



Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2021 – 10:10 AM EDT

PDB ID : 7KZ5
Title : Crystal structure of KabA from *Bacillus cereus* UW85 in complex with the plp external aldimine adduct with kanosamine-6-phosphate
Authors : Prasertanan, T.; Palmer, D.R.J.; Sanders, D.A.R.
Deposited on : 2020-12-10
Resolution : 1.60 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	:	2.18
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)	:	2.18

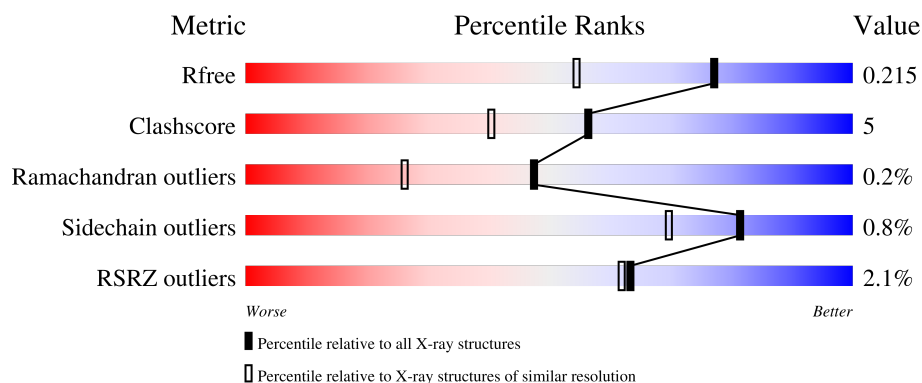
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	445	<div> <div>2%</div> <div>91%</div> <div>7%</div> </div>
1	B	445	<div> <div>2%</div> <div>86%</div> <div>12%</div> </div>
1	C	445	<div> <div>2%</div> <div>89%</div> <div>10%</div> </div>
1	D	445	<div> <div>3%</div> <div>87%</div> <div>12%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15818 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.

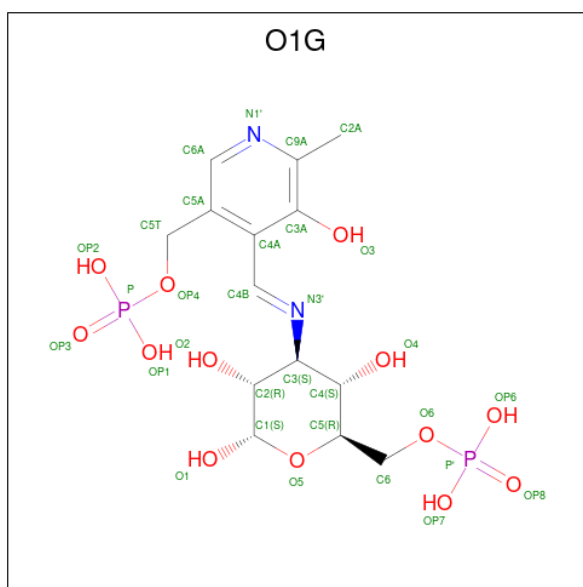
- Molecule 1 is a protein called Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	439	Total	C	N	O	S	0	3	0
			3563	2276	608	664	15			
1	B	439	Total	C	N	O	S	0	5	0
			3577	2286	611	665	15			
1	C	439	Total	C	N	O	S	0	5	0
			3572	2285	609	662	16			
1	D	439	Total	C	N	O	S	0	5	0
			3573	2283	608	667	15			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP C0JRF5
A	0	ALA	-	expression tag	UNP C0JRF5
A	1	MET	-	expression tag	UNP C0JRF5
A	2	ASP	-	expression tag	UNP C0JRF5
B	-1	GLY	-	expression tag	UNP C0JRF5
B	0	ALA	-	expression tag	UNP C0JRF5
B	1	MET	-	expression tag	UNP C0JRF5
B	2	ASP	-	expression tag	UNP C0JRF5
C	-1	GLY	-	expression tag	UNP C0JRF5
C	0	ALA	-	expression tag	UNP C0JRF5
C	1	MET	-	expression tag	UNP C0JRF5
C	2	ASP	-	expression tag	UNP C0JRF5
D	-1	GLY	-	expression tag	UNP C0JRF5
D	0	ALA	-	expression tag	UNP C0JRF5
D	1	MET	-	expression tag	UNP C0JRF5
D	2	ASP	-	expression tag	UNP C0JRF5

- Molecule 2 is 3-deoxy-3-[(E)-({3-hydroxy-2-methyl-5-[(phosphonooxy)methyl]pyridin-4-yl}methylidene)amino]-6-O-phosphono- α -D-glucopyranose (three-letter code: O1G) (formula: $C_{14}H_{22}N_2O_{13}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	14	2	13	2		
2	B	1	Total	C	N	O	P	0	0
			31	14	2	13	2		
2	C	1	Total	C	N	O	P	0	0
			31	14	2	13	2		
2	D	1	Total	C	N	O	P	0	0
			31	14	2	13	2		

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

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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	335	Total O 335 335	0	0
5	B	280	Total O 280 280	0	0
5	C	339	Total O 339 339	0	0

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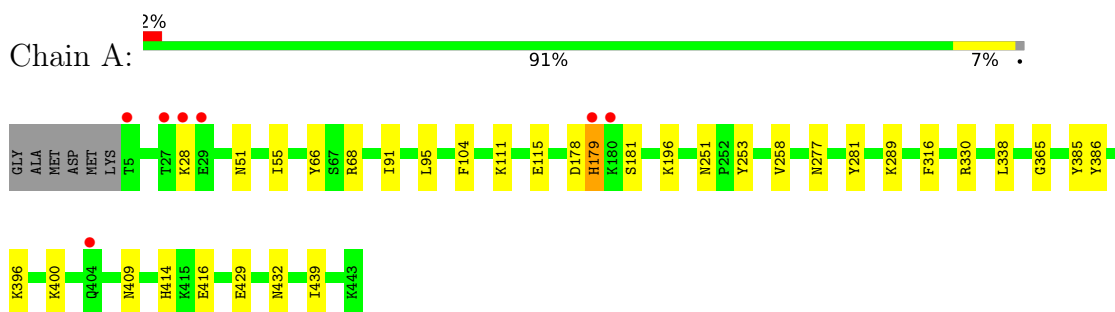
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	249	Total 249	O 249	0	0

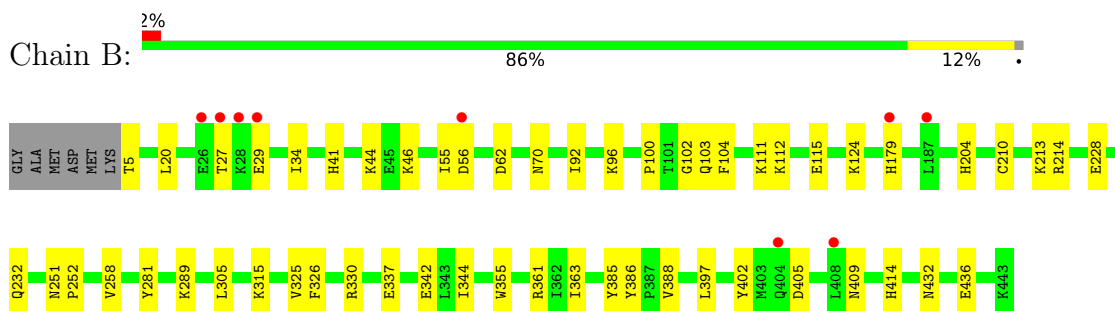
3 Residue-property plots [i](#)

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.

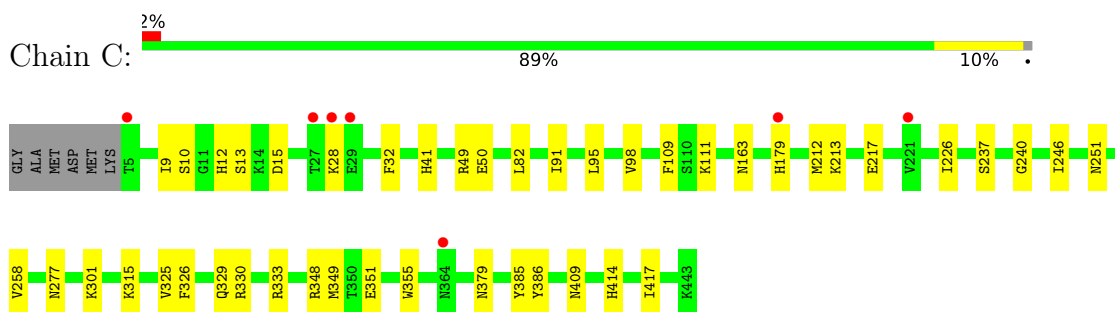
- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



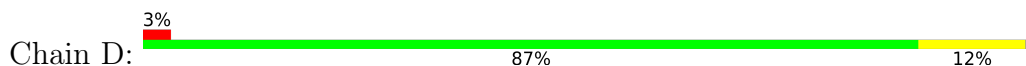
- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme

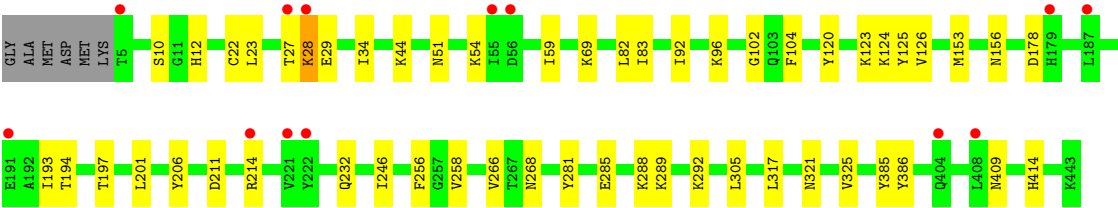


- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	66.67Å 66.60Å 111.78Å 77.35° 81.06° 87.91°	Depositor
Resolution (Å)	46.75 – 1.60 47.25 – 1.60	Depositor EDS
% Data completeness (in resolution range)	92.8 (46.75-1.60) 92.8 (47.25-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 1.60Å)	Xtriage
Refinement program	PHENIX dev_2398	Depositor
R, R_{free}	0.175 , 0.215 0.176 , 0.215	Depositor DCC
R_{free} test set	11353 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.089	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.035 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15818	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.80% 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: NA, EDO, O1G

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	0.85	0/3624	0.88	0/4885
1	B	0.80	0/3642	0.83	0/4910
1	C	0.84	0/3640	0.87	0/4906
1	D	0.76	0/3637	0.83	0/4903
All	All	0.81	0/14543	0.85	0/19604

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	3563	0	3644	32	0
1	B	3577	0	3659	41	0
1	C	3572	0	3664	32	0
1	D	3573	0	3655	41	0
2	A	31	0	8	1	0
2	B	31	0	8	1	0
2	C	31	0	8	2	0
2	D	31	0	8	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1	0	0	0	0
4	A	64	0	95	11	0
4	B	56	0	83	10	0
4	C	44	0	66	9	0
4	D	40	0	60	3	0
5	A	335	0	0	5	0
5	B	280	0	0	6	0
5	C	339	0	0	4	0
5	D	249	0	0	1	0
All	All	15818	0	14958	147	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 5.

All (147) 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:333:ARG:HH11	4:C:510:EDO:H11	1.29	0.94
1:A:68:ARG:HH22	1:D:54:LYS:NZ	1.75	0.83
1:A:68:ARG:HH22	1:D:54:LYS:HZ3	1.28	0.81
1:D:193:ILE:HG23	1:D:197:THR:HG21	1.61	0.81
1:B:204:HIS:HD2	1:B:210:CYS:H	1.28	0.79
1:D:126:VAL:HG22	1:D:266[B]:VAL:HG23	1.64	0.79
1:C:13:SER:HB3	1:C:351:GLU:O	1.83	0.79
1:B:388:VAL:HA	4:B:504:EDO:H12	1.65	0.76
1:A:330:ARG:HH22	1:A:432:ASN:HD21	1.35	0.73
1:B:204:HIS:CD2	1:B:210:CYS:H	2.07	0.72
1:B:325[A]:VAL:HG12	1:B:355:TRP:HD1	1.56	0.71
1:C:111:LYS:HE3	5:C:721:HOH:O	1.90	0.71
1:D:34:ILE:H	1:D:51:ASN:HD21	1.39	0.71
1:D:206:TYR:H	1:D:232:GLN:NE2	1.89	0.70
1:B:325[A]:VAL:HG12	1:B:355:TRP:CD1	2.26	0.70
1:B:388:VAL:HG13	4:B:504:EDO:H11	1.76	0.68
1:B:432:ASN:O	1:B:436:GLU:HG3	1.94	0.68
4:A:508:EDO:H11	1:B:100:PRO:HA	1.76	0.67
1:C:82:LEU:HD22	5:C:722:HOH:O	1.95	0.66
1:D:193:ILE:HG23	1:D:197:THR:CG2	2.27	0.65
1:C:333:ARG:NH1	4:C:510:EDO:H11	2.10	0.64
1:B:46:LYS:NZ	1:B:62:ASP:O	2.27	0.64
1:D:82:LEU:O	4:D:505:EDO:H21	1.98	0.63
1:C:349[A]:MET:HE1	1:C:355:TRP:HE1	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:ASN:ND2	1:A:253:TYR:H	1.97	0.62
1:D:123:LYS:HD3	4:D:502:EDO:H21	1.81	0.61
1:B:204:HIS:HE1	1:B:228:GLU:OE1	1.85	0.59
1:D:178:ASP:O	1:D:409[A]:ASN:ND2	2.35	0.59
1:C:213:LYS:O	1:C:217:GLU:HG3	2.04	0.58
1:B:213:LYS:HG2	5:B:621:HOH:O	2.02	0.57
1:C:13:SER:HB2	1:C:237:SER:H	1.68	0.57
1:C:348:ARG:HG3	4:C:507:EDO:H11	1.85	0.57
1:D:29:GLU:OE2	1:D:29:GLU:N	2.32	0.57
1:A:111:LYS:NZ	1:A:115:GLU:HG3	2.20	0.56
1:B:70:ASN:ND2	1:C:28:LYS:HE3	2.20	0.56
1:D:194:THR:O	1:D:197:THR:HG22	2.05	0.56
1:D:246:ILE:HB	1:D:266[B]:VAL:HG12	1.88	0.56
4:B:508:EDO:H21	5:B:866:HOH:O	2.05	0.55
1:D:10:SER:OG	1:D:12:HIS:HD2	1.89	0.55
1:B:92:ILE:HG22	1:B:96:LYS:HE3	1.88	0.55
1:B:179:HIS:NE2	1:B:409[A]:ASN:OD1	2.39	0.55
1:C:417[A]:ILE:HG21	4:C:508:EDO:H11	1.89	0.55
1:D:153:MET:CE	1:D:201:LEU:HD23	2.37	0.55
1:A:51:ASN:ND2	5:A:611:HOH:O	2.39	0.54
1:A:91:ILE:O	1:A:95:LEU:HG	2.06	0.54
1:A:396:LYS:HB3	1:A:400:LYS:HZ3	1.73	0.54
1:A:251:ASN:HD21	1:A:253:TYR:HD2	1.55	0.53
1:B:20:LEU:HD11	1:B:326:PHE:HD1	1.73	0.53
1:D:246:ILE:HB	1:D:266[B]:VAL:CG1	2.39	0.53
1:B:402:TYR:CE1	4:B:502:EDO:H21	2.43	0.53
1:D:124:LYS:H	1:D:268:ASN:HD22	1.56	0.53
1:A:68:ARG:NH2	1:D:54:LYS:HZ3	2.01	0.52
1:B:41[A]:HIS:HB3	1:B:315:LYS:HB3	1.92	0.52
1:C:163:ASN:HD21	1:D:292:LYS:NZ	2.08	0.52
1:A:365:GLY:HA2	4:A:505:EDO:H12	1.91	0.51
1:C:10:SER:OG	1:C:12:HIS:HD2	1.93	0.51
1:C:98:VAL:HG22	1:C:109:PHE:CE1	2.45	0.51
1:D:22:CYS:HA	1:D:27:THR:HG22	1.93	0.51
1:A:281:TYR:OH	1:A:289:LYS:HE3	2.11	0.51
1:B:281:TYR:OH	1:B:289:LYS:HE3	2.11	0.51
1:A:179:HIS:CE1	1:A:409[A]:ASN:HD21	2.29	0.51
1:B:5:THR:HG22	1:B:29:GLU:O	2.12	0.50
1:C:12:HIS:CE1	4:C:509:EDO:H11	2.45	0.50
1:C:41[A]:HIS:HB3	1:C:315:LYS:HB3	1.93	0.50
1:B:386:TYR:O	1:B:414:HIS:HE1	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:ARG:NH2	1:A:432:ASN:HD21	2.09	0.49
1:D:124:LYS:HD3	1:D:125:TYR:CZ	2.48	0.49
1:C:385:TYR:CZ	4:C:504:EDO:H21	2.47	0.49
1:C:325:VAL:HG21	5:C:719:HOH:O	2.12	0.49
1:D:386:TYR:O	1:D:414:HIS:HE1	1.96	0.49
1:D:211:ASP:OD2	1:D:214:ARG:NE	2.45	0.48
4:A:508:EDO:C1	5:A:607:HOH:O	2.60	0.48
1:C:179:HIS:CE1	1:C:409:ASN:OD1	2.67	0.48
4:C:508:EDO:H21	5:C:678:HOH:O	2.13	0.48
1:A:115:GLU:HA	4:A:511:EDO:H21	1.94	0.48
1:A:414:HIS:HB3	4:A:509:EDO:H11	1.96	0.48
1:D:124:LYS:H	1:D:268:ASN:ND2	2.12	0.48
2:A:501:O1G:H12	2:A:501:O1G:O4	2.14	0.47
1:C:91:ILE:O	1:C:95:LEU:HG	2.14	0.47
1:A:178:ASP:OD2	1:A:181:SER:OG	2.24	0.47
1:B:55:ILE:HD12	1:B:56:ASP:N	2.30	0.47
1:D:12:HIS:HE1	1:D:120:TYR:O	1.96	0.47
1:C:9:ILE:HD11	1:C:32:PHE:CD1	2.49	0.47
1:B:111:LYS:O	1:B:115:GLU:HG3	2.15	0.47
1:A:179:HIS:N	1:A:179:HIS:ND1	2.63	0.47
1:B:344:ILE:HA	1:B:361:ARG:O	2.15	0.47
1:A:396:LYS:HB3	1:A:400:LYS:NZ	2.29	0.47
4:A:505:EDO:O2	5:A:601:HOH:O	2.20	0.47
1:C:212:MET:HE1	1:C:226:ILE:HG21	1.96	0.46
1:B:124:LYS:O	4:B:506:EDO:H22	2.15	0.46
1:C:111:LYS:HA	1:C:111:LYS:HD2	1.52	0.46
1:D:281:TYR:OH	1:D:289:LYS:HE3	2.15	0.46
1:A:396:LYS:HE3	1:A:400:LYS:HZ1	1.81	0.46
1:D:321:ASN:O	1:D:325:VAL:HG23	2.16	0.46
2:C:501:O1G:O4	2:C:501:O1G:H12	2.16	0.46
1:D:28:LYS:HG2	1:D:29:GLU:OE2	2.16	0.46
1:D:156:ASN:HB2	4:D:511:EDO:H11	1.96	0.46
1:A:277:ASN:OD1	4:A:506:EDO:H12	2.16	0.45
2:C:501:O1G:N3'	2:C:501:O1G:O3	2.50	0.45
1:A:104:PHE:HB2	1:B:252:PRO:HG2	1.98	0.45
1:B:325[A]:VAL:HG11	5:B:695:HOH:O	2.15	0.45
1:A:55:ILE:O	1:D:69:LYS:HD3	2.16	0.45
1:C:329:GLN:OE1	4:C:510:EDO:H21	2.17	0.45
1:C:240:GLY:HA2	1:C:246:ILE:HD11	1.98	0.44
1:B:402:TYR:HE1	4:B:502:EDO:H21	1.83	0.44
4:A:508:EDO:H12	5:A:607:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:508:EDO:H12	5:B:756:HOH:O	2.18	0.44
1:B:397:LEU:HD23	4:B:513:EDO:C2	2.48	0.44
1:B:102:GLY:HA2	1:B:104:PHE:CZ	2.53	0.43
1:B:232:GLN:HG3	2:B:501:O1G:H7	1.99	0.43
1:D:92:ILE:HG22	1:D:96:LYS:HE3	2.00	0.43
1:C:326:PHE:CE2	1:C:330:ARG:HD2	2.54	0.43
1:B:55:ILE:H	1:B:55:ILE:HG13	1.66	0.43
1:D:83:ILE:O	5:D:601:HOH:O	2.21	0.43
1:D:285:GLU:HB3	1:D:288:LYS:HB2	2.00	0.43
1:A:416:GLU:HG2	4:A:505:EDO:H11	2.00	0.43
1:B:385:TYR:HA	1:B:386:TYR:HA	1.85	0.43
1:B:388:VAL:HA	4:B:504:EDO:C1	2.40	0.43
1:B:34:ILE:HG21	1:B:44:LYS:HG3	2.01	0.43
1:A:111:LYS:HZ1	1:A:115:GLU:HG3	1.82	0.42
1:B:44:LYS:HE2	5:B:854:HOH:O	2.19	0.42
1:D:34:ILE:H	1:D:51:ASN:ND2	2.11	0.42
1:B:213:LYS:HG3	1:B:214:ARG:N	2.34	0.42
1:B:337:GLU:HG2	5:B:801:HOH:O	2.18	0.42
1:C:49:ARG:HG2	1:C:50:GLU:HG3	2.02	0.42
1:D:385:TYR:HA	1:D:386:TYR:HA	1.89	0.42
1:D:256:PHE:HB2	1:D:317:LEU:HD13	2.00	0.42
1:C:349[A]:MET:HE3	1:C:355:TRP:HZ2	1.84	0.42
1:A:385:TYR:HA	1:A:386:TYR:HA	1.88	0.42
1:D:102:GLY:HA2	1:D:104:PHE:CZ	2.55	0.42
1:A:196:LYS:HG2	5:A:888:HOH:O	2.19	0.42
1:C:385:TYR:CE2	4:C:504:EDO:H21	2.55	0.42
1:D:206:TYR:H	1:D:232:GLN:HE21	1.65	0.42
1:B:103:GLN:HA	4:B:503:EDO:H11	2.02	0.42
1:B:179:HIS:CE1	1:B:409[A]:ASN:HD21	2.37	0.42
1:A:338:LEU:HD11	1:A:439:ILE:HD13	2.02	0.41
1:B:27:THR:HG23	1:B:29:GLU:HG2	2.02	0.41
1:C:348:ARG:HB2	1:C:348:ARG:HH21	1.85	0.41
1:A:316:PHE:HB2	4:A:515:EDO:H12	2.03	0.41
1:A:414:HIS:ND1	4:A:509:EDO:H21	2.36	0.41
1:D:23:LEU:HB3	1:D:59:ILE:HD11	2.03	0.41
1:B:342:GLU:O	1:B:363:ILE:HD11	2.21	0.40
1:C:277:ASN:OD1	1:C:301:LYS:NZ	2.43	0.40
1:A:66:TYR:CE1	1:A:429:GLU:HG3	2.56	0.40
1:A:68:ARG:HH22	1:D:54:LYS:HZ2	1.60	0.40
1:C:386:TYR:O	1:C:414:HIS:HE1	2.04	0.40
1:D:34:ILE:HG23	1:D:44:LYS:HE3	2.03	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	440/445 (99%)	429 (98%)	10 (2%)	1 (0%)	47	26
1	B	442/445 (99%)	425 (96%)	16 (4%)	1 (0%)	47	26
1	C	442/445 (99%)	431 (98%)	10 (2%)	1 (0%)	47	26
1	D	442/445 (99%)	431 (98%)	10 (2%)	1 (0%)	47	26
All	All	1766/1780 (99%)	1716 (97%)	46 (3%)	4 (0%)	47	26

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	VAL
1	B	258	VAL
1	C	258	VAL
1	D	258	VAL

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	400/401 (100%)	398 (100%)	2 (0%)	88	80
1	B	402/401 (100%)	397 (99%)	5 (1%)	71	54
1	C	402/401 (100%)	399 (99%)	3 (1%)	84	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	402/401 (100%)	400 (100%)	2 (0%)	88	80
All	All	1606/1604 (100%)	1594 (99%)	12 (1%)	81	73

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LYS
1	A	179	HIS
1	B	112	LYS
1	B	251	ASN
1	B	305	LEU
1	B	330	ARG
1	B	405	ASP
1	C	15	ASP
1	C	251	ASN
1	C	379	ASN
1	D	28	LYS
1	D	305	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	251	ASN
1	A	278	GLN
1	A	321	ASN
1	A	432	ASN
1	B	103	GLN
1	B	204	HIS
1	C	12	HIS
1	C	122	ASN
1	C	163	ASN
1	D	12	HIS
1	D	51	ASN
1	D	70	ASN
1	D	232	GLN
1	D	268	ASN
1	D	299	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 57 ligands modelled in this entry, 2 are monoatomic - leaving 55 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	EDO	D	502	-	3,3,3	0.27	0	2,2,2	0.62	0
4	EDO	C	512	-	3,3,3	0.39	0	2,2,2	0.27	0
4	EDO	D	509	-	3,3,3	0.39	0	2,2,2	0.29	0
2	O1G	A	501	-	32,32,32	1.88	9 (28%)	43,48,48	1.74	10 (23%)
4	EDO	A	503	-	3,3,3	0.34	0	2,2,2	0.30	0
4	EDO	A	517	-	3,3,3	0.37	0	2,2,2	0.46	0
4	EDO	A	504	-	3,3,3	0.57	0	2,2,2	0.30	0
4	EDO	A	515	-	3,3,3	0.52	0	2,2,2	0.23	0
4	EDO	D	505	-	3,3,3	0.26	0	2,2,2	0.75	0
4	EDO	A	506	-	3,3,3	0.60	0	2,2,2	0.29	0
4	EDO	B	514	-	3,3,3	0.38	0	2,2,2	0.81	0
4	EDO	B	515	-	3,3,3	0.39	0	2,2,2	0.92	0
4	EDO	D	503	-	3,3,3	0.71	0	2,2,2	0.66	0
4	EDO	D	504	-	3,3,3	0.11	0	2,2,2	1.31	0
4	EDO	B	505	-	3,3,3	0.35	0	2,2,2	0.56	0
4	EDO	A	513	-	3,3,3	0.27	0	2,2,2	0.50	0
4	EDO	A	514	-	3,3,3	0.55	0	2,2,2	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	A	518	-	3,3,3	0.36	0	2,2,2	0.15	0
4	EDO	A	509	-	3,3,3	0.63	0	2,2,2	0.45	0
4	EDO	C	513	-	3,3,3	0.25	0	2,2,2	0.84	0
4	EDO	B	503	-	3,3,3	0.24	0	2,2,2	0.77	0
4	EDO	B	513	-	3,3,3	0.35	0	2,2,2	0.97	0
4	EDO	C	505	-	3,3,3	0.22	0	2,2,2	0.83	0
4	EDO	A	505	-	3,3,3	0.39	0	2,2,2	0.62	0
4	EDO	C	506	-	3,3,3	0.62	0	2,2,2	0.39	0
2	O1G	C	501	-	32,32,32	1.76	7 (21%)	43,48,48	2.06	15 (34%)
4	EDO	C	508	-	3,3,3	0.46	0	2,2,2	1.01	0
4	EDO	C	511	-	3,3,3	0.47	0	2,2,2	1.18	0
4	EDO	A	507	-	3,3,3	0.77	0	2,2,2	1.06	0
4	EDO	A	510	-	3,3,3	0.57	0	2,2,2	0.54	0
4	EDO	B	509	-	3,3,3	0.72	0	2,2,2	0.60	0
4	EDO	C	509	-	3,3,3	0.48	0	2,2,2	0.48	0
4	EDO	B	502	-	3,3,3	0.58	0	2,2,2	0.26	0
4	EDO	D	507	-	3,3,3	0.30	0	2,2,2	0.68	0
4	EDO	B	512	-	3,3,3	0.37	0	2,2,2	0.21	0
4	EDO	D	510	-	3,3,3	0.43	0	2,2,2	0.87	0
4	EDO	A	508	-	3,3,3	0.76	0	2,2,2	0.46	0
2	O1G	D	501	-	32,32,32	1.82	9 (28%)	43,48,48	1.94	15 (34%)
4	EDO	B	511	-	3,3,3	0.17	0	2,2,2	0.81	0
4	EDO	A	512	-	3,3,3	0.47	0	2,2,2	0.27	0
4	EDO	C	510	-	3,3,3	0.44	0	2,2,2	0.18	0
4	EDO	B	510	-	3,3,3	0.50	0	2,2,2	0.71	0
4	EDO	C	504	-	3,3,3	0.34	0	2,2,2	0.47	0
4	EDO	D	506	-	3,3,3	0.41	0	2,2,2	0.65	0
4	EDO	B	504	-	3,3,3	0.89	0	2,2,2	1.23	0
4	EDO	D	508	-	3,3,3	0.39	0	2,2,2	0.19	0
4	EDO	D	511	-	3,3,3	0.54	0	2,2,2	0.64	0
4	EDO	A	511	-	3,3,3	0.30	0	2,2,2	0.42	0
4	EDO	B	508	-	3,3,3	0.32	0	2,2,2	0.34	0
2	O1G	B	501	-	32,32,32	1.74	7 (21%)	43,48,48	1.86	14 (32%)
4	EDO	A	516	-	3,3,3	0.32	0	2,2,2	0.16	0
4	EDO	B	506	-	3,3,3	0.24	0	2,2,2	0.36	0
4	EDO	B	507	-	3,3,3	0.20	0	2,2,2	0.87	0
4	EDO	C	507	-	3,3,3	0.33	0	2,2,2	0.97	0
4	EDO	C	503	-	3,3,3	0.38	0	2,2,2	0.52	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	D	502	-	-	1/1/1/1	-
4	EDO	C	512	-	-	0/1/1/1	-
4	EDO	D	509	-	-	0/1/1/1	-
2	O1G	A	501	-	-	3/17/37/37	0/2/2/2
4	EDO	A	503	-	-	0/1/1/1	-
4	EDO	A	517	-	-	0/1/1/1	-
4	EDO	A	504	-	-	0/1/1/1	-
4	EDO	A	515	-	-	0/1/1/1	-
4	EDO	D	505	-	-	1/1/1/1	-
4	EDO	A	506	-	-	1/1/1/1	-
4	EDO	B	514	-	-	0/1/1/1	-
4	EDO	B	515	-	-	1/1/1/1	-
4	EDO	D	503	-	-	1/1/1/1	-
4	EDO	D	504	-	-	1/1/1/1	-
4	EDO	B	505	-	-	0/1/1/1	-
4	EDO	A	513	-	-	0/1/1/1	-
4	EDO	A	514	-	-	1/1/1/1	-
4	EDO	A	518	-	-	1/1/1/1	-
4	EDO	A	509	-	-	1/1/1/1	-
4	EDO	C	513	-	-	0/1/1/1	-
4	EDO	B	503	-	-	1/1/1/1	-
4	EDO	B	513	-	-	1/1/1/1	-
4	EDO	C	505	-	-	1/1/1/1	-
4	EDO	A	505	-	-	1/1/1/1	-
4	EDO	C	506	-	-	0/1/1/1	-
2	O1G	C	501	-	-	4/17/37/37	0/2/2/2
4	EDO	C	508	-	-	0/1/1/1	-
4	EDO	C	511	-	-	0/1/1/1	-
4	EDO	A	507	-	-	0/1/1/1	-
4	EDO	A	510	-	-	1/1/1/1	-
4	EDO	B	509	-	-	0/1/1/1	-
4	EDO	C	509	-	-	0/1/1/1	-
4	EDO	B	502	-	-	1/1/1/1	-
4	EDO	D	507	-	-	1/1/1/1	-
4	EDO	B	512	-	-	1/1/1/1	-
4	EDO	D	510	-	-	0/1/1/1	-
4	EDO	A	508	-	-	0/1/1/1	-
2	O1G	D	501	-	-	4/17/37/37	0/2/2/2
4	EDO	B	511	-	-	1/1/1/1	-
4	EDO	A	512	-	-	1/1/1/1	-
4	EDO	C	510	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	510	-	-	0/1/1/1	-
4	EDO	C	504	-	-	1/1/1/1	-
4	EDO	D	506	-	-	1/1/1/1	-
4	EDO	B	504	-	-	1/1/1/1	-
4	EDO	D	508	-	-	0/1/1/1	-
4	EDO	D	511	-	-	1/1/1/1	-
4	EDO	A	511	-	-	0/1/1/1	-
4	EDO	B	508	-	-	0/1/1/1	-
2	O1G	B	501	-	-	1/17/37/37	0/2/2/2
4	EDO	A	516	-	-	0/1/1/1	-
4	EDO	B	506	-	-	0/1/1/1	-
4	EDO	B	507	-	-	0/1/1/1	-
4	EDO	C	507	-	-	0/1/1/1	-
4	EDO	C	503	-	-	0/1/1/1	-

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	O1G	C4A-C4B	5.44	1.57	1.46
2	C	501	O1G	C4A-C4B	5.09	1.56	1.46
2	D	501	O1G	C4A-C4B	4.89	1.55	1.46
2	A	501	O1G	C4A-C4B	4.78	1.55	1.46
2	A	501	O1G	C4B-N3'	4.31	1.35	1.27
2	D	501	O1G	C2A-C9A	4.15	1.57	1.50
2	C	501	O1G	C4B-N3'	3.51	1.33	1.27
2	B	501	O1G	C4B-N3'	3.44	1.33	1.27
2	A	501	O1G	C3-N3'	3.26	1.50	1.46
2	D	501	O1G	O5-C1	3.23	1.51	1.42
2	A	501	O1G	O5-C1	3.21	1.50	1.42
2	C	501	O1G	P'-O6	3.03	1.70	1.60
2	D	501	O1G	C4B-N3'	2.94	1.32	1.27
2	B	501	O1G	C2-C3	-2.93	1.48	1.53
2	B	501	O1G	O5-C1	2.82	1.49	1.42
2	C	501	O1G	C2-C3	-2.71	1.48	1.53
2	C	501	O1G	C3-N3'	2.67	1.50	1.46
2	A	501	O1G	C4A-C3A	2.61	1.44	1.40
2	D	501	O1G	C3-N3'	2.60	1.50	1.46
2	C	501	O1G	O5-C1	2.58	1.49	1.42
2	B	501	O1G	O2-C2	2.56	1.49	1.43
2	A	501	O1G	C2-C3	-2.42	1.49	1.53
2	A	501	O1G	C2A-C9A	2.42	1.54	1.50
2	B	501	O1G	C2A-C9A	2.39	1.54	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	O1G	P'-O6	2.37	1.67	1.60
2	D	501	O1G	P-OP4	2.30	1.67	1.60
2	C	501	O1G	O5-C5	2.20	1.49	1.44
2	D	501	O1G	P'-O6	2.16	1.67	1.60
2	D	501	O1G	C9A-N1'	-2.15	1.29	1.33
2	D	501	O1G	C2-C3	-2.13	1.49	1.53
2	A	501	O1G	O5-C5	2.11	1.49	1.44
2	B	501	O1G	O5-C5	2.00	1.49	1.44

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	O1G	C3-N3'-C4B	5.31	127.40	117.99
2	C	501	O1G	C4A-C3A-C9A	-5.03	117.07	120.19
2	C	501	O1G	C4A-C4B-N3'	-4.88	112.18	123.01
2	A	501	O1G	C4A-C4B-N3'	-4.77	112.40	123.01
2	D	501	O1G	C4A-C4B-N3'	-4.52	112.97	123.01
2	C	501	O1G	C5A-C6A-N1'	-4.07	117.03	123.82
2	D	501	O1G	C3-N3'-C4B	3.84	124.81	117.99
2	A	501	O1G	OP6-P'-O6	-3.82	96.57	106.73
2	D	501	O1G	O4-C4-C5	-3.81	99.84	109.30
2	C	501	O1G	C3A-C4A-C5A	3.69	121.09	118.26
2	D	501	O1G	C1-O5-C5	-3.65	106.78	113.66
2	A	501	O1G	C3-N3'-C4B	3.53	124.24	117.99
2	D	501	O1G	C5A-C6A-N1'	-3.43	118.11	123.82
2	B	501	O1G	C4A-C4B-N3'	-3.36	115.53	123.01
2	C	501	O1G	O6-C6-C5	-3.36	97.43	108.99
2	B	501	O1G	O1-C1-C2	3.20	118.03	109.03
2	A	501	O1G	OP6-P'-OP8	3.16	123.06	110.68
2	C	501	O1G	C3-N3'-C4B	3.09	123.48	117.99
2	C	501	O1G	C6A-N1'-C9A	3.08	124.87	119.17
2	B	501	O1G	C5A-C6A-N1'	-3.05	118.75	123.82
2	D	501	O1G	C3A-C4A-C5A	-2.95	116.00	118.26
2	D	501	O1G	C5T-C5A-C6A	-2.93	114.55	119.37
2	A	501	O1G	O6-C6-C5	-2.88	99.09	108.99
2	D	501	O1G	C5A-C4A-C4B	2.84	126.23	121.56
2	B	501	O1G	OP7-P'-OP6	2.78	118.26	107.64
2	A	501	O1G	O1-C1-C2	2.75	116.77	109.03
2	B	501	O1G	O5-C5-C6	-2.67	101.27	106.67
2	A	501	O1G	OP7-P'-O6	2.62	113.70	106.73
2	B	501	O1G	O2-C2-C1	2.53	115.03	109.16
2	D	501	O1G	C6-C5-C4	-2.51	106.85	112.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	O1G	O4-C4-C3	-2.48	103.46	109.14
2	B	501	O1G	C6A-N1'-C9A	2.42	123.64	119.17
2	C	501	O1G	OP6-P'-OP8	2.39	120.05	110.68
2	A	501	O1G	C2A-C9A-N1'	2.33	122.23	117.67
2	C	501	O1G	OP6-P'-O6	-2.33	100.53	106.73
2	B	501	O1G	OP7-P'-O6	-2.31	100.58	106.73
2	D	501	O1G	C6A-N1'-C9A	2.27	123.37	119.17
2	C	501	O1G	O3-C3A-C9A	2.26	122.41	117.49
2	C	501	O1G	OP1-P-OP4	2.25	112.72	106.73
2	B	501	O1G	O6-C6-C5	-2.23	101.31	108.99
2	A	501	O1G	C2A-C9A-C3A	-2.21	118.16	120.89
2	D	501	O1G	O6-P'-OP8	2.20	112.66	106.47
2	C	501	O1G	O6-P'-OP8	-2.20	100.30	106.47
2	D	501	O1G	O1-C1-C2	2.19	115.21	109.03
2	D	501	O1G	C6A-C5A-C4A	2.18	122.17	118.15
2	D	501	O1G	OP7-P'-OP6	2.18	115.98	107.64
2	C	501	O1G	OP2-P-OP4	-2.15	101.02	106.73
2	C	501	O1G	O1-C1-C2	2.13	115.04	109.03
2	C	501	O1G	O4-C4-C3	-2.12	104.30	109.14
2	A	501	O1G	O1-C1-O5	-2.08	104.15	110.38
2	D	501	O1G	O5-C1-C2	-2.05	106.62	110.28
2	B	501	O1G	OP6-P'-O6	-2.04	101.29	106.73
2	B	501	O1G	C6-C5-C4	-2.04	107.84	112.09
2	B	501	O1G	O4-C4-C5	-2.03	104.25	109.30

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	O1G	C4-C3-N3'-C4B
2	B	501	O1G	C4-C3-N3'-C4B
2	C	501	O1G	C5T-OP4-P-OP3
2	C	501	O1G	C4-C3-N3'-C4B
2	C	501	O1G	C4-C5-C6-O6
2	D	501	O1G	C2-C3-N3'-C4B
2	D	501	O1G	C4-C3-N3'-C4B
4	A	505	EDO	O1-C1-C2-O2
4	A	509	EDO	O1-C1-C2-O2
4	B	503	EDO	O1-C1-C2-O2
4	B	504	EDO	O1-C1-C2-O2
4	B	511	EDO	O1-C1-C2-O2
4	B	512	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
4	C	505	EDO	O1-C1-C2-O2
4	D	502	EDO	O1-C1-C2-O2
4	D	505	EDO	O1-C1-C2-O2
4	D	506	EDO	O1-C1-C2-O2
4	D	507	EDO	O1-C1-C2-O2
2	A	501	O1G	C5T-OP4-P-OP3
2	A	501	O1G	C4-C5-C6-O6
4	A	506	EDO	O1-C1-C2-O2
4	B	502	EDO	O1-C1-C2-O2
4	C	504	EDO	O1-C1-C2-O2
4	B	515	EDO	O1-C1-C2-O2
4	D	504	EDO	O1-C1-C2-O2
4	C	510	EDO	O1-C1-C2-O2
4	D	511	EDO	O1-C1-C2-O2
2	D	501	O1G	C5T-OP4-P-OP1
2	D	501	O1G	C3A-C4A-C4B-N3'
2	C	501	O1G	C2-C3-N3'-C4B
4	A	510	EDO	O1-C1-C2-O2
4	B	513	EDO	O1-C1-C2-O2
4	D	503	EDO	O1-C1-C2-O2
4	A	512	EDO	O1-C1-C2-O2
4	A	514	EDO	O1-C1-C2-O2
4	A	518	EDO	O1-C1-C2-O2

There are no ring outliers.

23 monomers are involved in 37 short contacts:

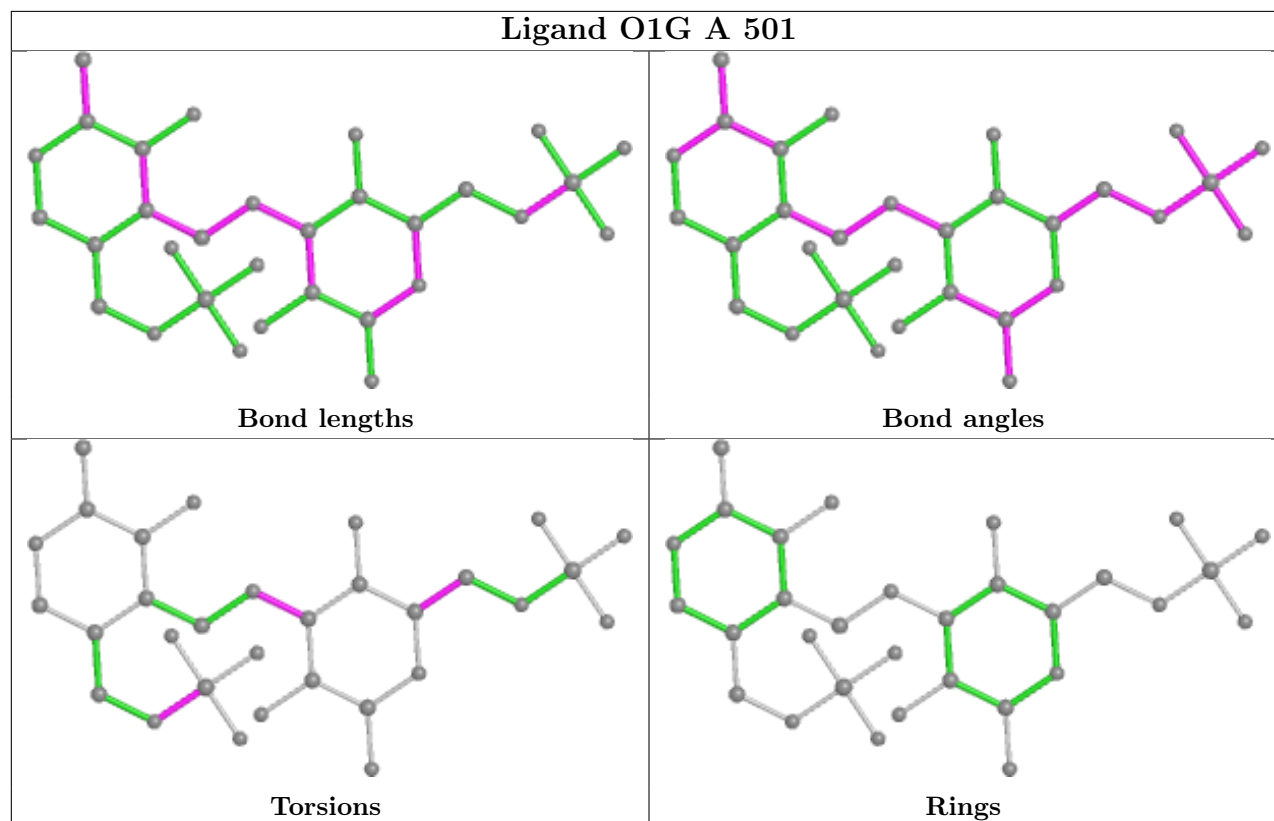
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	502	EDO	1	0
2	A	501	O1G	1	0
4	A	515	EDO	1	0
4	D	505	EDO	1	0
4	A	506	EDO	1	0
4	A	509	EDO	2	0
4	B	503	EDO	1	0
4	B	513	EDO	1	0
4	A	505	EDO	3	0
2	C	501	O1G	2	0
4	C	508	EDO	2	0
4	C	509	EDO	1	0
4	B	502	EDO	2	0
4	A	508	EDO	3	0

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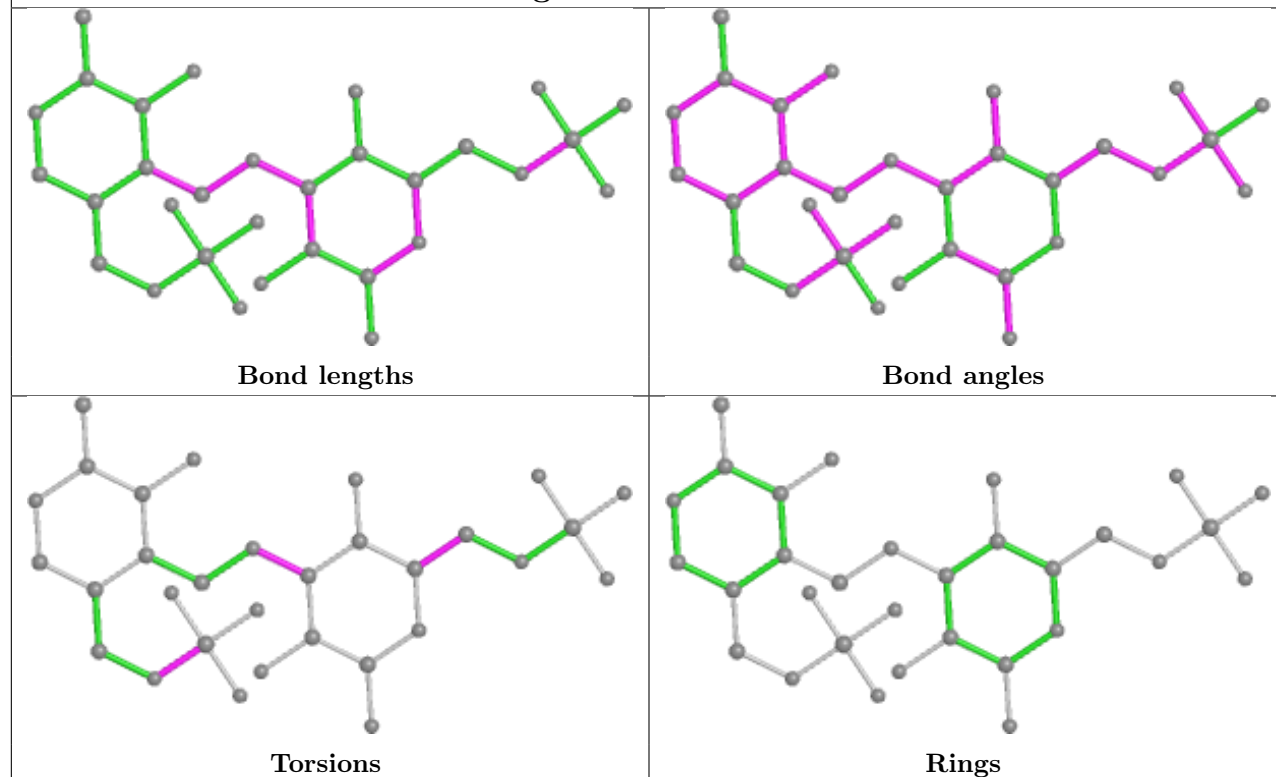
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	510	EDO	3	0
4	C	504	EDO	2	0
4	B	504	EDO	3	0
4	D	511	EDO	1	0
4	A	511	EDO	1	0
4	B	508	EDO	2	0
2	B	501	O1G	1	0
4	B	506	EDO	1	0
4	C	507	EDO	1	0

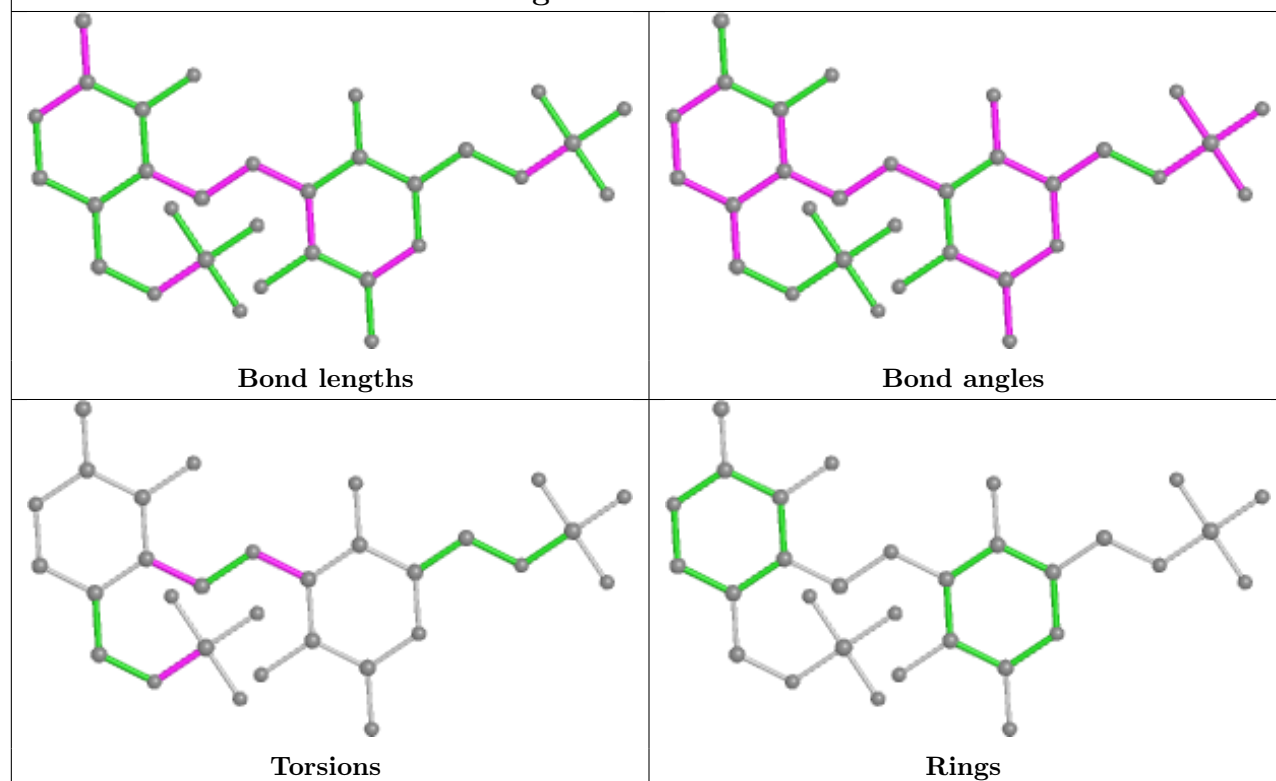
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

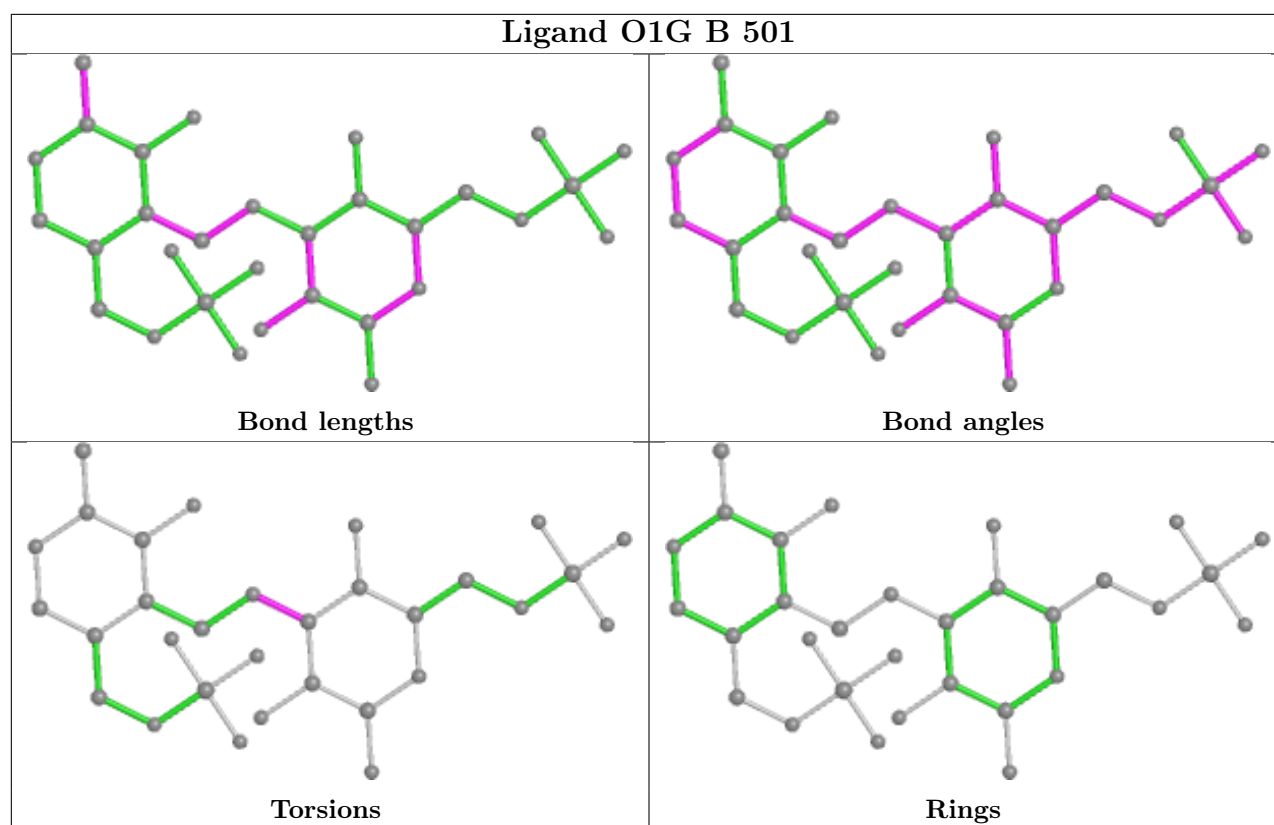


Ligand O1G C 501



Ligand O1G D 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, 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	439/445 (98%)	-0.20	7 (1%) 72 71	11, 20, 37, 95	1 (0%)
1	B	439/445 (98%)	-0.17	9 (2%) 63 62	12, 22, 41, 79	0
1	C	439/445 (98%)	-0.22	7 (1%) 72 71	11, 20, 36, 79	1 (0%)
1	D	439/445 (98%)	-0.07	13 (2%) 50 48	12, 24, 44, 71	1 (0%)
All	All	1756/1780 (98%)	-0.16	36 (2%) 63 62	11, 21, 40, 95	3 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	29	GLU	4.8
1	D	179	HIS	4.5
1	A	29	GLU	4.4
1	D	56	ASP	3.9
1	D	187	LEU	3.8
1	A	179	HIS	3.8
1	D	27	THR	3.5
1	D	222	TYR	3.3
1	B	179	HIS	3.3
1	D	221	VAL	3.2
1	C	28	LYS	3.2
1	A	27	THR	3.1
1	B	29	GLU	3.1
1	B	404[A]	GLN	3.0
1	D	214	ARG	2.9
1	C	5	THR	2.7
1	A	28	LYS	2.7
1	D	404	GLN	2.6
1	B	28	LYS	2.5
1	D	408	LEU	2.5
1	B	56	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	55	ILE	2.4
1	C	179	HIS	2.3
1	B	27	THR	2.3
1	D	28	LYS	2.3
1	A	180	LYS	2.3
1	A	404	GLN	2.2
1	C	27	THR	2.2
1	B	26	GLU	2.2
1	C	364	ASN	2.2
1	D	5	THR	2.2
1	A	5	THR	2.2
1	C	221	VAL	2.2
1	B	187	LEU	2.0
1	D	191	GLU	2.0
1	B	408	LEU	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	EDO	D	508	4/4	0.58	0.28	48,48,49,49	0
4	EDO	A	518	4/4	0.62	0.11	51,51,51,52	0
4	EDO	B	512	4/4	0.74	0.14	41,43,44,44	0
4	EDO	D	506	4/4	0.77	0.20	32,35,35,37	0
4	EDO	B	511	4/4	0.77	0.15	43,43,44,47	0
4	EDO	D	507	4/4	0.78	0.13	36,39,42,44	0
4	EDO	C	511	4/4	0.79	0.17	23,35,39,43	0
4	EDO	D	504	4/4	0.80	0.25	32,32,37,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	EDO	B	507	4/4	0.81	0.16	31,32,32,34	0
4	EDO	B	508	4/4	0.82	0.12	36,38,39,42	0
4	EDO	A	503	4/4	0.82	0.18	29,34,35,35	0
4	EDO	A	507	4/4	0.84	0.26	23,30,34,41	0
4	EDO	A	509	4/4	0.84	0.17	28,31,32,36	0
4	EDO	C	510	4/4	0.85	0.18	31,37,41,44	0
4	EDO	C	505	4/4	0.86	0.13	36,41,44,47	0
4	EDO	D	502	4/4	0.87	0.19	46,47,47,48	0
4	EDO	A	504	4/4	0.87	0.17	28,28,29,37	0
4	EDO	C	507	4/4	0.87	0.21	37,41,46,49	0
4	EDO	A	508	4/4	0.87	0.14	28,30,31,32	0
4	EDO	C	504	4/4	0.87	0.11	38,38,38,39	0
4	EDO	B	502	4/4	0.88	0.14	23,27,35,37	0
4	EDO	A	511	4/4	0.89	0.14	39,43,47,52	0
4	EDO	B	506	4/4	0.89	0.14	42,44,45,48	0
4	EDO	A	513	4/4	0.89	0.12	29,31,32,36	0
4	EDO	A	505	4/4	0.89	0.28	38,41,45,46	0
4	EDO	A	512	4/4	0.90	0.11	29,30,35,36	0
4	EDO	A	506	4/4	0.91	0.14	25,29,29,35	0
4	EDO	B	513	4/4	0.91	0.15	19,22,28,34	0
4	EDO	A	515	4/4	0.91	0.10	19,23,25,26	0
3	NA	A	502	1/1	0.92	0.12	46,46,46,46	0
4	EDO	B	509	4/4	0.92	0.09	21,22,24,25	0
4	EDO	B	503	4/4	0.92	0.13	27,34,39,44	0
4	EDO	B	514	4/4	0.93	0.11	26,27,32,34	0
4	EDO	C	508	4/4	0.93	0.11	25,29,30,32	0
4	EDO	A	514	4/4	0.93	0.09	25,26,26,27	0
4	EDO	A	510	4/4	0.93	0.12	21,22,26,27	0
4	EDO	C	513	4/4	0.93	0.09	31,32,32,32	0
4	EDO	B	504	4/4	0.94	0.11	23,26,31,34	0
4	EDO	B	505	4/4	0.94	0.20	40,40,42,42	0
4	EDO	C	509	4/4	0.94	0.09	28,31,37,40	0
3	NA	C	502	1/1	0.94	0.12	45,45,45,45	0
4	EDO	C	506	4/4	0.94	0.08	20,23,23,24	0
4	EDO	C	512	4/4	0.94	0.12	23,26,27,27	0
4	EDO	D	511	4/4	0.94	0.06	23,25,26,27	0
4	EDO	C	503	4/4	0.95	0.07	26,27,27,35	0
4	EDO	D	505	4/4	0.95	0.08	34,39,40,43	0
4	EDO	D	510	4/4	0.95	0.06	26,27,29,33	0
4	EDO	A	517	4/4	0.95	0.09	25,27,28,28	0
4	EDO	D	503	4/4	0.96	0.08	20,21,22,23	0
4	EDO	A	516	4/4	0.96	0.09	25,27,28,28	0

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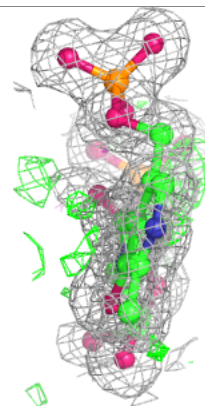
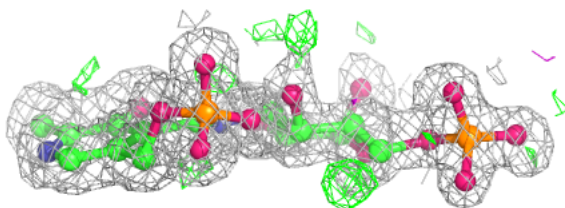
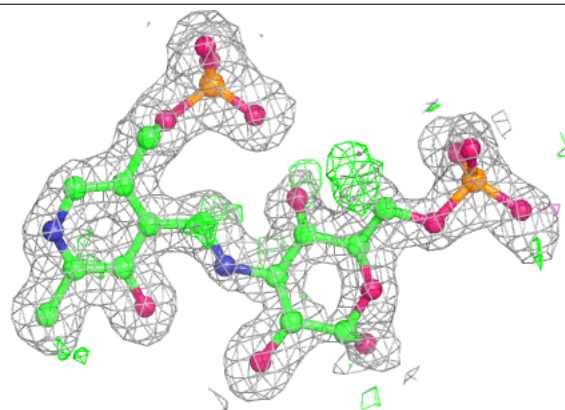
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	EDO	B	510	4/4	0.96	0.08	23,26,28,29	0
4	EDO	B	515	4/4	0.96	0.08	24,25,26,27	0
4	EDO	D	509	4/4	0.97	0.07	22,23,26,27	0
2	O1G	B	501	31/31	0.97	0.11	10,17,22,26	16
2	O1G	C	501	31/31	0.97	0.10	9,17,25,36	13
2	O1G	D	501	31/31	0.98	0.11	12,17,24,31	14
2	O1G	A	501	31/31	0.98	0.11	12,18,24,25	15

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.

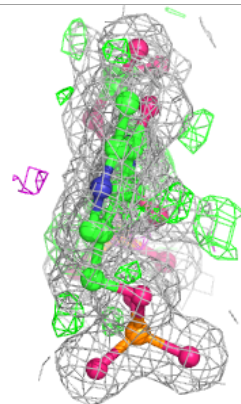
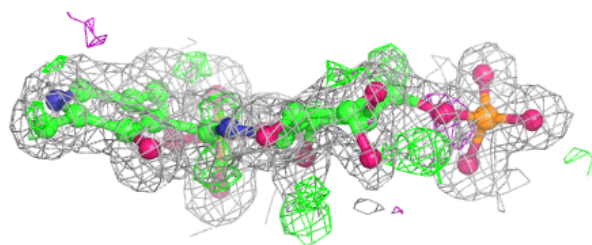
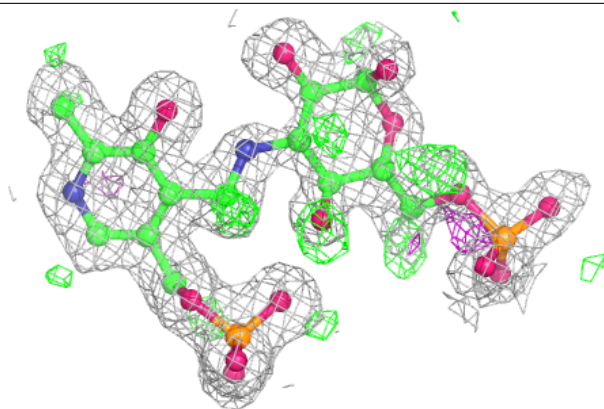
Electron density around O1G B 501:

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

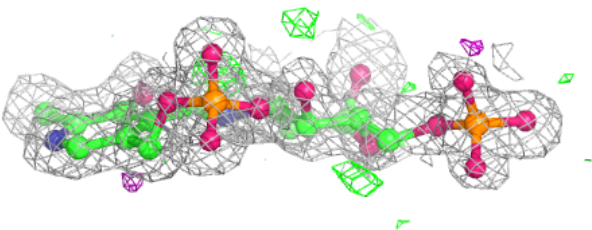
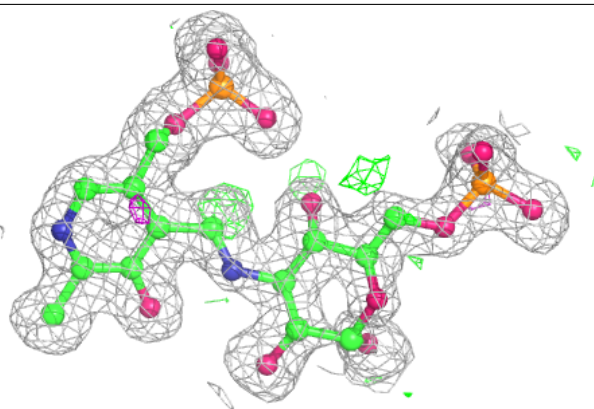


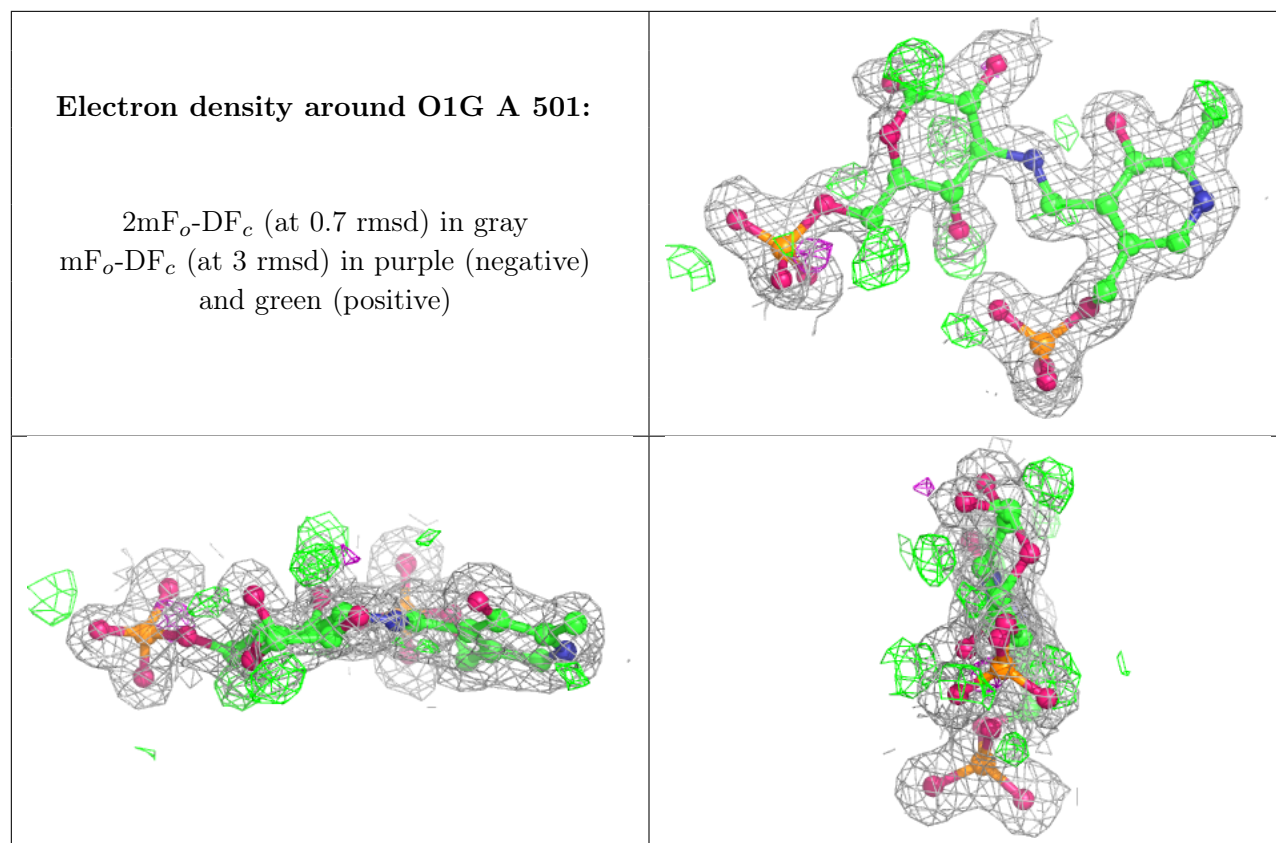
Electron density around O1G C 501:

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

**Electron density around O1G D 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.