647 L650 E654 L655 F656 I657 K664 Q665 K666 T667 E668 T671
V676 R681 F686 G689 F704 T705 P706 I707 N714 S736 Q737 G755 Q762 P772 L777 R778 N779 E780 F799 R804 R805 N806 ALA GLY ILE ALA TYR GLN LEU ASP ARG ASN GLY GLN GLN GLU ASN R822 E823 S824 S845 R849 A850 L851 S870
T871 S872 C875 F876 L877 N878 L882 R883 N884 I885 Y886

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	74.49Å 141.88Å 180.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.48 – 2.50 29.48 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.48-2.50) 100.0 (29.48-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.92 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.166 , 0.231 0.166 , 0.231	Depositor DCC
R_{free} test set	2000 reflections (2.98%)	wwPDB-VP
Wilson B-factor (Å^2)	40.0	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 44.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	14207	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GCB, CA, NA

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	0.44	0/6985	0.60	0/9483
1	B	0.46	0/7013	0.60	0/9520
All	All	0.45	0/13998	0.60	0/19003

There are no bond length outliers.

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	6802	0	6566	120	0
1	B	6830	0	6594	106	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	13	7	0	1	0
4	B	13	7	0	0	0
5	A	225	0	0	1	0
5	B	306	0	0	4	0
All	All	14193	14	13160	219	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 8.

All (219) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:830:ILE:HG12	1:A:866:ILE:HD13	1.44	0.99
1:B:290:THR:HG23	1:B:292:THR:H	1.41	0.85
1:B:112:LYS:HG3	1:B:155:ASN:HA	1.58	0.84
1:A:793:TYR:HD2	1:A:882:LEU:HD11	1.44	0.82
1:A:830:ILE:HG12	1:A:866:ILE:CD1	2.10	0.80
1:B:804:ARG:HH11	1:B:804:ARG:HG2	1.46	0.80
1:A:290:THR:OG1	1:A:292:THR:HG23	1.83	0.79
1:B:287:ASN:OD1	1:B:290:THR:HG22	1.82	0.79
1:B:510:LEU:HD13	1:B:567:ILE:HD11	1.69	0.75
1:A:442:LYS:HB3	1:A:443:PRO:HD3	1.69	0.74
1:B:885:ILE:HG22	1:B:886:TYR:CD2	2.24	0.73
1:A:94:LEU:HD11	1:A:141:LEU:HD13	1.71	0.72
1:B:539:HIS:ND1	5:B:1001:HOH:O	2.23	0.71
1:A:845:SER:HB2	1:A:851[A]:LEU:HD11	1.73	0.70
1:A:882:LEU:HD12	1:A:883:ARG:N	2.07	0.70
1:B:287:ASN:HB3	1:B:290:THR:CG2	2.22	0.69
1:B:83:ARG:HG2	1:B:155:ASN:O	1.94	0.67
1:B:307:ARG:HD2	5:B:1004:HOH:O	1.93	0.67
1:A:793:TYR:CD2	1:A:882:LEU:HD11	2.30	0.66
1:B:287:ASN:CG	1:B:290:THR:HG22	2.16	0.66
1:B:287:ASN:HB3	1:B:290:THR:HG22	1.81	0.63
1:A:435:TYR:HB2	1:A:441:LEU:HD23	1.78	0.63
1:B:510:LEU:HD13	1:B:567:ILE:CD1	2.29	0.63
1:B:883:ARG:NH2	1:B:885:ILE:HD11	2.14	0.63
1:A:435:TYR:HB2	1:A:441:LEU:CD2	2.30	0.62
1:A:32:GLN:HA	1:A:294:LEU:O	2.00	0.61
1:A:845:SER:HG	1:B:41:SER:HG	1.45	0.61
1:B:288:ARG:HG2	1:B:288:ARG:HH11	1.67	0.60
1:A:32:GLN:OE1	1:A:293:LEU:HD22	2.01	0.60
1:B:822:ARG:HH12	1:B:823:GLU:HB2	1.66	0.60
1:A:244:LEU:HD21	1:A:261:GLU:O	2.02	0.59
1:B:804:ARG:HG3	1:B:804:ARG:O	2.01	0.59
1:B:614:ARG:NH2	1:B:617:ILE:HD13	2.17	0.59
1:B:706:PRO:O	1:B:707:ILE:HD13	2.03	0.59
1:A:519:THR:O	1:A:521:PRO:HD3	2.03	0.58
1:B:325:GLY:HA2	1:B:358:PHE:O	2.03	0.58
1:A:112:LYS:HG3	1:A:155:ASN:HA	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:GLN:HG3	1:B:293:LEU:HD23	1.85	0.57
1:A:81:TYR:CE1	1:A:83[B]:ARG:HD3	2.39	0.57
1:A:440:GLU:O	1:A:443:PRO:HD2	2.04	0.57
1:A:851[B]:LEU:HD13	1:B:40:ASP:CB	2.35	0.57
1:B:444:VAL:HG22	1:B:447:ARG:HH22	1.70	0.57
1:B:519:THR:O	1:B:521:PRO:HD3	2.04	0.57
1:B:650:LEU:HD23	1:B:689:GLY:HA3	1.86	0.56
1:A:289:LYS:O	1:A:291:LYS:NZ	2.35	0.56
1:A:614:ARG:NH2	1:A:617:ILE:HD13	2.21	0.56
1:B:287:ASN:CB	1:B:290:THR:HG22	2.34	0.56
1:B:384:GLU:HA	1:B:422:GLY:O	2.06	0.56
1:A:564:THR:HB	1:A:567:ILE:HD12	1.87	0.56
1:B:822:ARG:NH1	1:B:823:GLU:HB2	2.21	0.56
1:A:325:GLY:HA2	1:A:358:PHE:O	2.06	0.56
1:A:325:GLY:HA3	1:A:571:THR:HG22	1.86	0.56
1:B:804:ARG:HG2	1:B:804:ARG:NH1	2.20	0.56
1:B:323:LEU:HB2	1:B:569:GLY:HA3	1.87	0.55
1:B:154:ASP:OD2	1:B:157:ARG:HG3	2.07	0.55
1:A:87:TRP:CH2	1:A:152:CYS:HB2	2.41	0.55
1:B:76:TYR:CE1	1:B:579:SER:HB2	2.42	0.55
1:B:822:ARG:HG3	1:B:872:SER:HB2	1.88	0.55
1:B:403:ARG:O	1:B:407:GLU:HG3	2.07	0.55
1:B:755:GLY:O	1:B:778:ARG:HD3	2.08	0.55
1:A:245:LYS:O	1:A:246:LYS:HD2	2.06	0.54
1:B:822:ARG:CG	1:B:872:SER:HB2	2.38	0.54
1:B:777:LEU:O	1:B:777:LEU:HD12	2.08	0.54
1:A:437:THR:HG23	1:A:440:GLU:OE1	2.06	0.54
1:A:124:GLY:HA3	1:A:135:PHE:CE2	2.42	0.53
1:A:185:GLN:HB2	1:A:294:LEU:HD13	1.91	0.53
1:B:706:PRO:C	1:B:707:ILE:HD13	2.29	0.52
1:B:288:ARG:HG2	1:B:288:ARG:NH1	2.24	0.52
1:A:223:GLU:HA	1:A:252:ALA:HB2	1.92	0.52
1:B:101:LYS:HB3	1:B:182:VAL:O	2.09	0.52
1:B:497:GLY:HA3	1:B:503:LEU:HD23	1.92	0.52
1:A:326:ILE:HD13	1:A:326:ILE:H	1.75	0.52
1:B:347:ASP:O	1:B:351:MET:HG3	2.09	0.51
1:A:73[A]:THR:HG22	1:A:334:PRO:O	2.09	0.51
1:B:599:ASP:O	1:B:600:ARG:HB2	2.10	0.51
1:A:708:PRO:HG2	1:A:711:LEU:HD23	1.93	0.51
1:A:387:ILE:HB	1:A:423:TYR:O	2.10	0.51
1:A:360:ARG:NH1	1:A:526:GLU:HG3	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:493:ASN:OD1	1:B:526:GLU:HG3	2.11	0.51
1:A:138:THR:HB	1:A:139:PRO:HD3	1.92	0.50
1:A:232:THR:HG22	1:A:240:LEU:HD23	1.94	0.50
1:A:110:ALA:O	1:A:129:GLY:HA2	2.12	0.50
1:A:390:ILE:HG13	1:A:431:THR:OG1	2.12	0.49
1:A:737:GLN:OE1	1:B:346:ARG:NH1	2.43	0.49
1:B:42:TRP:CE3	1:B:88:TYR:HB3	2.47	0.49
1:A:492:TRP:HB3	1:A:494:LEU:HD21	1.93	0.49
1:A:506:PHE:CE1	1:A:510:LEU:HD11	2.48	0.49
1:A:721:LEU:O	1:A:881:LYS:HA	2.12	0.49
1:A:50:THR:HG23	1:A:50:THR:O	2.12	0.49
1:A:526:GLU:OE1	4:A:903:GCB:C1	2.61	0.49
1:A:496:GLN:HB3	1:A:502:ASP:O	2.12	0.48
1:A:223:GLU:CA	1:A:252:ALA:HB2	2.44	0.48
1:A:457:ARG:HG3	1:A:457:ARG:HH11	1.78	0.48
1:B:52:ALA:O	1:B:89:ARG:NH1	2.45	0.48
1:A:730:PHE:CD2	1:A:742:PRO:HD3	2.48	0.48
1:B:762:GLN:OE1	1:B:762:GLN:HA	2.15	0.47
1:A:145:THR:HB	1:A:146:PRO:HD2	1.96	0.47
1:A:599:ASP:O	1:A:600:ARG:HB2	2.13	0.47
1:B:50:THR:O	1:B:51:ALA:HB3	2.14	0.47
1:A:384:GLU:HA	1:A:422:GLY:O	2.14	0.47
1:A:403:ARG:O	1:A:407:GLU:HG3	2.14	0.47
1:A:435:TYR:O	1:A:441:LEU:HD21	2.15	0.47
1:A:735:GLU:HG2	1:B:343:MET:HE2	1.96	0.47
1:A:111:GLY:HA2	1:A:112:LYS:HA	1.70	0.47
1:A:64:SER:HB3	1:B:849:ARG:CZ	2.45	0.46
1:A:230:THR:HG23	1:A:245:LYS:HG2	1.96	0.46
1:A:231:HIS:CE1	1:A:283:THR:HG23	2.51	0.46
1:A:323:LEU:HB2	1:A:569:GLY:HA3	1.96	0.46
1:A:654:GLU:O	1:A:687:ALA:HA	2.14	0.46
1:B:467:ILE:HD13	1:B:467:ILE:HA	1.71	0.46
1:A:76:TYR:CE1	1:A:579:SER:HB2	2.49	0.46
1:B:597:TYR:CZ	1:B:603:LYS:HG2	2.51	0.46
1:B:434:LYS:HE2	1:B:435:TYR:CZ	2.51	0.46
1:A:278:LEU:HD11	1:A:376:LYS:HA	1.98	0.46
1:A:700:LEU:N	1:A:700:LEU:HD23	2.30	0.46
1:A:851[B]:LEU:HD11	1:A:853:LYS:CG	2.46	0.46
1:A:575:PHE:O	1:A:596:VAL:HG22	2.16	0.46
1:A:287:ASN:O	1:A:291:LYS:HA	2.15	0.46
1:B:93:THR:HG23	1:B:146:PRO:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:614:ARG:CZ	1:B:617:ILE:HD13	2.46	0.46
1:B:736:SER:O	1:B:737:GLN:HB2	2.14	0.45
1:A:107:LEU:N	1:A:107:LEU:HD12	2.31	0.45
1:A:130:TYR:CG	1:A:364:TYR:HB2	2.52	0.45
1:A:617:ILE:HB	1:A:618:PRO:HD2	1.97	0.45
1:B:641:TYR:HA	1:B:676:VAL:O	2.17	0.45
1:A:777:LEU:C	1:A:777:LEU:HD12	2.37	0.45
1:B:245:LYS:O	1:B:246:LYS:HD3	2.16	0.45
1:B:681:ARG:HD3	1:B:704:PHE:O	2.17	0.45
1:A:574:ASN:O	1:A:594:GLY:HA2	2.17	0.45
1:A:360:ARG:HH12	1:A:526:GLU:HG3	1.82	0.45
1:A:746:TYR:CE2	1:A:785:ARG:HG2	2.52	0.45
1:A:457:ARG:HG3	1:A:457:ARG:NH1	2.32	0.45
1:A:106:ARG:NH1	1:A:108:ASP:OD1	2.50	0.45
1:A:175:ARG:NH2	1:A:366:GLN:HA	2.31	0.45
1:B:218:LYS:NZ	1:B:220:ASP:OD2	2.46	0.45
1:B:681:ARG:HD3	1:B:681:ARG:HA	1.79	0.45
1:A:182:VAL:HB	1:A:183:PRO:HD2	1.99	0.44
1:A:611:ALA:O	1:A:615:LYS:NZ	2.50	0.44
1:A:380:LEU:HD22	1:A:419:ILE:HD13	1.99	0.44
1:B:95:PRO:HD2	1:B:98:TRP:CE3	2.53	0.44
1:B:608:TYR:HB2	1:B:647:TYR:CE1	2.53	0.44
1:B:573:TRP:HA	1:B:574:ASN:HA	1.72	0.44
1:B:655:LEU:HD12	1:B:686:PHE:O	2.17	0.44
1:A:51:ALA:HB1	1:A:54:ASP:OD2	2.17	0.44
1:B:772:PRO:HD3	5:B:1239:HOH:O	2.17	0.44
1:B:882:LEU:C	1:B:882:LEU:HD23	2.38	0.44
1:B:824:SER:HA	1:B:872:SER:O	2.17	0.44
1:A:755:GLY:O	1:A:778:ARG:HD3	2.18	0.43
1:B:472:PHE:CE2	1:B:490:VAL:HB	2.53	0.43
1:B:777:LEU:HD22	1:B:876:PHE:HB2	1.99	0.43
1:A:95:PRO:HG2	1:A:98:TRP:CG	2.53	0.43
1:B:226:THR:HG22	1:B:288:ARG:CB	2.48	0.43
1:B:646:ILE:O	1:B:671:THR:HA	2.18	0.43
1:A:646:ILE:O	1:A:671:THR:HA	2.18	0.43
1:A:845:SER:CB	1:A:851[A]:LEU:HD11	2.45	0.43
1:A:851[B]:LEU:HD13	1:B:40:ASP:HB2	2.00	0.43
1:A:67:LEU:HB3	1:A:68:PRO:HA	2.00	0.43
1:A:467:ILE:HD13	1:A:467:ILE:HA	1.76	0.43
1:A:843:LYS:HD2	1:A:843:LYS:O	2.17	0.43
1:B:232:THR:HG23	1:B:240:LEU:HD13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:THR:HG22	1:A:288:ARG:CB	2.49	0.43
1:B:94:LEU:HD11	1:B:141:LEU:HD22	2.01	0.43
1:A:290:THR:CB	1:A:292:THR:HG23	2.49	0.43
1:A:735:GLU:HG2	1:B:343:MET:CE	2.47	0.43
1:B:111:GLY:HA2	1:B:112:LYS:HA	1.60	0.43
1:B:777:LEU:HD12	1:B:777:LEU:C	2.39	0.43
1:A:824:SER:HA	1:A:873:GLY:HA3	1.99	0.43
1:B:88:TYR:O	1:B:150:ALA:HA	2.19	0.43
1:A:249:GLN:C	1:A:250:LEU:HD12	2.40	0.42
1:A:84:GLY:HA2	1:A:156:ALA:HB2	2.01	0.42
1:B:242:GLN:HG2	1:B:244:LEU:HD12	2.00	0.42
1:B:804:ARG:NH1	1:B:804:ARG:CG	2.81	0.42
1:B:882:LEU:HD23	1:B:883:ARG:N	2.34	0.42
1:A:183:PRO:HG3	1:A:294:LEU:O	2.19	0.42
1:B:535:LEU:HD13	1:B:544:ASP:O	2.19	0.42
1:B:845:SER:HB2	1:B:851:LEU:HD11	2.00	0.42
1:A:42:TRP:CD2	1:A:90:ARG:HB2	2.54	0.42
1:A:231:HIS:ND1	1:A:283:THR:HG23	2.34	0.42
1:A:163:ILE:HD12	1:A:164:SER:HB3	2.01	0.42
1:A:683:PRO:HD2	1:A:702:ILE:O	2.19	0.42
1:B:360:ARG:HB2	1:B:571:THR:HG21	2.01	0.42
1:A:388:ILE:HG22	1:A:389:ASP:N	2.35	0.42
1:A:536:HIS:O	1:A:729:CYS:HA	2.20	0.42
1:A:765:ILE:HD12	1:A:800:THR:HG21	2.01	0.42
1:B:666:LYS:NZ	1:B:668:GLU:OE2	2.42	0.42
1:B:644:VAL:CG2	1:B:676:VAL:HG21	2.50	0.42
1:A:103:ILE:N	1:A:103:ILE:HD13	2.35	0.42
1:B:780:GLU:N	1:B:875:CYS:O	2.52	0.42
1:A:491:GLY:HA2	1:A:523:ILE:O	2.20	0.42
1:A:691:TYR:C	1:A:692:GLN:HG2	2.40	0.42
1:A:857:ILE:HD13	1:A:857:ILE:HA	1.87	0.41
1:A:277:ARG:HA	1:A:277:ARG:HD3	1.83	0.41
1:B:444:VAL:HG22	1:B:447:ARG:NH2	2.33	0.41
1:A:195:ASP:OD2	1:A:299:HIS:HE1	2.03	0.41
1:A:867:ARG:NH1	5:A:1002:HOH:O	2.24	0.41
1:B:224:LYS:NZ	1:B:249:GLN:HG2	2.36	0.41
1:B:300:TYR:OH	1:B:376:LYS:HE3	2.21	0.41
1:B:536:HIS:CE1	1:B:626:ASP:O	2.74	0.41
1:B:654:GLU:HB2	1:B:664:LYS:HG2	2.02	0.41
1:B:885:ILE:HD13	1:B:885:ILE:HA	1.82	0.41
1:A:232:THR:HG22	1:A:240:LEU:CD2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:ASN:OD1	1:A:288:ARG:NE	2.54	0.41
1:B:420:THR:HA	1:B:467:ILE:O	2.20	0.41
1:B:532:ASP:OD2	1:B:534:ARG:HB2	2.21	0.41
1:B:644:VAL:HA	5:B:1219:HOH:O	2.21	0.41
1:B:657:ILE:HG23	1:B:657:ILE:O	2.21	0.41
1:A:685:LEU:O	1:A:699:GLY:HA2	2.20	0.41
1:B:460:LYS:HA	1:B:460:LYS:HD3	1.85	0.40
1:B:799:PHE:HA	1:B:878:ASN:OD1	2.21	0.40
1:A:129:GLY:O	1:A:174[B]:TYR:OH	2.31	0.40
1:A:179:LEU:HD23	1:A:179:LEU:HA	1.94	0.40
1:A:434:LYS:HE3	1:A:435:TYR:CZ	2.56	0.40
1:A:615:LYS:HA	1:A:615:LYS:HD3	1.87	0.40
1:B:170:PHE:CD2	1:B:337:PRO:HD3	2.56	0.40
1:B:650:LEU:CD2	1:B:689:GLY:HA3	2.52	0.40
1:B:230:THR:HA	1:B:244:LEU:O	2.21	0.40
1:A:32:GLN:HB2	1:A:33:ARG:H	1.78	0.40
1:A:612:ALA:HA	1:A:670:TYR:CE1	2.57	0.40
1:B:187:PHE:CZ	1:B:283:THR:HG22	2.57	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	837/886 (94%)	806 (96%)	31 (4%)	0	100	100
1	B	840/886 (95%)	812 (97%)	28 (3%)	0	100	100
All	All	1677/1772 (95%)	1618 (96%)	59 (4%)	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	730/767 (95%)	716 (98%)	14 (2%)	57	80
1	B	733/767 (96%)	720 (98%)	13 (2%)	59	81
All	All	1463/1534 (95%)	1436 (98%)	27 (2%)	57	81

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	140	PHE
1	A	164	SER
1	A	232	THR
1	A	277	ARG
1	A	304	ARG
1	A	326	ILE
1	A	382	TRP
1	A	653	VAL
1	A	664	LYS
1	A	695	THR
1	A	798	LEU
1	A	823	GLU
1	A	843	LYS
1	A	870	SER
1	B	50	THR
1	B	55	SER
1	B	93	THR
1	B	125	GLU
1	B	304	ARG
1	B	382	TRP
1	B	438	GLU
1	B	586	SER
1	B	777	LEU
1	B	804	ARG
1	B	822	ARG
1	B	870	SER
1	B	878	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	299	HIS

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 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\#\|Z\| > 2$
4	GCB	A	903	-	13,13,13	3.65	7 (53%)	16,19,19	1.83	4 (25%)
4	GCB	B	903	-	13,13,13	3.69	8 (61%)	16,19,19	2.62	7 (43%)

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
4	GCB	A	903	-	-	0/4/24/24	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GCB	B	903	-	-	0/4/24/24	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	903	GCB	O5-C1	9.78	1.49	1.34
4	B	903	GCB	O5-C1	8.78	1.47	1.34
4	B	903	GCB	C3-C2	-6.60	1.43	1.53
4	A	903	GCB	C3-C2	-5.74	1.44	1.53
4	B	903	GCB	O5-C5	4.46	1.52	1.46
4	A	903	GCB	O5-C5	2.82	1.50	1.46
4	B	903	GCB	C4-C5	-2.63	1.48	1.53
4	B	903	GCB	C4-C3	-2.46	1.46	1.52
4	A	903	GCB	O4-C4	-2.45	1.37	1.43
4	A	903	GCB	C2-C1	2.42	1.57	1.52
4	A	903	GCB	C4-C3	-2.39	1.46	1.52
4	B	903	GCB	O2-C2	2.31	1.46	1.42
4	A	903	GCB	O2-C2	2.29	1.46	1.42
4	B	903	GCB	C2-C1	2.25	1.57	1.52
4	B	903	GCB	O4-C4	-2.21	1.37	1.43

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	903	GCB	O5-C1-C2	-5.78	110.57	119.20
4	A	903	GCB	O5-C1-C2	-4.82	112.00	119.20
4	B	903	GCB	O5-C5-C4	4.71	119.98	110.26
4	B	903	GCB	C3-C4-C5	-4.20	102.06	109.25
4	B	903	GCB	O4-C4-C5	-3.02	102.97	109.74
4	B	903	GCB	O5-C1-O1	2.97	122.81	118.47
4	A	903	GCB	O5-C5-C4	2.50	115.43	110.26
4	A	903	GCB	O5-C1-O1	2.50	122.12	118.47
4	B	903	GCB	O6B-C6-O6A	-2.19	119.11	124.09
4	B	903	GCB	O3-C3-C4	-2.09	105.53	110.35
4	A	903	GCB	O4-C4-C3	2.02	115.01	110.35

There are no chirality outliers.

There are no torsion outliers.

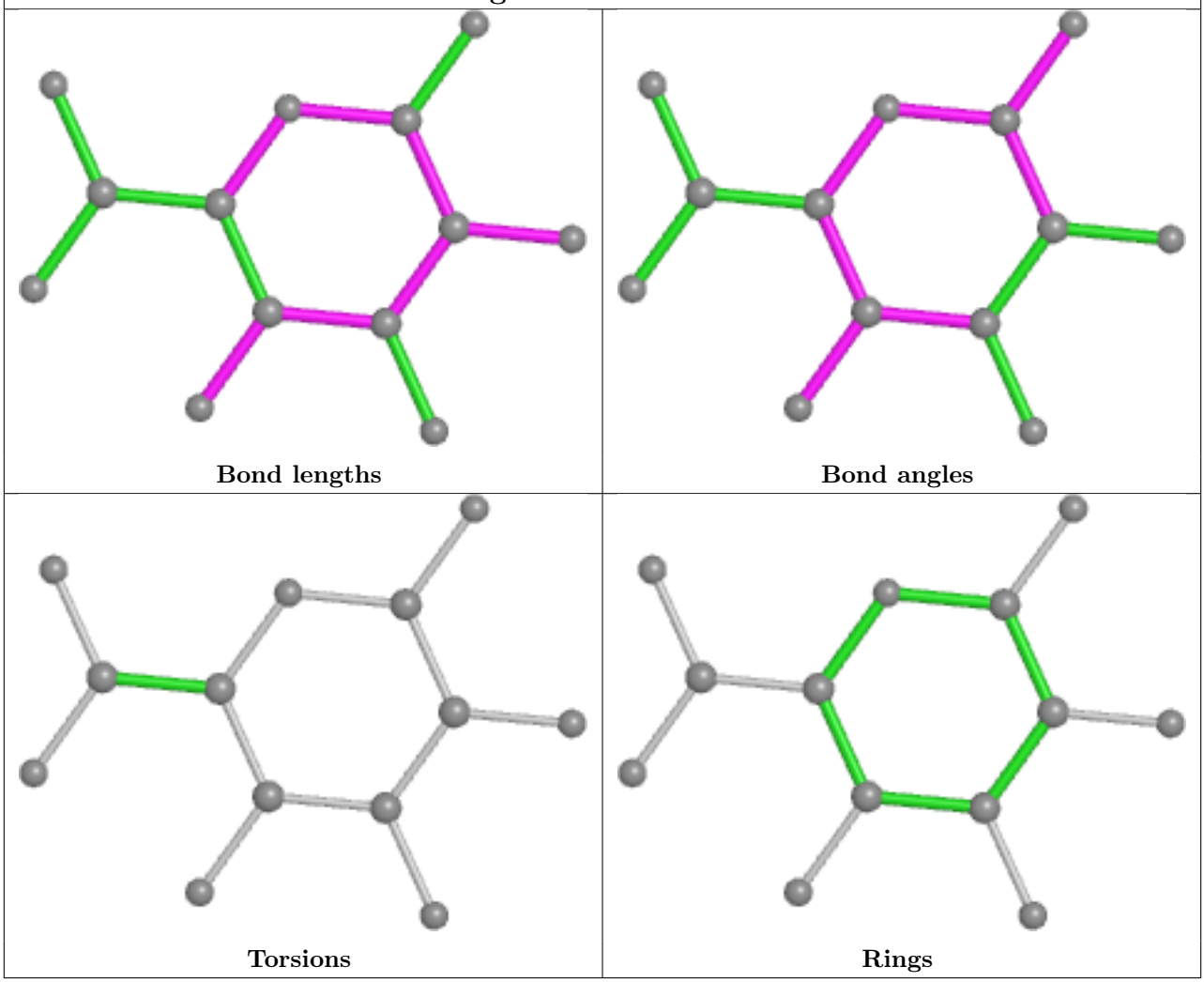
There are no ring outliers.

1 monomer is involved in 1 short contact:

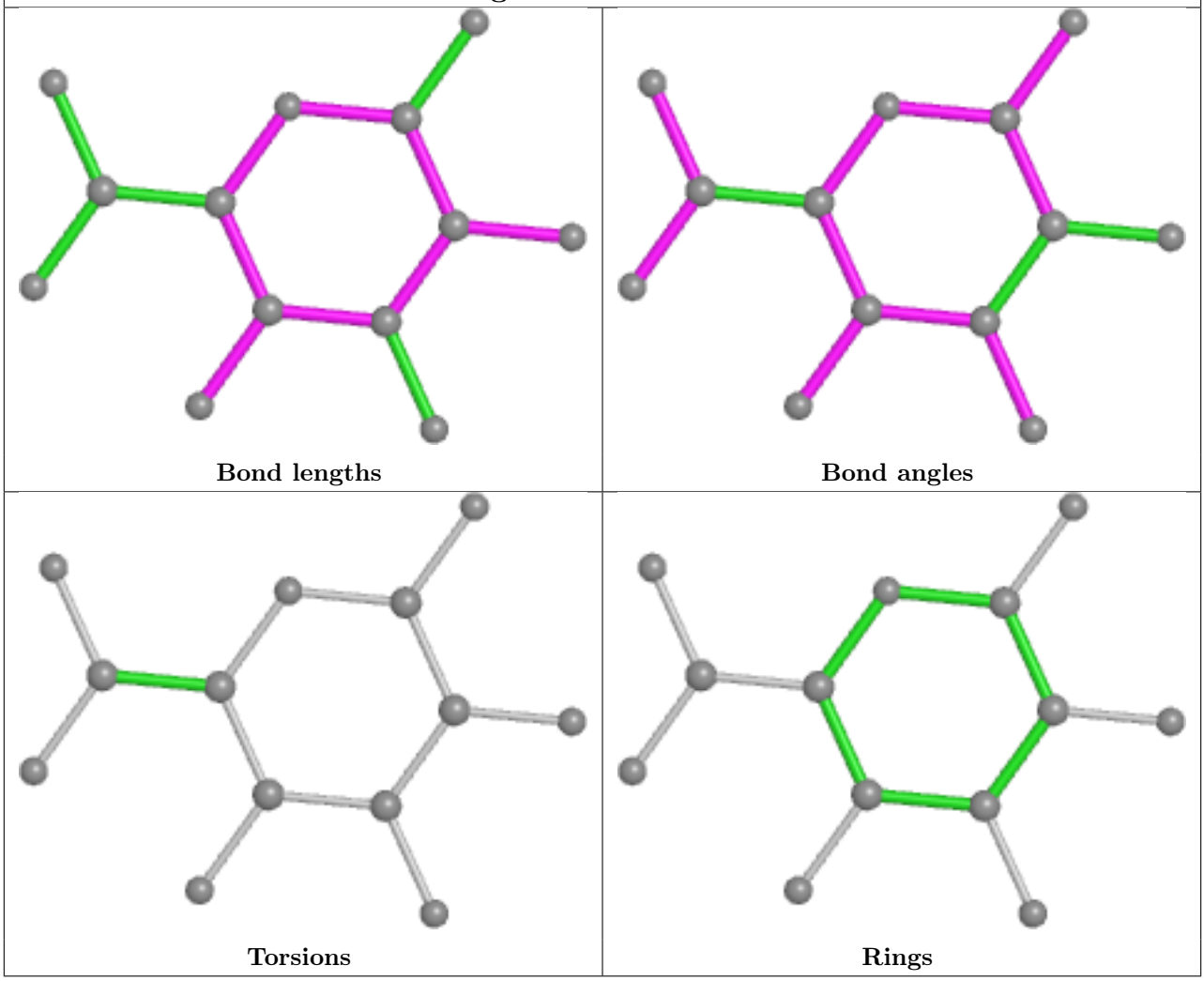
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	903	GCB	1	0

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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand GCB A 903



Ligand GCB B 903



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	837/886 (94%)	-0.10	28 (3%) 46 50	28, 41, 60, 81	0
1	B	840/886 (94%)	-0.27	6 (0%) 87 89	26, 37, 53, 81	0
All	All	1677/1772 (94%)	-0.19	34 (2%) 65 68	26, 39, 58, 81	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	805	ARG	4.9
1	A	50	THR	3.4
1	A	437	THR	3.3
1	A	290	THR	3.1
1	A	439	ALA	3.0
1	B	806	ASN	2.9
1	A	289	LYS	2.9
1	A	144	ASP	2.9
1	A	691	TYR	2.9
1	A	291	LYS	2.7
1	A	441	LEU	2.6
1	A	327	CYS	2.6
1	A	99	LYS	2.6
1	B	822	ARG	2.5
1	A	145	THR	2.5
1	A	382	TRP	2.4
1	B	50	THR	2.4
1	B	714	ASN	2.3
1	A	360	ARG	2.3
1	B	804	ARG	2.3
1	A	435	TYR	2.3
1	A	471	ALA	2.2
1	A	288	ARG	2.2
1	A	693	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	57	PHE	2.2
1	A	615	LYS	2.2
1	A	694	LYS	2.2
1	A	525	SER	2.1
1	A	573	TRP	2.1
1	A	362	SER	2.1
1	A	207	GLU	2.1
1	A	714	ASN	2.1
1	A	361	ILE	2.1
1	A	692	GLN	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)

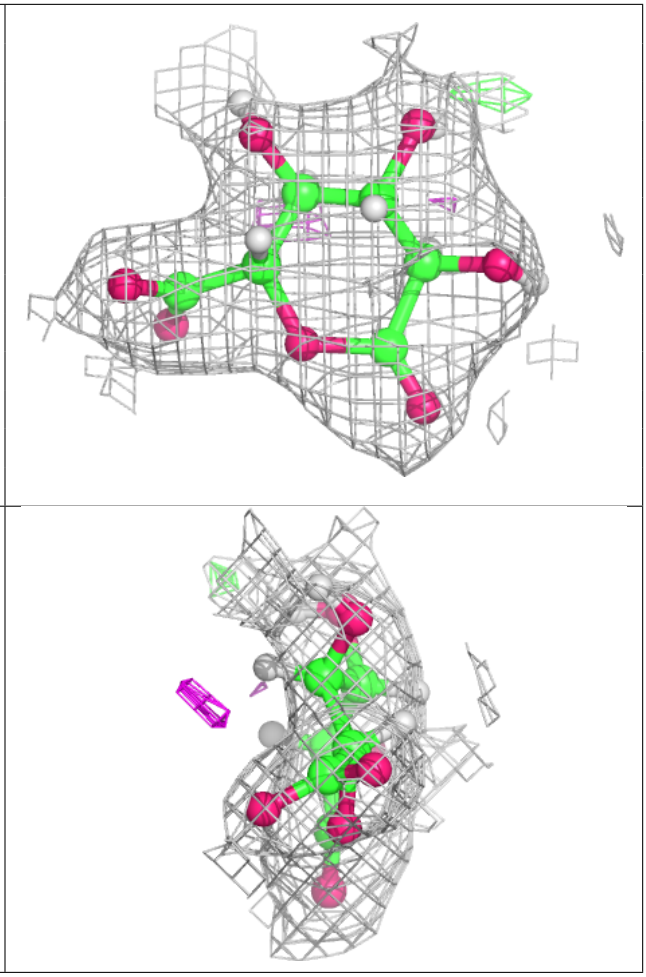
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	GCB	B	903	13/13	0.95	0.12	30,38,44,46	0
4	GCB	A	903	13/13	0.96	0.12	35,42,46,50	0
3	NA	B	902	1/1	0.98	0.22	26,26,26,26	0
3	NA	A	902	1/1	0.99	0.23	25,25,25,25	0
2	CA	A	901	1/1	0.99	0.07	37,37,37,37	0
2	CA	B	901	1/1	1.00	0.07	28,28,28,28	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.

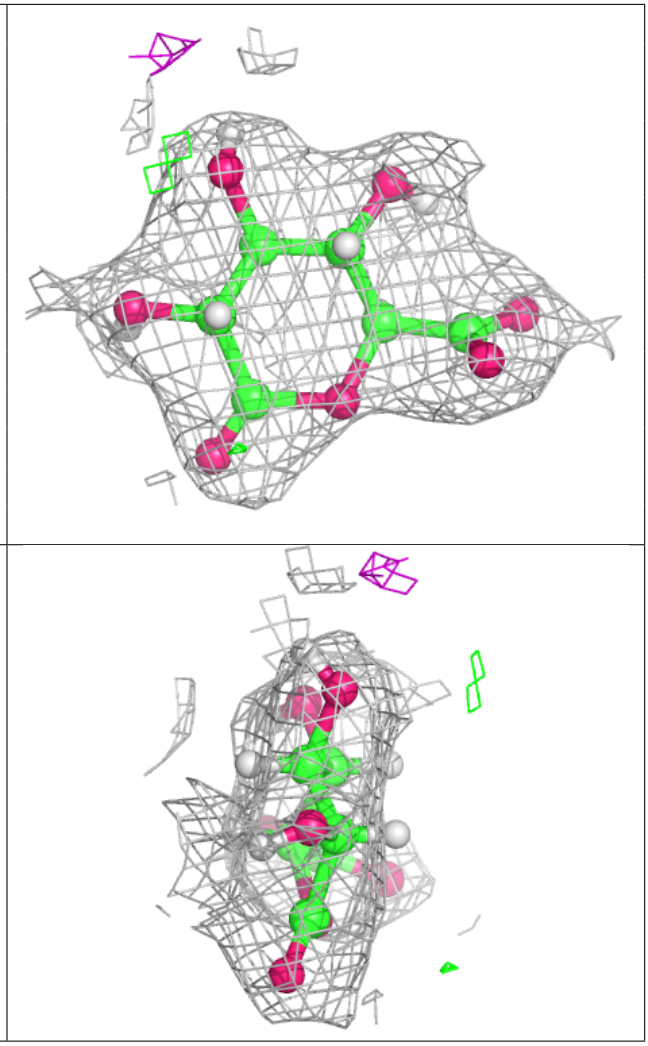
Electron density around GCB B 903:

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



Electron density around GCB A 903:

$2mF_o$-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.