



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 12:43 AM UTC

PDB ID : 1NC4 / pdb_00001nc4
Title : Crystal Structure of Monoclonal Antibody 2D12.5 Fab Complexed with Gd-DOTA
Authors : Corneillie, T.M.; Fisher, A.J.; Meares, C.F.
Deposited on : 2002-12-04
Resolution : 2.25 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

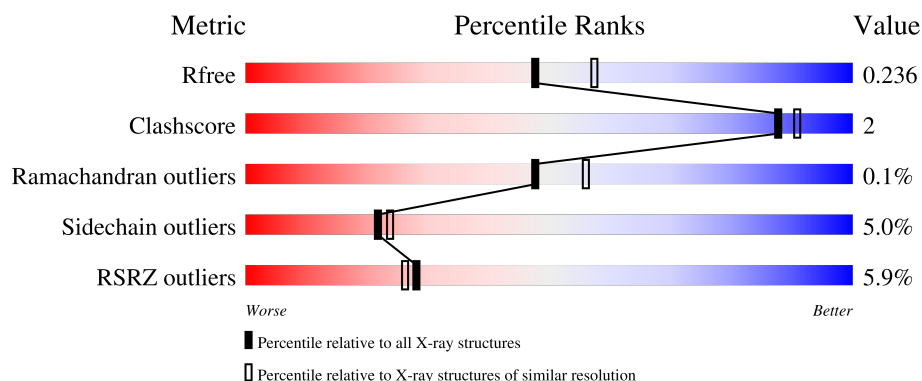
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	215	12% 85% 12% ..
1	C	215	5% 93% 7% .
2	B	221	6% 86% 11% ..
3	D	221	% 88% 9% .

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6944 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 MONOCLONAL ANTIBODY 2D12.5, LAMBDA LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1600	1001	270	323	6			
1	C	215	Total	C	N	O	S	0	0	0
			1627	1015	274	331	7			

- Molecule 2 is a protein called MONOCLONAL ANTIBODY 2D12.5, IGG1 GAMMA HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	219	Total	C	N	O	S	0	0	0
			1658	1055	273	323	7			

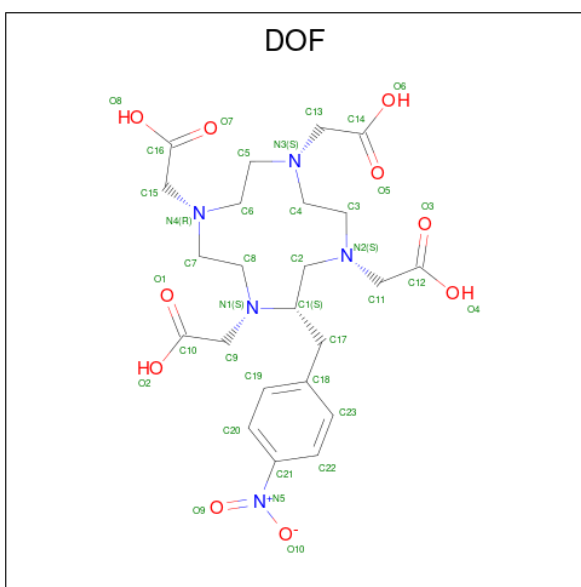
- Molecule 3 is a protein called MONOCLONAL ANTIBODY 2D12.5, IGG1 GAMMA HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	215	Total	C	N	O	S	0	0	0
			1633	1040	268	318	7			

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 5 is (S)-2-(4-NITROBENZYL)-1,4,7,10-TETRAAZACYCLODODECANE-N,N',N'',N'''-TETRAACETATE (CCD ID: DOF) (formula: C₂₃H₃₃N₅O₁₀).

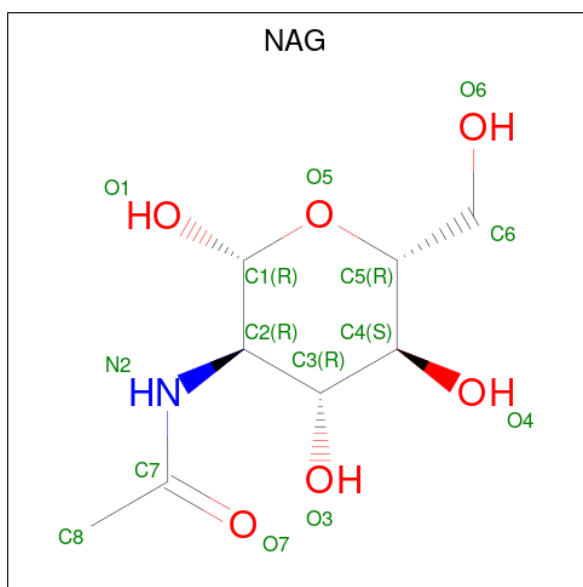


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			38	23	5	10		
5	D	1	Total	C	N	O	0	0
			38	23	5	10		

- Molecule 6 is GADOLINIUM ION (CCD ID: GD3) (formula: Gd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Gd	0	0
			1	1		
6	D	1	Total	Gd	0	0
			1	1		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	D	1	Total	C	N	O	0	0
			14	8	1	5		

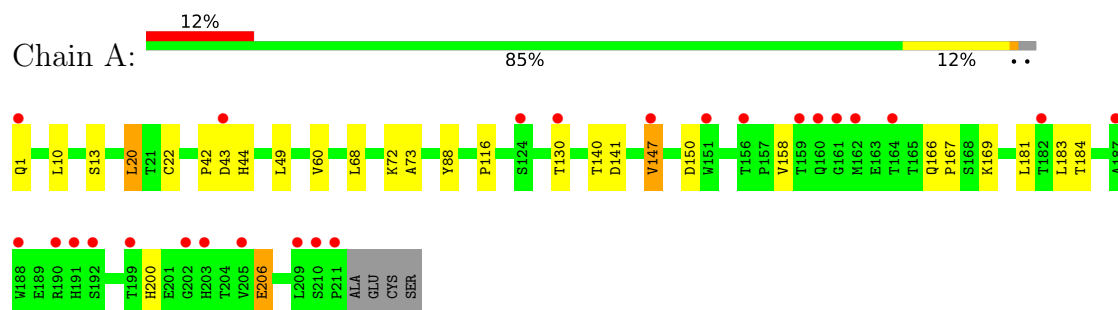
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	79	Total	O	0	0
			79	79		
8	B	77	Total	O	0	0
			77	77		
8	C	85	Total	O	0	0
			85	85		
8	D	91	Total	O	0	0
			91	91		

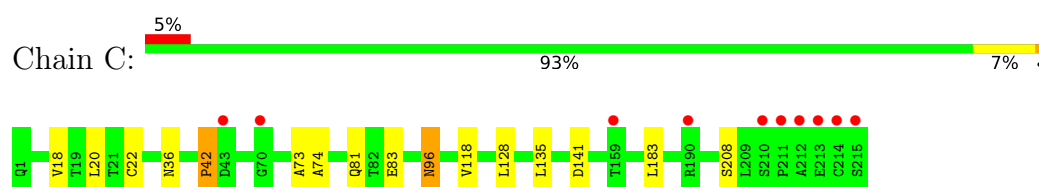
3 Residue-property plots [i](#)

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.

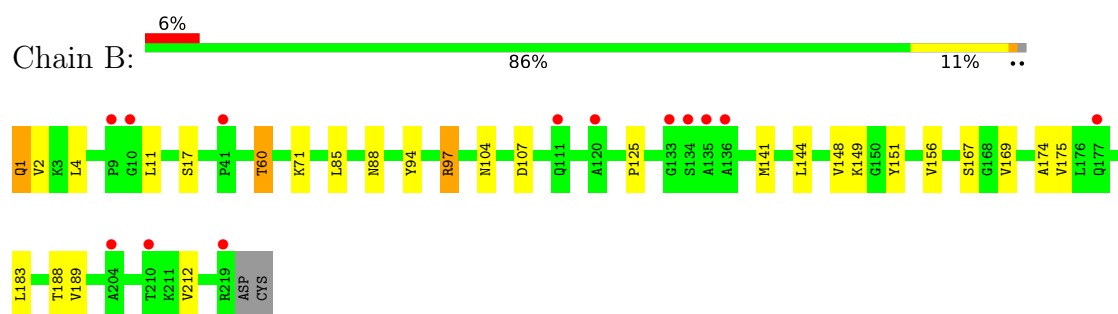
- Molecule 1: MONOCLONAL ANTIBODY 2D12.5, LAMBDA LIGHT CHAIN



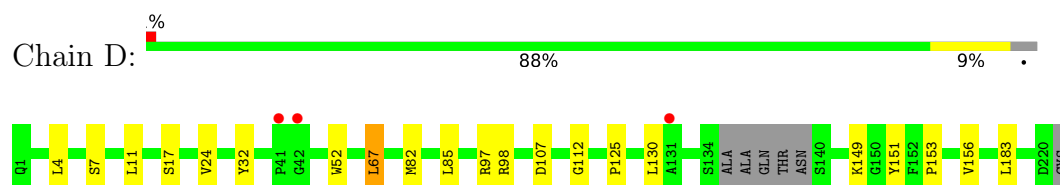
- Molecule 1: MONOCLONAL ANTIBODY 2D12.5, LAMBDA LIGHT CHAIN



- Molecule 2: MONOCLONAL ANTIBODY 2D12.5, IGG1 GAMMA HEAVY CHAIN



- Molecule 3: MONOCLONAL ANTIBODY 2D12.5, IGG1 GAMMA HEAVY CHAIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.57Å 82.24Å 160.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.39 – 2.25 29.39 – 2.25	Depositor EDS
% Data completeness (in resolution range)	91.5 (29.39-2.25) 91.5 (29.39-2.25)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.02 (at 2.24Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.208 , 0.248 0.199 , 0.236	Depositor DCC
R_{free} test set	2014 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6944	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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: PCA, NAG, DOF, CL, GD3

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	0.74	0/1640	1.03	9/2245 (0.4%)
1	C	0.72	0/1667	1.03	3/2280 (0.1%)
2	B	0.75	1/1698 (0.1%)	1.09	9/2328 (0.4%)
3	D	0.77	1/1679 (0.1%)	1.06	7/2298 (0.3%)
All	All	0.74	2/6684 (0.0%)	1.05	28/9151 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	82	MET	SD-CE	-5.47	1.65	1.79
2	B	189	VAL	CA-CB	5.33	1.61	1.55

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	67	LEU	N-CA-C	7.34	121.99	110.17
3	D	107	ASP	N-CA-C	7.19	119.02	111.03
1	A	141	ASP	N-CA-C	6.90	120.91	111.39
1	A	166	GLN	N-CA-C	-6.26	102.19	110.07
2	B	107	ASP	N-CA-C	6.21	117.93	111.03
3	D	32	TYR	N-CA-C	6.19	119.58	109.24
1	A	169	LYS	N-CA-C	6.18	119.43	110.28
2	B	167	SER	N-CA-C	6.17	118.00	111.28
1	C	208	SER	N-CA-C	6.11	118.25	109.14
3	D	112	GLY	N-CA-C	6.04	120.28	112.18
2	B	17	SER	N-CA-C	5.97	119.30	110.46
3	D	149	LYS	N-CA-C	5.93	119.48	109.76
2	B	174	ALA	N-CA-C	5.90	118.51	110.55
1	A	13	SER	N-CA-C	-5.79	101.70	110.50
2	B	60	THR	N-CA-C	-5.72	101.39	109.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	149	LYS	N-CA-C	5.71	119.55	109.96
1	C	141	ASP	N-CA-C	5.56	119.10	111.54
3	D	17	SER	N-CA-C	5.50	119.07	110.32
2	B	71	LYS	N-CA-C	5.46	117.10	108.96
2	B	169	VAL	N-CA-C	5.39	116.74	108.71
2	B	104	ASN	N-CA-C	5.32	117.85	109.39
1	A	130	THR	N-CA-C	-5.28	106.07	112.88
1	A	60	VAL	N-CA-C	-5.23	104.14	108.63
1	A	116	PRO	N-CA-C	5.16	119.59	111.38
3	D	156	VAL	N-CA-C	-5.12	100.50	108.95
1	A	167	PRO	N-CA-C	5.12	118.91	111.03
1	A	206	GLU	N-CA-C	5.08	117.52	109.50
1	C	74	ALA	N-CA-C	5.01	117.58	109.06

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	1600	0	1540	8	0
1	C	1627	0	1561	6	0
2	B	1658	0	1609	8	0
3	D	1633	0	1584	6	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	B	38	0	29	0	0
5	D	38	0	29	4	0
6	B	1	0	0	0	0
6	D	1	0	0	0	0
7	D	14	0	13	0	0
8	A	79	0	0	0	0
8	B	77	0	0	2	0
8	C	85	0	0	1	0
8	D	91	0	0	1	0
All	All	6944	0	6365	26	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 2.

All (26) 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:96:ASN:HD22	1:C:96:ASN:H	1.47	0.62
5:D:501:DOF:O7	5:D:501:DOF:H132	1.99	0.62
1:A:1:GLN:HE22	2:B:60:THR:HG23	1.68	0.58
1:C:96:ASN:HD22	1:C:96:ASN:N	2.00	0.56
1:A:1:GLN:NE2	2:B:60:THR:HG23	2.22	0.54
5:D:501:DOF:O7	5:D:501:DOF:HC51	2.09	0.52
1:C:36:ASN:HD21	3:D:98:ARG:HH21	1.60	0.49
1:A:10:LEU:HD12	1:A:20:LEU:HD22	1.95	0.49
3:D:7:SER:HB3	8:D:555:HOH:O	2.13	0.48
1:C:42:PRO:HD3	8:C:514:HOH:O	2.12	0.48
1:A:49:LEU:HD11	1:A:88:TYR:HE2	1.79	0.47
1:C:22:CYS:HB3	1:C:73:ALA:HB3	1.96	0.47
3:D:11:LEU:HB2	3:D:153:PRO:HG3	1.97	0.46
1:A:158:VAL:HG11	1:A:181:LEU:HD11	1.98	0.45
5:D:501:DOF:O7	5:D:501:DOF:C5	2.64	0.45
1:A:44:HIS:HB3	2:B:94:TYR:OH	2.17	0.45
1:A:22:CYS:HB3	1:A:73:ALA:HB3	1.99	0.45
1:C:36:ASN:HD21	3:D:98:ARG:NH2	2.15	0.44
1:A:147:VAL:HG23	1:A:200:HIS:HB2	2.00	0.44
2:B:1:PCA:HB2	8:B:539:HOH:O	2.19	0.43
2:B:148:VAL:HB	2:B:183:LEU:HD23	1.99	0.43
2:B:125:PRO:HB3	2:B:151:TYR:HB3	2.02	0.42
3:D:125:PRO:HB3	3:D:151:TYR:HB3	2.02	0.42
3:D:52:TRP:NE1	5:D:501:DOF:O7	2.53	0.41
2:B:141:MET:HE3	2:B:188:THR:HG22	2.03	0.40
2:B:97:ARG:NH2	8:B:562:HOH:O	2.54	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	209/215 (97%)	201 (96%)	7 (3%)	1 (0%)	24	24
1	C	213/215 (99%)	202 (95%)	11 (5%)	0	100	100
2	B	217/221 (98%)	207 (95%)	10 (5%)	0	100	100
3	D	211/221 (96%)	208 (99%)	3 (1%)	0	100	100
All	All	850/872 (98%)	818 (96%)	31 (4%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	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	180/183 (98%)	170 (94%)	10 (6%)	19	20
1	C	183/183 (100%)	173 (94%)	10 (6%)	19	20
2	B	187/189 (99%)	177 (95%)	10 (5%)	20	22
3	D	186/190 (98%)	179 (96%)	7 (4%)	29	36
All	All	736/745 (99%)	699 (95%)	37 (5%)	22	24

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	42	PRO
1	A	68	LEU
1	A	72	LYS
1	A	140	THR
1	A	147	VAL
1	A	150	ASP

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Mol	Chain	Res	Type
1	A	183	LEU
1	A	184	THR
1	A	206	GLU
2	B	2	VAL
2	B	4	LEU
2	B	11	LEU
2	B	85	LEU
2	B	88	ASN
2	B	97	ARG
2	B	144	LEU
2	B	156	VAL
2	B	175	VAL
2	B	212	VAL
1	C	18	VAL
1	C	20	LEU
1	C	42	PRO
1	C	81	GLN
1	C	83	GLU
1	C	96	ASN
1	C	118	VAL
1	C	128	LEU
1	C	135	LEU
1	C	183	LEU
3	D	4	LEU
3	D	24	VAL
3	D	67	LEU
3	D	85	LEU
3	D	97	ARG
3	D	130	LEU
3	D	183	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
1	A	44	HIS
1	A	173	ASN
2	B	86	GLN
2	B	88	ASN
2	B	111	GLN
2	B	177	GLN
1	C	96	ASN

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Mol	Chain	Res	Type
1	C	191	HIS
1	C	200	HIS
3	D	77	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
2	PCA	B	1	2	7,8,9	2.72	2 (28%)	9,10,12	1.99	4 (44%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PCA	B	1	2	-	0/0/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	PCA	CD-N	5.59	1.48	1.34
2	B	1	PCA	CA-N	4.26	1.51	1.46

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	PCA	CA-N-CD	-3.17	102.74	113.58
2	B	1	PCA	OE-CD-CG	-3.02	121.32	126.72
2	B	1	PCA	CG-CD-N	2.33	114.08	108.39
2	B	1	PCA	CB-CA-N	2.18	109.25	103.24

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	PCA	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\# Z > 2$
5	DOF	D	501	6	39,39,39	1.53	7 (17%)	48,52,52	1.39	9 (18%)
7	NAG	D	400	3	14,14,15	2.04	4 (28%)	17,19,21	2.11	5 (29%)
5	DOF	B	500	6	39,39,39	2.00	11 (28%)	48,52,52	1.52	10 (20%)

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
5	DOF	D	501	6	-	5/45/47/47	0/2/2/2
7	NAG	D	400	3	-	0/6/23/26	0/1/1/1
5	DOF	B	500	6	-	8/45/47/47	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	500	DOF	C17-C1	5.15	1.61	1.53
7	D	400	NAG	C1-C2	5.01	1.59	1.52
5	B	500	DOF	C8-N1	4.09	1.54	1.47
5	B	500	DOF	C20-C21	3.87	1.46	1.38
7	D	400	NAG	C4-C5	3.54	1.60	1.53
5	D	501	DOF	C20-C21	3.30	1.44	1.38
5	B	500	DOF	C22-C21	3.28	1.44	1.38
5	D	501	DOF	C17-C1	3.18	1.58	1.53
5	D	501	DOF	C21-N5	-3.11	1.37	1.45
5	D	501	DOF	C22-C21	3.00	1.44	1.38
5	B	500	DOF	C15-N4	2.90	1.53	1.47
5	B	500	DOF	C20-C19	2.90	1.43	1.38
5	B	500	DOF	C23-C18	2.66	1.44	1.38
7	D	400	NAG	O7-C7	-2.57	1.17	1.23
5	D	501	DOF	C23-C18	2.43	1.43	1.38
5	B	500	DOF	C19-C18	2.39	1.43	1.38
5	B	500	DOF	C13-C14	2.37	1.56	1.51
5	D	501	DOF	C8-N1	2.22	1.50	1.47
5	D	501	DOF	C9-C10	2.22	1.55	1.51
7	D	400	NAG	O5-C5	2.15	1.47	1.43
5	B	500	DOF	C23-C22	2.13	1.42	1.38
5	B	500	DOF	C21-N5	-2.09	1.40	1.45

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	400	NAG	C2-N2-C7	5.56	130.35	122.90
5	B	500	DOF	C20-C21-N5	3.52	122.41	119.34
5	D	501	DOF	C22-C21-N5	3.11	122.05	119.34
5	B	500	DOF	C1-C2-N2	2.99	117.26	113.40
5	B	500	DOF	C9-N1-C1	-2.93	108.00	113.00
7	D	400	NAG	O7-C7-N2	2.89	127.08	121.98
7	D	400	NAG	C4-C3-C2	-2.88	106.80	111.02
5	B	500	DOF	C15-N4-C7	-2.83	105.10	111.91
5	D	501	DOF	C9-N1-C1	-2.81	108.19	113.00
5	D	501	DOF	C11-N2-C3	-2.79	105.19	111.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	500	DOF	C3-C4-N3	2.76	119.54	112.98
5	B	500	DOF	C22-C21-N5	2.57	121.58	119.34
5	D	501	DOF	C7-C8-N1	2.56	117.94	112.72
5	B	500	DOF	C5-C6-N4	2.47	118.83	112.98
5	B	500	DOF	C5-N3-C4	-2.45	105.61	111.44
5	D	501	DOF	C20-C21-N5	2.45	121.47	119.34
5	B	500	DOF	C22-C21-C20	-2.41	116.00	119.88
5	B	500	DOF	C13-N3-C5	-2.39	106.15	111.91
5	D	501	DOF	C1-C2-N2	2.36	116.44	113.40
5	D	501	DOF	C13-N3-C5	-2.29	106.39	111.91
7	D	400	NAG	C1-C2-N2	-2.20	106.96	110.43
5	D	501	DOF	C22-C21-C20	-2.11	116.47	119.88
5	D	501	DOF	C5-C6-N4	2.11	117.99	112.98
7	D	400	NAG	O4-C4-C3	-2.01	105.64	110.38

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	500	DOF	N2-C11-C12-O3
5	B	500	DOF	N2-C11-C12-O4
5	D	501	DOF	N2-C11-C12-O4
5	D	501	DOF	N3-C13-C14-O6
5	D	501	DOF	N3-C13-C14-O5
5	B	500	DOF	C14-C13-N3-C5
5	D	501	DOF	N2-C11-C12-O3
5	D	501	DOF	N1-C1-C2-N2
5	B	500	DOF	N1-C1-C2-N2
5	B	500	DOF	C6-C5-N3-C4
5	B	500	DOF	N3-C13-C14-O5
5	B	500	DOF	N2-C3-C4-N3
5	B	500	DOF	C1-C17-C18-C23

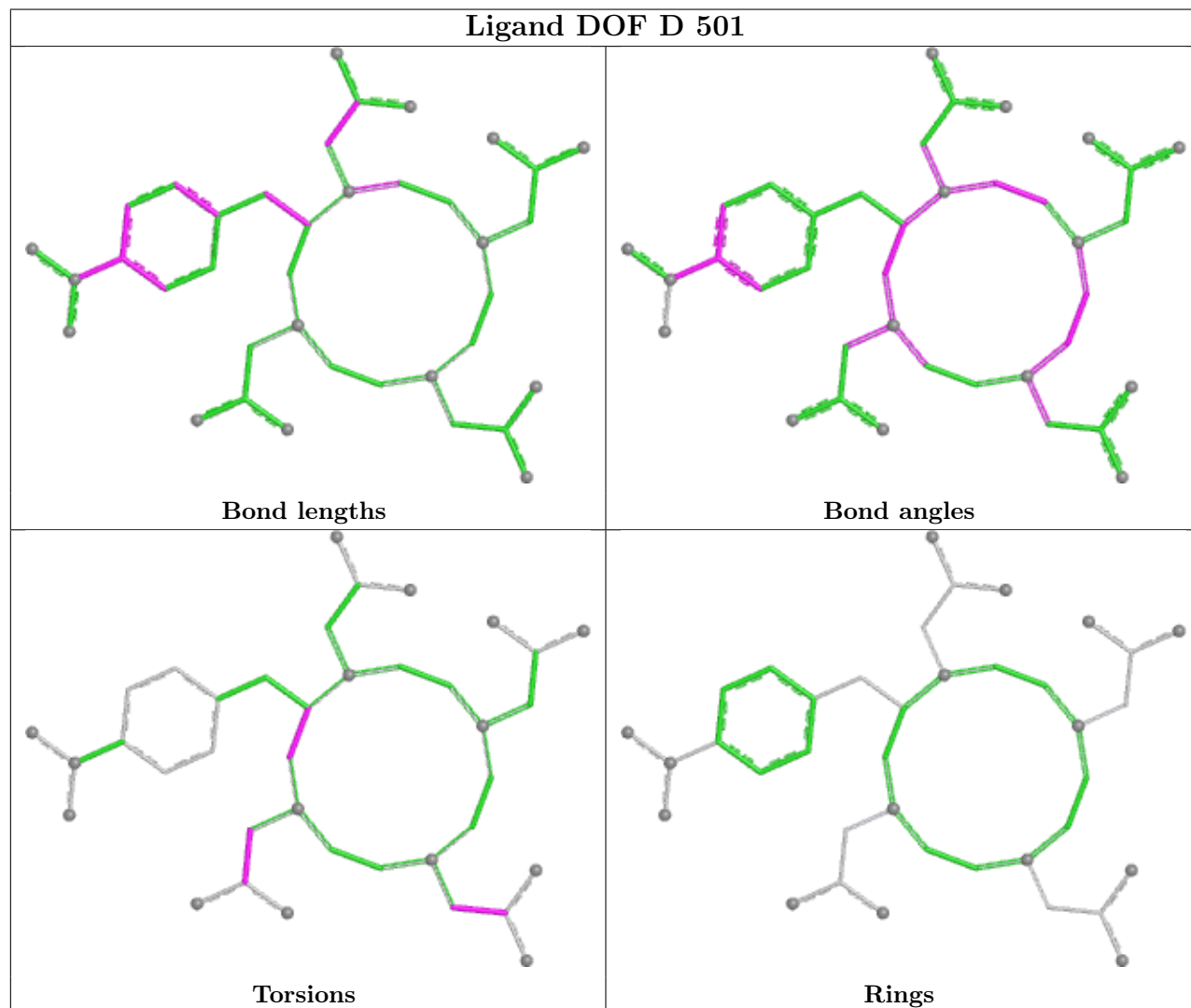
There are no ring outliers.

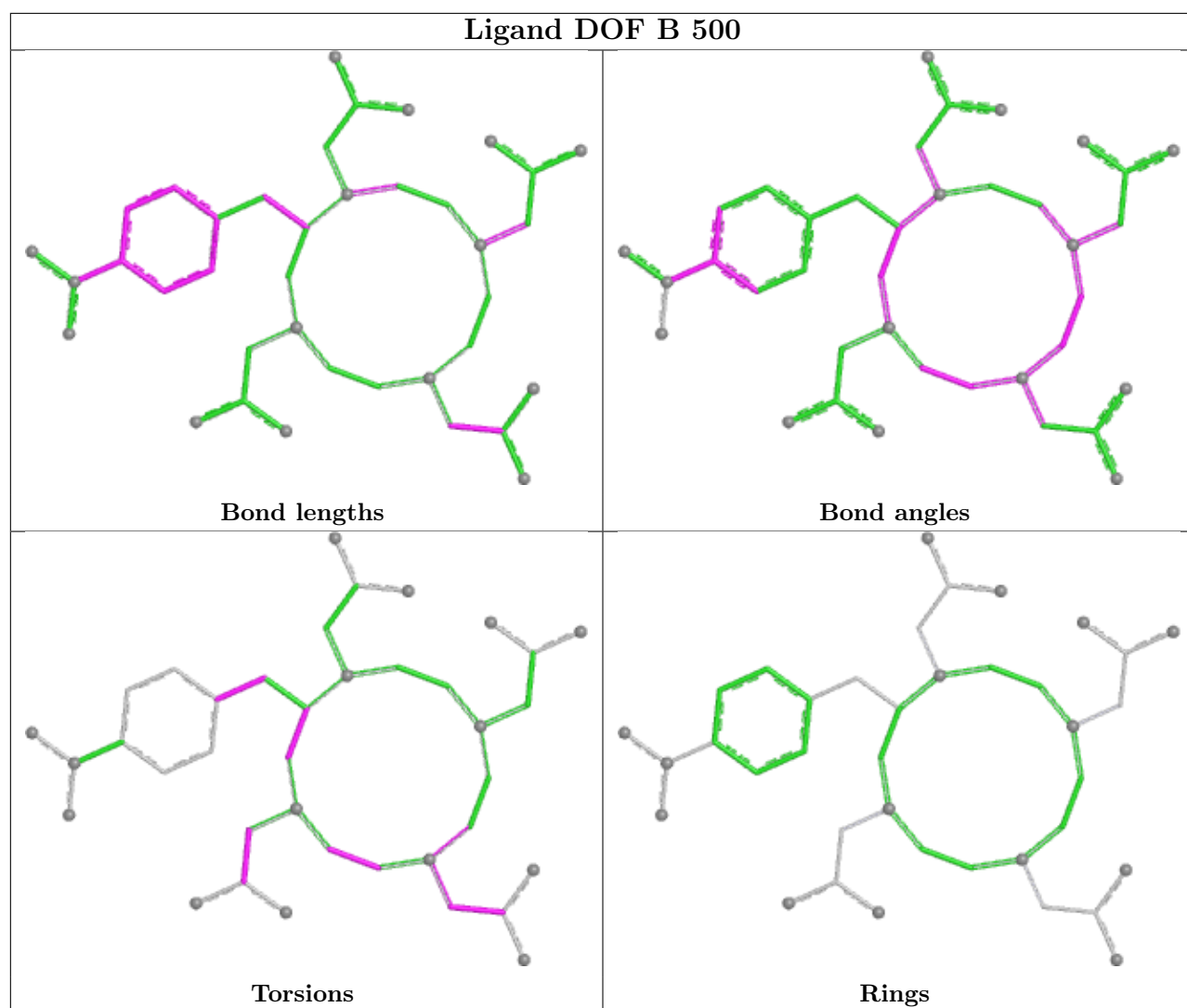
1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	501	DOF	4	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	211/215 (98%)	0.41	25 (11%) 9 8	10, 28, 68, 82	0
1	C	215/215 (100%)	0.05	10 (4%) 36 35	14, 27, 45, 83	0
2	B	218/221 (98%)	0.15	13 (5%) 27 25	11, 27, 54, 61	0
3	D	215/221 (97%)	0.02	3 (1%) 73 74	12, 27, 52, 66	0
All	All	859/872 (98%)	0.16	51 (5%) 28 26	10, 27, 60, 83	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	212	ALA	4.0
1	A	211	PRO	3.9
1	C	211	PRO	3.8
2	B	9	PRO	3.7
1	A	164	THR	3.6
1	A	192	SER	3.4
1	C	43	ASP	3.4
1	A	162	MET	3.2
1	A	156	THR	3.1
3	D	42	GLY	3.0
1	A	130	THR	2.9
1	A	1	GLN	2.8
1	A	190	ARG	2.8
2	B	219	ARG	2.8
1	A	191	HIS	2.8
1	A	187	ALA	2.8
2	B	135	ALA	2.8
2	B	177	GLN	2.7
1	C	214	CYS	2.7
2	B	41	PRO	2.7
1	A	159	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	210	SER	2.7
2	B	111	GLN	2.6
2	B	133	GLY	2.6
1	A	209	LEU	2.6
1	C	213	GLU	2.6
3	D	131	ALA	2.5
1	C	190	ARG	2.4
2	B	210	THR	2.4
1	A	205	VAL	2.4
1	C	159	THR	2.4
1	A	199	THR	2.4
2	B	134	SER	2.4
1	A	202	GLY	2.4
2	B	136	ALA	2.4
1	A	124	SER	2.4
1	A	161	GLY	2.3
1	C	210	SER	2.3
1	A	203	HIS	2.3
1	C	215	SER	2.3
1	C	70	GLY	2.2
1	A	151	TRP	2.2
1	A	147	VAL	2.2
2	B	120	ALA	2.2
1	A	43	ASP	2.2
3	D	41	PRO	2.2
2	B	204	ALA	2.2
1	A	160	GLN	2.1
1	A	188	TRP	2.1
2	B	10	GLY	2.1
1	A	182	THR	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PCA	B	1	8/9	0.85	0.11	31,33,33,36	0

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

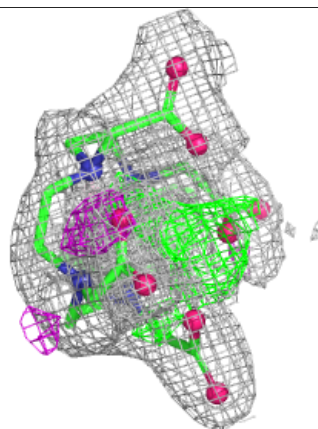
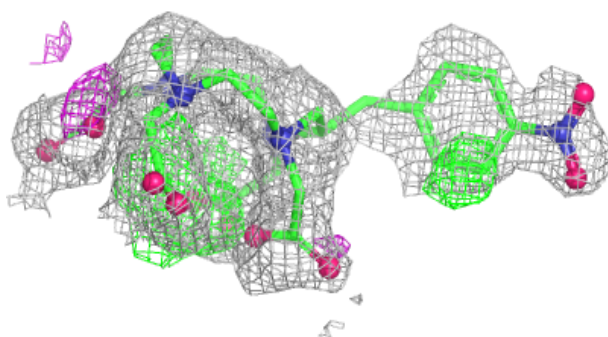
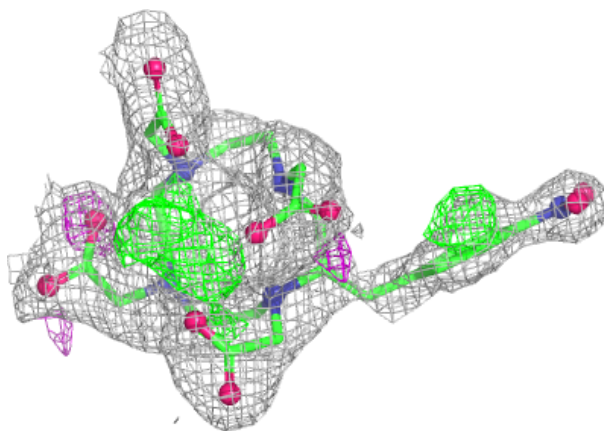
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

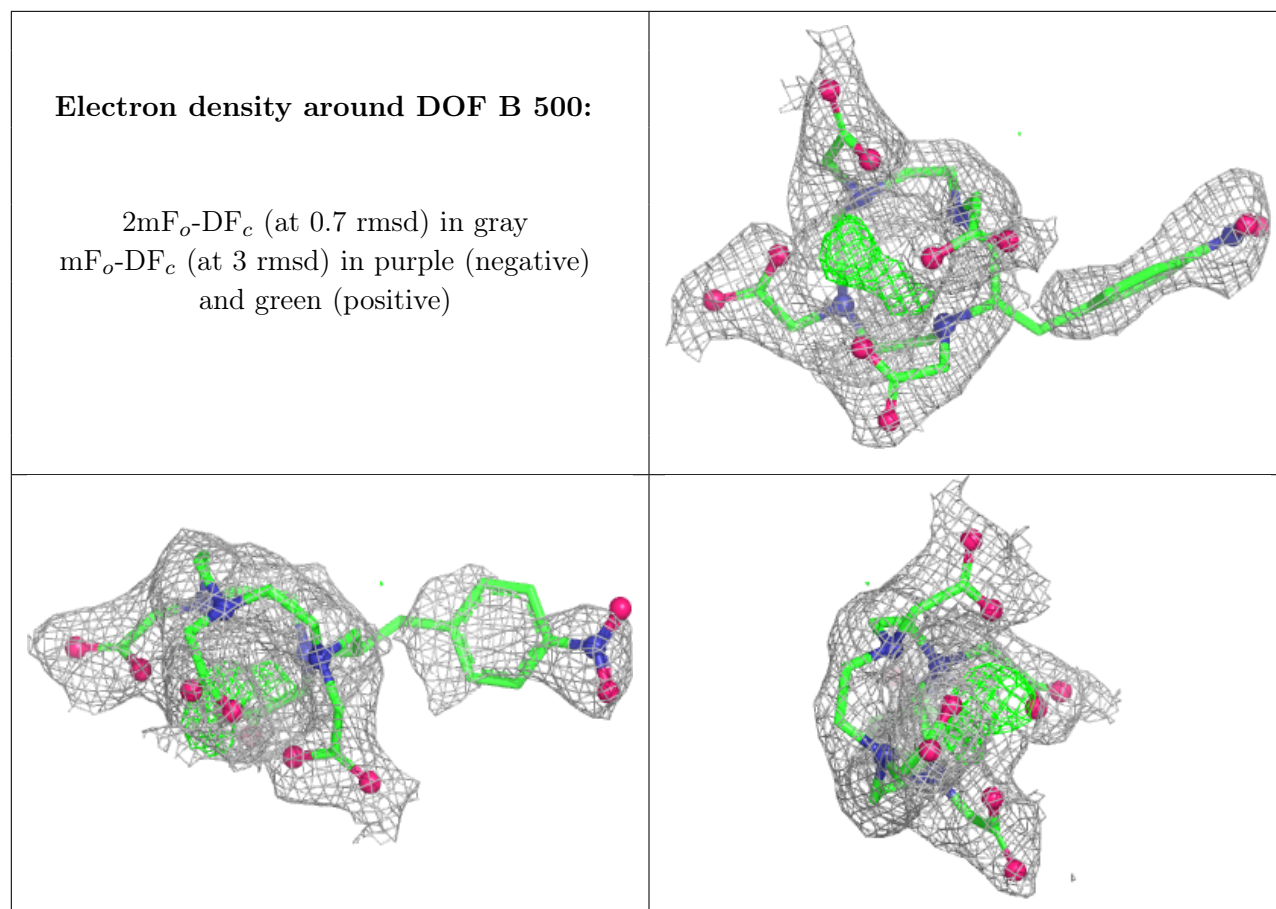
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	NAG	D	400	14/15	0.70	0.15	45,47,52,53	0
4	CL	A	504	1/1	0.89	0.11	54,54,54,54	0
5	DOF	D	501	38/38	0.93	0.12	13,21,30,31	10
4	CL	C	505	1/1	0.93	0.11	48,48,48,48	0
5	DOF	B	500	38/38	0.95	0.09	16,24,56,58	0
6	GD3	D	503	1/1	1.00	0.03	18,18,18,18	0
6	GD3	B	502	1/1	1.00	0.03	19,19,19,19	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 DOF 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.