



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 25, 2020 – 01:08 pm BST

PDB ID : 2BWO  
Title : 5-Aminolevulinate Synthase from *Rhodobacter capsulatus* in complex with succinyl-CoA  
Authors : Astner, I.; Schulze, J.O.; van den Heuvel, J.J.; Jahn, D.; Schubert, W.-D.; Heinz, D.W.  
Deposited on : 2005-07-15  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

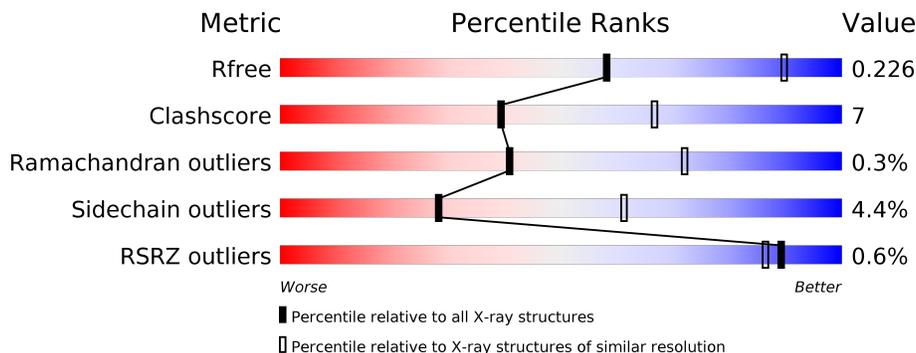
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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

Mol	Chain	Length	Quality of chain
1	A	401	 81% 16% ..
1	B	401	 84% 14% .
1	D	401	 79% 18% ..
1	E	401	 82% 17% .

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12602 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 5-AMINOLEVULINATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	398	3048	1936	539	558	15	0	0	0
1	B	399	3059	1942	543	559	15	0	0	0
1	D	397	3043	1933	538	557	15	0	0	0
1	E	399	3059	1942	543	559	15	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

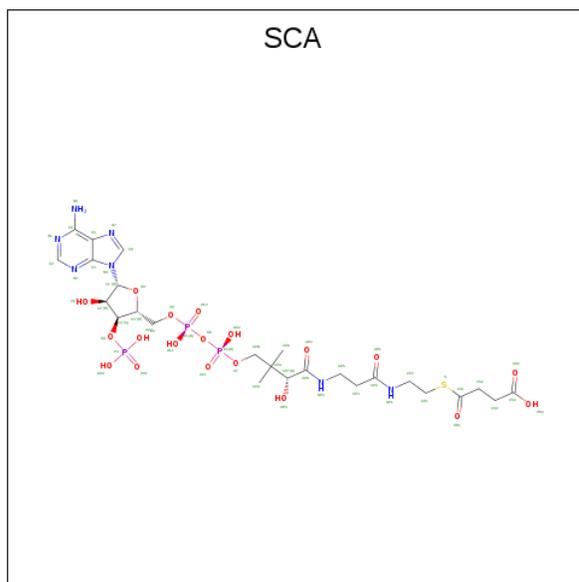
Chain	Residue	Modelled	Actual	Comment	Reference
A	102	GLY	ASP	variant	UNP P18079
A	105	GLN	GLY	variant	UNP P18079
A	117	ASN	ILE	variant	UNP P18079
A	128	VAL	LEU	variant	UNP P18079
A	205	GLU	ASP	variant	UNP P18079
A	262	ARG	LYS	variant	UNP P18079
B	102	GLY	ASP	variant	UNP P18079
B	105	GLN	GLY	variant	UNP P18079
B	117	ASN	ILE	variant	UNP P18079
B	128	VAL	LEU	variant	UNP P18079
B	205	GLU	ASP	variant	UNP P18079
B	262	ARG	LYS	variant	UNP P18079
D	102	GLY	ASP	variant	UNP P18079
D	105	GLN	GLY	variant	UNP P18079
D	117	ASN	ILE	variant	UNP P18079
D	128	VAL	LEU	variant	UNP P18079
D	205	GLU	ASP	variant	UNP P18079
D	262	ARG	LYS	variant	UNP P18079
E	102	GLY	ASP	variant	UNP P18079
E	105	GLN	GLY	variant	UNP P18079
E	117	ASN	ILE	variant	UNP P18079

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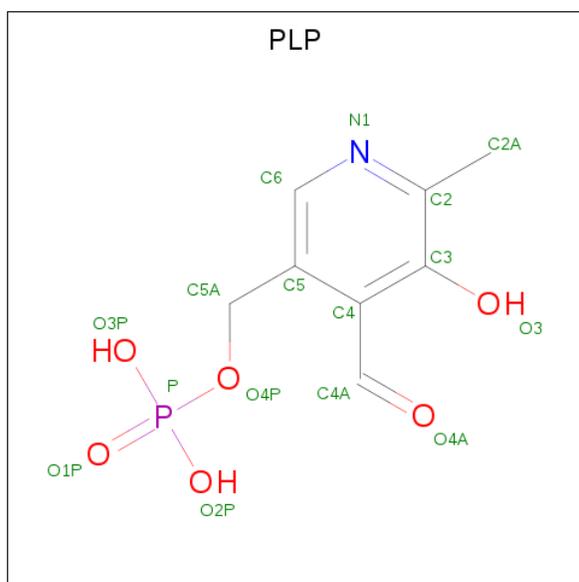
Chain	Residue	Modelled	Actual	Comment	Reference
E	128	VAL	LEU	variant	UNP P18079
E	205	GLU	ASP	variant	UNP P18079
E	262	ARG	LYS	variant	UNP P18079

- Molecule 2 is SUCCINYL-COENZYME A (three-letter code: SCA) (formula:  $C_{25}H_{40}N_7O_{19}P_3S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			S
2	A	1	Total	C	N	O	P	S	30	0
			55	25	7	19	3	1		
2	B	1	Total	C	N	O	P	S	25	0
			55	25	7	19	3	1		
2	D	1	Total	C	N	O	P	S	20	0
			55	25	7	19	3	1		
2	E	1	Total	C	N	O	P	S	24	0
			55	25	7	19	3	1		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula:  $C_8H_{10}NO_6P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	D	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	E	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

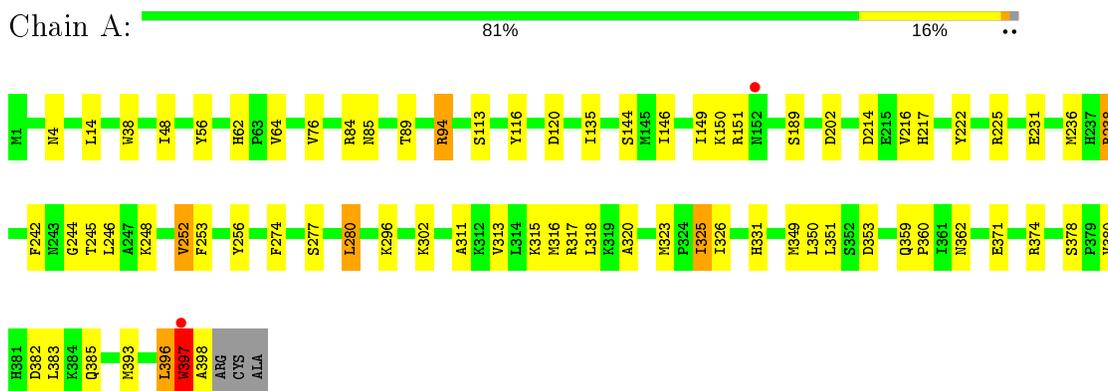
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	23	Total	O	0	0
			23	23		
4	D	30	Total	O	0	0
			30	30		
4	E	25	Total	O	0	0
			25	25		

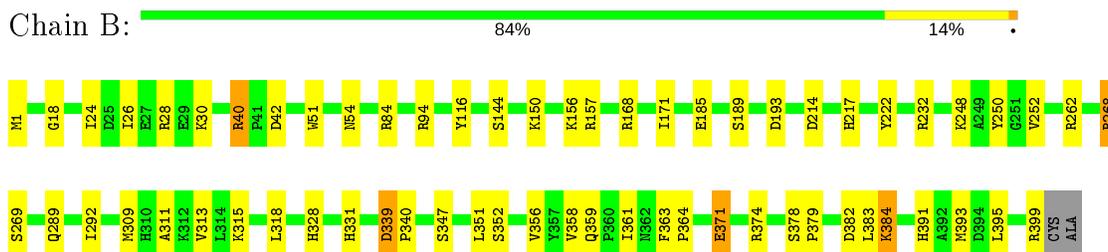
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

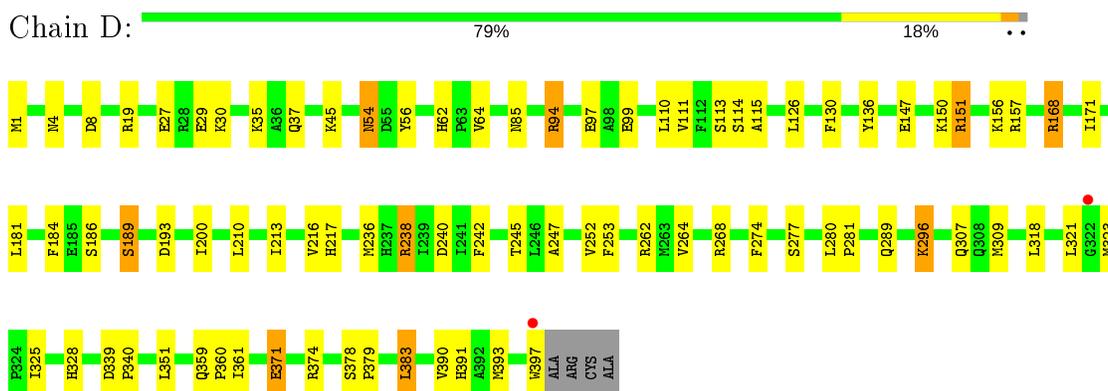
- Molecule 1: 5-AMINOLEVULINATE SYNTHASE



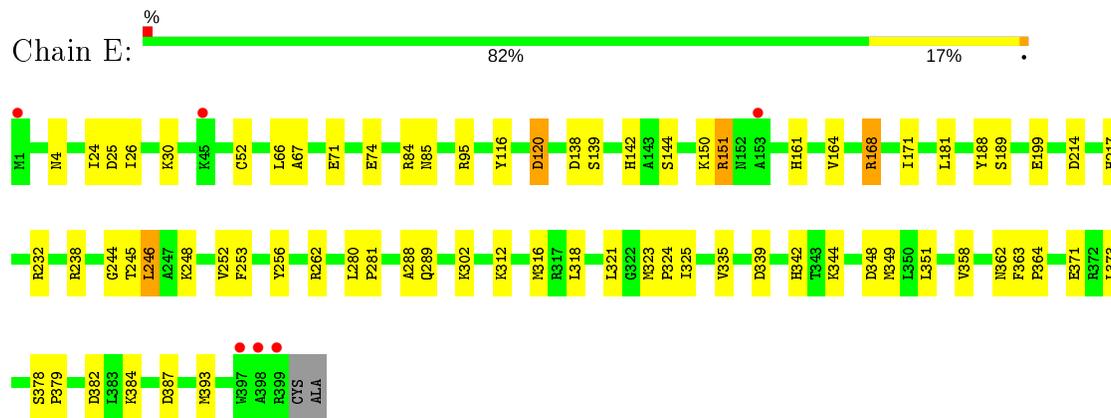
- Molecule 1: 5-AMINOLEVULINATE SYNTHASE



- Molecule 1: 5-AMINOLEVULINATE SYNTHASE



- Molecule 1: 5-AMINOLEVULINATE SYNTHASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.94Å 91.36Å 247.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	119.52 – 2.80 29.78 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.3 (119.52-2.80) 99.5 (29.78-2.80)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.24 (at 2.80Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.160 , 0.224 0.164 , 0.226	Depositor DCC
$R_{free}$ test set	1876 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.9	Xtrriage
Anisotropy	0.138	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 45.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	12602	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SCA, PLP

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	0.86	1/3118 (0.0%)	0.79	6/4225 (0.1%)
1	B	0.85	0/3129	0.78	3/4239 (0.1%)
1	D	0.88	2/3113 (0.1%)	0.81	2/4218 (0.0%)
1	E	0.86	0/3129	0.78	2/4239 (0.0%)
All	All	0.86	3/12489 (0.0%)	0.79	13/16921 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	27	GLU	CG-CD	7.30	1.62	1.51
1	A	248	LYS	CE-NZ	5.24	1.62	1.49
1	D	130	PHE	CE1-CZ	5.13	1.47	1.37

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	398	ALA	N-CA-C	6.16	127.64	111.00
1	D	19	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	A	374	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	D	168	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	E	120	ASP	CB-CG-OD1	5.62	123.36	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	396	LEU	Peptide
1	A	397	TRP	Peptide

## 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	3048	0	3028	44	0
1	B	3059	0	3043	42	0
1	D	3043	0	3025	54	0
1	E	3059	0	3043	56	0
2	A	55	0	31	0	0
2	B	55	0	31	1	0
2	D	55	0	35	1	0
2	E	55	0	31	0	0
3	A	15	0	6	0	0
3	B	16	0	8	3	0
3	D	16	0	8	3	0
3	E	16	0	8	1	0
4	A	32	0	0	0	0
4	B	23	0	0	0	0
4	D	30	0	0	1	0
4	E	25	0	0	2	0
All	All	12602	0	12297	183	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 7.

The worst 5 of 183 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:D:94:ARG:HG2	1:D:94:ARG:HH11	1.21	1.03
1:D:383:LEU:H	1:D:383:LEU:HD23	1.21	1.03
1:A:94:ARG:HH11	1:A:94:ARG:CG	1.73	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:ARG:HH11	1:A:94:ARG:HG3	1.34	0.91
1:E:151:ARG:HH11	1:E:151:ARG:CB	1.85	0.90

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	396/401 (99%)	380 (96%)	15 (4%)	1 (0%)	41 72
1	B	397/401 (99%)	384 (97%)	12 (3%)	1 (0%)	41 72
1	D	395/401 (98%)	380 (96%)	14 (4%)	1 (0%)	41 72
1	E	397/401 (99%)	382 (96%)	14 (4%)	1 (0%)	41 72
All	All	1585/1604 (99%)	1526 (96%)	55 (4%)	4 (0%)	41 72

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	252	VAL
1	D	252	VAL
1	E	252	VAL
1	B	252	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	311/313 (99%)	298 (96%)	13 (4%)	30	63
1	B	312/313 (100%)	300 (96%)	12 (4%)	33	67
1	D	311/313 (99%)	292 (94%)	19 (6%)	18	48
1	E	312/313 (100%)	301 (96%)	11 (4%)	36	70
All	All	1246/1252 (100%)	1191 (96%)	55 (4%)	28	61

5 of 55 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	399	ARG
1	D	189	SER
1	E	246	LEU
1	D	1	MET
1	D	94	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	307	GLN
1	B	331	HIS
1	D	391	HIS
1	A	385	GLN
1	D	37	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry

8 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	PLP	A	600	1	15,15,16	2.15	3 (20%)	20,22,23	2.26	7 (35%)
2	SCA	B	500	-	46,57,57	1.49	4 (8%)	58,84,84	1.98	12 (20%)
3	PLP	E	600	2	16,16,16	1.67	3 (18%)	20,23,23	1.05	0
2	SCA	D	500	-	46,57,57	1.33	5 (10%)	58,84,84	1.61	12 (20%)
3	PLP	B	600	-	16,16,16	1.49	2 (12%)	20,23,23	0.80	0
2	SCA	A	500	-	46,57,57	1.80	6 (13%)	58,84,84	3.44	13 (22%)
3	PLP	D	600	-	16,16,16	1.47	2 (12%)	20,23,23	0.94	1 (5%)
2	SCA	E	500	3	46,57,57	1.45	4 (8%)	58,84,84	2.27	9 (15%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PLP	A	600	1	-	2/6/6/8	0/1/1/1
2	SCA	B	500	-	-	28/50/72/72	0/3/3/3
3	PLP	E	600	2	-	2/8/8/8	0/1/1/1
2	SCA	D	500	-	-	26/50/72/72	0/3/3/3
3	PLP	B	600	-	-	4/8/8/8	0/1/1/1
2	SCA	A	500	-	-	15/50/72/72	0/3/3/3
3	PLP	D	600	-	-	4/8/8/8	0/1/1/1
2	SCA	E	500	3	-	21/50/72/72	0/3/3/3

The worst 5 of 29 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	SCA	CS3-CS2	-9.43	1.05	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	500	SCA	CS3-CS2	-7.19	1.16	1.52
2	B	500	SCA	CS3-CS2	-7.06	1.17	1.52
3	A	600	PLP	O3-C3	-5.37	1.24	1.37
2	D	500	SCA	O7-CPB	5.07	1.60	1.43

The worst 5 of 54 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	SCA	CS2-CS3-CS4	20.26	146.66	112.67
2	E	500	SCA	CS2-CS3-CS4	11.88	132.60	112.67
2	A	500	SCA	CS3-CS2-CS1	10.44	128.10	112.41
2	B	500	SCA	CS2-CS3-CS4	9.45	128.53	112.67
2	E	500	SCA	CS3-CS2-CS1	6.86	122.73	112.41

There are no chirality outliers.

5 of 102 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	500	SCA	C3'-O3'-P3-O31
2	B	500	SCA	C5'-O5'-P1-O12
2	B	500	SCA	CPB-O7-P2-O6
2	B	500	SCA	CPA-CPB-O7-P2
2	B	500	SCA	CP7-CPA-CPB-O7

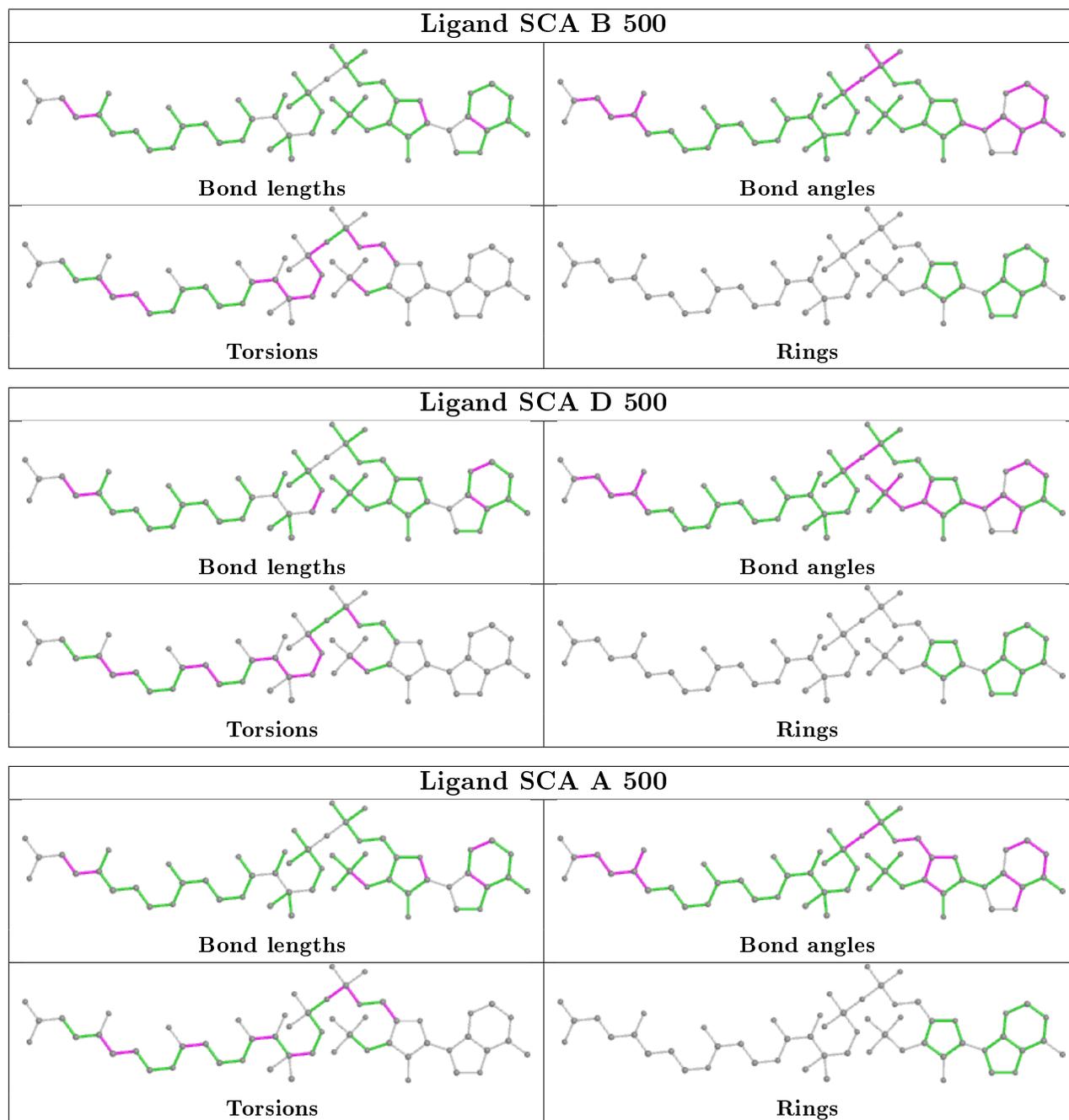
There are no ring outliers.

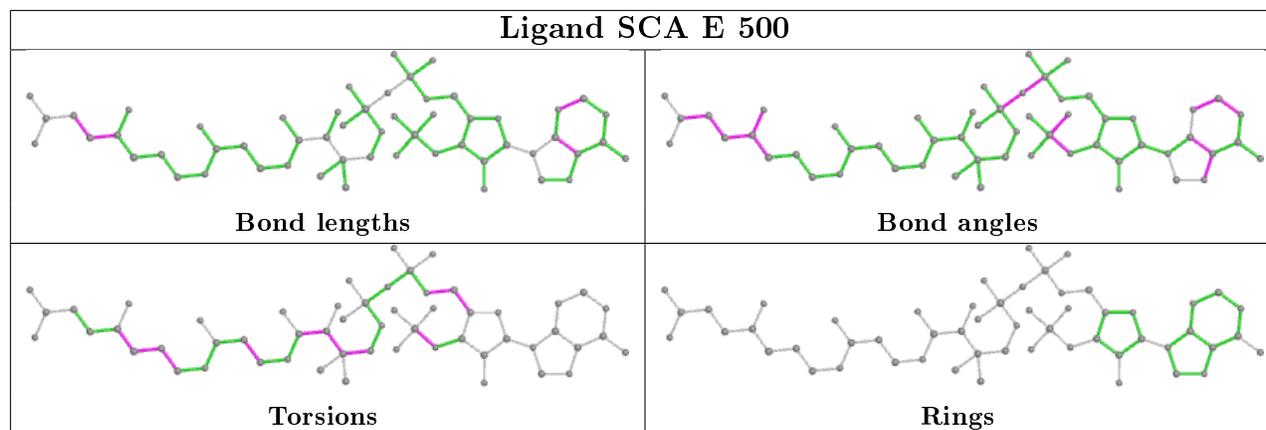
5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	SCA	1	0
3	E	600	PLP	1	0
2	D	500	SCA	1	0
3	B	600	PLP	3	0
3	D	600	PLP	3	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å <sup>2</sup> )	Q < 0.9
1	A	398/401 (99%)	-0.49	2 (0%) 91 88	13, 24, 45, 77	0
1	B	399/401 (99%)	-0.39	0 100 100	13, 24, 46, 85	0
1	D	397/401 (99%)	-0.54	2 (0%) 91 88	13, 24, 46, 77	0
1	E	399/401 (99%)	-0.45	6 (1%) 73 68	13, 24, 45, 91	0
All	All	1593/1604 (99%)	-0.47	10 (0%) 89 86	13, 24, 46, 91	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	1	MET	2.7
1	E	398	ALA	2.7
1	A	397	TRP	2.7
1	E	153	ALA	2.6
1	E	397	TRP	2.5

### 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](#)

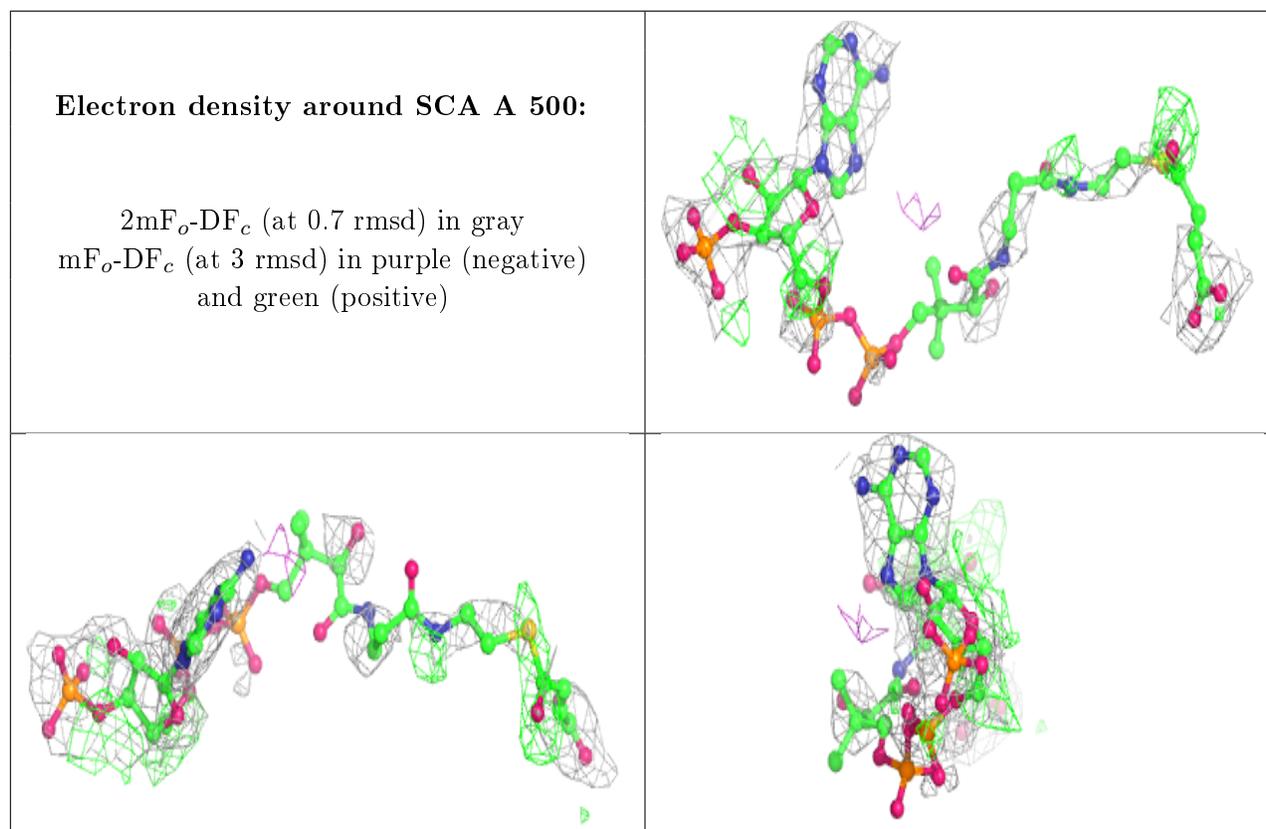
There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

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

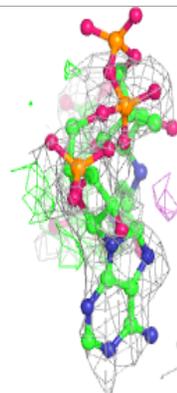
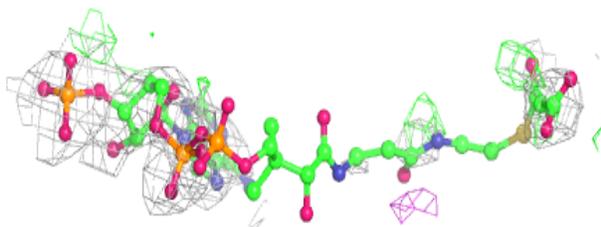
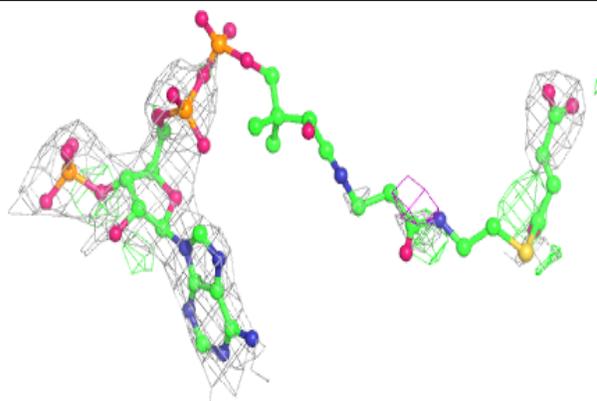
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SCA	A	500	55/55	0.89	0.20	18,80,121,122	55
2	SCA	E	500	55/55	0.89	0.20	33,79,130,131	55
2	SCA	B	500	55/55	0.90	0.19	28,71,113,114	55
2	SCA	D	500	55/55	0.91	0.18	29,62,100,101	55
3	PLP	E	600	16/16	0.96	0.16	29,46,60,67	0
3	PLP	D	600	16/16	0.97	0.12	33,39,59,71	0
3	PLP	B	600	16/16	0.98	0.16	38,47,67,73	0
3	PLP	A	600	15/16	0.98	0.11	32,34,38,40	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.

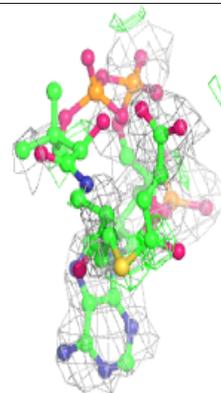
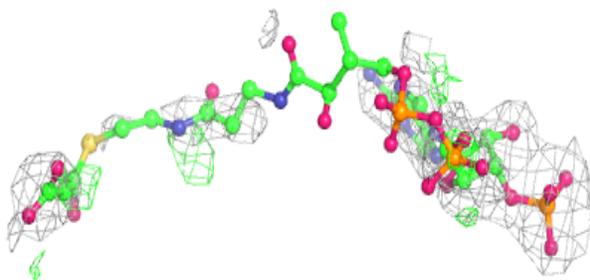
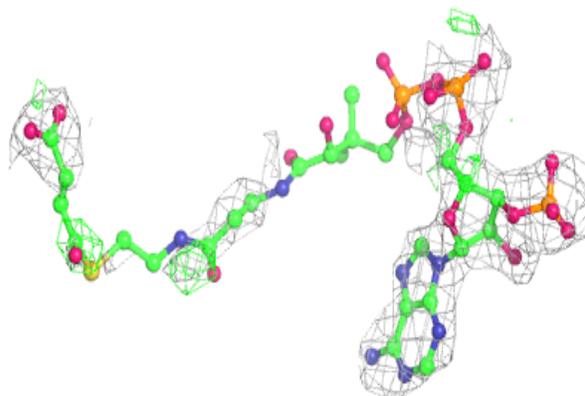


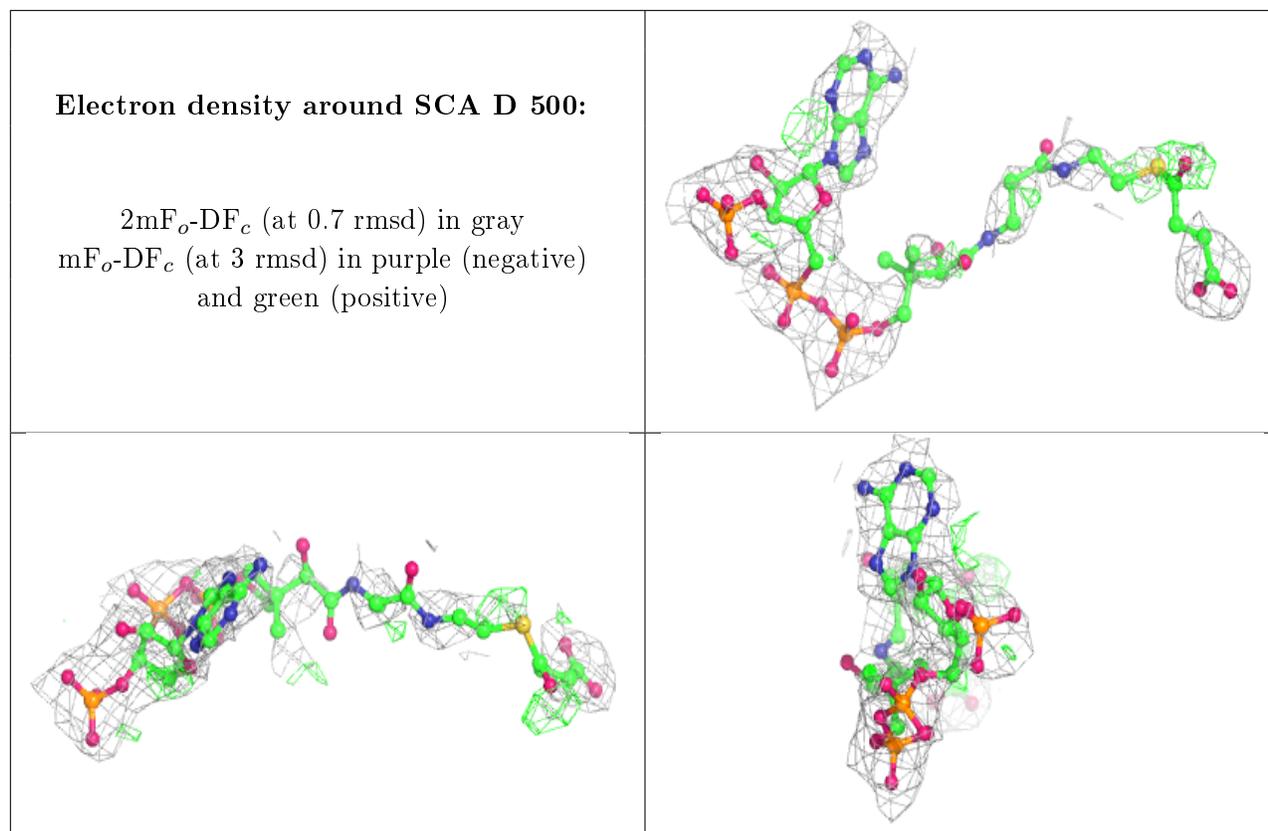
**Electron density around SCA E 500:**

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

**Electron density around SCA 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 [i](#)

There are no such residues in this entry.