



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:40 PM UTC

PDB ID : 1NC2 / pdb_00001nc2
Title : Crystal Structure of Monoclonal Antibody 2D12.5 Fab Complexed with Y-DOTA
Authors : Corneillie, T.M.; Fisher, A.J.; Meares, C.F.
Deposited on : 2002-12-04
Resolution : 2.10 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

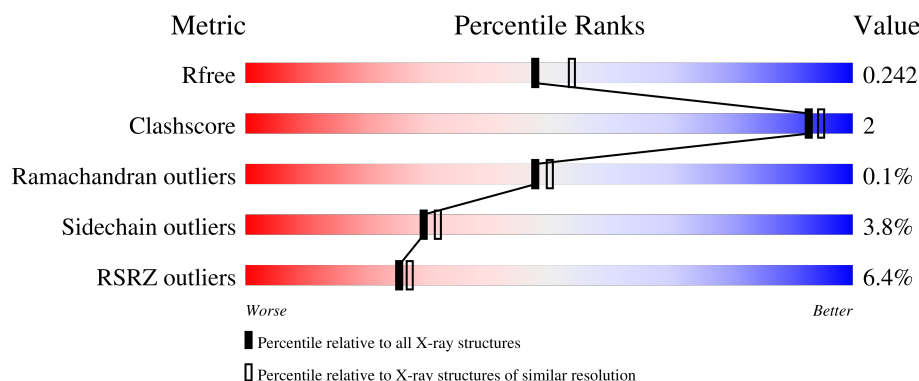
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	10% 88% 10% .
1	C	215	5% 93% 7%
2	B	221	7% 87% 11% ..
3	D	221	4% 88% 9% .

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6985 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
			1634	1040	268	318	8			

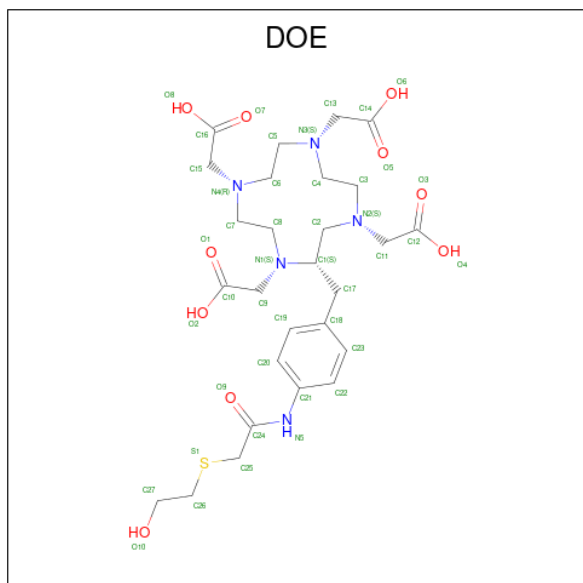
- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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 YTTRIUM (III) ION (CCD ID: YT3) (formula: Y).

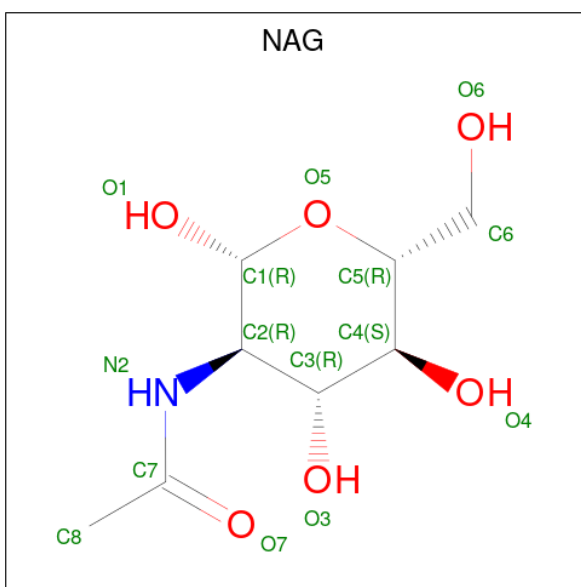
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Y 1 1	0	0
5	D	1	Total Y 1 1	0	0

- Molecule 6 is (S)-2-(4-(2-(2-HYDROXYETHYLTHIO)-ACETAMIDO)-BENZYL)-1,4,7,10-TETRAAZACYCLODODECANE-N,N',N'',N'''-TETRAACETATE (CCD ID: DOE) (formula: C₂₇H₄₁N₅O₁₀S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	S	0	0
			43	27	5	10	1		
6	D	1	Total	C	N	O	S	0	0
			43	27	5	10	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		

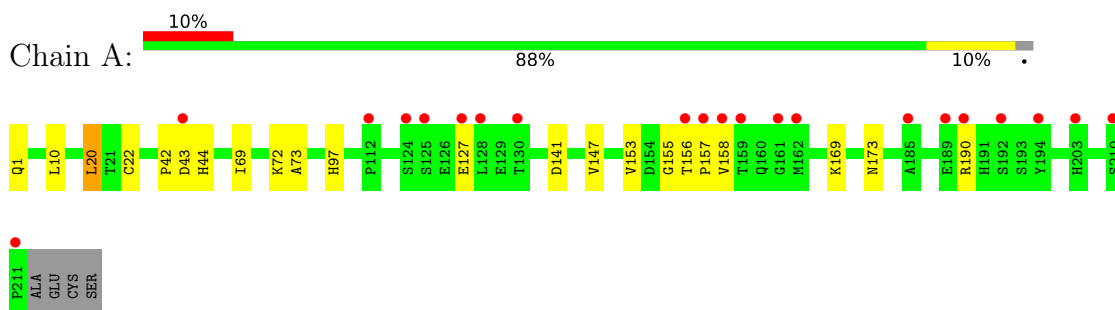
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	74	Total	O	0	0
			74	74		
8	B	88	Total	O	0	0
			88	88		
8	C	102	Total	O	0	0
			102	102		
8	D	98	Total	O	0	0
			98	98		

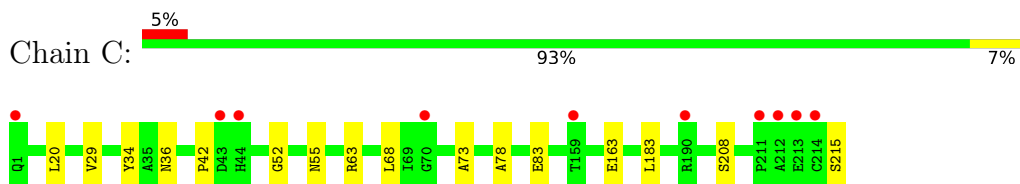
3 Residue-property plots [i](#)

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.

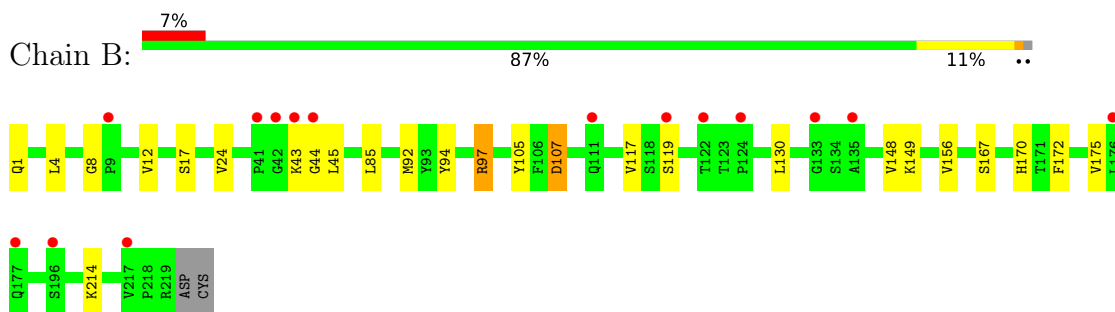
- Molecule 1: MONOCLONAL ANTIBODY 2D12.5, LAMBDA LIGHT CHAIN



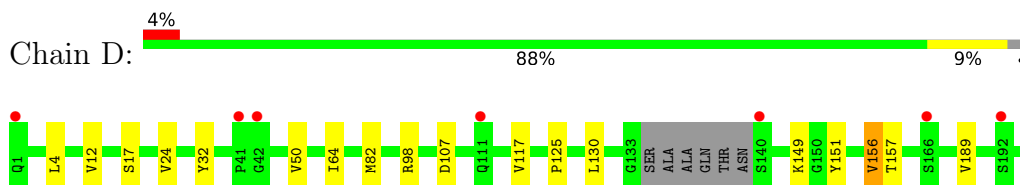
- Molecule 1: MONOCLONAL ANTIBODY 2D12.5, LAMBDA LIGHT CHAIN



- 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	67.33Å 81.30Å 160.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.20 – 2.10 29.20 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.9 (29.20-2.10) 96.0 (29.20-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 2.10Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.209 , 0.247 0.203 , 0.242	Depositor DCC
R_{free} test set	2558 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6985	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% 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: YT3, NAG, PCA, DOE, CL

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	0.75	0/1640	1.01	2/2245 (0.1%)
1	C	0.73	0/1667	1.02	2/2280 (0.1%)
2	B	0.78	0/1698	1.12	9/2328 (0.4%)
3	D	0.80	3/1680 (0.2%)	1.05	5/2298 (0.2%)
All	All	0.76	3/6685 (0.0%)	1.05	18/9151 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	82	MET	SD-CE	-6.89	1.62	1.79
3	D	50	VAL	CA-CB	5.33	1.60	1.55
3	D	189	VAL	CA-CB	5.33	1.59	1.55

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	107	ASP	N-CA-C	8.40	120.21	111.14
2	B	8	GLY	CA-C-N	8.38	128.44	119.90
2	B	8	GLY	C-N-CA	8.38	128.44	119.90
3	D	107	ASP	N-CA-C	6.79	118.47	111.14
1	A	141	ASP	N-CA-C	6.76	120.73	111.54
1	C	208	SER	N-CA-C	6.62	119.01	109.14
3	D	149	LYS	N-CA-C	6.55	120.13	109.59
2	B	172	PHE	N-CA-C	6.27	119.40	110.24
1	A	169	LYS	N-CA-C	6.22	119.28	110.50
2	B	149	LYS	N-CA-C	6.03	119.65	109.76
2	B	17	SER	N-CA-C	5.67	118.82	110.59
3	D	156	VAL	N-CA-C	-5.58	100.20	108.45
3	D	32	TYR	N-CA-C	5.41	118.22	109.40
2	B	167	SER	N-CA-C	5.37	117.83	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	119	SER	N-CA-C	-5.36	106.33	112.92
3	D	17	SER	N-CA-C	5.29	118.74	110.32
2	B	45	LEU	N-CA-C	5.20	118.41	110.20
1	C	52	GLY	N-CA-C	5.06	120.47	113.99

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	7	0
1	C	1627	0	1560	5	0
2	B	1658	0	1609	6	0
3	D	1634	0	1583	5	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
6	B	43	0	37	0	0
6	D	43	0	37	1	0
7	D	14	0	13	0	0
8	A	74	0	0	1	0
8	B	88	0	0	0	0
8	C	102	0	0	0	0
8	D	98	0	0	0	0
All	All	6985	0	6379	20	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 (20) 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)
2:B:43:LYS:HG2	2:B:44:GLY:H	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:CYS:HB3	1:A:73:ALA:HB3	1.86	0.57
1:A:10:LEU:HD12	1:A:20:LEU:HD22	1.88	0.56
1:A:1:GLN:HE22	1:A:97:HIS:CD2	2.28	0.51
2:B:97:ARG:HD3	2:B:107:ASP:OD2	2.13	0.48
1:C:36:ASN:HD21	3:D:98:ARG:HH21	1.60	0.47
1:C:215:SER:HA	3:D:219:ARG:HH11	1.80	0.47
1:A:44:HIS:HB3	2:B:94:TYR:OH	2.15	0.46
2:B:148:VAL:HG11	2:B:156:VAL:HG11	1.98	0.45
3:D:125:PRO:HB3	3:D:151:TYR:HB3	1.99	0.45
1:A:69:ILE:O	1:A:72:LYS:HG2	2.17	0.45
1:C:63:ARG:HB2	1:C:78:ALA:O	2.17	0.45
1:A:153:VAL:C	1:A:155:GLY:H	2.27	0.42
1:C:29:VAL:HG11	1:C:73:ALA:HB2	2.01	0.42
3:D:12:VAL:O	3:D:117:VAL:HA	2.20	0.41
3:D:157:THR:OG1	3:D:204:ALA:HB3	2.20	0.41
2:B:12:VAL:O	2:B:117:VAL:HA	2.20	0.41
8:A:531:HOH:O	2:B:170:HIS:HD2	2.02	0.41
1:C:34:TYR:CE1	6:D:501:DOE:HC62	2.55	0.41
1:A:156:THR:HA	1:A:157:PRO:HD3	1.97	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	22
1	C	213/215 (99%)	206 (97%)	7 (3%)	0	100	100
2	B	217/221 (98%)	205 (94%)	12 (6%)	0	100	100
3	D	211/221 (96%)	204 (97%)	7 (3%)	0	100	100
All	All	850/872 (98%)	816 (96%)	33 (4%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	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	180/183 (98%)	173 (96%)	7 (4%)	28	31
1	C	183/183 (100%)	176 (96%)	7 (4%)	29	32
2	B	187/189 (99%)	178 (95%)	9 (5%)	23	23
3	D	186/190 (98%)	181 (97%)	5 (3%)	39	45
All	All	736/745 (99%)	708 (96%)	28 (4%)	29	32

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	42	PRO
1	A	127	GLU
1	A	147	VAL
1	A	158	VAL
1	A	173	ASN
1	A	190	ARG
2	B	4	LEU
2	B	24	VAL
2	B	85	LEU
2	B	92	MET
2	B	97	ARG
2	B	105	TYR
2	B	130	LEU
2	B	175	VAL
2	B	214	LYS
1	C	20	LEU
1	C	42	PRO
1	C	55	ASN

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Mol	Chain	Res	Type
1	C	68	LEU
1	C	83	GLU
1	C	163	GLU
1	C	183	LEU
3	D	4	LEU
3	D	24	VAL
3	D	64	ILE
3	D	130	LEU
3	D	156	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
2	B	77	GLN
3	D	77	GLN
3	D	83	ASN
3	D	170	HIS

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	2.50	2 (28%)	9,10,12	2.14	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	5.63	1.48	1.34
2	B	1	PCA	CA-N	3.24	1.50	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	-3.32	102.19	113.58
2	B	1	PCA	OE-CD-CG	-3.10	121.19	126.72
2	B	1	PCA	CB-CA-N	2.54	110.24	103.24
2	B	1	PCA	CG-CD-N	2.28	113.96	108.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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
7	NAG	D	400	3	14,14,15	1.97	3 (21%)	17,19,21	1.98	4 (23%)
6	DOE	B	500	-	44,44,44	1.82	16 (36%)	54,57,57	1.40	6 (11%)
6	DOE	D	501	-	44,44,44	2.11	19 (43%)	54,57,57	1.36	6 (11%)

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
7	NAG	D	400	3	-	0/6/23/26	0/1/1/1
6	DOE	B	500	-	-	4/52/52/52	0/2/2/2
6	DOE	D	501	-	-	2/52/52/52	0/2/2/2

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	501	DOE	C17-C1	5.03	1.60	1.53
6	B	500	DOE	C17-C1	4.82	1.60	1.53
7	D	400	NAG	C1-C2	4.73	1.58	1.52
6	D	501	DOE	C6-N4	3.43	1.55	1.47
6	D	501	DOE	C2-N2	3.27	1.53	1.47
6	D	501	DOE	C23-C18	3.19	1.45	1.38
6	D	501	DOE	C20-C21	3.07	1.44	1.39
6	D	501	DOE	C20-C19	3.07	1.43	1.38
7	D	400	NAG	C4-C5	3.03	1.59	1.53
6	B	500	DOE	C4-N3	3.01	1.54	1.47
6	D	501	DOE	C22-C21	2.98	1.44	1.39
6	B	500	DOE	C23-C18	2.96	1.44	1.38
6	D	501	DOE	C8-N1	2.94	1.52	1.47
7	D	400	NAG	O7-C7	-2.91	1.16	1.23
6	D	501	DOE	C13-C14	2.89	1.57	1.51
6	D	501	DOE	C15-C16	2.84	1.57	1.51
6	B	500	DOE	C20-C21	2.83	1.44	1.39
6	D	501	DOE	O5-C14	2.76	1.31	1.22
6	B	500	DOE	C8-N1	2.69	1.51	1.47
6	B	500	DOE	C15-N4	2.67	1.52	1.47
6	D	501	DOE	C1-N1	2.60	1.55	1.48
6	D	501	DOE	C23-C22	2.58	1.43	1.38
6	B	500	DOE	C22-C21	2.54	1.43	1.39
6	B	500	DOE	C2-N2	2.54	1.52	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	500	DOE	C20-C19	2.49	1.42	1.38
6	B	500	DOE	O3-C12	2.36	1.29	1.22
6	D	501	DOE	O2-C10	-2.34	1.23	1.30
6	B	500	DOE	C9-C10	2.33	1.56	1.51
6	B	500	DOE	C11-N2	2.30	1.51	1.47
6	B	500	DOE	C21-N5	-2.30	1.37	1.41
6	D	501	DOE	C19-C18	2.25	1.43	1.38
6	D	501	DOE	O1-C10	2.20	1.29	1.22
6	D	501	DOE	C15-N4	2.17	1.51	1.47
6	B	500	DOE	O1-C10	2.16	1.29	1.22
6	D	501	DOE	O8-C16	-2.15	1.23	1.30
6	B	500	DOE	C23-C22	2.14	1.42	1.38
6	D	501	DOE	O7-C16	2.09	1.28	1.22
6	B	500	DOE	C13-N3	2.08	1.51	1.47

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	400	NAG	C2-N2-C7	5.25	129.94	122.90
6	B	500	DOE	C21-N5-C24	5.17	136.67	127.52
6	D	501	DOE	C21-N5-C24	4.09	134.76	127.52
6	B	500	DOE	C1-C2-N2	3.46	117.86	113.40
6	B	500	DOE	C9-N1-C1	-3.19	107.55	113.00
7	D	400	NAG	C4-C3-C2	-3.05	106.54	111.02
6	D	501	DOE	C5-N3-C4	-3.04	104.19	111.44
6	D	501	DOE	C9-N1-C1	-2.71	108.36	113.00
6	B	500	DOE	C7-N4-C6	-2.62	105.19	111.44
7	D	400	NAG	O7-C7-N2	2.62	126.61	121.98
6	D	501	DOE	C1-C2-N2	-2.61	110.04	113.40
7	D	400	NAG	C1-C2-N2	-2.27	106.86	110.43
6	B	500	DOE	C3-C4-N3	2.25	118.31	112.98
6	D	501	DOE	O9-C24-C25	-2.13	117.83	121.44
6	D	501	DOE	C13-N3-C5	-2.10	106.85	111.91
6	B	500	DOE	O3-C12-C11	-2.01	113.99	122.38

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	501	DOE	C24-C25-S1-C26
6	B	500	DOE	O9-C24-C25-S1
6	B	500	DOE	O2-C10-C9-N1

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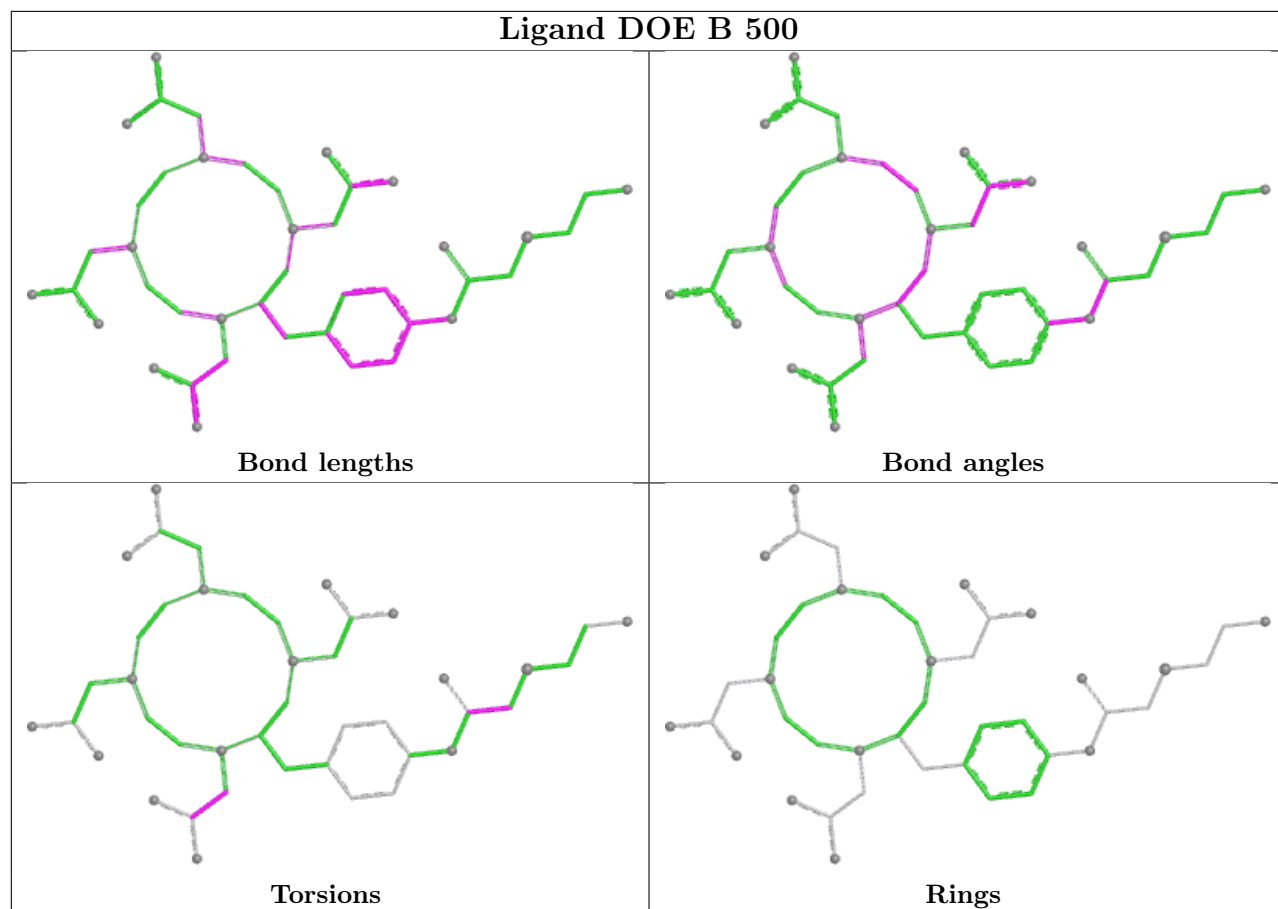
Mol	Chain	Res	Type	Atoms
6	B	500	DOE	N5-C24-C25-S1
6	D	501	DOE	N2-C11-C12-O4
6	B	500	DOE	O1-C10-C9-N1

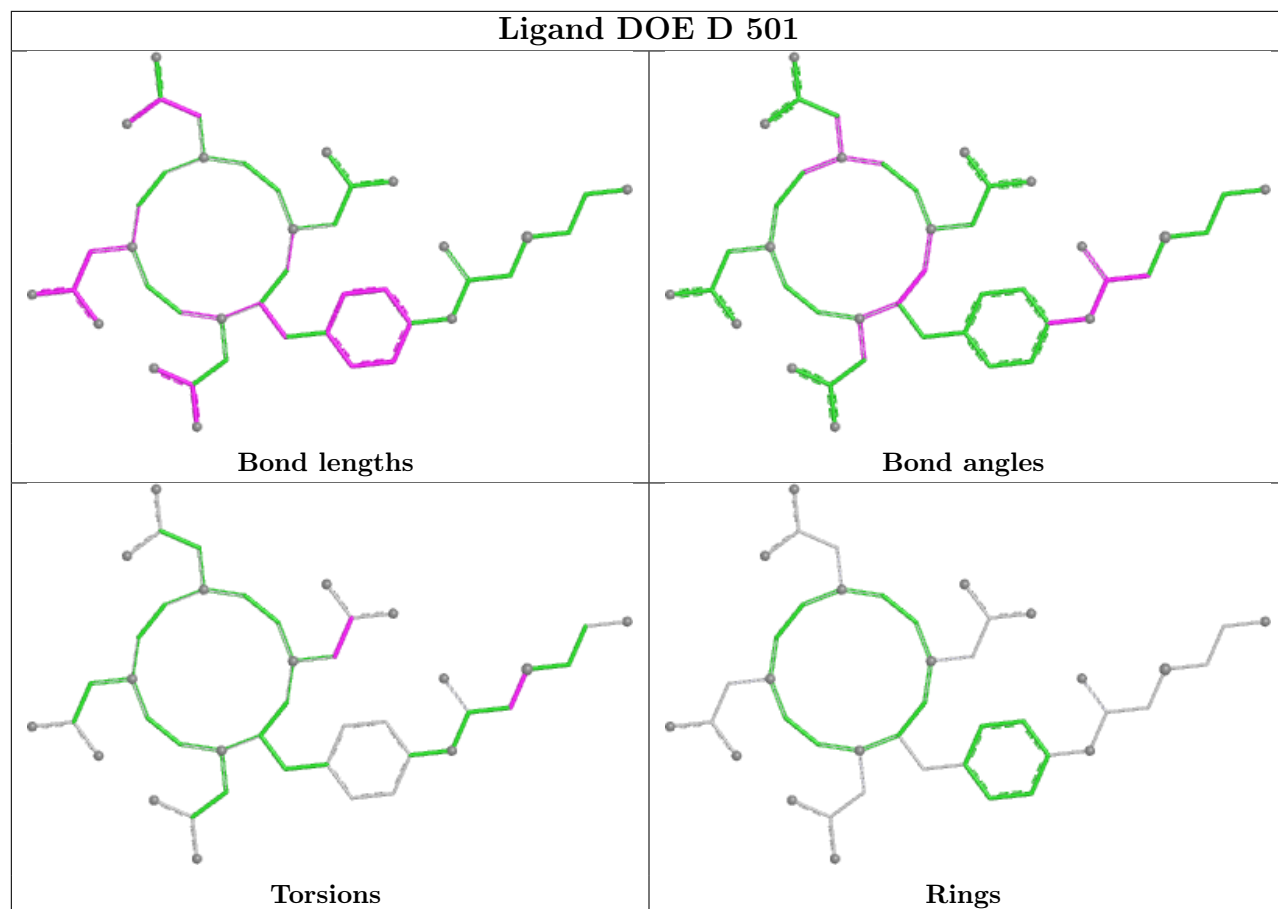
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	501	DOE	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	211/215 (98%)	0.55	21 (9%) 12 13	13, 29, 60, 72	0
1	C	215/215 (100%)	0.19	10 (4%) 36 38	15, 26, 42, 65	0
2	B	218/221 (98%)	0.35	15 (6%) 23 24	14, 26, 53, 58	0
3	D	215/221 (97%)	0.28	9 (4%) 40 42	14, 27, 49, 59	0
All	All	859/872 (98%)	0.34	55 (6%) 25 27	13, 27, 54, 72	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	158	VAL	5.3
1	A	161	GLY	5.1
1	C	212	ALA	4.2
2	B	41	PRO	4.1
1	A	156	THR	4.1
1	A	157	PRO	3.7
1	A	162	MET	3.7
1	A	211	PRO	3.6
1	A	159	THR	3.1
2	B	42	GLY	2.9
1	C	43	ASP	2.9
2	B	122	THR	2.8
1	A	185	ALA	2.8
3	D	42	GLY	2.8
2	B	176	LEU	2.8
2	B	111	GLN	2.8
2	B	133	GLY	2.8
1	C	159	THR	2.8
2	B	43	LYS	2.8
1	A	190	ARG	2.7
1	A	192	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	213	GLU	2.6
1	C	214	CYS	2.6
1	C	211	PRO	2.6
3	D	1	GLN	2.6
3	D	41	PRO	2.6
3	D	111	GLN	2.5
2	B	135	ALA	2.5
1	C	1	GLN	2.5
2	B	44	GLY	2.5
1	A	125	SER	2.5
2	B	9	PRO	2.5
1	A	194	TYR	2.4
1	A	130	THR	2.4
3	D	212	VAL	2.4
2	B	177	GLN	2.3
1	A	112	PRO	2.3
2	B	119	SER	2.3
1	A	128	LEU	2.3
1	C	190	ARG	2.3
3	D	140	SER	2.2
1	A	43	ASP	2.2
1	A	127	GLU	2.2
1	C	44	HIS	2.2
1	A	124	SER	2.2
1	A	189	GLU	2.2
2	B	124	PRO	2.1
2	B	217	VAL	2.1
1	A	203	HIS	2.1
2	B	196	SER	2.1
3	D	166	SER	2.1
3	D	198	THR	2.1
1	A	210	SER	2.0
3	D	192	SER	2.0
1	C	70	GLY	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($\text{\AA}^2$)	Q<0.9
2	PCA	B	1	8/9	0.88	0.09	23,24,27,29	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

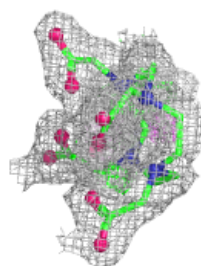
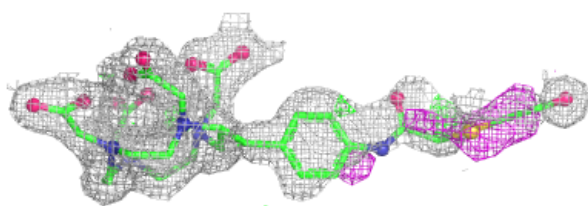
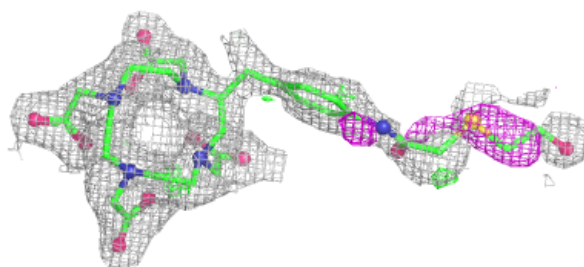
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
7	NAG	D	400	14/15	0.59	0.16	50,54,56,58	0
6	DOE	D	501	43/43	0.89	0.13	15,22,51,55	0
4	CL	A	504	1/1	0.89	0.14	65,65,65,65	0
6	DOE	B	500	43/43	0.91	0.12	14,23,65,68	0
4	CL	C	505	1/1	0.96	0.05	45,45,45,45	0
5	YT3	D	503	1/1	1.00	0.04	21,21,21,21	0
5	YT3	B	502	1/1	1.00	0.03	23,23,23,23	0

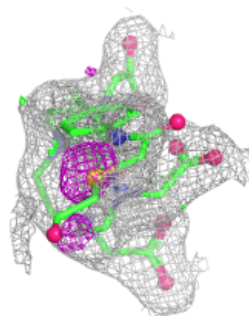
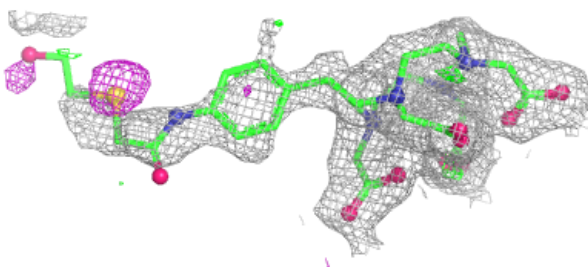
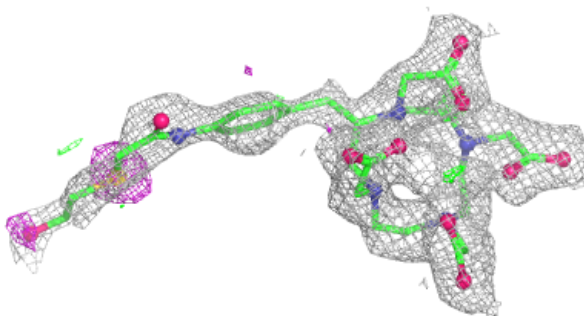
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around DOE 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)

**Electron density around DOE B 500:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.