Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2023 – 06:46 PM EDT

PDB ID	:	1OTW
Title	:	Crystal structure of PqqC in complex with PQQ and a putative H2O2
Authors	:	Magnusson, O.T.; Toyama, H.; Saeki, M.; Rojas, A.; Reed, J.C.; Liddington, R.C.; Klinman, J.P.; Schwarzenbacher, R.
Deposited on	:	2003-03-23
Resolution	:	2.30 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:

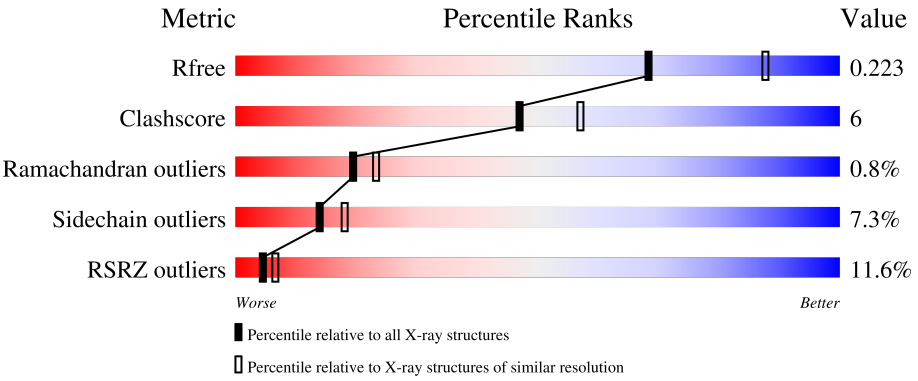
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 <=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	255	4% 88% 11%
1	B	255	20% 70% 24% 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	PEO	A	600	-	X	-	-
3	PEO	B	601	-	X	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4335 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.

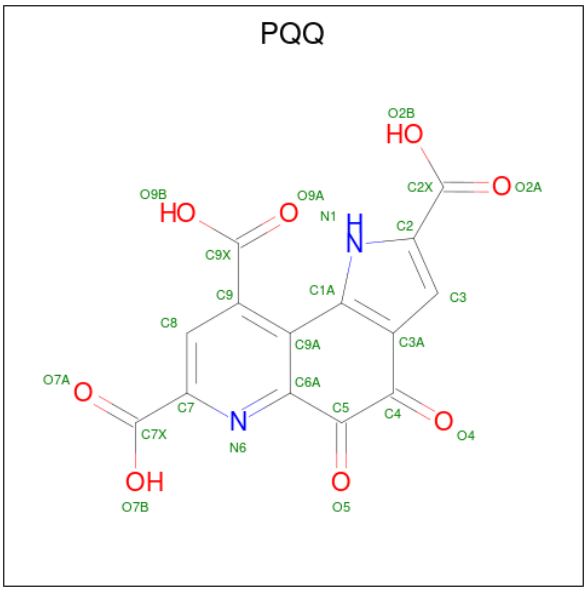
- Molecule 1 is a protein called Coenzyme PQQ synthesis protein C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	24	0	0
			2086	1323	379	373	11			
1	B	253	Total	C	N	O	S	85	0	0
			2068	1312	375	371	10			

There are 10 discrepancies between the modelled and reference sequences:

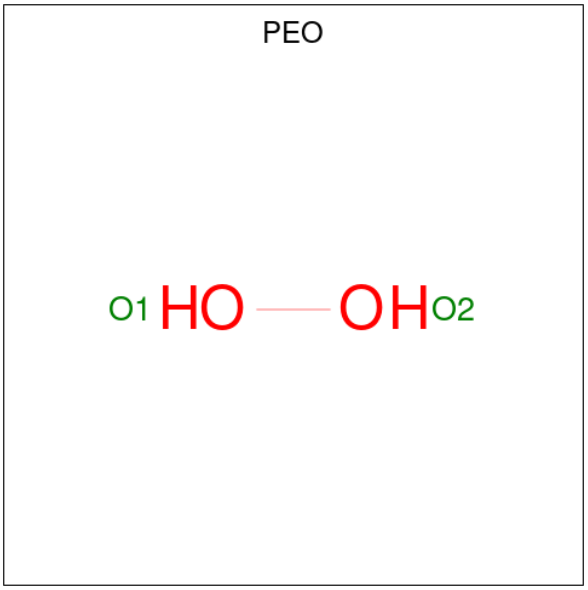
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	cloning artifact	UNP P27505
A	21	ASP	ALA	engineered mutation	UNP P27505
A	252	LEU	-	expression tag	UNP P27505
A	253	GLU	-	expression tag	UNP P27505
A	254	HIS	-	expression tag	UNP P27505
B	0	HIS	-	cloning artifact	UNP P27505
B	21	ASP	ALA	engineered mutation	UNP P27505
B	252	LEU	-	expression tag	UNP P27505
B	253	GLU	-	expression tag	UNP P27505
B	254	HIS	-	expression tag	UNP P27505

- Molecule 2 is PYRROLOQUINOLINE QUINONE (three-letter code: PQQ) (formula: $C_{14}H_6N_2O_8$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			24	14	2	8		
2	B	1	Total	C	N	O	0	0
			24	14	2	8		

- Molecule 3 is HYDROGEN PEROXIDE (three-letter code: PEO) (formula: H_2O_2).



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

- Molecule 4 is water.

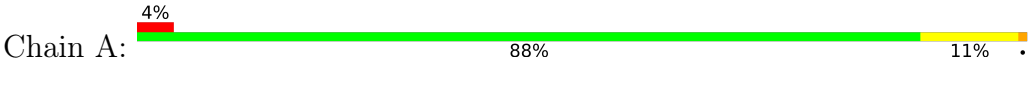
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	94	Total O 94 94	0	0
4	B	35	Total O 35 35	0	0

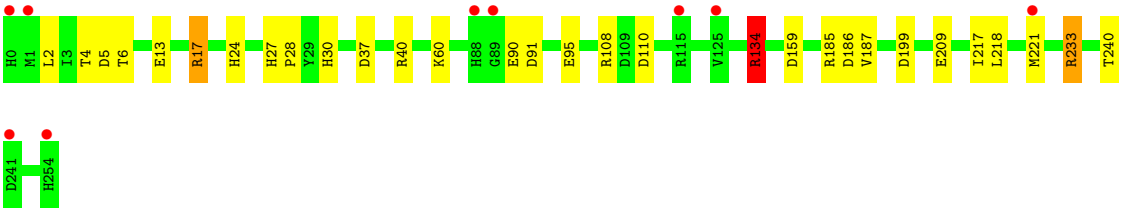
3 Residue-property plots (i)

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

• Molecule 1: Coenzyme PQQ synthesis protein C

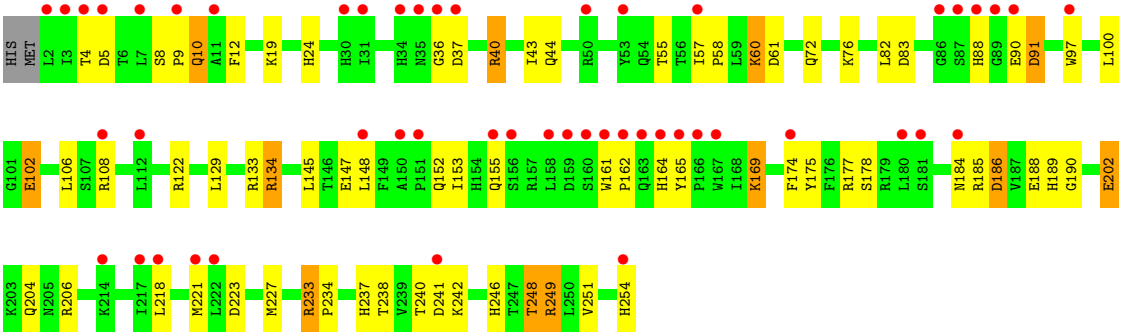
Chain A:





• Molecule 1: Coenzyme PQQ synthesis protein C

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	73.50Å 118.42Å 70.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.71 – 2.30 27.88 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.4 (70.71-2.30) 93.5 (27.88-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$< I/\sigma(I) >$ [1]	5.95 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.205 , 0.248 0.219 , 0.223	Depositor DCC
R_{free} test set	1311 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	36.6	Xtriage
Anisotropy	0.988	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 42.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	0.010 for l,-k,h	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4335	wwPDB-VP
Average B, all atoms (Å^2)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: PQQ, PEO

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.85	0/2146	0.90	10/2911 (0.3%)
1	B	0.67	0/2127	0.87	9/2886 (0.3%)
All	All	0.77	0/4273	0.89	19/5797 (0.3%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	17	ARG	NE-CZ-NH2	-10.69	114.95	120.30
1	A	17	ARG	NE-CZ-NH1	10.08	125.34	120.30
1	B	88	HIS	N-CA-C	7.65	131.67	111.00
1	B	223	ASP	CB-CG-OD2	7.62	125.16	118.30
1	B	186	ASP	CB-CG-OD2	7.12	124.71	118.30
1	A	37	ASP	CB-CG-OD2	6.31	123.98	118.30
1	B	83	ASP	CB-CG-OD2	6.14	123.83	118.30
1	A	91	ASP	CB-CG-OD2	5.76	123.48	118.30
1	B	88	HIS	CA-C-N	5.73	127.66	116.20
1	B	5	ASP	CB-CG-OD2	5.61	123.35	118.30
1	B	61	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	186	ASP	CB-CG-OD2	5.43	123.18	118.30
1	A	134	ARG	CG-CD-NE	5.41	123.17	111.80
1	B	37	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	17	ARG	CD-NE-CZ	5.25	130.95	123.60
1	A	110	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	159	ASP	CB-CG-OD2	5.04	122.83	118.30
1	B	108	ARG	NE-CZ-NH2	-5.02	117.79	120.30
1	A	199	ASP	CB-CG-OD1	5.02	122.82	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2086	0	2008	13	0
1	B	2068	0	1989	38	0
2	A	24	0	3	0	0
2	B	24	0	3	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	94	0	0	5	0
4	B	35	0	0	3	0
All	All	4335	0	4003	49	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 6.

All (49) 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:202:GLU:OE1	1:B:202:GLU:HA	1.73	0.87
1:B:240:THR:HG22	1:B:242:LYS:H	1.50	0.75
1:B:251:VAL:HG22	4:B:624:HOH:O	1.90	0.70
1:A:95:GLU:OE2	1:A:108:ARG:NH2	2.30	0.65
1:B:134:ARG:HG2	1:B:134:ARG:HH11	1.62	0.64
1:B:97:TRP:CE2	1:B:100:LEU:HD23	2.32	0.64
1:A:40:ARG:HD3	4:A:675:HOH:O	1.99	0.62
1:A:30:HIS:HD2	4:A:637:HOH:O	1.83	0.61
1:B:12:PHE:CD1	1:B:204:GLN:HG2	2.38	0.59
1:B:240:THR:HG22	1:B:241:ASP:N	2.19	0.58
1:B:76:LYS:HD3	1:B:189:HIS:CE1	2.38	0.58
1:B:134:ARG:HH11	1:B:134:ARG:CG	2.17	0.56
1:B:40:ARG:O	1:B:44:GLN:HG3	2.06	0.55
1:A:209:GLU:OE2	1:B:248:THR:CG2	2.55	0.55
1:B:91:ASP:OD1	1:B:91:ASP:N	2.40	0.55
1:B:233:ARG:N	1:B:234:PRO:CD	2.71	0.54
1:A:17:ARG:NH2	4:A:639:HOH:O	2.42	0.52
1:B:184:ASN:O	1:B:188:GLU:HG3	2.10	0.52
1:B:57:ILE:N	1:B:58:PRO:CD	2.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:SER:O	1:B:10:GLN:N	2.43	0.51
1:B:19:LYS:NZ	4:B:622:HOH:O	2.42	0.50
1:A:13:GLU:O	1:A:17:ARG:HG3	2.11	0.50
1:A:108:ARG:HG3	4:A:608:HOH:O	2.11	0.49
1:B:218:LEU:HD23	1:B:221:MET:HE1	1.96	0.48
1:B:145:LEU:O	1:B:148:LEU:HD12	2.14	0.47
1:B:202:GLU:OE1	1:B:202:GLU:CA	2.54	0.47
1:A:233:ARG:HD2	4:A:644:HOH:O	2.14	0.46
1:A:217:ILE:O	1:A:221:MET:HG3	2.15	0.46
1:A:218:LEU:HA	1:A:221:MET:HE3	1.97	0.46
1:B:147:GLU:OE1	1:B:186:ASP:OD2	2.34	0.46
1:B:227:MET:O	1:B:233:ARG:HB2	2.16	0.45
1:B:174:PHE:O	1:B:177:ARG:N	2.51	0.44
1:A:27:HIS:ND1	1:A:28:PRO:HD2	2.32	0.44
1:B:161:TRP:N	1:B:162:PRO:CD	2.80	0.44
1:B:82:LEU:HD23	1:B:82:LEU:HA	1.89	0.43
1:A:134:ARG:HD3	4:B:629:HOH:O	2.18	0.43
1:B:174:PHE:O	1:B:175:TYR:C	2.56	0.43
1:B:60:LYS:C	1:B:60:LYS:HD2	2.40	0.42
1:B:177:ARG:O	1:B:178:SER:C	2.59	0.41
1:B:43:ILE:HD11	1:B:165:TYR:CE2	2.55	0.41
1:B:129:LEU:O	1:B:133:ARG:HG3	2.21	0.41
1:B:237:HIS:CE1	1:B:238:THR:HG23	2.56	0.41
1:B:240:THR:CG2	1:B:241:ASP:N	2.84	0.41
1:B:36:GLY:HA3	1:B:164:HIS:HB3	2.02	0.41
1:B:189:HIS:O	1:B:190:GLY:C	2.58	0.41
1:B:55:THR:OG1	1:B:249:ARG:NH2	2.53	0.40
1:B:122:ARG:HD2	1:B:246:HIS:NE2	2.36	0.40
1:A:209:GLU:OE2	1:B:248:THR:HG23	2.20	0.40
1:B:102:GLU:HA	1:B:106:LEU:O	2.21	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	253/255 (99%)	247 (98%)	5 (2%)	1 (0%)	34 42
1	B	251/255 (98%)	231 (92%)	17 (7%)	3 (1%)	13 14
All	All	504/510 (99%)	478 (95%)	22 (4%)	4 (1%)	19 23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	169	LYS
1	B	249	ARG
1	A	6	THR
1	B	9	PRO

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	214/214 (100%)	203 (95%)	11 (5%)	24 33
1	B	212/214 (99%)	192 (91%)	20 (9%)	8 10
All	All	426/428 (100%)	395 (93%)	31 (7%)	14 18

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	4	THR
1	A	5	ASP
1	A	24	HIS
1	A	60	LYS
1	A	90	GLU
1	A	134	ARG
1	A	185	ARG
1	A	187	VAL

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Mol	Chain	Res	Type
1	A	233	ARG
1	A	240	THR
1	B	4	THR
1	B	10	GLN
1	B	24	HIS
1	B	40	ARG
1	B	60	LYS
1	B	72	GLN
1	B	90	GLU
1	B	91	ASP
1	B	102	GLU
1	B	134	ARG
1	B	152	GLN
1	B	153	ILE
1	B	155	GLN
1	B	169	LYS
1	B	185	ARG
1	B	202	GLU
1	B	206	ARG
1	B	233	ARG
1	B	248	THR
1	B	254	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	HIS
1	A	254	HIS
1	B	189	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

4 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	PEO	B	601	-	1,1,1	3.18	1 (100%)	-		
3	PEO	A	600	-	1,1,1	3.57	1 (100%)	-		
2	PQQ	B	501	-	23,26,26	1.86	7 (30%)	29,40,40	1.67	9 (31%)
2	PQQ	A	500	-	23,26,26	1.60	4 (17%)	29,40,40	2.28	9 (31%)

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	PQQ	B	501	-	-	0/10/28/28	0/3/3/3
2	PQQ	A	500	-	-	0/10/28/28	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	PQQ	C6A-C9A	5.06	1.48	1.42
2	A	500	PQQ	C6A-C9A	4.46	1.47	1.42
2	B	501	PQQ	C9-C9A	3.68	1.49	1.41
3	A	600	PEO	O2-O1	-3.57	1.06	1.42
3	B	601	PEO	O2-O1	-3.18	1.10	1.42
2	B	501	PQQ	C9A-C1A	2.89	1.48	1.42
2	A	500	PQQ	C9-C9A	2.89	1.47	1.41
2	A	500	PQQ	C9A-C1A	2.84	1.48	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	PQQ	C2-C2X	-2.54	1.47	1.50
2	B	501	PQQ	C6A-C5	-2.39	1.47	1.50
2	A	500	PQQ	O5-C5	2.29	1.28	1.23
2	B	501	PQQ	C7-C7X	-2.29	1.47	1.50
2	B	501	PQQ	C5-C4	-2.10	1.46	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	PQQ	O5-C5-C6A	-5.04	116.50	121.84
2	A	500	PQQ	O2B-C2X-C2	4.65	125.28	114.69
2	A	500	PQQ	C9-C9A-C6A	-4.21	116.05	121.68
2	A	500	PQQ	C9-C9A-C1A	4.05	126.20	122.88
2	A	500	PQQ	O2B-C2X-O2A	-3.93	114.63	123.35
2	A	500	PQQ	O5-C5-C4	3.83	125.95	119.31
2	B	501	PQQ	C9-C9A-C1A	3.46	125.72	122.88
2	B	501	PQQ	O7A-C7X-C7	-3.10	114.96	121.24
2	A	500	PQQ	C6A-N6-C7	3.02	122.99	117.91
2	B	501	PQQ	C9-C9A-C6A	-2.69	118.08	121.68
2	B	501	PQQ	O7B-C7X-C7	2.68	120.79	114.69
2	B	501	PQQ	O5-C5-C6A	-2.63	119.05	121.84
2	B	501	PQQ	O9B-C9X-C9	2.39	121.16	114.39
2	B	501	PQQ	C8-C9-C9A	-2.28	116.96	120.06
2	B	501	PQQ	C2-C3-C3A	-2.12	103.17	105.98
2	A	500	PQQ	O7B-C7X-C7	2.07	119.40	114.69
2	B	501	PQQ	C3A-C4-C5	-2.06	116.94	118.14
2	A	500	PQQ	C3A-C4-C5	2.02	119.31	118.14

There are no chirality outliers.

There are no torsion outliers.

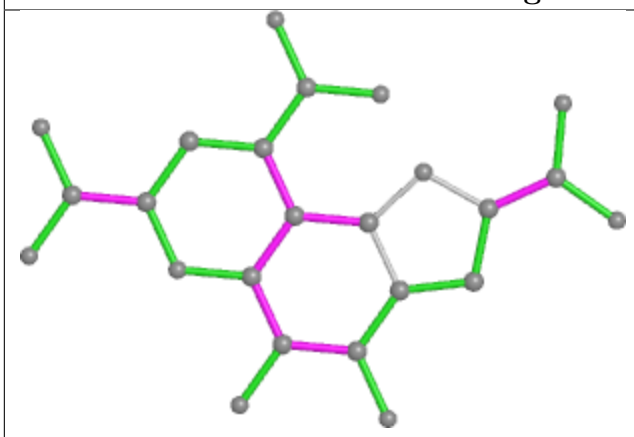
There are no ring outliers.

No monomer is involved in short contacts.

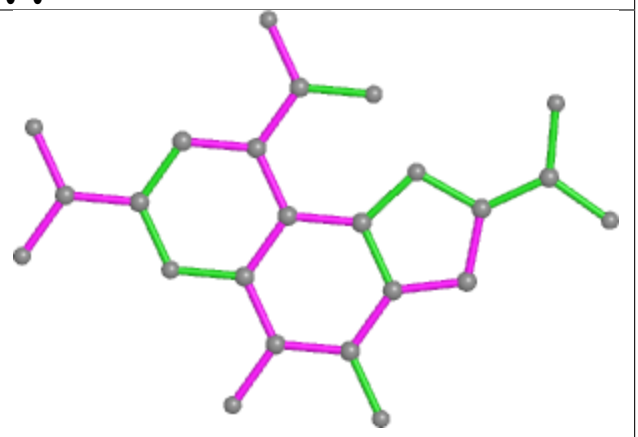
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

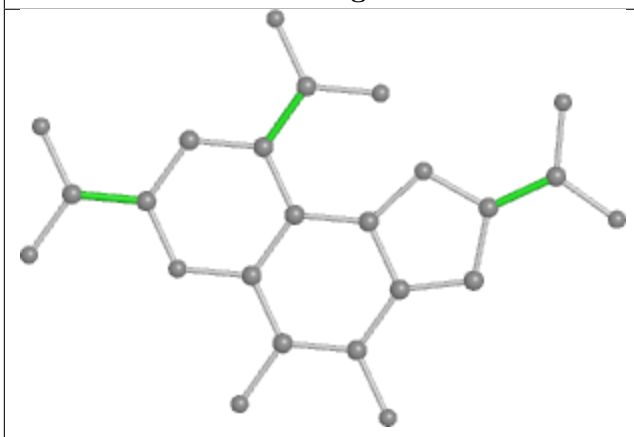
Ligand PQQ B 501



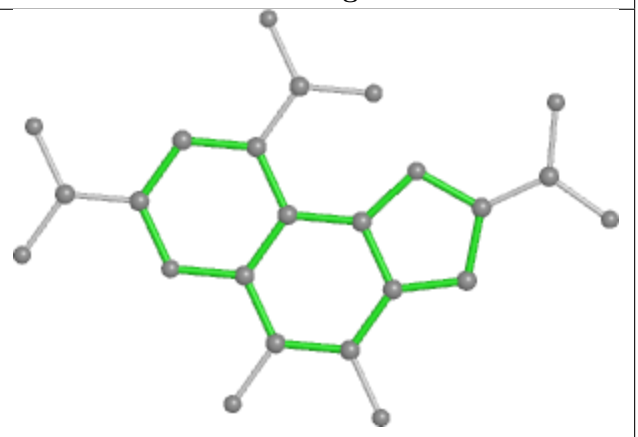
Bond lengths



Bond angles

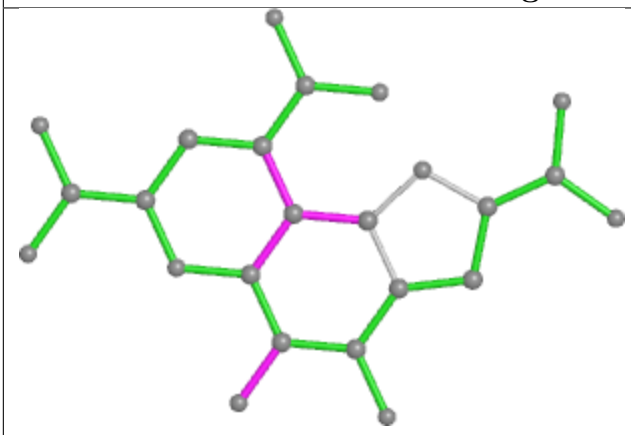


Torsions

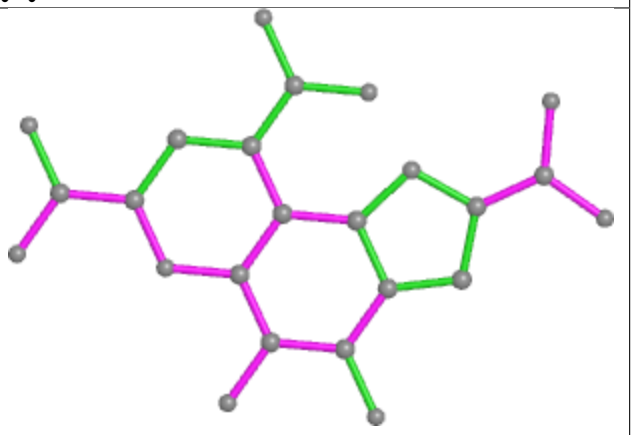


Rings

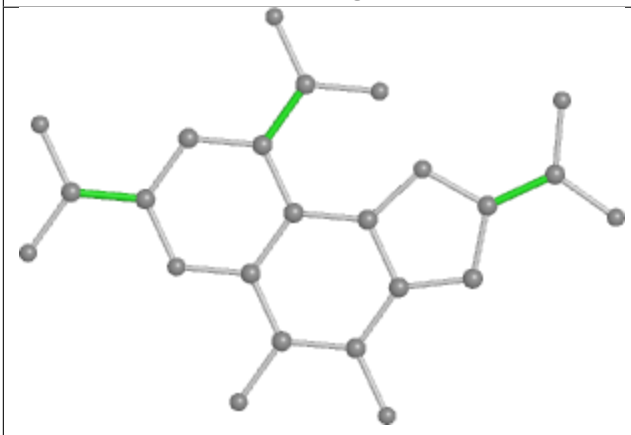
Ligand PQQ A 500



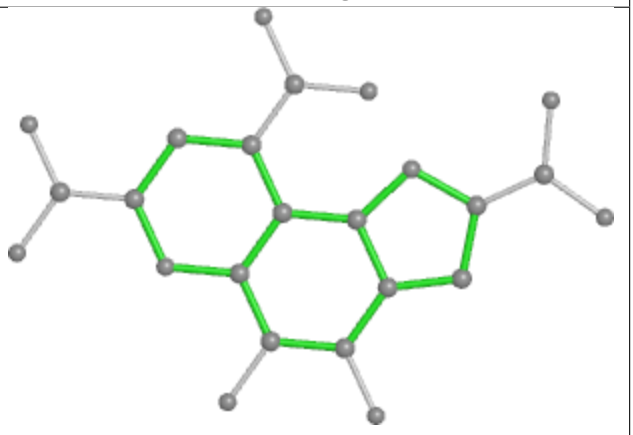
Bond lengths



Bond angles



Torsions



Rings

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, 95^{th} 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	255/255 (100%)	0.03	9 (3%) 44 51	24, 34, 54, 77	5 (1%)
1	B	253/255 (99%)	1.00	50 (19%) 1 1	39, 57, 78, 94	17 (6%)
All	All	508/510 (99%)	0.51	59 (11%) 4 6	24, 44, 75, 94	22 (4%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	89	GLY	6.3
1	B	160	SER	5.8
1	A	88	HIS	5.4
1	B	88	HIS	5.2
1	B	254	HIS	5.1
1	A	89	GLY	4.8
1	B	53	TYR	4.8
1	B	87	SER	4.7
1	B	161	TRP	4.3
1	B	35	ASN	4.2
1	B	156	SER	4.0
1	A	1	MET	3.8
1	B	97	TRP	3.8
1	B	218	LEU	3.7
1	B	166	PRO	3.7
1	A	0	HIS	3.5
1	B	180	LEU	3.4
1	B	3	ILE	3.4
1	B	50	ARG	3.3
1	B	34	HIS	3.3
1	B	159	ASP	3.2
1	B	36	GLY	3.1
1	B	221	MET	3.1
1	B	174	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	181	SER	3.1
1	B	184	ASN	3.0
1	B	4	THR	3.0
1	B	163	GLN	3.0
1	B	151	PRO	3.0
1	A	254	HIS	3.0
1	B	155	GLN	3.0
1	B	31	ILE	2.9
1	B	165	TYR	2.9
1	B	57	ILE	2.8
1	B	217	ILE	2.8
1	B	222	LEU	2.7
1	A	241	ASP	2.7
1	A	115	ARG	2.6
1	B	90	GLU	2.6
1	B	5	ASP	2.5
1	B	167	TRP	2.5
1	B	148	LEU	2.4
1	B	158	LEU	2.4
1	A	221	MET	2.4
1	B	86	GLY	2.4
1	B	7	LEU	2.3
1	B	112	LEU	2.3
1	B	108	ARG	2.3
1	B	164	HIS	2.3
1	B	9	PRO	2.2
1	B	2	LEU	2.2
1	B	162	PRO	2.1
1	B	214	LYS	2.1
1	A	125	VAL	2.1
1	B	150	ALA	2.1
1	B	241	ASP	2.1
1	B	11	ALA	2.1
1	B	37	ASP	2.0
1	B	30	HIS	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

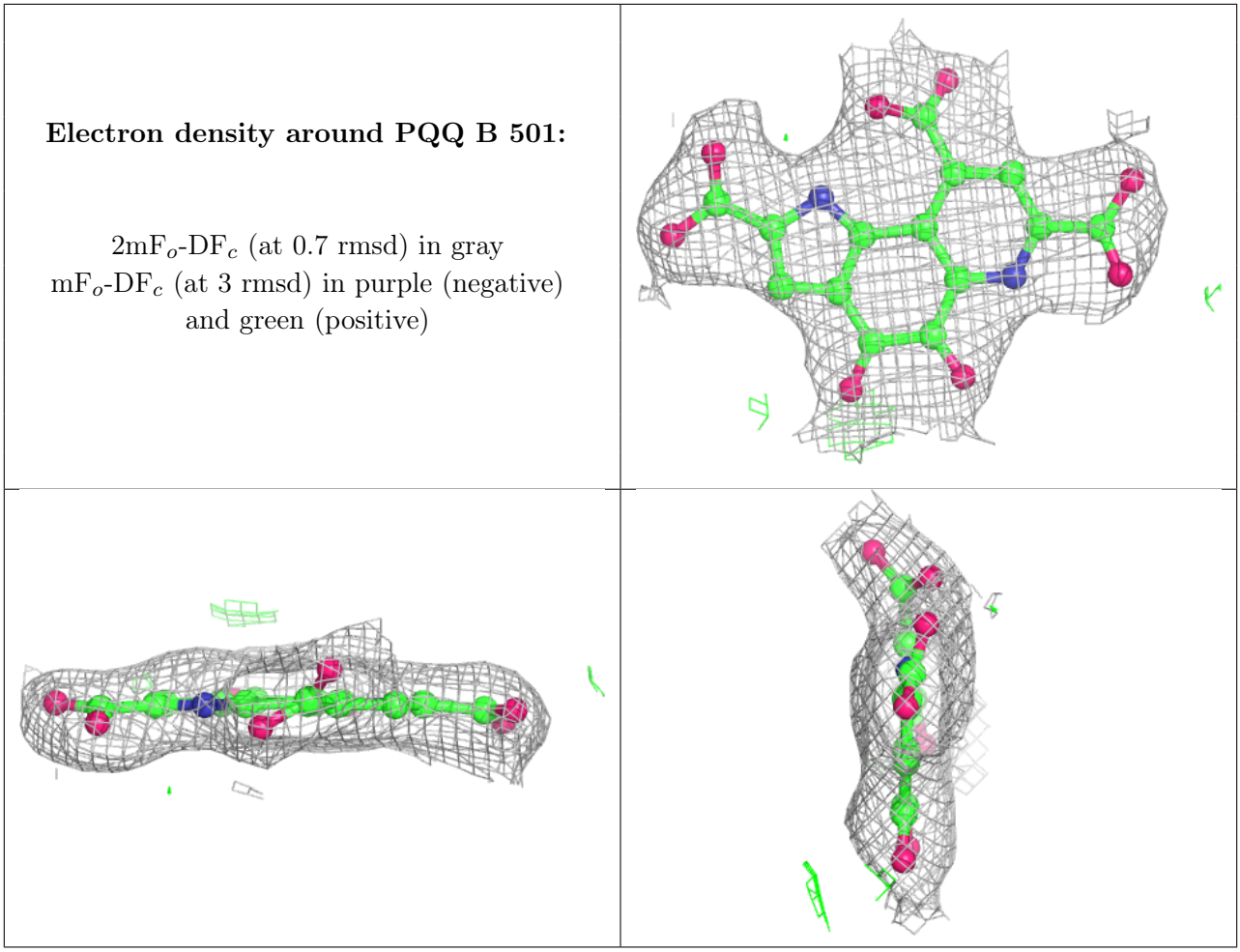
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
3	PEO	B	601	2/2	0.82	0.27	52,52,52,53	0
2	PQQ	B	501	24/24	0.93	0.20	56,58,60,60	0
2	PQQ	A	500	24/24	0.97	0.13	27,30,32,33	0
3	PEO	A	600	2/2	0.99	0.05	27,27,27,27	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.

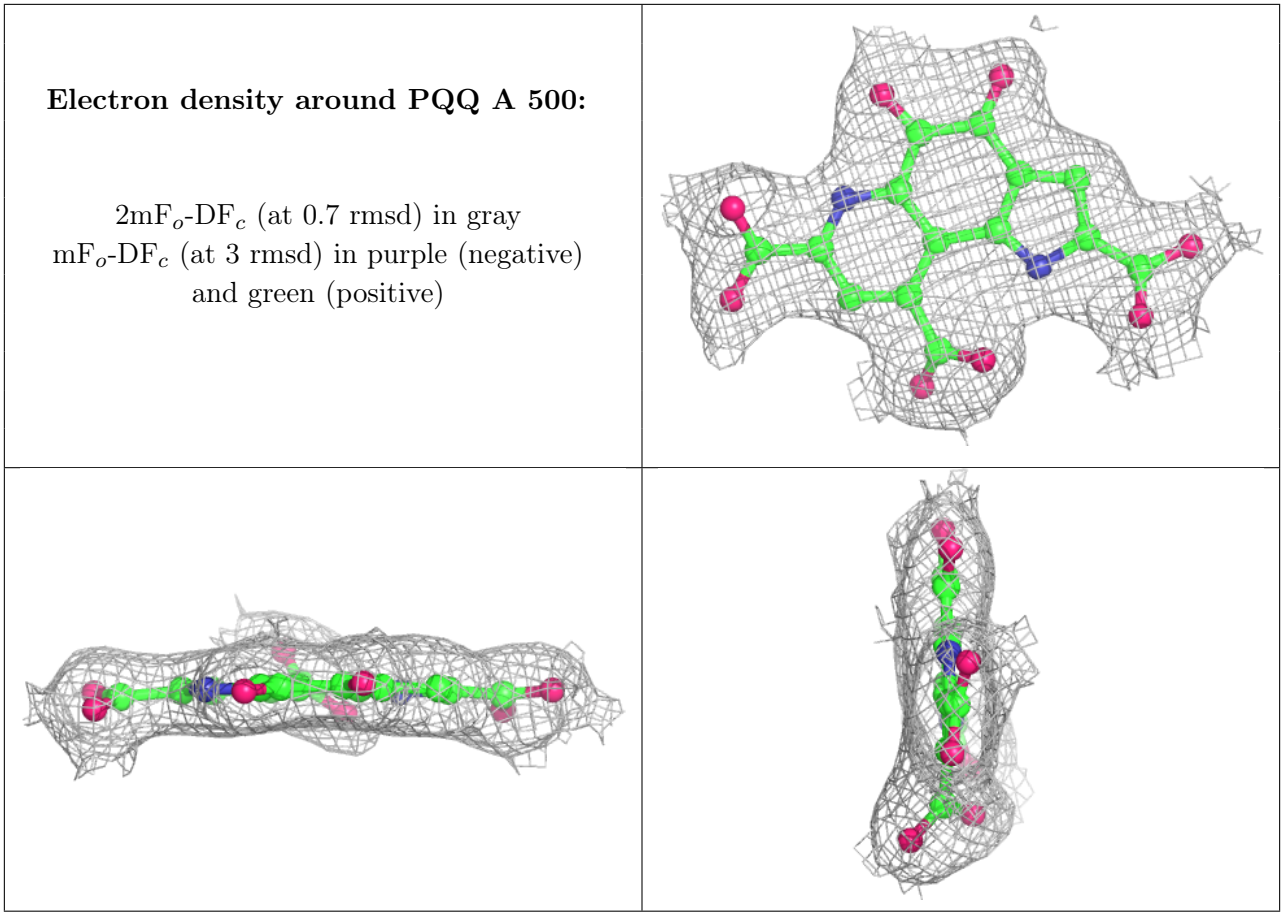
Electron density around PQQ B 501:

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



Electron density around PQQ A 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.