



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2022 – 05:03 PM EDT

PDB ID : 8CSC  
Title : WbbB D232N-Kdo adduct  
Authors : Forrester, T.J.B.; Kimber, M.S.  
Deposited on : 2022-05-12  
Resolution : 1.90 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2.31.2
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)	:	2.31.2

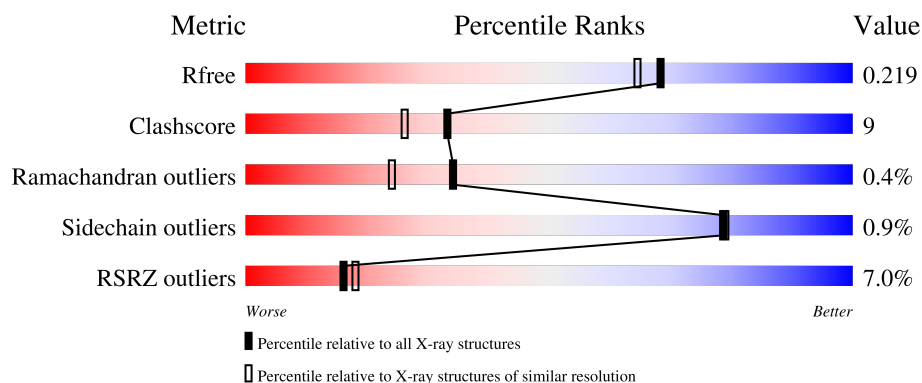
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	410	<div> <div>4%</div> <div> <div></div> <div>80%</div> <div>15%</div> <div>• •</div> </div> </div>
1	B	410	<div> <div>9%</div> <div> <div></div> <div>80%</div> <div>16%</div> <div>•</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6871 atoms, of which 48 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.

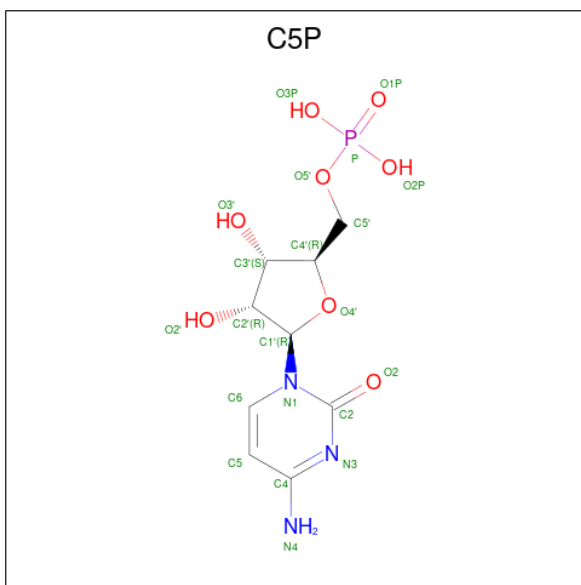
- Molecule 1 is a protein called N-acetyl glucosaminyl transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	8	0
			3193	2042	523	614	14			
1	B	393	Total	C	N	O	S	0	3	0
			3137	2011	510	604	12			

There are 22 discrepancies between the modelled and reference sequences:

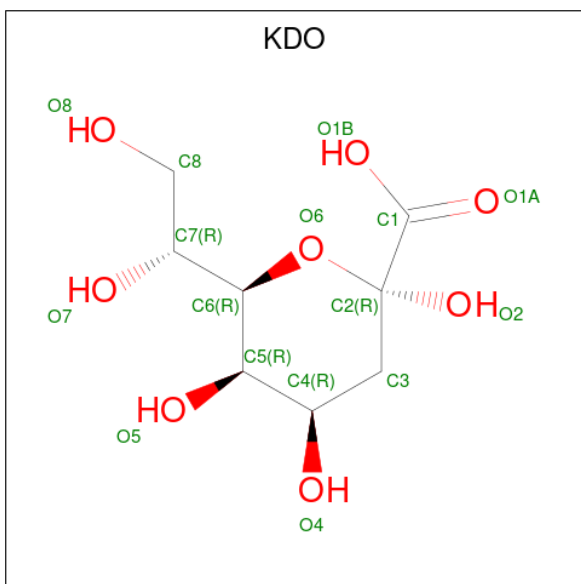
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP Q6U8B0
A	1	GLY	-	expression tag	UNP Q6U8B0
A	232	ASN	ASP	engineered mutation	UNP Q6U8B0
A	402	LEU	-	expression tag	UNP Q6U8B0
A	403	GLU	-	expression tag	UNP Q6U8B0
A	404	HIS	-	expression tag	UNP Q6U8B0
A	405	HIS	-	expression tag	UNP Q6U8B0
A	406	HIS	-	expression tag	UNP Q6U8B0
A	407	HIS	-	expression tag	UNP Q6U8B0
A	408	HIS	-	expression tag	UNP Q6U8B0
A	409	HIS	-	expression tag	UNP Q6U8B0
B	0	MET	-	expression tag	UNP Q6U8B0
B	1	GLY	-	expression tag	UNP Q6U8B0
B	232	ASN	ASP	engineered mutation	UNP Q6U8B0
B	402	LEU	-	expression tag	UNP Q6U8B0
B	403	GLU	-	expression tag	UNP Q6U8B0
B	404	HIS	-	expression tag	UNP Q6U8B0
B	405	HIS	-	expression tag	UNP Q6U8B0
B	406	HIS	-	expression tag	UNP Q6U8B0
B	407	HIS	-	expression tag	UNP Q6U8B0
B	408	HIS	-	expression tag	UNP Q6U8B0
B	409	HIS	-	expression tag	UNP Q6U8B0

- Molecule 2 is CYTIDINE-5'-MONOPHOSPHATE (three-letter code: C5P) (formula: C<sub>9</sub>H<sub>14</sub>N<sub>3</sub>O<sub>8</sub>P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0	0
			33	9	12	3	8	1		
2	B	1	Total	C	H	N	O	P	0	0
			33	9	12	3	8	1		

- Molecule 3 is 3-deoxy- $\alpha$ -D-manno-oct-2-ulopyranosonic acid (three-letter code: KDO) (formula:  $C_8H_{14}O_8$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			27	8	12	7		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	O	0	0
			27	8	12	7		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Na	0	0
			2	2		
4	B	2	Total	Na	0	0
			2	2		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Cl	0	0
			1	1		

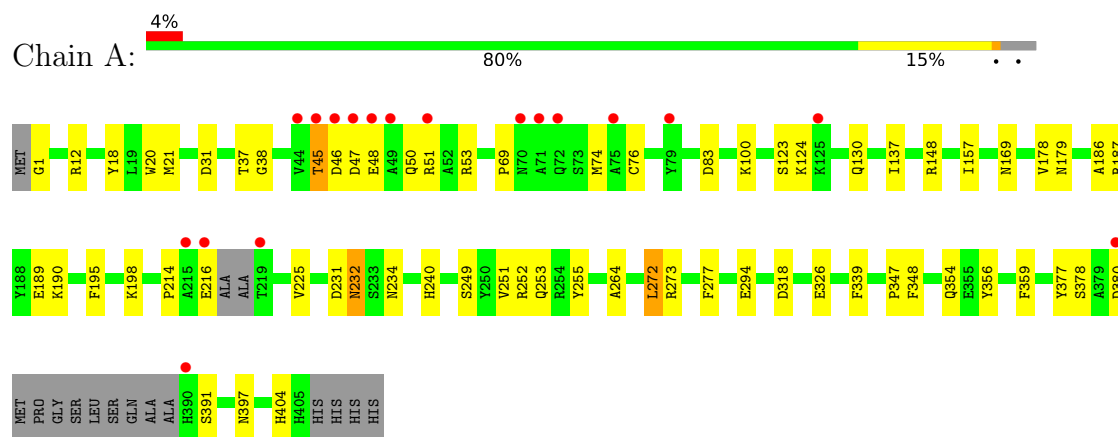
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	208	Total	O	0	0
			208	208		
6	B	208	Total	O	0	0
			208	208		

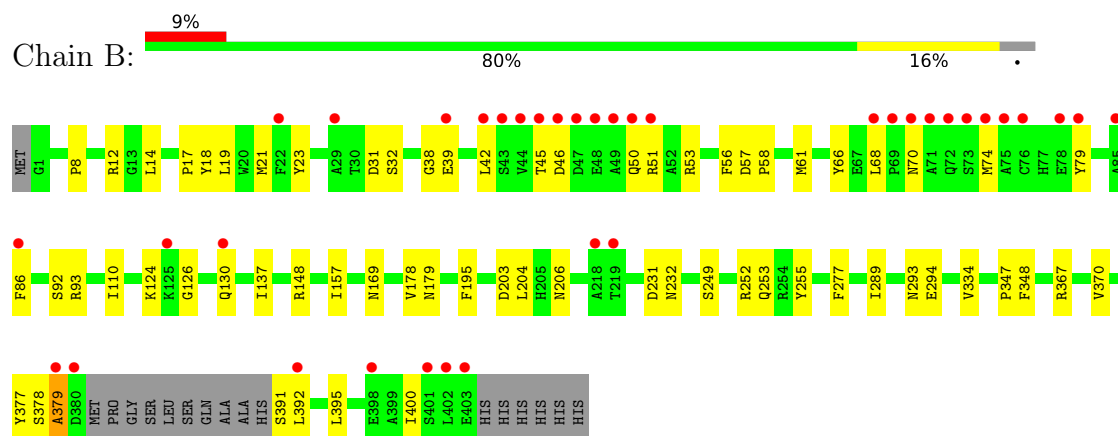
### 3 Residue-property plots [i](#)

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: N-acetyl glucosaminyl transferase



- Molecule 1: N-acetyl glucosaminyl transferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.60Å 158.22Å 117.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.96 – 1.90 47.35 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.96-1.90) 87.2 (47.35-1.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.40 (at 1.75Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, $R_{free}$	0.187 , 0.219 0.186 , 0.219	Depositor DCC
$R_{free}$ test set	4346 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.1	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.017 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	6871	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CSO, C5P, CL, NA, KDO

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	0.26	0/3259	0.45	0/4424
1	B	0.26	0/3201	0.46	0/4348
All	All	0.26	0/6460	0.46	0/8772

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

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	3193	0	3051	56	0
1	B	3137	0	3016	53	0
2	A	21	12	12	0	0
2	B	21	12	12	0	0
3	A	15	12	12	1	0
3	B	15	12	12	2	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	B	1	0	0	1	0
6	A	208	0	0	15	0
6	B	208	0	0	9	0
All	All	6823	48	6115	108	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 9.

All (108) 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:18[B]:TYR:OH	6:A:601:HOH:O	1.97	0.83
1:A:255:TYR:OH	1:A:294:GLU:OE2	1.96	0.83
1:A:214:PRO:O	6:A:602:HOH:O	1.97	0.82
1:A:76[A]:CYS:SG	6:A:803:HOH:O	2.37	0.82
1:A:234[B]:ASN:HD22	1:A:234[B]:ASN:H	1.27	0.81
1:B:206:ASN:ND2	6:B:604:HOH:O	2.13	0.80
1:B:204:LEU:HD23	1:B:400:ILE:HD11	1.62	0.79
1:A:190:LYS:NZ	6:A:608:HOH:O	2.15	0.77
1:A:354:GLN:OE1	6:A:603:HOH:O	2.02	0.76
1:B:92:SER:O	6:B:601:HOH:O	2.05	0.72
1:B:18[B]:TYR:OH	6:B:602:HOH:O	2.07	0.72
1:A:378:SER:OG	6:A:604:HOH:O	2.09	0.69
1:A:83:ASP:OD1	6:A:607:HOH:O	2.10	0.69
1:A:189[B]:GLU:OE1	6:A:605:HOH:O	2.09	0.69
1:A:326:GLU:OE1	6:A:606:HOH:O	2.10	0.68
1:B:58:PRO:O	6:B:603:HOH:O	2.12	0.66
1:B:203:ASP:OD1	6:B:605:HOH:O	2.14	0.66
1:B:45:THR:O	1:B:51:ARG:NH2	2.28	0.65
1:B:137:ILE:HG12	1:B:137:ILE:O	1.96	0.64
1:A:124:LYS:NZ	6:A:610:HOH:O	2.30	0.63
1:B:46:ASP:HA	1:B:51:ARG:HH22	1.63	0.63
1:A:18[A]:TYR:CZ	1:A:137:ILE:HD11	2.35	0.62
1:A:21[A]:MET:SD	1:A:137:ILE:HG22	2.40	0.62
1:A:277:PHE:N	6:A:611:HOH:O	2.33	0.62
1:B:169:ASN:HB3	1:B:348:PHE:HB2	1.83	0.61
1:B:255:TYR:OH	1:B:294:GLU:OE2	2.13	0.61
1:B:232:ASN:ND2	3:B:502:KDO:O6	2.31	0.61
1:A:1:GLY:N	1:A:130:GLN:O	2.34	0.60
1:A:123:SER:O	1:A:124:LYS:HG3	2.02	0.59
1:B:249:SER:O	1:B:253:GLN:HG2	2.02	0.59
1:A:216:GLU:HG3	6:A:602:HOH:O	2.03	0.58
1:B:334:VAL:HG23	6:B:683:HOH:O	2.01	0.58
1:A:179:ASN:HB3	1:A:234[A]:ASN:OD1	2.04	0.58
1:A:264:ALA:HB2	1:A:272:LEU:HD12	1.87	0.57
1:B:169:ASN:HB3	1:B:348:PHE:CB	2.35	0.56
1:B:252:ARG:HD2	1:B:277:PHE:HZ	1.71	0.56
1:A:249:SER:O	1:A:253:GLN:HG2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:PHE:HD1	1:B:68:LEU:HD21	1.71	0.56
1:B:391:SER:O	1:B:395:LEU:HG	2.06	0.55
1:B:74:MET:HB2	1:B:79:TYR:OH	2.07	0.54
1:B:18[A]:TYR:CD1	1:B:137:ILE:HD13	2.43	0.54
1:A:234[B]:ASN:HD22	1:A:234[B]:ASN:N	1.98	0.53
1:A:198:LYS:HE2	1:A:397:ASN:ND2	2.23	0.53
1:A:169:ASN:HB3	1:A:348:PHE:HB2	1.89	0.52
1:B:378:SER:O	1:B:379:ALA:CB	2.57	0.52
1:A:12:ARG:HD3	1:A:231:ASP:O	2.10	0.52
1:B:203:ASP:HB3	1:B:400:ILE:HG21	1.91	0.52
1:A:31:ASP:H	1:A:354:GLN:NE2	2.08	0.52
1:A:48:GLU:HA	1:A:51:ARG:HD3	1.92	0.52
1:A:31:ASP:H	1:A:354:GLN:HE22	1.58	0.51
1:B:12:ARG:HD3	1:B:231:ASP:O	2.10	0.51
1:A:169:ASN:HB3	1:A:348:PHE:CB	2.41	0.51
1:B:14:LEU:HD11	1:B:17:PRO:HA	1.93	0.50
1:A:21[A]:MET:SD	1:A:137:ILE:CG2	2.99	0.50
1:A:232:ASN:ND2	3:A:502:KDO:O6	2.43	0.50
1:B:57:ASP:O	1:B:61:MET:HG3	2.11	0.49
1:B:70:ASN:O	1:B:74:MET:HG2	2.13	0.49
1:A:234[B]:ASN:H	1:A:234[B]:ASN:ND2	2.06	0.48
1:B:21:MET:SD	1:B:137:ILE:HD12	2.52	0.48
1:B:347:PRO:HG3	1:B:377:TYR:CZ	2.48	0.48
1:A:45:THR:C	1:A:47:ASP:H	2.16	0.48
1:B:126:GLY:O	1:B:130:GLN:HG2	2.13	0.47
1:A:69:PRO:HG2	1:A:74:MET:CE	2.45	0.47
1:A:50:GLN:OE1	1:A:53:ARG:HD2	2.14	0.47
1:A:148:ARG:HG3	1:B:148:ARG:NH2	2.29	0.47
1:B:86:PHE:CZ	1:B:110:ILE:HG21	2.50	0.47
1:B:14:LEU:C	1:B:14:LEU:HD12	2.36	0.46
1:A:186:ALA:O	1:A:190:LYS:HG3	2.16	0.46
1:B:61:MET:HB3	1:B:66:TYR:O	2.15	0.46
1:A:20:TRP:CZ2	1:A:21[B]:MET:HG2	2.51	0.46
1:A:137:ILE:C	1:A:137:ILE:HD12	2.36	0.46
1:B:50:GLN:OE1	1:B:53:ARG:HD2	2.16	0.46
1:A:354:GLN:HG3	6:A:624:HOH:O	2.16	0.45
1:B:148:ARG:NH1	6:B:612:HOH:O	2.44	0.45
1:A:178:VAL:O	1:A:179:ASN:HB2	2.17	0.45
1:B:14:LEU:HD12	1:B:14:LEU:O	2.17	0.45
1:A:37:THR:OG1	1:A:38:GLY:N	2.50	0.44
1:A:45:THR:O	1:A:47:ASP:N	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:LYS:HB2	1:B:124:LYS:HE3	1.87	0.44
1:A:240:HIS:NE2	1:A:318[A]:ASP:OD2	2.50	0.44
1:B:378:SER:O	1:B:379:ALA:HB2	2.17	0.44
1:A:69:PRO:HG2	1:A:74:MET:HE1	2.00	0.44
1:B:367:ARG:HB3	1:B:367:ARG:NH1	2.33	0.43
1:B:39:GLU:HA	1:B:42:LEU:HD12	2.00	0.43
1:B:392:LEU:HD23	1:B:392:LEU:O	2.18	0.43
1:A:195:PHE:CZ	1:A:339:PHE:HB2	2.53	0.43
1:B:289:ILE:HG22	6:B:708:HOH:O	2.17	0.43
1:A:148:ARG:NH1	6:A:622:HOH:O	2.51	0.43
1:A:252:ARG:HD2	1:A:277:PHE:CZ	2.54	0.43
1:B:178:VAL:O	1:B:179:ASN:HB2	2.18	0.42
1:A:225:VAL:HG22	1:A:251:VAL:HG21	2.01	0.42
1:A:347:PRO:HG3	1:A:377:TYR:CZ	2.54	0.42
1:A:391:SER:HB2	6:A:715:HOH:O	2.20	0.41
1:A:137:ILE:HA	1:A:157:ILE:O	2.20	0.41
1:B:31[B]:ASP:OD1	1:B:32:SER:N	2.50	0.41
1:B:204:LEU:CD2	1:B:400:ILE:HD11	2.42	0.41
1:A:356:TYR:O	1:A:359:PHE:HB3	2.20	0.41
1:B:93:ARG:NE	6:B:615:HOH:O	2.53	0.41
1:B:195:PHE:CZ	1:B:392:LEU:HD22	2.55	0.41
1:B:232:ASN:HD21	3:B:502:KDO:H82	1.86	0.41
1:A:187:ARG:HB3	1:A:339:PHE:HZ	1.85	0.41
1:B:370:VAL:HG21	1:B:392:LEU:HD21	2.02	0.41
1:B:8:PRO:HB3	1:B:38:GLY:HA3	2.03	0.41
1:B:19:LEU:HD11	1:B:23:TYR:CZ	2.56	0.41
1:B:137:ILE:HA	1:B:157:ILE:O	2.21	0.40
1:A:123:SER:C	1:A:124:LYS:HG3	2.42	0.40
1:B:179:ASN:HA	5:B:505:CL:CL	2.58	0.40
1:A:100:LYS:HA	1:A:100:LYS:HD3	1.84	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	394/410 (96%)	381 (97%)	11 (3%)	2 (0%)	29	18
1	B	390/410 (95%)	380 (97%)	9 (2%)	1 (0%)	41	31
All	All	784/820 (96%)	761 (97%)	20 (3%)	3 (0%)	34	24

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	379	ALA
1	A	45	THR
1	A	46	ASP

### 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	343/347 (99%)	338 (98%)	5 (2%)	65	62
1	B	335/347 (96%)	334 (100%)	1 (0%)	92	93
All	All	678/694 (98%)	672 (99%)	6 (1%)	78	79

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	232	ASN
1	A	272	LEU
1	A	273	ARG
1	A	380	ASP
1	A	404	HIS
1	B	293	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	A	354	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
1	CSO	B	28	1	3,6,7	0.60	0	0,6,8	-	-
1	CSO	A	191	1	3,6,7	0.56	0	0,6,8	-	-
1	CSO	A	28	1	3,6,7	0.57	0	0,6,8	-	-
1	CSO	B	191	1	3,6,7	0.62	0	0,6,8	-	-

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
1	CSO	B	28	1	-	1/1/5/7	-
1	CSO	A	191	1	-	0/1/5/7	-
1	CSO	A	28	1	-	0/1/5/7	-
1	CSO	B	191	1	-	0/1/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	28	CSO	N-CA-CB-SG

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 5 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$
3	KDO	B	502	-	15,15,16	2.02	4 (26%)	19,21,24	1.40	4 (21%)
2	C5P	B	501	-	22,22,22	2.80	10 (45%)	33,33,33	1.51	9 (27%)
2	C5P	A	501	-	22,22,22	2.73	9 (40%)	33,33,33	1.61	9 (27%)
3	KDO	A	502	-	15,15,16	1.93	4 (26%)	19,21,24	1.50	3 (15%)

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	KDO	B	502	-	-	3/10/26/30	0/1/1/1
2	C5P	B	501	-	-	0/10/26/26	0/2/2/2
2	C5P	A	501	-	-	0/10/26/26	0/2/2/2
3	KDO	A	502	-	-	3/10/26/30	0/1/1/1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	C5P	P-O5'	8.09	1.86	1.60
2	A	501	C5P	P-O5'	7.50	1.84	1.60
3	B	502	KDO	O6-C2	5.59	1.51	1.43
3	A	502	KDO	O6-C2	5.33	1.50	1.43
2	A	501	C5P	C4-N4	4.85	1.45	1.33
2	B	501	C5P	C2-N3	4.80	1.46	1.36
2	A	501	C5P	C2-N3	4.70	1.45	1.36
2	B	501	C5P	C4-N4	4.69	1.44	1.33
2	A	501	C5P	C2-N1	3.30	1.47	1.40
2	A	501	C5P	C5-C4	3.02	1.49	1.42
2	B	501	C5P	C5-C4	2.93	1.49	1.42
2	B	501	C5P	C1'-N1	2.91	1.56	1.47
2	B	501	C5P	O5'-C5'	-2.85	1.33	1.44
2	A	501	C5P	O5'-C5'	-2.81	1.34	1.44
2	A	501	C5P	C1'-N1	2.70	1.55	1.47
2	A	501	C5P	C5'-C4'	2.66	1.59	1.51
2	B	501	C5P	C5'-C4'	2.63	1.59	1.51
3	B	502	KDO	C3-C4	-2.62	1.48	1.52
3	B	502	KDO	C2-C1	2.61	1.54	1.52
3	A	502	KDO	C2-C1	2.54	1.54	1.52
3	A	502	KDO	C3-C4	-2.53	1.48	1.52
2	B	501	C5P	C2-N1	2.39	1.45	1.40
2	B	501	C5P	O3'-C3'	-2.23	1.37	1.43
3	B	502	KDO	O6-C6	2.19	1.47	1.44
2	A	501	C5P	O3'-C3'	-2.19	1.37	1.43
2	B	501	C5P	C6-C5	2.03	1.39	1.35
3	A	502	KDO	O6-C6	2.01	1.47	1.44

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	KDO	O6-C6-C5	3.90	113.35	107.87
2	A	501	C5P	C5'-C4'-C3'	-3.07	103.66	115.18
2	A	501	C5P	O3'-C3'-C4'	3.01	119.75	111.05
2	A	501	C5P	O2P-P-O5'	-2.78	99.34	106.73
2	B	501	C5P	O2'-C2'-C3'	-2.76	102.90	111.82
2	B	501	C5P	C5'-C4'-C3'	-2.71	105.01	115.18
2	A	501	C5P	O5'-P-O1P	-2.65	99.05	106.47
2	B	501	C5P	O2P-P-O5'	-2.62	99.77	106.73
3	A	502	KDO	C7-C6-C5	-2.60	109.71	114.03
3	B	502	KDO	C4-C3-C2	2.54	114.37	109.81
2	B	501	C5P	C4-N3-C2	2.50	124.30	120.25
3	B	502	KDO	O6-C6-C5	2.46	111.32	107.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	C5P	O3'-C3'-C2'	-2.44	103.92	111.82
2	A	501	C5P	P-O5'-C5'	2.41	124.94	118.30
3	B	502	KDO	O1B-C1-C2	2.40	119.89	113.03
2	B	501	C5P	P-O5'-C5'	2.37	124.82	118.30
2	A	501	C5P	O2'-C2'-C3'	-2.35	104.21	111.82
2	A	501	C5P	C4-N3-C2	2.35	124.05	120.25
3	B	502	KDO	O6-C2-C3	2.28	113.60	110.46
3	A	502	KDO	O1B-C1-C2	2.24	119.43	113.03
2	B	501	C5P	O3'-C3'-C4'	2.24	117.51	111.05
2	B	501	C5P	O5'-P-O1P	-2.15	100.43	106.47
2	B	501	C5P	O2P-P-O3P	2.08	115.60	107.64
2	A	501	C5P	N1-C2-N3	-2.05	115.07	118.81
2	B	501	C5P	O2-C2-N3	2.02	125.63	122.33

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	502	KDO	O1A-C1-C2-O6
3	A	502	KDO	O1A-C1-C2-O6
3	A	502	KDO	O1A-C1-C2-C3
3	A	502	KDO	O1B-C1-C2-C3
3	B	502	KDO	O1A-C1-C2-C3
3	B	502	KDO	O1B-C1-C2-C3

There are no ring outliers.

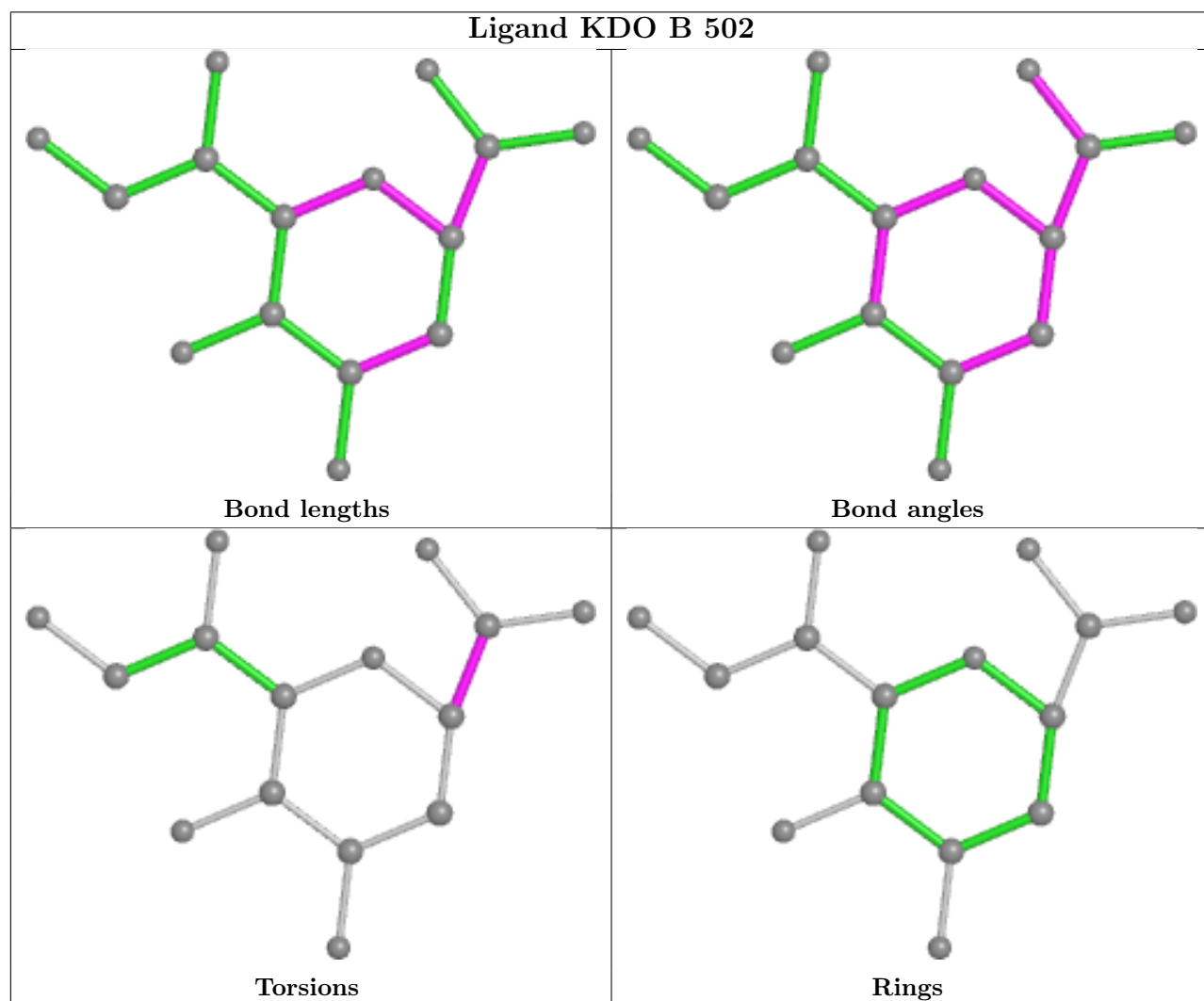
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	KDO	2	0
3	A	502	KDO	1	0

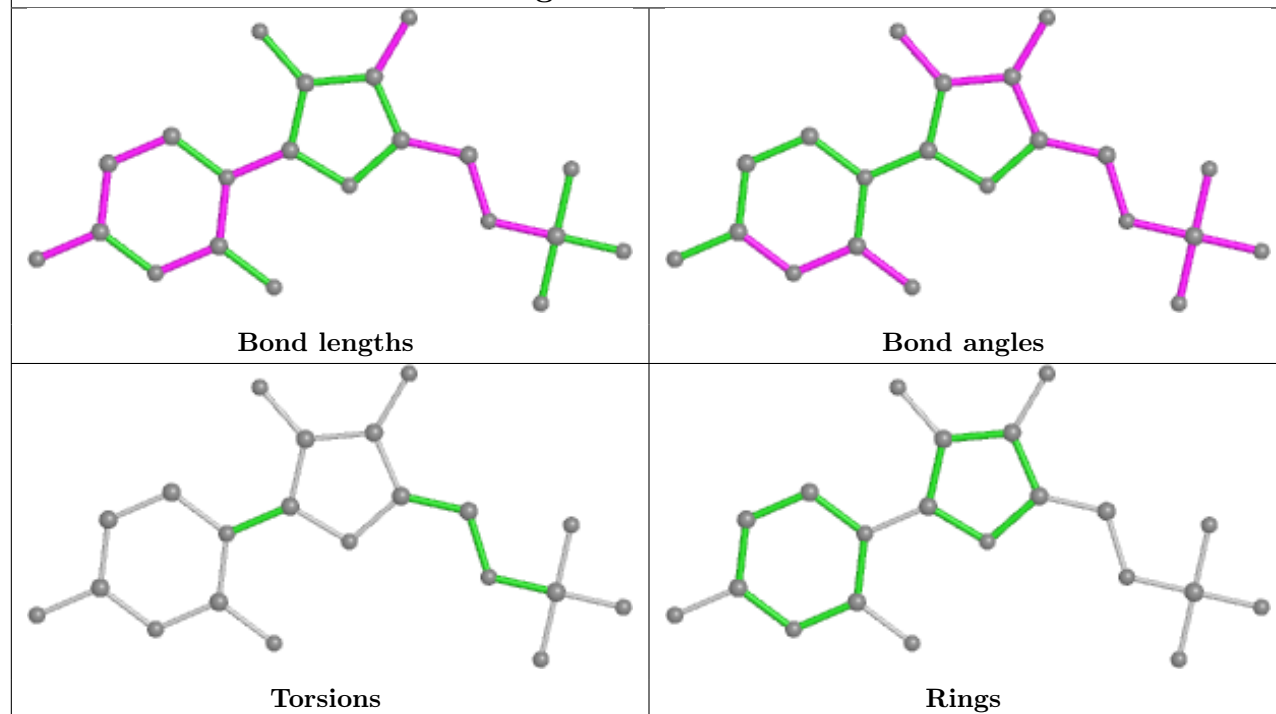
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



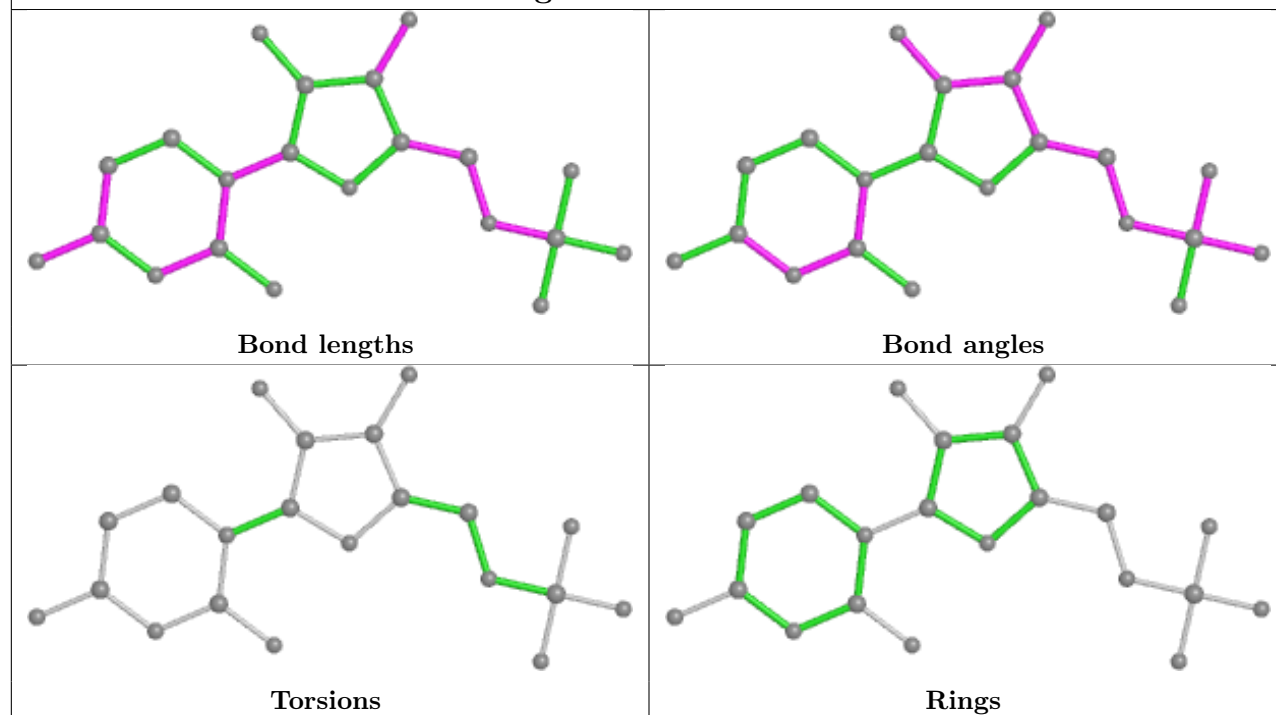
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

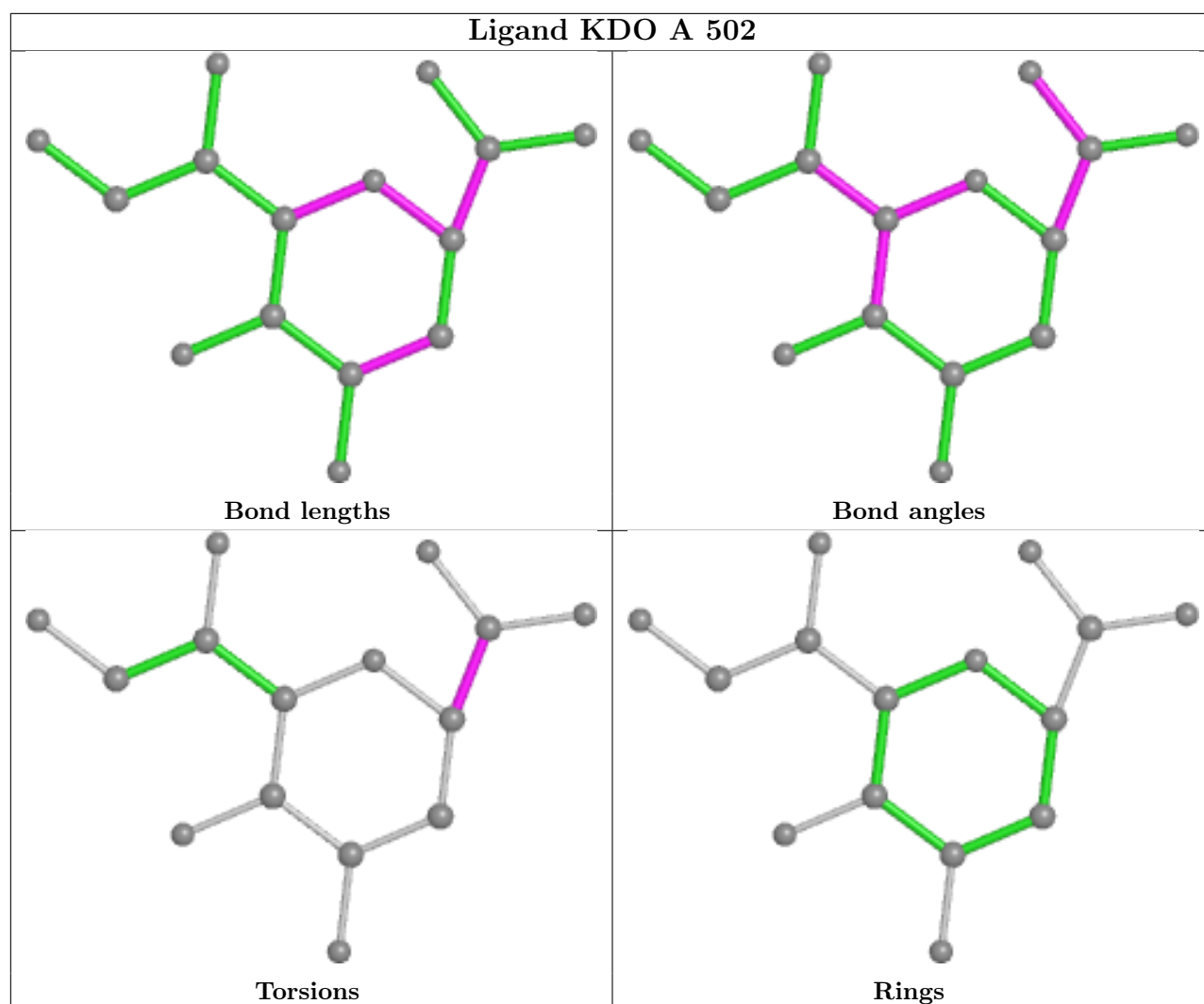


## Ligand C5P B 501



## Ligand C5P A 501





## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	392/410 (95%)	-0.00	18 (4%) 32 35	27, 44, 81, 116	0
1	B	391/410 (95%)	0.38	37 (9%) 8 9	26, 47, 93, 133	0
All	All	783/820 (95%)	0.19	55 (7%) 16 18	26, 45, 87, 133	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	45	THR	7.9
1	B	44	VAL	7.5
1	B	50	GLN	7.4
1	B	49	ALA	6.2
1	B	46	ASP	6.0
1	B	51	ARG	5.5
1	B	72	GLN	5.0
1	B	71	ALA	5.0
1	A	219	THR	4.9
1	B	48	GLU	4.5
1	B	70	ASN	4.3
1	B	380	ASP	4.2
1	A	51	ARG	3.8
1	B	39	GLU	3.8
1	A	46	ASP	3.7
1	A	380	ASP	3.5
1	B	42	LEU	3.4
1	B	75	ALA	3.3
1	B	47	ASP	3.3
1	B	74	MET	3.2
1	B	401	SER	3.1
1	B	69	PRO	3.0
1	B	86	PHE	3.0
1	A	49	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	29	ALA	2.9
1	A	390	HIS	2.9
1	A	48	GLU	2.8
1	B	43	SER	2.8
1	B	73	SER	2.8
1	A	71	ALA	2.8
1	B	22	PHE	2.7
1	A	125	LYS	2.7
1	B	402	LEU	2.7
1	A	72	GLN	2.7
1	A	44	VAL	2.7
1	B	125	LYS	2.6
1	A	45	THR	2.6
1	B	79	TYR	2.6
1	A	75	ALA	2.6
1	B	85	ALA	2.6
1	B	130	GLN	2.6
1	A	215	ALA	2.6
1	B	218	ALA	2.4
1	B	219	THR	2.4
1	B	379	ALA	2.3
1	B	403	GLU	2.3
1	B	76	CYS	2.2
1	A	70	ASN	2.1
1	B	78	GLU	2.1
1	A	216	GLU	2.1
1	B	68	LEU	2.1
1	A	79	TYR	2.1
1	B	398	GLU	2.1
1	B	392	LEU	2.1
1	A	47	ASP	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	CSO	A	191	7/8	0.86	0.17	46,51,59,63	0
1	CSO	B	28	7/8	0.87	0.11	66,69,73,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
1	CSO	B	191	7/8	0.92	0.15	56,65,71,102	0
1	CSO	A	28	7/8	0.93	0.07	52,54,58,61	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

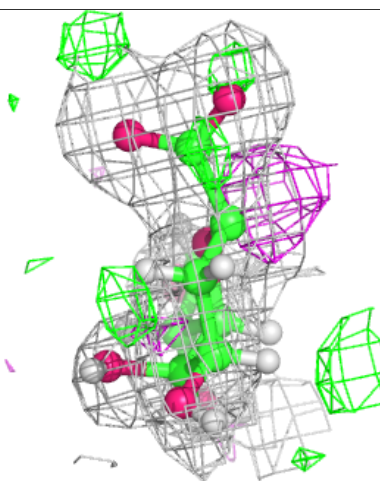
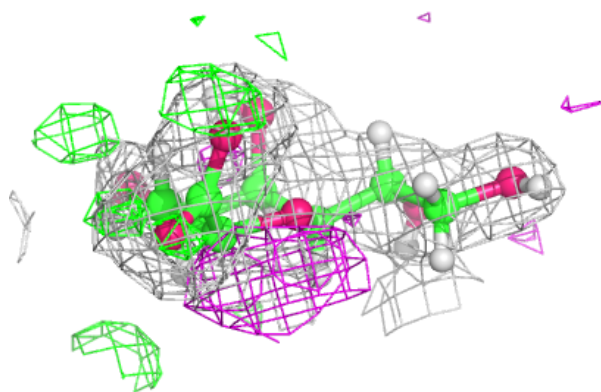
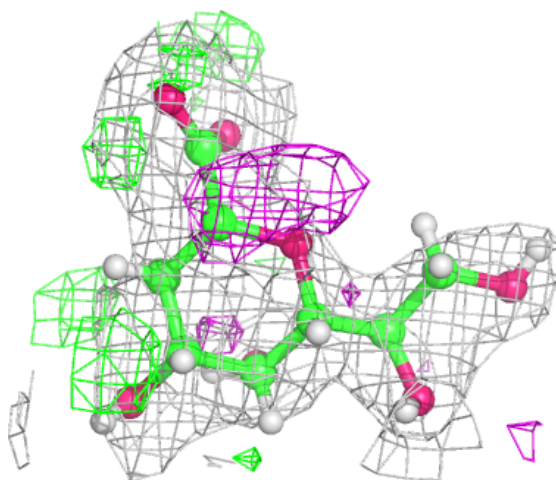
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	KDO	A	502	15/16	0.64	0.18	40,54,66,67	27
4	NA	B	504	1/1	0.69	0.09	52,52,52,52	0
3	KDO	B	502	15/16	0.74	0.19	40,51,60,62	27
4	NA	B	503	1/1	0.87	0.13	69,69,69,69	0
4	NA	A	504	1/1	0.95	0.14	44,44,44,44	0
2	C5P	A	501	21/21	0.96	0.12	28,34,45,49	0
5	CL	B	505	1/1	0.97	0.05	43,43,43,43	0
2	C5P	B	501	21/21	0.98	0.10	27,33,43,44	0
4	NA	A	503	1/1	0.99	0.11	35,35,35,35	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.

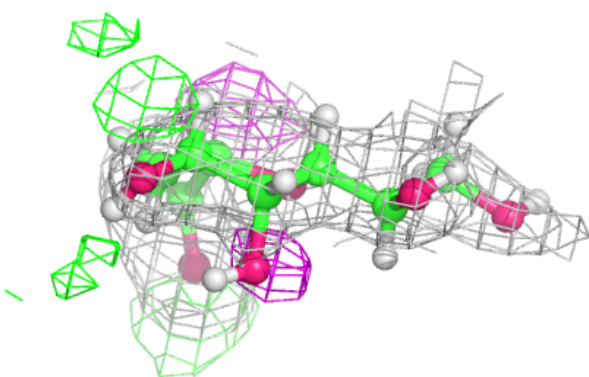
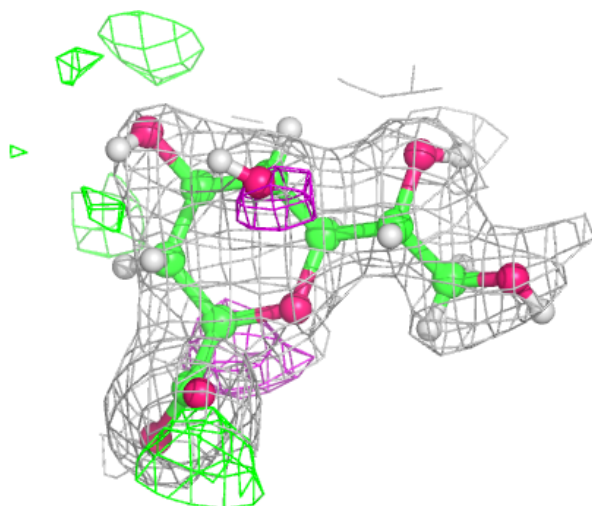
**Electron density around KDO A 502:**

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



**Electron density around KDO B 502:**

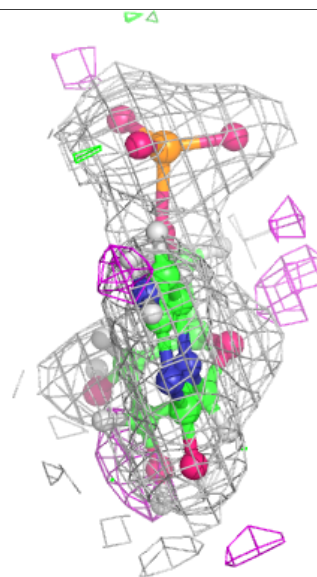
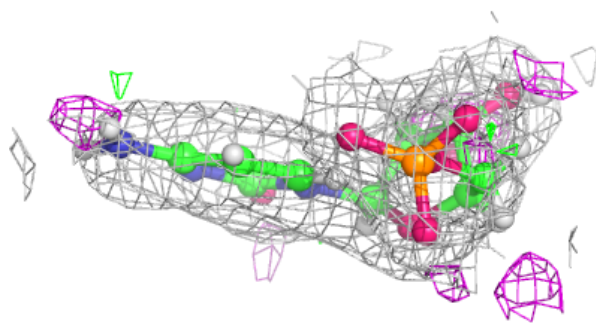
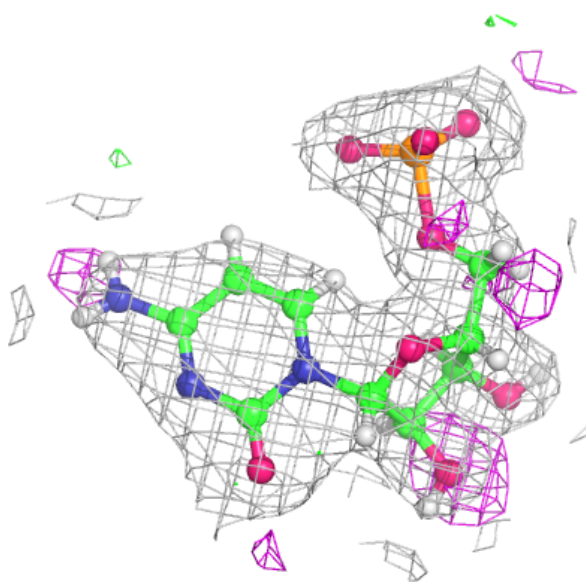
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





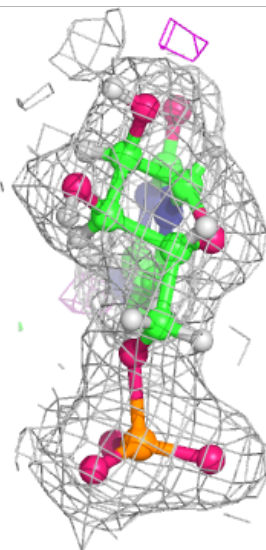
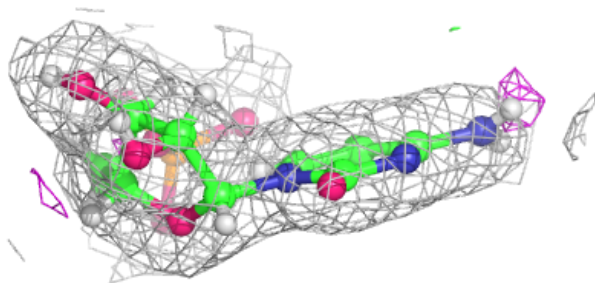
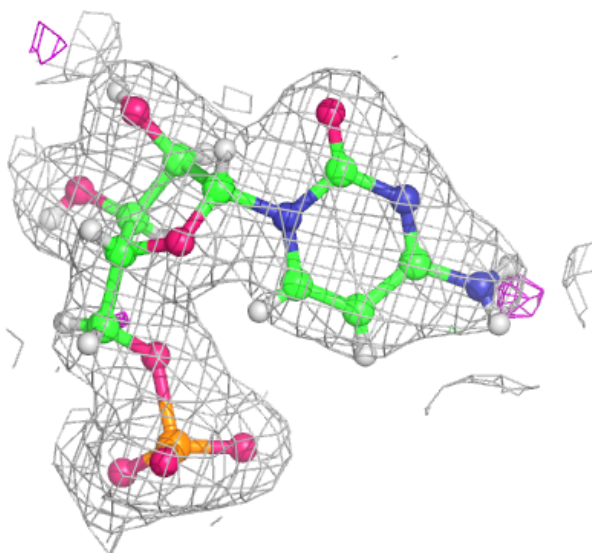
**Electron density around C5P A 501:**

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



**Electron density around C5P B 501:**

$2mF_o-DF_c$  (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.