



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:04 PM UTC

PDB ID : 1KV9 / pdb_00001kv9
Title : Structure at 1.9 Å Resolution of a Quinohemoprotein Alcohol Dehydrogenase from *Pseudomonas putida* HK5
Authors : Chen, Z.-W.; Matsushita, K.; Yamashita, T.; Fujii, T.; Toyama, H.; Adachi, O.; Bellamy, H.D.; Mathews, F.S.
Deposited on : 2002-01-25
Resolution : 1.90 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

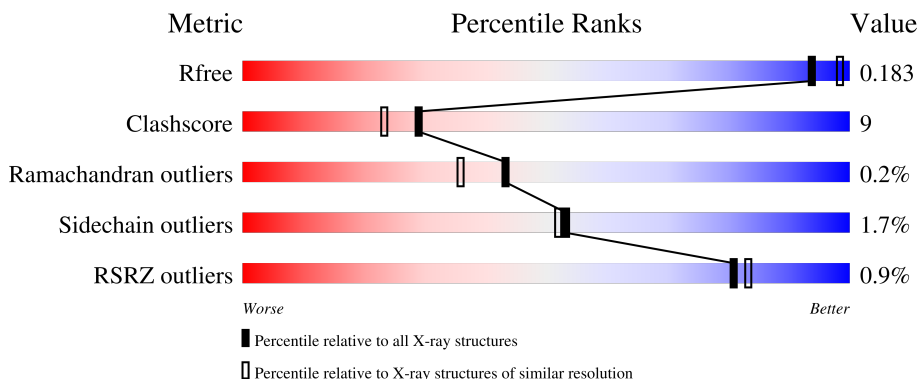
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	668	80% 18% ..

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.

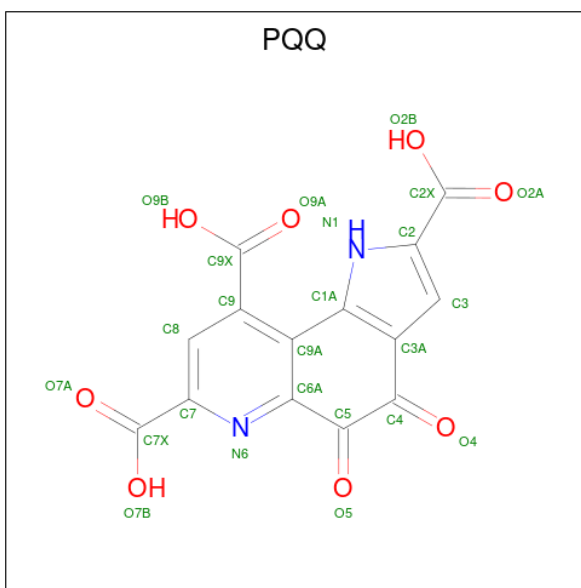
- Molecule 1 is a protein called TYPE II QUINOHENOPROTEIN ALCOHOL DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	664	Total	C	N	O	S	0	0	0
			5103	3251	882	951	19			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

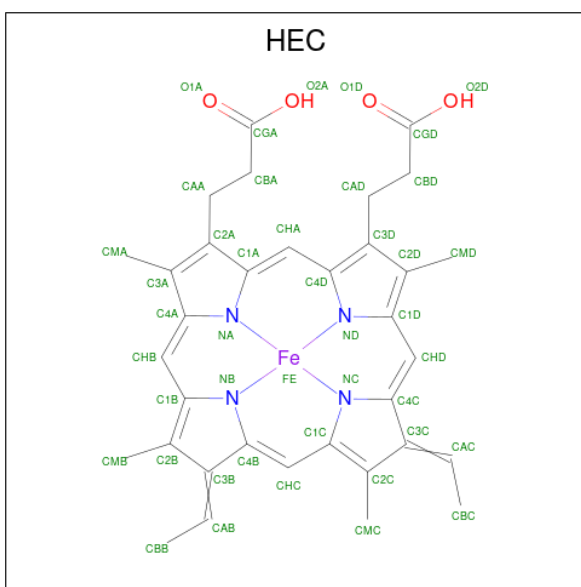
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0

- Molecule 3 is PYRROLOQUINOLINE QUINONE (CCD ID: PQQ) (formula: $C_{14}H_6N_2O_8$).



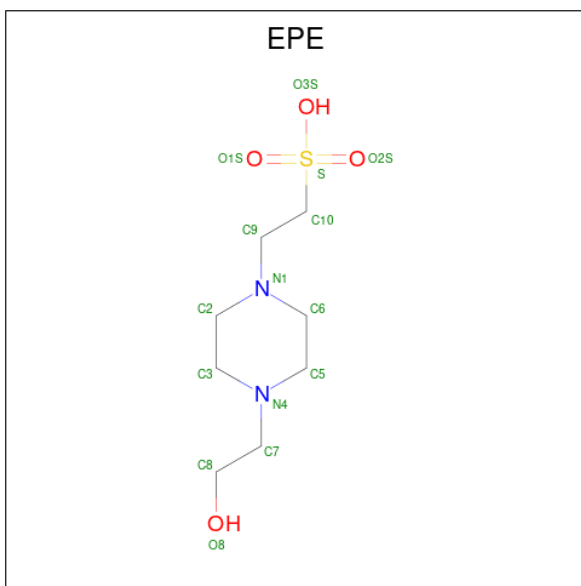
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			24	14	2	8		

- Molecule 4 is HEME C (CCD ID: HEC) (formula: $\text{C}_{34}\text{H}_{34}\text{FeN}_4\text{O}_4$).



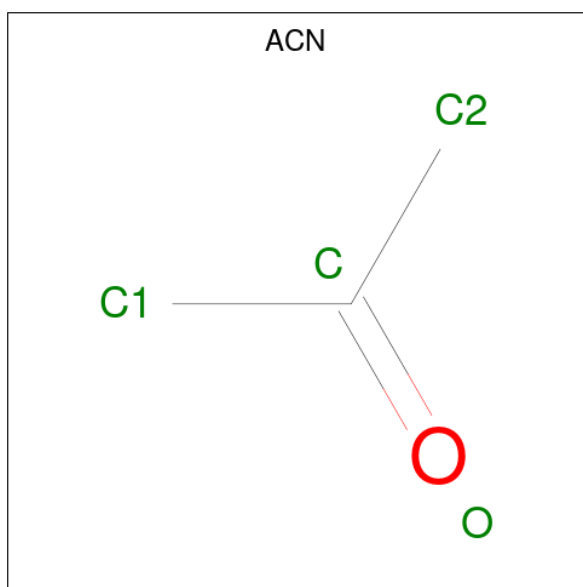
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



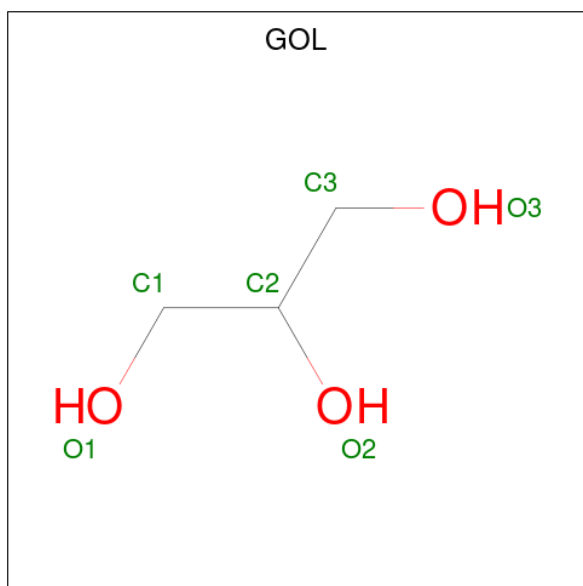
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S		
			15	8	2	4	1	0	0

- Molecule 6 is ACETONE (CCD ID: ACN) (formula: C_3H_6O).



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

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		

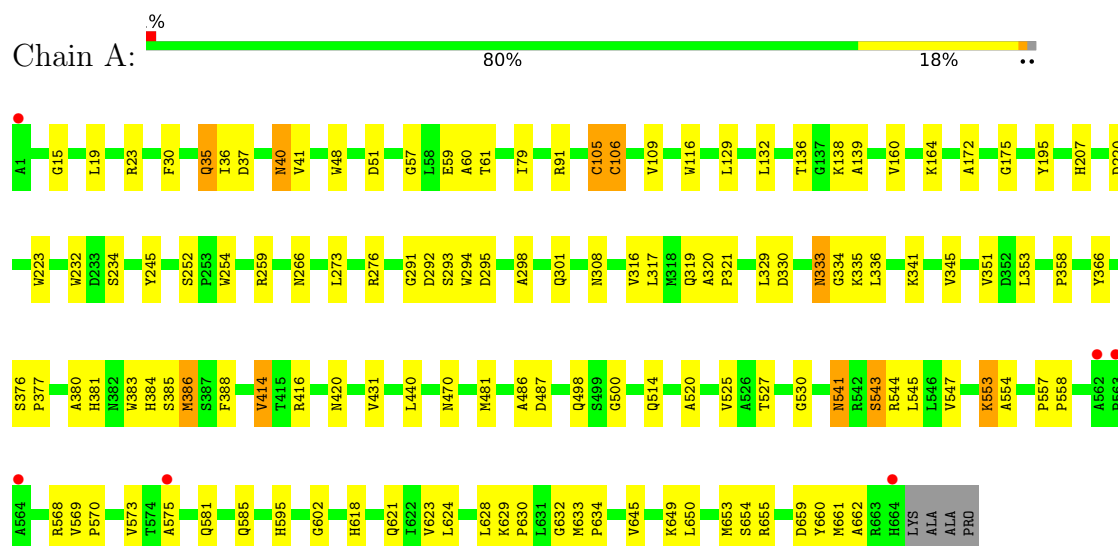
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	732	Total 732	O 732	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: TYPE II QUINOHENOPROTEIN ALCOHOL DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.81Å 57.44Å 67.52Å 89.65° 69.34° 68.39°	Depositor
Resolution (Å)	30.00 – 1.90 30.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	87.1 (30.00-1.90) 87.1 (30.00-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 1.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.154 , 0.189 0.149 , 0.183	Depositor DCC
R_{free} test set	5189 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.565	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.9	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	0.96	EDS
Total number of atoms	5928	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.28% 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: ACN, CA, PQQ, GOL, EPE, HEC

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	0.39	2/5245 (0.0%)	0.96	31/7142 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	595	HIS	CE1-NE2	11.79	1.44	1.32
1	A	595	HIS	ND1-CE1	5.81	1.38	1.32

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	595	HIS	ND1-CG-CD2	13.81	119.91	106.10
1	A	106	CYS	N-CA-C	10.13	122.60	110.44
1	A	595	HIS	CB-CG-CD2	-9.24	119.18	131.20
1	A	105	CYS	N-CA-C	8.23	121.29	111.33
1	A	234	SER	N-CA-C	7.63	121.27	110.50
1	A	500	GLY	N-CA-C	-6.63	103.59	112.14
1	A	623	VAL	N-CA-C	6.30	117.08	110.72
1	A	385	SER	N-CA-C	6.28	118.89	110.35
1	A	595	HIS	CG-CD2-NE2	-6.17	101.03	107.20
1	A	380	ALA	N-CA-C	-6.14	105.05	112.54
1	A	61	THR	N-CA-C	-6.04	99.86	109.04
1	A	316	VAL	N-CA-C	5.96	117.75	109.29
1	A	109	VAL	N-CA-C	5.91	118.28	109.17
1	A	19	LEU	N-CA-C	5.83	120.56	113.38
1	A	386	MET	N-CA-C	-5.76	101.67	110.14
1	A	481	MET	N-CA-C	-5.68	99.93	109.07
1	A	440	LEU	N-CA-C	-5.61	99.37	108.41
1	A	545	LEU	N-CA-C	-5.61	99.58	108.73
1	A	60	ALA	N-CA-C	5.56	118.34	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	431	VAL	CA-C-N	5.55	125.57	119.90
1	A	431	VAL	C-N-CA	5.55	125.57	119.90
1	A	298	ALA	N-CA-C	-5.45	103.91	110.88
1	A	569	VAL	CA-C-N	5.41	126.25	119.98
1	A	569	VAL	C-N-CA	5.41	126.25	119.98
1	A	543	SER	N-CA-C	5.39	117.99	109.96
1	A	35	GLN	N-CA-C	-5.25	105.64	111.36
1	A	232	TRP	N-CA-C	5.25	119.78	113.17
1	A	530	GLY	N-CA-C	5.22	120.73	111.02
1	A	57	GLY	N-CA-C	5.15	117.39	111.36
1	A	129	LEU	N-CA-C	-5.07	100.15	108.41
1	A	595	HIS	CG-ND1-CE1	-5.03	100.75	109.30

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	5103	0	4974	89	0
2	A	1	0	0	0	0
3	A	24	0	3	2	0
4	A	43	0	30	0	0
5	A	15	0	17	0	0
6	A	4	0	6	0	0
7	A	6	0	8	1	0
8	A	732	0	0	1	0
All	All	5928	0	5038	89	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 9.

All (89) 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:A:581:GLN:HE21	1:A:585:GLN:HE21	1.07	0.94
1:A:416:ARG:H	1:A:420:ASN:HD21	1.11	0.92
1:A:40:ASN:C	1:A:40:ASN:HD22	1.94	0.74
1:A:195:TYR:H	1:A:207:HIS:HE1	1.40	0.69
1:A:15:GLY:HA2	1:A:23:ARG:NH2	2.09	0.68
1:A:660:TYR:HD1	1:A:661:MET:HE2	1.60	0.66
1:A:381:HIS:HE1	1:A:384:HIS:O	1.78	0.66
1:A:106:CYS:HB3	1:A:525:VAL:HB	1.77	0.66
1:A:470:ASN:HD22	1:A:486:ALA:HB3	1.61	0.65
1:A:319:GLN:HG2	1:A:321:PRO:HD3	1.79	0.64
1:A:581:GLN:HE21	1:A:585:GLN:NE2	1.90	0.63
1:A:573:VAL:HG11	1:A:654:SER:HA	1.80	0.62
1:A:649:LYS:HG2	1:A:653:MET:HE2	1.80	0.62
1:A:498:GLN:HE21	1:A:544:ARG:HH12	1.47	0.61
1:A:320:ALA:O	1:A:381:HIS:HD2	1.83	0.61
1:A:259:ARG:HH21	1:A:266:ASN:HD22	1.49	0.60
1:A:333:ASN:CG	1:A:335:LYS:HG3	2.27	0.60
1:A:553:LYS:H	1:A:553:LYS:HE2	1.67	0.59
1:A:553:LYS:H	1:A:553:LYS:CE	2.16	0.59
1:A:568:ARG:O	1:A:570:PRO:HD3	2.02	0.58
1:A:254:TRP:O	1:A:293:SER:HA	2.03	0.58
1:A:655:ARG:HH12	1:A:659:ASP:CG	2.12	0.58
1:A:553:LYS:H	1:A:553:LYS:CD	2.18	0.57
1:A:498:GLN:NE2	1:A:544:ARG:HH12	2.02	0.56
1:A:37:ASP:H	1:A:40:ASN:HD21	1.52	0.55
1:A:40:ASN:HD22	1:A:41:VAL:N	2.04	0.55
1:A:195:TYR:H	1:A:207:HIS:CE1	2.22	0.55
1:A:301:GLN:NE2	1:A:386:MET:H	2.05	0.55
1:A:581:GLN:NE2	1:A:585:GLN:HE21	1.91	0.54
1:A:629:LYS:N	1:A:630:PRO:HD2	2.22	0.54
1:A:333:ASN:ND2	1:A:335:LYS:H	2.06	0.54
1:A:164:LYS:HE2	8:A:1201:HOH:O	2.07	0.54
1:A:333:ASN:ND2	1:A:335:LYS:HG3	2.23	0.54
1:A:575:ALA:HB1	1:A:650:LEU:HD21	1.90	0.54
1:A:172:ALA:HB3	3:A:801:PQQ:O7A	2.09	0.53
1:A:470:ASN:HD21	1:A:487:ASP:H	1.54	0.53
1:A:416:ARG:H	1:A:420:ASN:ND2	1.94	0.53
1:A:470:ASN:ND2	1:A:486:ALA:HB3	2.23	0.53
1:A:333:ASN:C	1:A:333:ASN:HD22	2.17	0.53
1:A:333:ASN:HD22	1:A:334:GLY:N	2.07	0.52
1:A:660:TYR:CD1	1:A:661:MET:HE2	2.44	0.51
1:A:553:LYS:H	1:A:553:LYS:HD3	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ASP:H	1:A:40:ASN:ND2	2.07	0.51
1:A:633:MET:HE3	1:A:634:PRO:O	2.11	0.50
1:A:320:ALA:O	1:A:381:HIS:CD2	2.64	0.50
1:A:330:ASP:HB3	1:A:333:ASN:HD21	1.76	0.49
1:A:341:LYS:HD3	1:A:345:VAL:HB	1.95	0.49
1:A:618:HIS:HA	1:A:621:GLN:OE1	2.13	0.48
1:A:655:ARG:HB3	1:A:655:ARG:NH1	2.28	0.48
1:A:40:ASN:C	1:A:40:ASN:ND2	2.64	0.48
1:A:381:HIS:CE1	1:A:384:HIS:O	2.64	0.47
1:A:79:ILE:HG22	1:A:91:ARG:CB	2.45	0.47
1:A:59:GLU:HG2	1:A:383:TRP:CH2	2.51	0.46
1:A:527:THR:O	1:A:602:GLY:HA3	2.16	0.46
1:A:51:ASP:HB3	7:A:805:GOL:H32	1.97	0.46
1:A:351:VAL:HA	1:A:358:PRO:HA	1.98	0.46
1:A:573:VAL:HG13	1:A:573:VAL:O	2.15	0.46
1:A:317:LEU:C	1:A:317:LEU:HD23	2.40	0.45
1:A:132:LEU:HD23	1:A:139:ALA:HA	1.99	0.45
1:A:175:GLY:HA2	1:A:252:SER:HB3	1.98	0.45
1:A:30:PHE:HB2	1:A:388:PHE:O	2.18	0.44
1:A:136:THR:OG1	1:A:138:LYS:HG2	2.18	0.44
1:A:220:ASP:HA	1:A:223:TRP:CH2	2.53	0.43
1:A:79:ILE:HG22	1:A:91:ARG:HB2	2.01	0.43
1:A:376:SER:HB2	1:A:377:PRO:CD	2.48	0.43
1:A:553:LYS:CD	1:A:553:LYS:N	2.81	0.43
1:A:48:TRP:CZ2	1:A:547:VAL:HG21	2.54	0.43
1:A:632:GLY:O	1:A:634:PRO:HD3	2.19	0.43
1:A:553:LYS:HD3	1:A:553:LYS:N	2.33	0.43
1:A:35:GLN:HB3	1:A:514:GLN:HE21	1.83	0.42
1:A:105:CYS:SG	1:A:106:CYS:N	2.92	0.42
1:A:659:ASP:O	1:A:662:ALA:HB3	2.19	0.42
1:A:294:TRP:O	1:A:295:ASP:HB3	2.20	0.42
1:A:59:GLU:OE1	3:A:801:PQQ:O2B	2.38	0.42
1:A:520:ALA:O	1:A:543:SER:HA	2.20	0.41
1:A:557:PRO:HA	1:A:558:PRO:HD3	1.92	0.41
1:A:36:ILE:H	1:A:514:GLN:NE2	2.18	0.41
1:A:116:TRP:CG	1:A:160:VAL:HG11	2.56	0.41
1:A:245:TYR:HA	1:A:273:LEU:O	2.20	0.41
1:A:575:ALA:HB2	1:A:650:LEU:HD11	2.02	0.41
1:A:329:LEU:HD23	1:A:336:LEU:HD12	2.02	0.41
1:A:414:VAL:HG13	1:A:416:ARG:NH1	2.36	0.41
1:A:541:ASN:ND2	1:A:541:ASN:C	2.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:ASP:HB3	1:A:333:ASN:ND2	2.36	0.41
1:A:553:LYS:HE2	1:A:554:ALA:N	2.36	0.41
1:A:291:GLY:O	1:A:292:ASP:C	2.65	0.40
1:A:575:ALA:CB	1:A:650:LEU:HD21	2.51	0.40
1:A:624:LEU:HG	1:A:645:VAL:HG21	2.03	0.40
1:A:541:ASN:C	1:A:541:ASN:HD22	2.29	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	662/668 (99%)	630 (95%)	31 (5%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	366	TYR

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	519/521 (100%)	510 (98%)	9 (2%)	53	52

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	276	ARG
1	A	308	ASN
1	A	333	ASN
1	A	353	LEU
1	A	414	VAL
1	A	541	ASN
1	A	553	LYS
1	A	628	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	40	ASN
1	A	65	HIS
1	A	143	GLN
1	A	207	HIS
1	A	221	GLN
1	A	255	ASN
1	A	266	ASN
1	A	300	GLN
1	A	301	GLN
1	A	381	HIS
1	A	420	ASN
1	A	470	ASN
1	A	494	GLN
1	A	498	GLN
1	A	514	GLN
1	A	535	ASN
1	A	541	ASN
1	A	555	GLN
1	A	579	GLN
1	A	581	GLN
1	A	589	GLN
1	A	618	HIS
1	A	620	GLN
1	A	664	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	$\# Z > 2$
6	ACN	A	803	-	3,3,3	0.43	0	3,3,3	0.59	0
4	HEC	A	901	1	46,50,50	1.64	6 (13%)	58,82,82	1.45	4 (6%)
3	PQQ	A	801	2	26,26,26	2.31	10 (38%)	34,40,40	1.70	9 (26%)
5	EPE	A	804	-	15,15,15	1.78	3 (20%)	19,20,20	1.51	4 (21%)
7	GOL	A	805	-	5,5,5	1.17	0	5,5,5	0.84	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PQQ	A	801	2	-	0/12/28/28	0/3/3/3
4	HEC	A	901	1	-	8/14/54/54	-
7	GOL	A	805	-	-	2/4/4/4	-
5	EPE	A	804	-	-	1/9/19/19	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	PQQ	C9A-C1A	6.25	1.53	1.46
4	A	901	HEC	CAC-C3C	5.74	1.53	1.35
4	A	901	HEC	CAB-C3B	5.63	1.53	1.35
3	A	801	PQQ	C9-C9A	4.23	1.47	1.41
5	A	804	EPE	C2-N1	3.58	1.56	1.46
5	A	804	EPE	C3-N4	3.47	1.56	1.46
3	A	801	PQQ	C8-C9	3.40	1.45	1.39
5	A	804	EPE	C6-N1	3.25	1.55	1.46
3	A	801	PQQ	C9A-C6A	3.09	1.45	1.41
3	A	801	PQQ	C5-C4	3.08	1.59	1.54
3	A	801	PQQ	O7A-C7X	2.98	1.31	1.22
4	A	901	HEC	CMC-C2C	2.62	1.56	1.50
3	A	801	PQQ	C6A-C5	2.61	1.53	1.50
3	A	801	PQQ	C1A-N1	-2.48	1.33	1.37
3	A	801	PQQ	C9-C9X	2.44	1.54	1.49
4	A	901	HEC	C3C-C2C	-2.20	1.33	1.41
3	A	801	PQQ	O2B-C2X	-2.15	1.24	1.30
4	A	901	HEC	CMD-C2D	2.07	1.55	1.50
4	A	901	HEC	O2A-CGA	-2.06	1.24	1.30

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	HEC	CBB-CAB-C3B	-6.64	114.16	127.43
4	A	901	HEC	CBC-CAC-C3C	-5.37	116.70	127.43
3	A	801	PQQ	C9A-C6A-C5	4.06	125.36	121.59
3	A	801	PQQ	O7A-C7X-C7	-3.35	114.56	121.30
5	A	804	EPE	C9-N1-C6	-3.26	102.54	111.24
5	A	804	EPE	C7-N4-C5	-2.89	103.55	111.24
3	A	801	PQQ	C5-C6A-N6	-2.89	110.25	115.06
3	A	801	PQQ	C8-C9-C9A	-2.77	116.78	120.33
3	A	801	PQQ	C3A-C4-C5	-2.65	114.04	117.03
4	A	901	HEC	CBA-CAA-C2A	-2.52	105.55	112.53
4	A	901	HEC	O1A-CGA-CBA	-2.45	115.31	123.09
5	A	804	EPE	C3-C2-N1	-2.42	105.77	110.65
3	A	801	PQQ	C1A-C3A-C4	2.35	124.31	122.33
3	A	801	PQQ	O4-C4-C3A	2.19	127.29	121.92
3	A	801	PQQ	C3-C2-C2X	-2.09	127.71	132.53
3	A	801	PQQ	O7B-C7X-O7A	2.03	127.72	123.35
5	A	804	EPE	O1S-S-C10	-2.03	103.67	106.73

There are no chirality outliers.

All (11) torsion outliers are listed below:

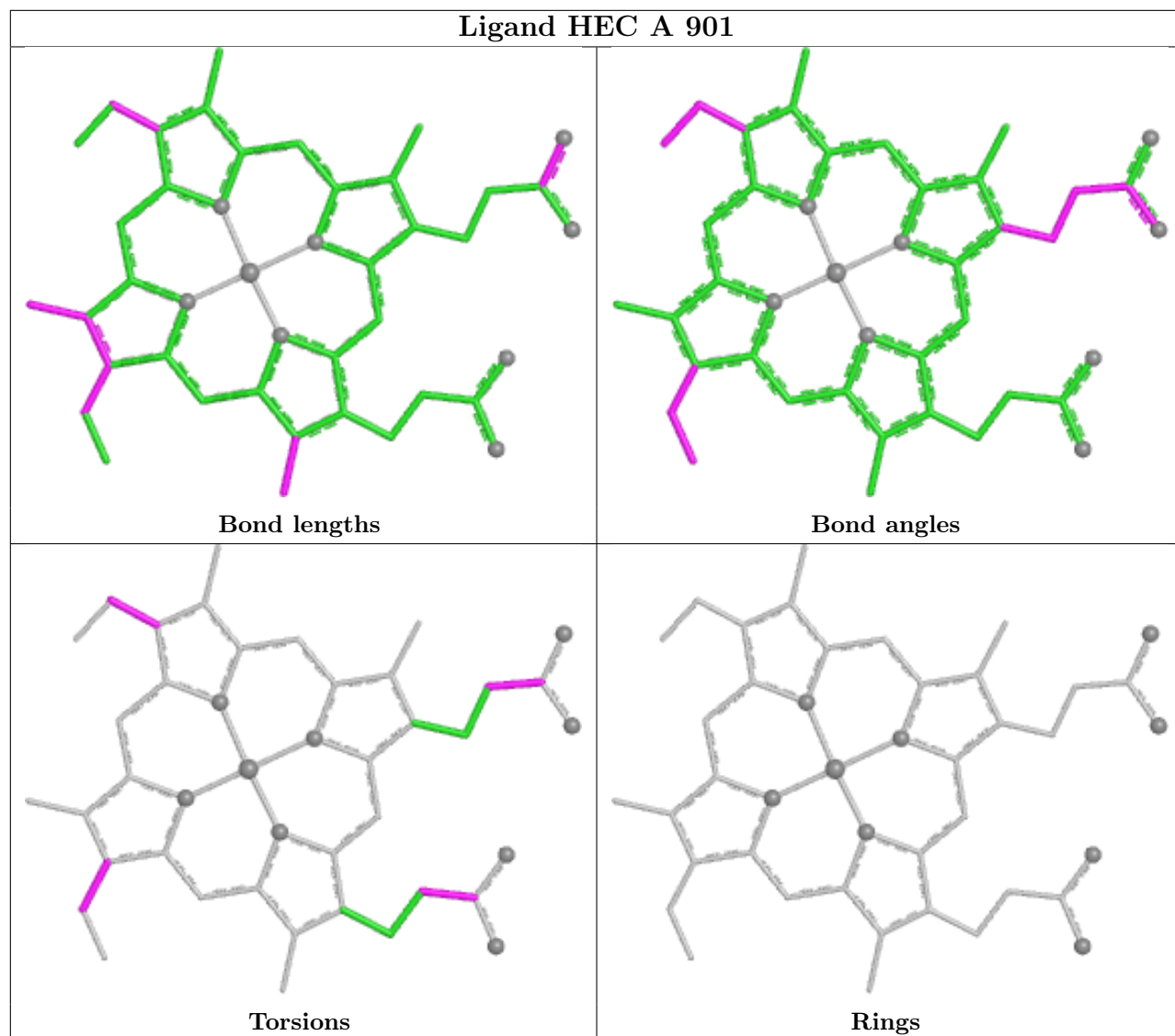
Mol	Chain	Res	Type	Atoms
4	A	901	HEC	C2B-C3B-CAB-CBB
4	A	901	HEC	C4B-C3B-CAB-CBB
4	A	901	HEC	C2C-C3C-CAC-CBC
4	A	901	HEC	C4C-C3C-CAC-CBC
5	A	804	EPE	N4-C7-C8-O8
7	A	805	GOL	C1-C2-C3-O3
7	A	805	GOL	O2-C2-C3-O3
4	A	901	HEC	CAD-CBD-CGD-O2D
4	A	901	HEC	CAD-CBD-CGD-O1D
4	A	901	HEC	CAA-CBA-CGA-O2A
4	A	901	HEC	CAA-CBA-CGA-O1A

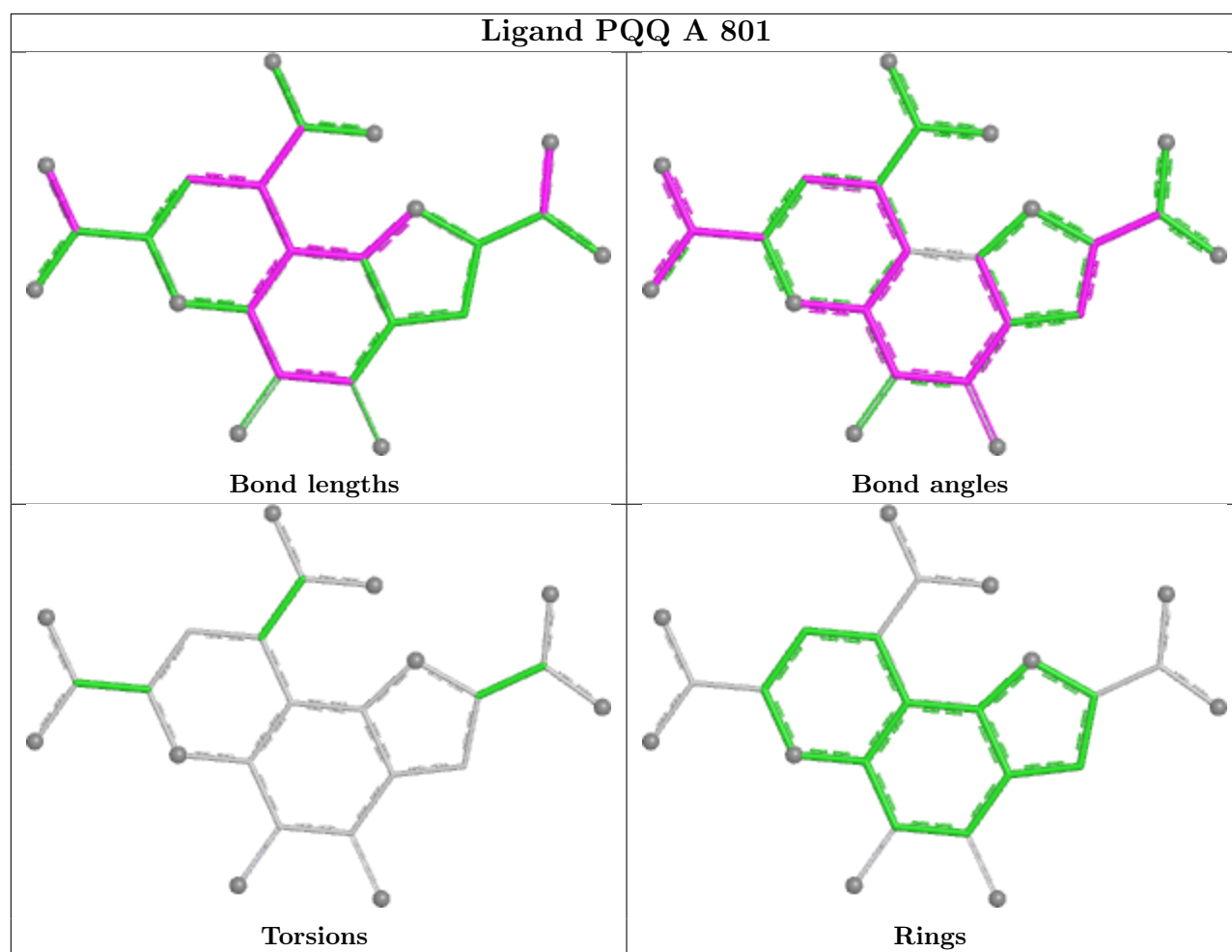
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	PQQ	2	0
7	A	805	GOL	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 [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, 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	664/668 (99%)	-0.57	6 (0%) 81 83	12, 18, 34, 62	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	ALA	6.3
1	A	564	ALA	2.8
1	A	563	PRO	2.5
1	A	562	ALA	2.2
1	A	575	ALA	2.1
1	A	664	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 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.

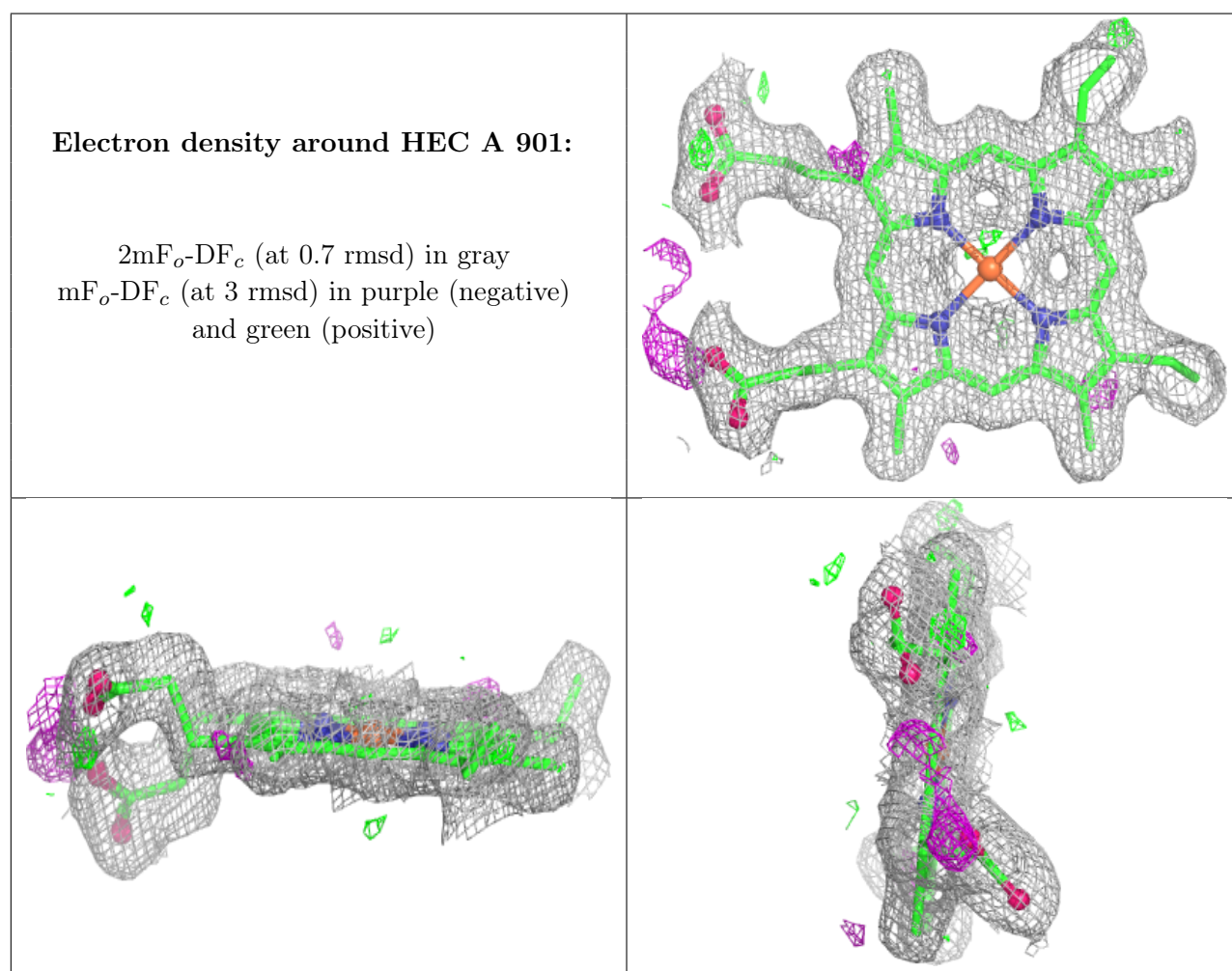
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	GOL	A	805	6/6	0.84	0.12	44,45,46,46	0

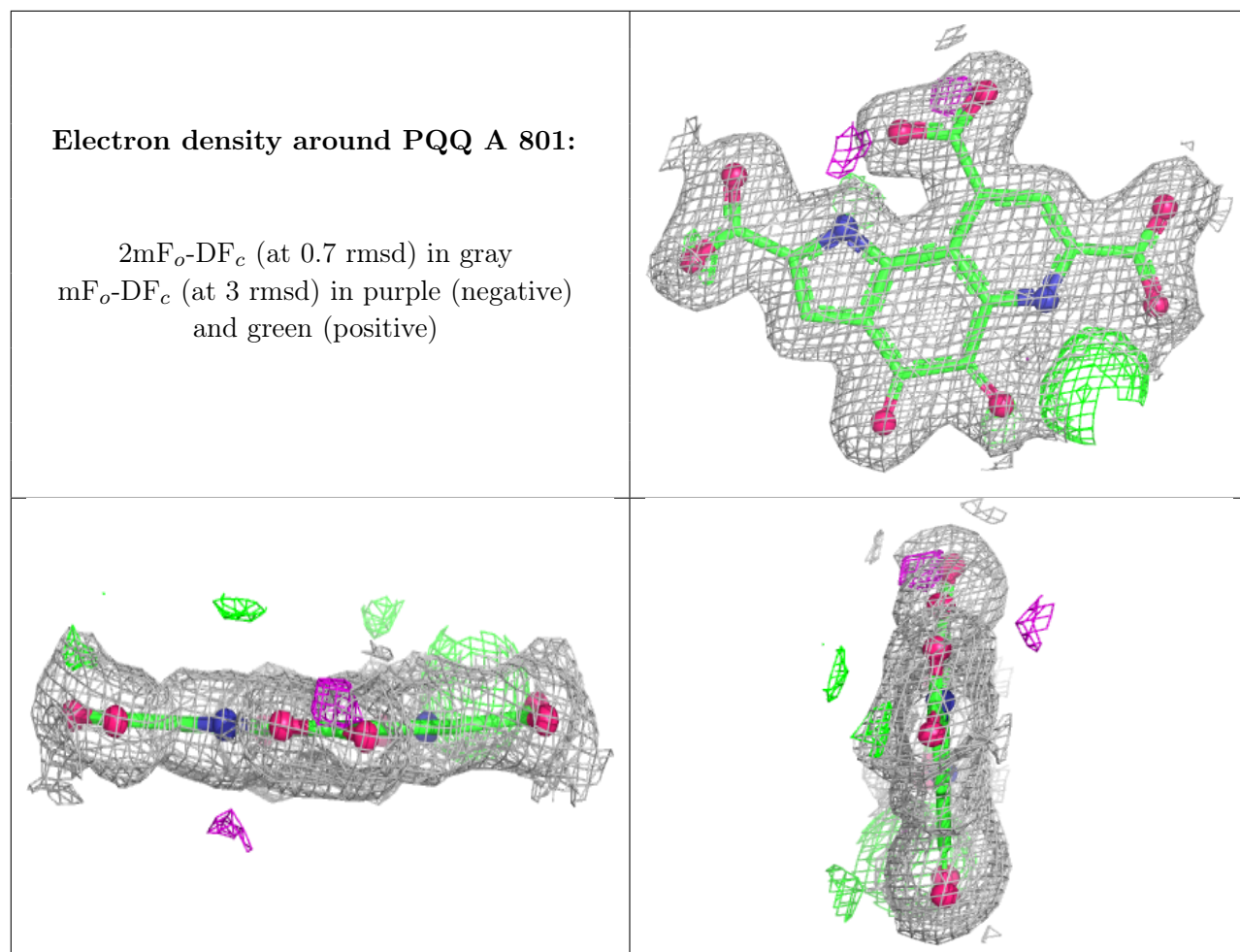
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EPE	A	804	15/15	0.85	0.12	41,47,56,57	0
6	ACN	A	803	4/4	0.90	0.10	30,30,31,31	0
4	HEC	A	901	43/43	0.97	0.06	13,19,23,28	0
3	PQQ	A	801	24/24	0.97	0.05	13,17,18,21	0
2	CA	A	802	1/1	1.00	0.21	20,20,20,20	0

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





6.5 Other polymers [i](#)

There are no such residues in this entry.