



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:47 AM UTC

PDB ID : 7CB8 / pdb_00007cb8
Title : Structure of a FIC-domain protein from Mycobacterium marinum in complex with CDP
Authors : Kumar, S.; Singh, A.; Penmatsa, A.; Surolia, A.
Deposited on : 2020-06-11
Resolution : 2.60 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 types of validation reports are described at

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

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

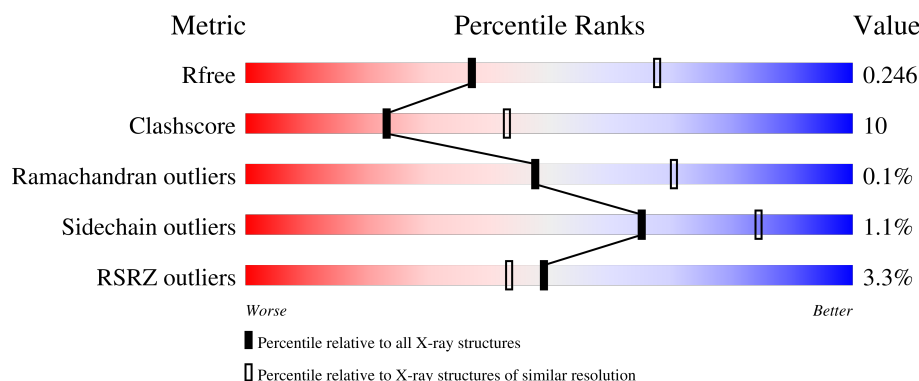
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	412	 2% 75% 20% • •
1	B	412	 4% 77% 17% • 5%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	A	504	-	-	X	-
3	EDO	A	505	-	-	X	-
5	ACT	A	509	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6280 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.

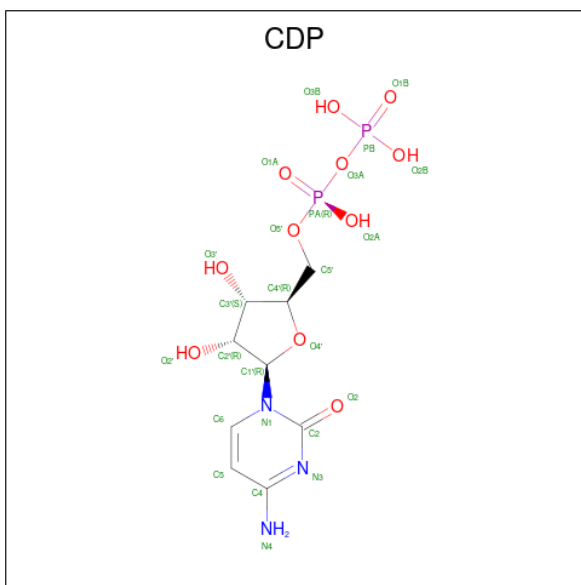
- Molecule 1 is a protein called Fic-domain containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	0	0
			2977	1853	553	564	7			
1	B	393	Total	C	N	O	S	0	0	0
			2942	1831	541	563	7			

There are 16 discrepancies between the modelled and reference sequences:

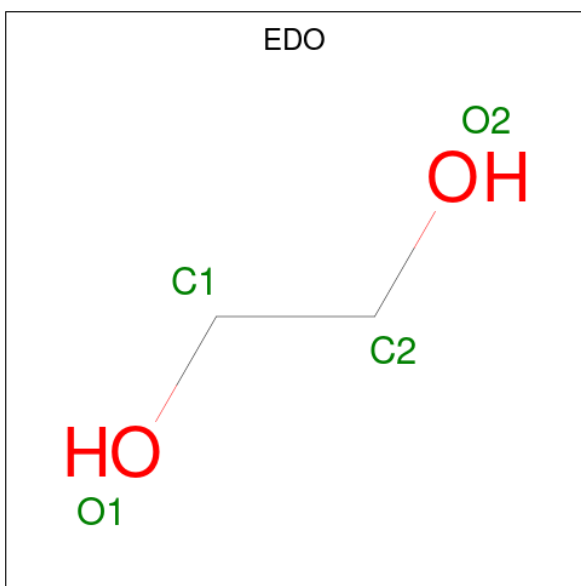
Chain	Residue	Modelled	Actual	Comment	Reference
A	405	LEU	-	expression tag	UNP B2HP08
A	406	GLU	-	expression tag	UNP B2HP08
A	407	HIS	-	expression tag	UNP B2HP08
A	408	HIS	-	expression tag	UNP B2HP08
A	409	HIS	-	expression tag	UNP B2HP08
A	410	HIS	-	expression tag	UNP B2HP08
A	411	HIS	-	expression tag	UNP B2HP08
A	412	HIS	-	expression tag	UNP B2HP08
B	405	LEU	-	expression tag	UNP B2HP08
B	406	GLU	-	expression tag	UNP B2HP08
B	407	HIS	-	expression tag	UNP B2HP08
B	408	HIS	-	expression tag	UNP B2HP08
B	409	HIS	-	expression tag	UNP B2HP08
B	410	HIS	-	expression tag	UNP B2HP08
B	411	HIS	-	expression tag	UNP B2HP08
B	412	HIS	-	expression tag	UNP B2HP08

- Molecule 2 is CYTIDINE-5'-DIPHOSPHATE (CCD ID: CDP) (formula: $\text{C}_9\text{H}_{15}\text{N}_3\text{O}_{11}\text{P}_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 25	C 9	N 3	O 11	P 2	0	0
2	B	1	Total 25	C 9	N 3	O 11	P 2	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



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

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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

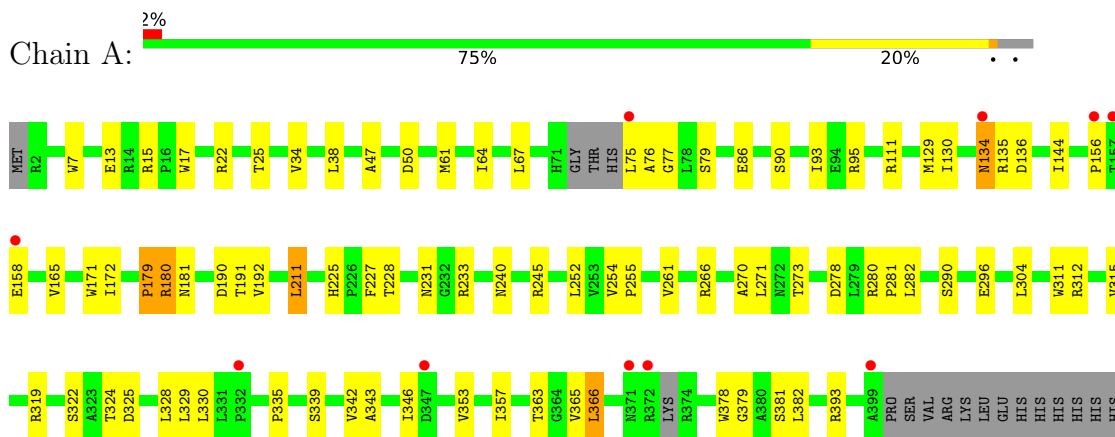
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	118	Total	O	0	0
			118	118		
6	B	122	Total	O	0	0
			122	122		

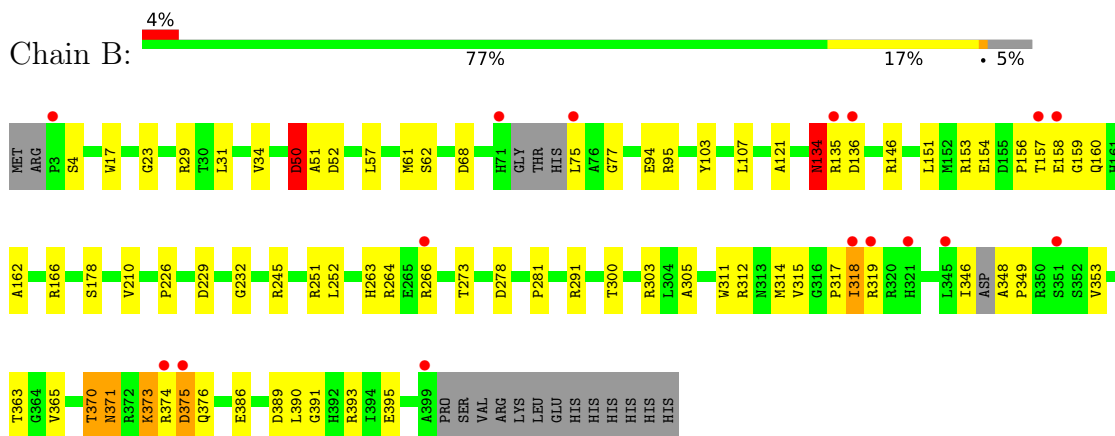
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Fic-domain containing protein



• Molecule 1: Fic-domain containing protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.07Å 107.90Å 110.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.49 – 2.60 48.49 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.3 (48.49-2.60) 97.2 (48.49-2.60)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.61Å)	Xtriage
Refinement program	PHENIX (1.11.1-2575-000)	Depositor
R, R_{free}	0.195 , 0.248 0.196 , 0.246	Depositor DCC
R_{free} test set	1334 reflections (4.51%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.451	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6280	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, EDO, CDP, SO4

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	0.45	1/3031 (0.0%)	0.71	3/4132 (0.1%)
1	B	0.51	1/2994 (0.0%)	0.95	22/4085 (0.5%)
All	All	0.48	2/6025 (0.0%)	0.84	25/8217 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	179	PRO	C-O	-6.86	1.15	1.24
1	B	178	SER	C-O	-5.28	1.18	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	375	ASP	N-CA-C	-17.74	83.52	108.54
1	A	135	ARG	N-CA-C	-13.54	93.84	112.25
1	B	374	ARG	CB-CA-C	12.72	135.74	110.42
1	B	375	ASP	N-CA-CB	9.54	127.76	110.60
1	B	157	THR	N-CA-C	-8.62	102.51	113.72
1	B	135	ARG	N-CA-C	-7.83	100.11	111.30
1	B	318	ILE	CB-CA-C	-7.30	99.32	111.29
1	B	135	ARG	CA-C-N	7.00	132.82	122.74
1	B	135	ARG	C-N-CA	7.00	132.82	122.74
1	B	319	ARG	N-CA-C	-6.91	102.12	111.87
1	B	134	ASN	CA-CB-CG	-6.75	105.86	112.60
1	B	319	ARG	N-CA-CB	6.28	121.19	111.39
1	B	136	ASP	N-CA-C	-6.11	105.13	112.58
1	B	23	GLY	N-CA-C	5.92	121.08	112.55
1	B	50	ASP	CA-C-N	5.83	128.90	120.79
1	B	50	ASP	C-N-CA	5.83	128.90	120.79
1	A	134	ASN	N-CA-CB	5.74	120.20	110.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	ARG	CB-CA-C	-5.74	101.23	109.84
1	B	134	ASN	N-CA-C	5.31	118.95	111.52
1	B	51	ALA	N-CA-C	5.27	117.45	111.02
1	B	156	PRO	CB-CA-C	-5.16	103.04	111.56
1	B	134	ASN	N-CA-CB	5.13	119.71	111.91
1	B	50	ASP	N-CA-CB	5.11	119.13	110.49
1	B	371	ASN	CB-CA-C	-5.07	101.58	109.84
1	B	134	ASN	O-C-N	5.02	128.61	122.58

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	2977	0	2950	71	0
1	B	2942	0	2893	52	0
2	A	25	0	11	1	0
2	B	25	0	11	1	0
3	A	16	0	24	15	0
3	B	20	0	30	6	0
4	A	5	0	0	1	0
4	B	10	0	0	1	0
5	A	12	0	9	3	0
5	B	8	0	6	1	0
6	A	118	0	0	2	0
6	B	122	0	0	4	0
All	All	6280	0	5934	123	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 10.

All (123) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:SER:HB2	1:B:50:ASP:HB2	1.41	1.01
1:A:325:ASP:O	1:A:329:LEU:HD13	1.71	0.89
1:A:342:VAL:HG11	1:A:357:ILE:HD11	1.62	0.81
1:A:130:ILE:HD11	1:A:240:ASN:HB3	1.64	0.79
1:B:4:SER:CB	1:B:50:ASP:HB2	2.12	0.78
1:A:13:GLU:OE2	1:A:15:ARG:HD3	1.84	0.76
1:A:319:ARG:HG3	1:A:322:SER:HB3	1.69	0.74
1:B:75:LEU:HD11	1:B:305:ALA:HB2	1.68	0.74
1:A:111:ARG:HH22	3:B:503:EDO:H22	1.54	0.73
1:B:61:MET:HE1	1:B:210:VAL:HG13	1.70	0.72
1:B:273:THR:HG21	1:B:281:PRO:HG2	1.74	0.70
1:B:264:ARG:H	5:B:509:ACT:H3	1.57	0.69
1:B:346:ILE:HD11	1:B:353:VAL:HG22	1.75	0.69
1:A:134:ASN:HB2	1:A:245:ARG:HA	1.76	0.67
1:A:273:THR:HG21	1:A:281:PRO:HG2	1.77	0.65
1:A:228:THR:HB	3:A:504:EDO:H22	1.77	0.65
1:B:134:ASN:HB2	1:B:245:ARG:HA	1.77	0.65
1:A:381:SER:HB2	3:A:505:EDO:H21	1.79	0.65
1:B:263:HIS:ND1	3:B:506:EDO:H22	2.12	0.64
1:B:348:ALA:N	1:B:349:PRO:HD3	2.11	0.64
1:B:77:GLY:H	3:B:502:EDO:H21	1.61	0.64
1:A:134:ASN:C	1:A:136:ASP:H	2.05	0.63
1:A:315:VAL:HA	1:A:363:THR:HG21	1.80	0.63
1:B:146:ARG:NH1	6:B:602:HOH:O	2.30	0.62
1:B:370:THR:CG2	1:B:373:LYS:HB3	2.31	0.61
1:B:61:MET:CE	1:B:210:VAL:HG13	2.33	0.59
1:A:233:ARG:O	5:A:509:ACT:H1	2.03	0.59
1:A:111:ARG:NH2	3:B:503:EDO:H22	2.17	0.58
1:A:181:ASN:ND2	6:A:601:HOH:O	2.29	0.58
1:A:366:LEU:HG	1:A:378:TRP:HB3	1.84	0.58
1:A:365:VAL:O	3:A:505:EDO:H22	2.04	0.58
1:A:312:ARG:HH22	1:A:329:LEU:CD1	2.17	0.57
1:A:312:ARG:HH22	1:A:329:LEU:HD12	1.71	0.55
1:A:382:LEU:HG	3:A:505:EDO:H12	1.87	0.55
1:B:62:SER:HA	3:B:505:EDO:H22	1.88	0.55
1:A:95:ARG:HG2	1:B:95:ARG:HB3	1.88	0.54
1:A:342:VAL:HG22	1:A:346:ILE:HD13	1.90	0.54
1:B:303:ARG:NH1	1:B:389:ASP:OD2	2.41	0.53
1:A:382:LEU:H	3:A:505:EDO:H21	1.72	0.53
1:A:90:SER:OG	5:A:509:ACT:H3	2.09	0.53
1:A:129:MET:HE3	1:A:144:ILE:HG12	1.91	0.52
1:B:370:THR:HG21	1:B:373:LYS:HB3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:ILE:HD11	1:A:225:HIS:CE1	2.45	0.51
1:A:76:ALA:HA	3:A:502:EDO:O2	2.11	0.51
1:A:339:SER:O	1:A:342:VAL:HG12	2.11	0.51
1:B:121:ALA:HB1	1:B:151:LEU:HA	1.93	0.51
1:B:391:GLY:O	1:B:395:GLU:HG3	2.11	0.50
1:B:17:TRP:HB2	1:B:34:VAL:HG12	1.92	0.50
1:B:153:ARG:HG2	1:B:154:GLU:HG3	1.92	0.50
1:A:93:ILE:HG12	1:A:261:VAL:HG23	1.94	0.50
1:A:134:ASN:HB2	1:A:245:ARG:O	2.12	0.49
1:A:13:GLU:OE1	1:A:38:LEU:HD11	2.12	0.49
1:B:386:GLU:OE2	1:B:393:ARG:NH2	2.46	0.48
1:A:64:ILE:HD11	1:A:290:SER:HB3	1.95	0.48
1:B:291:ARG:NH1	6:B:608:HOH:O	2.47	0.48
1:B:312:ARG:HA	1:B:315:VAL:HG22	1.96	0.48
1:B:346:ILE:HD11	1:B:353:VAL:CG2	2.43	0.47
1:A:77:GLY:CA	3:A:502:EDO:H12	2.44	0.47
1:A:271:LEU:HD23	2:A:501:CDP:C2	2.49	0.47
1:B:315:VAL:HG12	1:B:363:THR:HG21	1.95	0.47
1:A:296:GLU:OE1	1:A:393:ARG:HD2	2.14	0.47
1:A:319:ARG:HG3	1:A:322:SER:CB	2.40	0.47
1:B:373:LYS:C	1:B:375:ASP:H	2.22	0.47
1:A:273:THR:CG2	1:A:278:ASP:HB3	2.44	0.47
1:A:311:TRP:NE1	3:A:505:EDO:O2	2.33	0.47
1:B:158:GLU:C	1:B:160:GLN:H	2.22	0.47
1:B:68:ASP:HA	1:B:252:LEU:HD12	1.97	0.47
1:B:159:GLY:HA2	1:B:162:ALA:HB2	1.97	0.46
1:A:50:ASP:OD2	1:A:280:ARG:NH2	2.47	0.46
1:A:79:SER:HB3	3:A:502:EDO:O1	2.15	0.46
1:A:17:TRP:HB2	1:A:34:VAL:HG12	1.98	0.46
2:B:501:CDP:O1B	6:B:601:HOH:O	2.20	0.46
1:A:61:MET:SD	1:A:211:LEU:HD22	2.56	0.46
1:A:312:ARG:NH2	1:A:329:LEU:HD12	2.31	0.46
1:A:324:THR:O	1:A:328:LEU:HB2	2.16	0.46
1:A:227:PHE:HB3	3:A:504:EDO:H11	1.99	0.45
1:B:103:TYR:CE2	1:B:107:LEU:HD11	2.52	0.45
1:B:370:THR:HG23	1:B:373:LYS:CB	2.47	0.45
1:A:165:VAL:CG2	1:A:192:VAL:HG11	2.47	0.45
1:A:22:ARG:HH22	3:A:503:EDO:H12	1.81	0.44
1:A:67:LEU:HG	1:A:252:LEU:HD11	2.00	0.44
1:A:180:ARG:NH1	4:B:507:SO4:O3	2.45	0.44
1:B:29:ARG:NH1	3:B:503:EDO:O2	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:PRO:C	1:A:158:GLU:H	2.25	0.44
1:A:382:LEU:H	3:A:505:EDO:C2	2.30	0.44
1:A:330:LEU:HD21	1:A:346:ILE:HD11	2.00	0.44
1:A:75:LEU:HD23	1:A:76:ALA:O	2.18	0.44
1:B:166:ARG:HG2	1:B:226:PRO:O	2.18	0.43
1:B:153:ARG:NH2	6:B:605:HOH:O	2.42	0.43
1:B:346:ILE:HG13	1:B:349:PRO:HD2	2.00	0.43
1:A:64:ILE:HG12	1:A:254:VAL:HG21	2.01	0.43
1:B:314:MET:HE3	1:B:365:VAL:HG22	2.00	0.43
1:B:370:THR:HG23	1:B:373:LYS:HB3	2.01	0.43
1:A:342:VAL:CG1	1:A:357:ILE:HD11	2.41	0.43
1:A:228:THR:HB	3:A:504:EDO:C2	2.47	0.43
1:A:335:PRO:O	1:A:379:GLY:HA2	2.19	0.43
1:B:17:TRP:CE3	1:B:31:LEU:HD23	2.54	0.43
1:B:160:GLN:H	1:B:160:GLN:HG2	1.57	0.43
1:A:270:ALA:HB1	1:A:282:LEU:HA	2.01	0.43
1:A:304:LEU:HD23	1:A:304:LEU:HA	1.89	0.42
1:B:311:TRP:O	1:B:315:VAL:HG13	2.19	0.42
1:B:317:PRO:O	1:B:318:ILE:C	2.62	0.42
1:A:86:GLU:OE1	1:A:255:PRO:HA	2.20	0.41
1:A:225:HIS:CD2	1:A:231:ASN:ND2	2.88	0.41
1:A:273:THR:HG23	1:A:278:ASP:HB3	2.02	0.41
1:B:158:GLU:HG2	1:B:229:ASP:OD2	2.20	0.41
1:A:365:VAL:HA	3:A:505:EDO:H11	2.02	0.41
1:A:90:SER:CB	5:A:509:ACT:H3	2.50	0.41
1:A:343:ALA:HA	1:A:353:VAL:HG11	2.03	0.41
1:B:266:ARG:HE	1:B:266:ARG:HB3	1.64	0.41
1:B:300:THR:OG1	1:B:390:LEU:HD13	2.20	0.41
1:B:94:GLU:OE2	1:B:232:GLY:HA3	2.21	0.41
1:B:251:ARG:HG3	1:B:252:LEU:HG	2.03	0.41
1:B:370:THR:O	1:B:376:GLN:NE2	2.48	0.41
1:A:7:TRP:CZ3	1:A:47:ALA:HB2	2.56	0.40
1:A:190:ASP:OD1	1:A:191:THR:HG23	2.21	0.40
1:A:25:THR:HG21	4:A:506:SO4:O3	2.21	0.40
1:B:273:THR:CG2	1:B:278:ASP:HB3	2.51	0.40
1:A:266:ARG:NH1	6:A:605:HOH:O	2.45	0.40
1:B:52:ASP:OD1	1:B:52:ASP:N	2.55	0.40
1:A:171:TRP:CE3	1:A:179:PRO:HG3	2.56	0.40
1:A:228:THR:H	3:A:504:EDO:C1	2.35	0.40
1:B:311:TRP:HA	1:B:314:MET:HE2	2.04	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	388/412 (94%)	377 (97%)	11 (3%)	0	100	100
1	B	387/412 (94%)	365 (94%)	21 (5%)	1 (0%)	36	58
All	All	775/824 (94%)	742 (96%)	32 (4%)	1 (0%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	50	ASP

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	309/339 (91%)	307 (99%)	2 (1%)	78	91
1	B	302/339 (89%)	297 (98%)	5 (2%)	53	78
All	All	611/678 (90%)	604 (99%)	7 (1%)	65	84

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	211	LEU
1	A	366	LEU
1	B	57	LEU
1	B	134	ASN
1	B	370	THR

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Mol	Chain	Res	Type
1	B	371	ASN
1	B	373	LYS

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	181	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

19 ligands 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	$\# Z > 2$
3	EDO	A	503	-	3,3,3	0.48	0	2,2,2	0.38	0
3	EDO	B	503	-	3,3,3	0.40	0	2,2,2	0.59	0
3	EDO	B	504	-	3,3,3	0.39	0	2,2,2	0.48	0
3	EDO	B	502	-	3,3,3	0.51	0	2,2,2	0.30	0
2	CDP	B	501	-	25,26,26	3.61	14 (56%)	38,40,40	1.04	2 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	B	506	-	3,3,3	0.63	0	2,2,2	0.11	0
2	CDP	A	501	-	25,26,26	3.67	14 (56%)	38,40,40	1.01	2 (5%)
3	EDO	A	505	-	3,3,3	0.46	0	2,2,2	0.32	0
3	EDO	A	502	-	3,3,3	0.48	0	2,2,2	0.17	0
5	ACT	A	509	-	3,3,3	0.86	0	3,3,3	1.53	0
3	EDO	A	504	-	3,3,3	0.56	0	2,2,2	0.16	0
3	EDO	B	505	-	3,3,3	0.82	0	2,2,2	0.43	0
4	SO4	A	506	-	4,4,4	0.28	0	6,6,6	0.46	0
4	SO4	B	508	-	4,4,4	0.25	0	6,6,6	0.27	0
5	ACT	B	510	-	3,3,3	1.10	0	3,3,3	1.01	0
4	SO4	B	507	-	4,4,4	0.25	0	6,6,6	0.19	0
5	ACT	A	508	-	3,3,3	0.78	0	3,3,3	1.73	2 (66%)
5	ACT	A	507	-	3,3,3	0.89	0	3,3,3	1.21	0
5	ACT	B	509	-	3,3,3	0.83	0	3,3,3	1.28	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
3	EDO	A	505	-	-	0/1/1/1	-
3	EDO	A	503	-	-	0/1/1/1	-
3	EDO	B	503	-	-	0/1/1/1	-
3	EDO	B	504	-	-	0/1/1/1	-
3	EDO	B	502	-	-	0/1/1/1	-
2	CDP	B	501	-	-	4/16/32/32	0/2/2/2
3	EDO	A	502	-	-	0/1/1/1	-
3	EDO	A	504	-	-	1/1/1/1	-
3	EDO	B	506	-	-	1/1/1/1	-
2	CDP	A	501	-	-	3/16/32/32	0/2/2/2
3	EDO	B	505	-	-	0/1/1/1	-

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	CDP	C3'-C4'	-8.88	1.30	1.53
2	A	501	CDP	C3'-C4'	-8.65	1.31	1.53
2	A	501	CDP	O4'-C4'	7.11	1.60	1.45
2	B	501	CDP	O4'-C4'	6.87	1.60	1.45
2	B	501	CDP	C6-C5	5.96	1.48	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	CDP	C6-C5	5.95	1.48	1.35
2	B	501	CDP	C2-N3	5.90	1.48	1.36
2	A	501	CDP	PA-O3A	5.82	1.65	1.59
2	A	501	CDP	C2-N3	5.25	1.46	1.36
2	B	501	CDP	O4'-C1'	-4.82	1.30	1.42
2	B	501	CDP	C4-N3	4.72	1.43	1.34
2	A	501	CDP	O4'-C1'	-4.66	1.31	1.42
2	A	501	CDP	C4-N3	4.46	1.43	1.34
2	B	501	CDP	C4-N4	4.37	1.44	1.33
2	A	501	CDP	C4-N4	4.03	1.43	1.33
2	A	501	CDP	C2-N1	3.95	1.48	1.40
2	B	501	CDP	C2-N1	3.65	1.47	1.40
2	B	501	CDP	PA-O3A	3.27	1.63	1.59
2	B	501	CDP	O2-C2	-3.24	1.17	1.23
2	A	501	CDP	O2'-C2'	-3.00	1.35	1.43
2	B	501	CDP	C6-N1	2.88	1.45	1.38
2	A	501	CDP	C6-N1	2.85	1.44	1.38
2	A	501	CDP	O3'-C3'	2.77	1.49	1.43
2	B	501	CDP	O3'-C3'	2.54	1.49	1.43
2	B	501	CDP	O2'-C2'	-2.47	1.36	1.43
2	A	501	CDP	O2-C2	-2.47	1.19	1.23
2	B	501	CDP	C5-C4	2.40	1.48	1.42
2	A	501	CDP	C5-C4	2.29	1.48	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	CDP	C2'-C1'-N1	-3.14	104.52	113.25
2	A	501	CDP	O4'-C1'-N1	2.44	113.89	108.36
2	A	501	CDP	C4'-O4'-C1'	-2.26	104.48	109.47
2	B	501	CDP	O4'-C1'-N1	2.21	113.37	108.36
5	A	508	ACT	OXT-C-O	-2.17	113.96	122.03
5	A	508	ACT	OXT-C-CH3	2.07	123.74	115.05

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	CDP	C3'-C4'-C5'-O5'
2	B	501	CDP	O4'-C4'-C5'-O5'
2	A	501	CDP	C3'-C4'-C5'-O5'
2	A	501	CDP	O4'-C4'-C5'-O5'

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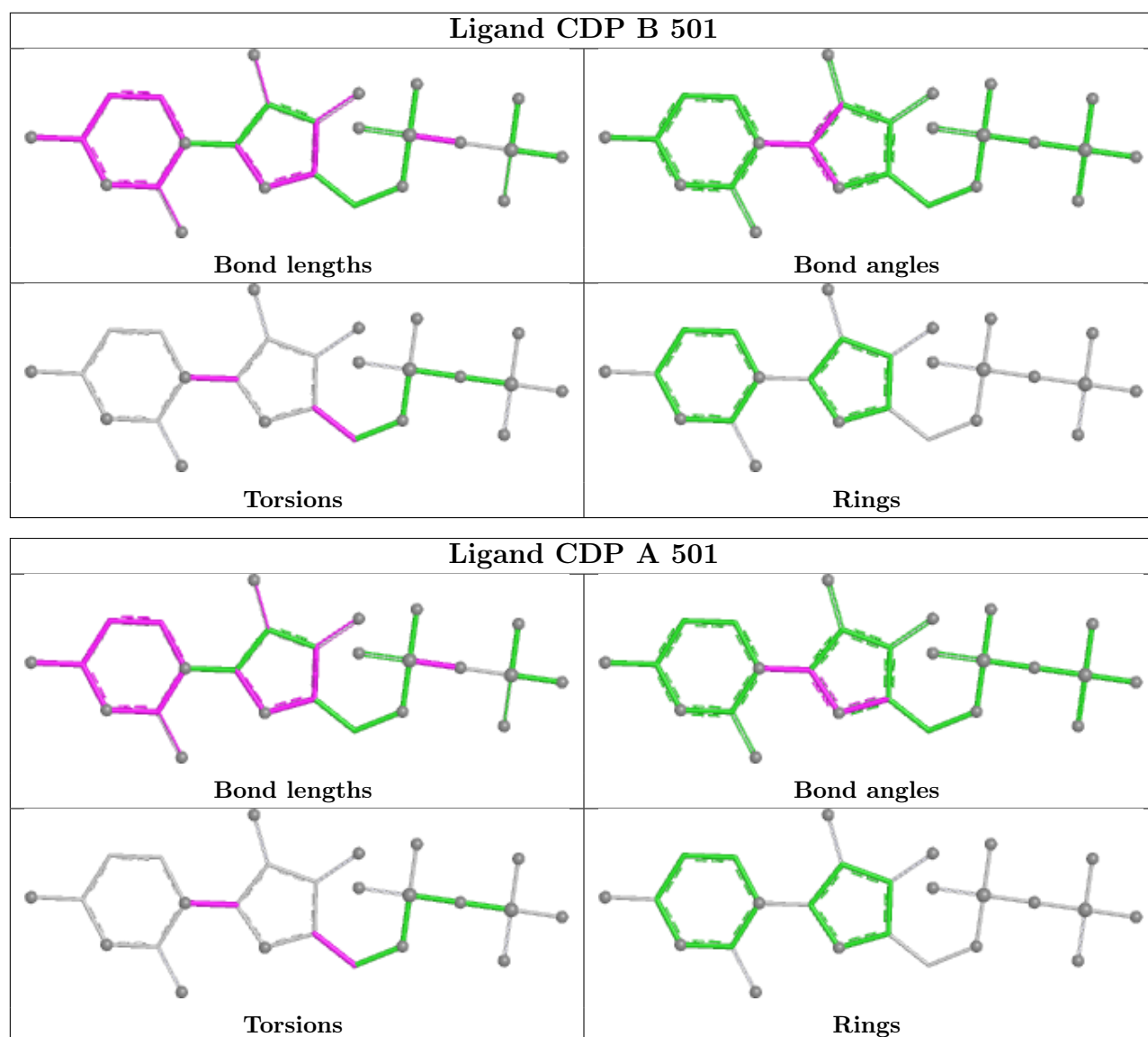
Mol	Chain	Res	Type	Atoms
3	A	504	EDO	O1-C1-C2-O2
3	B	506	EDO	O1-C1-C2-O2
2	B	501	CDP	C2'-C1'-N1-C6
2	B	501	CDP	O4'-C1'-N1-C6
2	A	501	CDP	O4'-C1'-N1-C6

There are no ring outliers.

14 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	EDO	1	0
3	B	503	EDO	3	0
3	B	502	EDO	1	0
2	B	501	CDP	1	0
3	B	506	EDO	1	0
2	A	501	CDP	1	0
3	A	505	EDO	7	0
3	A	502	EDO	3	0
5	A	509	ACT	3	0
3	A	504	EDO	4	0
3	B	505	EDO	1	0
4	A	506	SO4	1	0
4	B	507	SO4	1	0
5	B	509	ACT	1	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	394/412 (95%)	-0.18	10 (2%) 58 52	20, 32, 62, 80	0
1	B	393/412 (95%)	-0.13	16 (4%) 41 36	17, 30, 67, 79	0
All	All	787/824 (95%)	-0.16	26 (3%) 49 43	17, 31, 66, 80	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	157	THR	4.5
1	A	399	ALA	4.0
1	B	3	PRO	3.9
1	B	318	ILE	3.8
1	B	319	ARG	3.5
1	B	399	ALA	3.3
1	B	135	ARG	3.0
1	B	71	HIS	2.8
1	B	351	SER	2.8
1	A	158	GLU	2.8
1	B	374	ARG	2.8
1	B	321	HIS	2.6
1	A	371	ASN	2.6
1	A	372	ARG	2.4
1	A	75	LEU	2.4
1	A	134	ASN	2.4
1	A	332	PRO	2.4
1	A	347	ASP	2.4
1	A	156	PRO	2.4
1	B	158	GLU	2.2
1	B	375	ASP	2.2
1	B	136	ASP	2.2
1	B	157	THR	2.2
1	B	75	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	266	ARG	2.0
1	B	345	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 oligosaccharides in this entry.

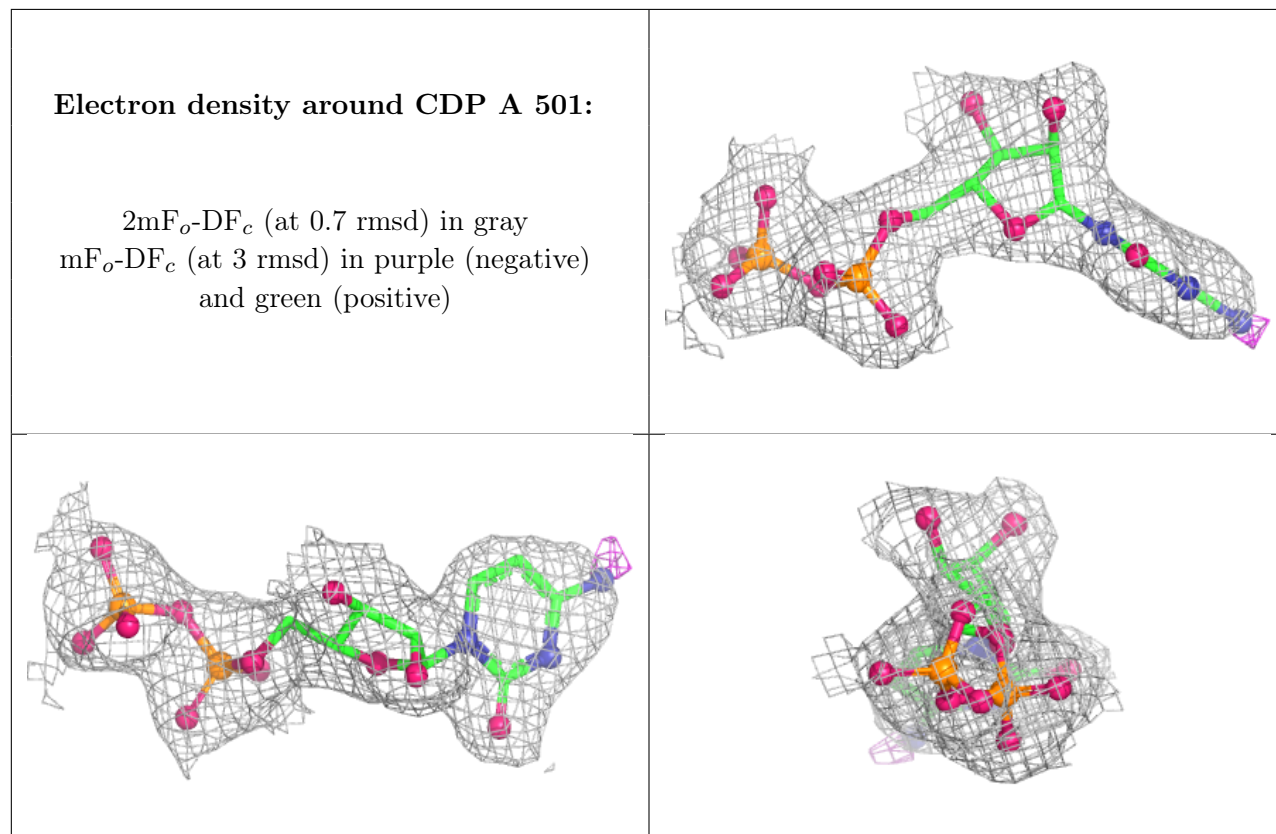
6.4 Ligands [i](#)

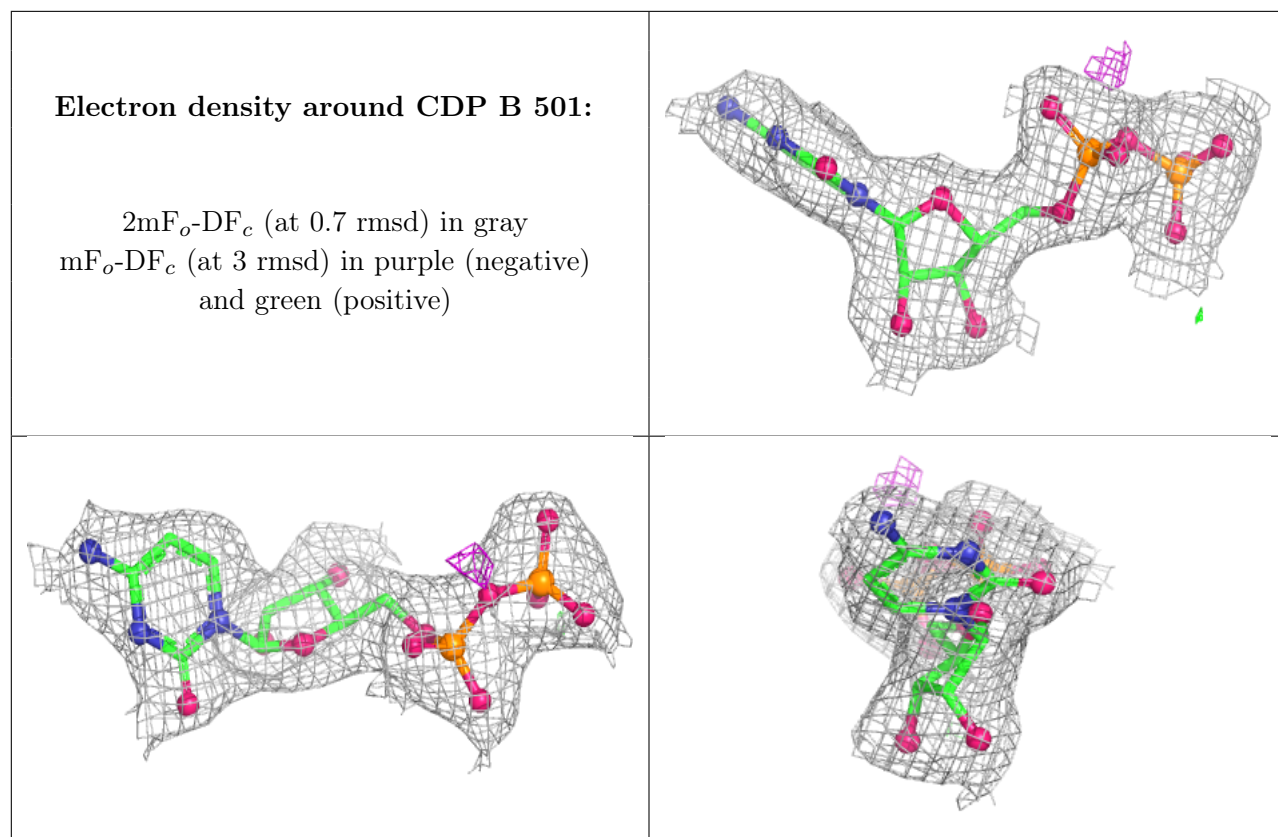
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	B	504	4/4	0.75	0.21	56,58,58,59	0
3	EDO	A	503	4/4	0.76	0.15	44,48,51,52	0
3	EDO	A	502	4/4	0.80	0.16	44,45,46,55	0
3	EDO	B	503	4/4	0.82	0.19	49,50,53,54	0
3	EDO	B	506	4/4	0.83	0.17	26,39,43,44	0
3	EDO	B	502	4/4	0.84	0.17	28,37,44,47	0
5	ACT	A	507	4/4	0.84	0.15	36,41,45,48	0
3	EDO	B	505	4/4	0.85	0.20	29,33,34,34	0
5	ACT	A	509	4/4	0.85	0.26	20,40,41,42	0
3	EDO	A	505	4/4	0.88	0.19	38,42,43,47	0
5	ACT	A	508	4/4	0.88	0.21	40,41,42,44	0
3	EDO	A	504	4/4	0.88	0.17	35,39,40,47	0
5	ACT	B	509	4/4	0.90	0.15	27,29,35,37	0
4	SO4	B	507	5/5	0.94	0.10	47,47,60,61	0
2	CDP	A	501	25/25	0.96	0.07	19,24,37,42	0
5	ACT	B	510	4/4	0.96	0.11	22,22,27,34	0
4	SO4	B	508	5/5	0.97	0.07	43,46,49,53	0
4	SO4	A	506	5/5	0.98	0.05	32,34,34,37	0
2	CDP	B	501	25/25	0.98	0.05	18,22,30,38	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.