



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 06:43 am BST

PDB ID : 6DJX
Title : Crystal Structure of pParkin-pUb-UbcH7 complex
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Deposited on : 2018-05-27
Resolution : 4.80 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 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

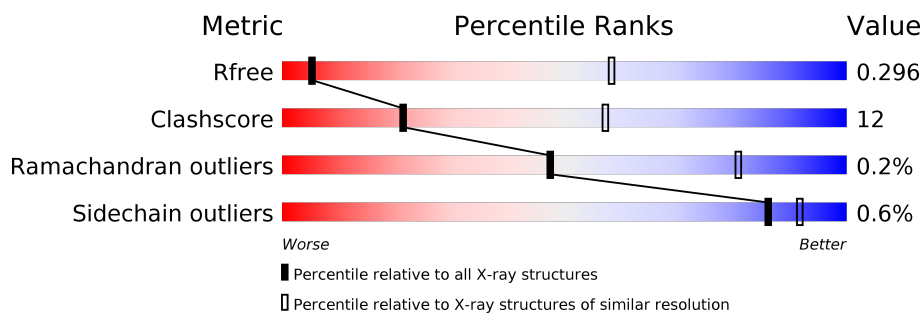
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.80 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	1096 (5.80-3.80)
Clashscore	141614	1170 (5.80-3.80)
Ramachandran outliers	138981	1105 (5.80-3.80)
Sidechain outliers	138945	1085 (5.80-3.80)

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 ≥ 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	432	
2	B	76	
3	C	159	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4346 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 RBR-type E3 ubiquitin transferase,RBR-type E3 ubiquitin transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	P	S	0	0	0
			2492	1569	429	465	1	28			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	GLY	-	expression tag	UNP A0A034W4L8
A	21	GLY	-	expression tag	UNP A0A034W4L8
A	22	GLU	-	expression tag	UNP A0A034W4L8
A	23	ASN	-	expression tag	UNP A0A034W4L8
A	24	LEU	-	expression tag	UNP A0A034W4L8
A	25	TYR	-	expression tag	UNP A0A034W4L8
A	26	LEU	-	expression tag	UNP A0A034W4L8
A	27	GLY	-	expression tag	UNP A0A034W4L8
A	28	GLY	-	expression tag	UNP A0A034W4L8
A	463	ALA	CYS	engineered mutation	UNP A0A034W4L8

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	74	Total	C	N	O	P	S	0	0	0
			597	374	103	118	1	1			

- Molecule 3 is a protein called Ubiquitin-conjugating enzyme E2 L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	153	Total	C	N	O	S	0	0	0
			1251	801	217	228	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	GLY	-	expression tag	UNP P68036
C	-3	PRO	-	expression tag	UNP P68036
C	-2	LEU	-	expression tag	UNP P68036
C	-1	GLY	-	expression tag	UNP P68036
C	0	SER	-	expression tag	UNP P68036
C	86	LYS	CYS	engineered mutation	UNP P68036

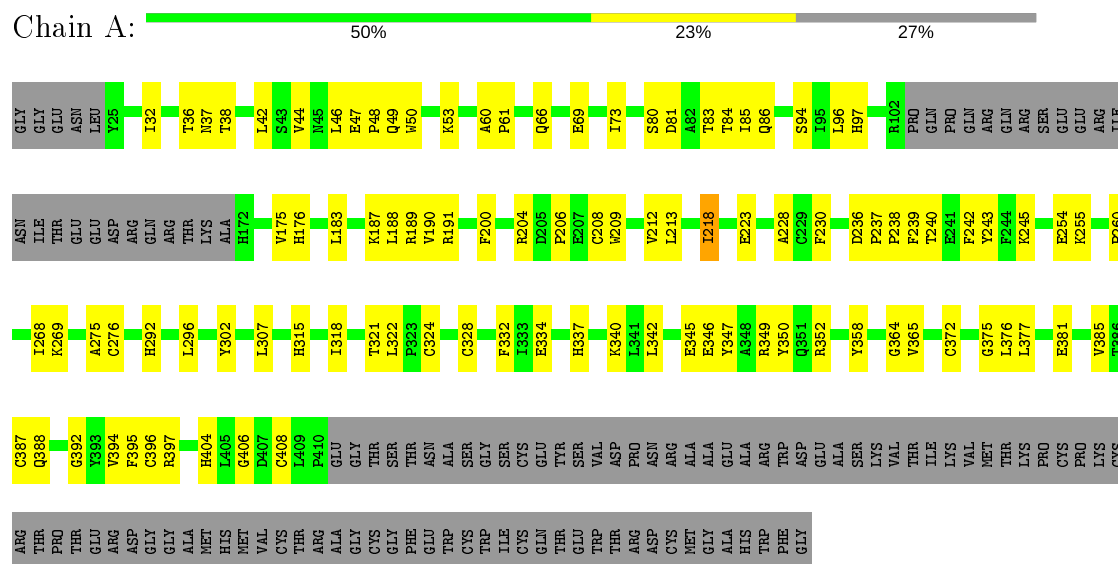
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by author).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	Zn	0	0
			6	6		

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. 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. 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: RBR-type E3 ubiquitin transferase,RBR-type E3 ubiquitin transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	135.67Å 135.67Å 87.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.86 – 4.80 48.86 – 4.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.86-4.80) 99.8 (48.86-4.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.19 (at 4.86Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.257 , 0.290 0.263 , 0.296	Depositor DCC
R_{free} test set	236 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	235.0	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 247.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4346	wwPDB-VP
Average B, all atoms (Å ²)	299.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.24	0/2538	0.42	0/3429
2	B	0.21	0/592	0.43	0/795
3	C	0.22	0/1282	0.40	0/1730
All	All	0.23	0/4412	0.42	0/5954

There are no bond length outliers.

There are no bond angle outliers.

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	2492	0	2408	77	0
2	B	597	0	622	24	0
3	C	1251	0	1270	16	0
4	A	6	0	0	0	0
All	All	4346	0	4300	104	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 12.

All (104) 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:212:VAL:HG13	1:A:213:LEU:HG	1.65	0.78
1:A:318:ILE:HG23	1:A:334:GLU:HA	1.73	0.71
3:C:87:LEU:HD12	3:C:88:PRO:HD2	1.76	0.66
1:A:376:LEU:HD22	2:B:71:LEU:HB2	1.77	0.65
1:A:377:LEU:O	2:B:72:ARG:NH1	2.30	0.65
1:A:377:LEU:HB2	2:B:72:ARG:HD2	1.80	0.63
3:C:87:LEU:HD23	3:C:90:ILE:HG13	1.81	0.61
1:A:342:LEU:HB3	1:A:346:GLU:HB2	1.82	0.60
1:A:53:LYS:HB2	1:A:81:ASP:HB3	1.84	0.59
1:A:204:ARG:HG3	1:A:206:PRO:HD3	1.85	0.58
1:A:315:HIS:HB3	1:A:318:ILE:HB	1.86	0.58
1:A:404:HIS:ND1	1:A:406:GLY:O	2.35	0.58
1:A:387:CYS:HB3	1:A:392:GLY:H	1.69	0.57
1:A:254:GLU:HG3	1:A:255:LYS:HG3	1.85	0.57
1:A:321:THR:OG1	1:A:322:LEU:N	2.36	0.57
1:A:73:ILE:HG22	1:A:97:HIS:HB2	1.88	0.56
1:A:268:ILE:HG23	1:A:269:LYS:HG3	1.87	0.56
1:A:324:CYS:HB3	1:A:328:CYS:H	1.71	0.56
1:A:188:LEU:HD21	1:A:242:PHE:HD2	1.70	0.56
1:A:94:SEP:O3P	1:A:245:LYS:NZ	2.36	0.55
1:A:321:THR:HG21	1:A:332:PHE:HB3	1.89	0.54
1:A:276:CYS:HB3	3:C:2:ALA:HB1	1.88	0.54
1:A:376:LEU:CD2	2:B:71:LEU:HB2	2.37	0.54
1:A:376:LEU:HD13	2:B:73:LEU:HD11	1.89	0.53
1:A:385:VAL:HG11	1:A:395:PHE:CZ	2.43	0.53
2:B:56:LEU:HD22	2:B:61:ILE:HD12	1.89	0.53
3:C:86:LYS:HE2	3:C:119:HIS:HB3	1.89	0.53
1:A:36:THR:OG1	1:A:37:ASN:N	2.43	0.52
1:A:404:HIS:NE2	1:A:408:CYS:HB3	2.24	0.52
3:C:7:LEU:HD22	3:C:32:LEU:HG	1.91	0.52
1:A:387:CYS:HB3	1:A:392:GLY:N	2.25	0.51
1:A:372:CYS:HA	2:B:9:THR:HG22	1.92	0.51
1:A:292:HIS:HE1	1:A:328:CYS:SG	2.33	0.51
1:A:190:VAL:HB	1:A:200:PHE:CZ	2.46	0.51
1:A:204:ARG:HG2	1:A:218:ILE:HD13	1.92	0.51
1:A:376:LEU:HD11	1:A:388:GLN:H	1.76	0.51
3:C:76:HIS:HB3	3:C:79:ILE:HD12	1.94	0.50
2:B:1:MET:HE1	2:B:63:LYS:HA	1.94	0.50
1:A:42:LEU:HG	1:A:44:VAL:HG23	1.94	0.49
1:A:364:GLY:HA3	1:A:377:LEU:HD23	1.95	0.48
2:B:61:ILE:HG23	2:B:65:SEP:HB3	1.94	0.48
1:A:200:PHE:CE1	1:A:240:THR:HB	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:VAL:HG22	1:A:213:LEU:H	1.79	0.48
2:B:8:LEU:HD22	2:B:8:LEU:H	1.79	0.48
1:A:188:LEU:HD21	1:A:242:PHE:CD2	2.49	0.48
1:A:175:VAL:HG12	1:A:260:PRO:HA	1.95	0.48
1:A:365:VAL:HG21	1:A:397:ARG:CD	2.44	0.47
3:C:22:PHE:HD1	3:C:39:ILE:HG12	1.80	0.47
1:A:381:GLU:N	1:A:381:GLU:OE1	2.49	0.46
2:B:31:GLN:HB2	2:B:36:ILE:O	2.15	0.46
1:A:200:PHE:CZ	1:A:240:THR:HB	2.50	0.46
1:A:337:HIS:NE2	2:B:60:ASN:O	2.49	0.46
1:A:365:VAL:HG21	1:A:397:ARG:HD2	1.97	0.46
1:A:32:ILE:HG21	1:A:96:LEU:HD13	1.97	0.46
1:A:191:ARG:HG3	1:A:243:TYR:CE1	2.51	0.46
1:A:237:PRO:HB2	1:A:239:PHE:CE1	2.50	0.45
1:A:66:GLN:HB2	1:A:69:GLU:HG2	1.98	0.45
2:B:36:ILE:H	2:B:36:ILE:HG12	1.58	0.45
1:A:49:GLN:HA	1:A:84:THR:HG23	1.99	0.45
1:A:189:ARG:HB2	1:A:243:TYR:CZ	2.52	0.45
1:A:352:ARG:HE	2:B:6:LYS:HE2	1.82	0.44
1:A:46:LEU:HD22	1:A:50:TRP:CZ3	2.52	0.44
1:A:176:HIS:CE1	1:A:340:LYS:HE3	2.52	0.44
1:A:347:TYR:O	1:A:350:TYR:HB3	2.18	0.44
1:A:48:PRO:HB2	1:A:86:GLN:HB2	1.98	0.44
1:A:37:ASN:ND2	1:A:223:GLU:HB3	2.33	0.44
1:A:47:GLU:HB3	1:A:50:TRP:NE1	2.33	0.44
1:A:375:GLY:O	2:B:70:VAL:HG13	2.17	0.43
1:A:183:LEU:HD21	1:A:296:LEU:HD21	1.99	0.43
3:C:26:GLN:HB3	3:C:36:GLN:HB2	2.01	0.43
1:A:228:ALA:HA	1:A:230:PHE:CE2	2.53	0.43
1:A:385:VAL:HG11	1:A:395:PHE:CE1	2.53	0.43
3:C:46:TYR:CD1	3:C:144:THR:HG21	2.53	0.43
1:A:53:LYS:N	1:A:81:ASP:O	2.50	0.43
1:A:345:GLU:OE2	1:A:349:ARG:NH2	2.52	0.43
1:A:307:LEU:HD11	1:A:349:ARG:HG3	2.00	0.43
1:A:345:GLU:HG2	1:A:349:ARG:HH21	1.83	0.43
2:B:31:GLN:HE22	2:B:38:PRO:HG3	1.84	0.43
1:A:377:LEU:CB	2:B:72:ARG:HD2	2.49	0.43
1:A:302:TYR:CE2	1:A:322:LEU:HD13	2.54	0.42
3:C:11:LEU:HB2	3:C:35:TRP:CZ2	2.55	0.42
3:C:28:ASP:N	3:C:28:ASP:OD1	2.51	0.42
1:A:60:ALA:HB3	1:A:61:PRO:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:ASP:O	1:A:238:PRO:HD3	2.20	0.42
1:A:80:SER:O	1:A:83:THR:HG22	2.18	0.42
2:B:28:ALA:HA	2:B:31:GLN:HE21	1.83	0.42
3:C:75:TYR:CZ	3:C:125:LEU:HG	2.55	0.42
1:A:204:ARG:HG2	1:A:218:ILE:CD1	2.49	0.41
3:C:86:LYS:HB2	3:C:121:LEU:HD11	2.00	0.41
1:A:94:SEP:O2P	1:A:187:LYS:NZ	2.51	0.41
2:B:26:VAL:HG21	2:B:56:LEU:HD21	2.02	0.41
3:C:102:ASP:OD1	3:C:103:GLN:N	2.54	0.41
3:C:64:LYS:HA	3:C:65:PRO:HD3	1.92	0.41
2:B:5:VAL:HG23	2:B:13:ILE:HB	2.03	0.41
1:A:50:TRP:O	1:A:85:ILE:HG13	2.20	0.41
2:B:28:ALA:HA	2:B:31:GLN:HG2	2.02	0.41
2:B:45:PHE:O	2:B:48:LYS:HG2	2.20	0.41
1:A:385:VAL:O	1:A:394:VAL:HA	2.21	0.41
1:A:307:LEU:HD21	1:A:349:ARG:HD2	2.03	0.41
1:A:358:TYR:OH	2:B:49:GLN:NE2	2.52	0.41
2:B:37:PRO:HA	2:B:38:PRO:HD3	1.90	0.41
1:A:208:CYS:SG	1:A:209:TRP:N	2.94	0.40
1:A:275:ALA:O	3:C:6:ARG:HD2	2.20	0.40
1:A:396:CYS:O	1:A:397:ARG:HB3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	312/432 (72%)	291 (93%)	21 (7%)	0	100	100
2	B	71/76 (93%)	70 (99%)	0	1 (1%)	11	46
3	C	151/159 (95%)	143 (95%)	8 (5%)	0	100	100
All	All	534/667 (80%)	504 (94%)	29 (5%)	1 (0%)	47	81

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	11	LYS

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	277/373 (74%)	275 (99%)	2 (1%)	84	90
2	B	67/67 (100%)	66 (98%)	1 (2%)	65	80
3	C	137/141 (97%)	137 (100%)	0	100	100
All	All	481/581 (83%)	478 (99%)	3 (1%)	86	92

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	218	ILE
2	B	36	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
2	SEP	B	65	2	8,9,10	1.38	1 (12%)	8,12,14	1.56	2 (25%)
1	SEP	A	94	1	8,9,10	1.38	1 (12%)	8,12,14	1.40	2 (25%)

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	SEP	B	65	2	-	4/5/8/10	-
1	SEP	A	94	1	-	3/5/8/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	SEP	P-O1P	3.01	1.60	1.50
2	B	65	SEP	P-O1P	2.98	1.60	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	65	SEP	OG-CB-CA	3.26	111.32	108.14
1	A	94	SEP	P-OG-CB	-2.95	110.17	118.30
2	B	65	SEP	P-OG-CB	-2.40	111.69	118.30
1	A	94	SEP	OG-CB-CA	2.21	110.30	108.14

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	65	SEP	CB-OG-P-O1P
2	B	65	SEP	CB-OG-P-O2P
2	B	65	SEP	CB-OG-P-O3P
1	A	94	SEP	N-CA-CB-OG
1	A	94	SEP	CA-CB-OG-P
2	B	65	SEP	CA-CB-OG-P
1	A	94	SEP	CB-OG-P-O2P

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	65	SEP	1	0
1	A	94	SEP	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

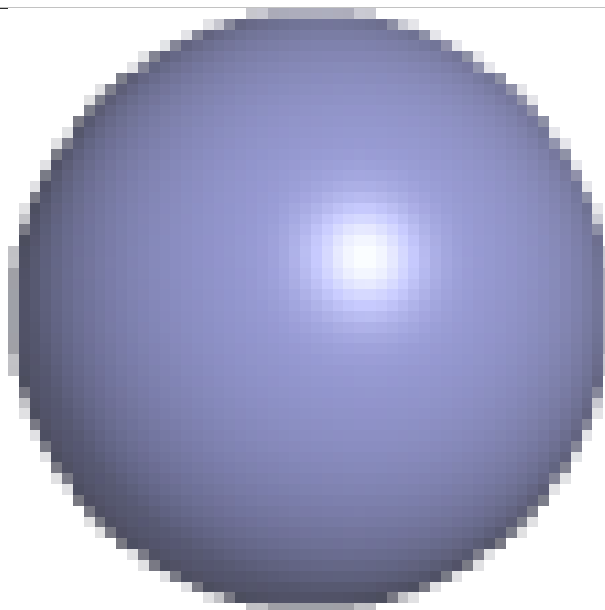
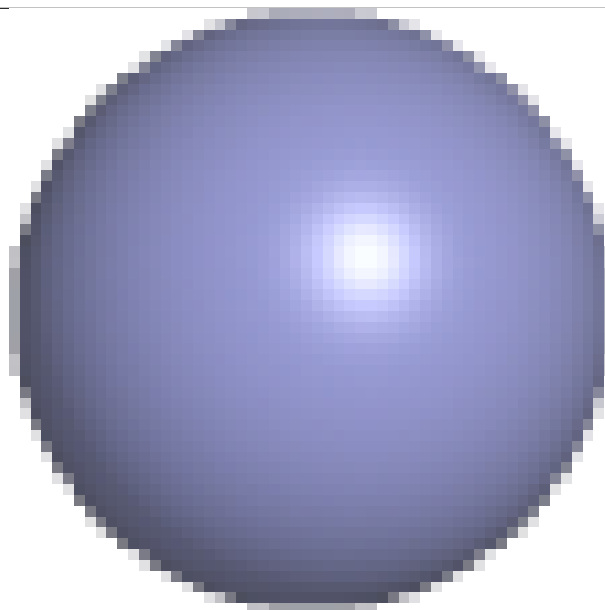
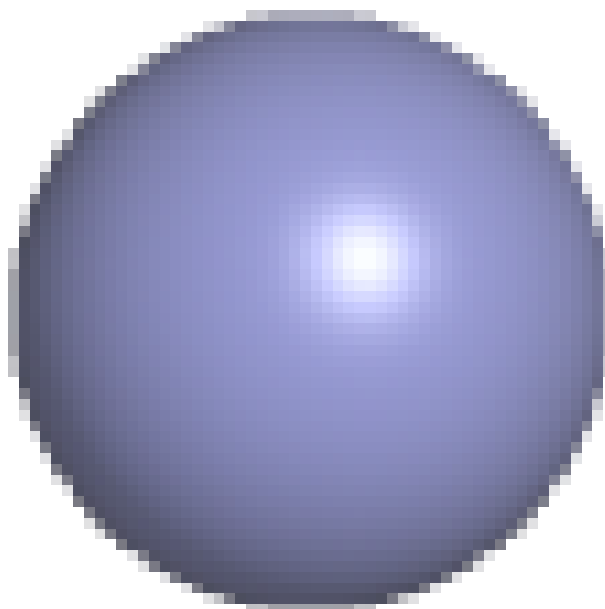
6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

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.

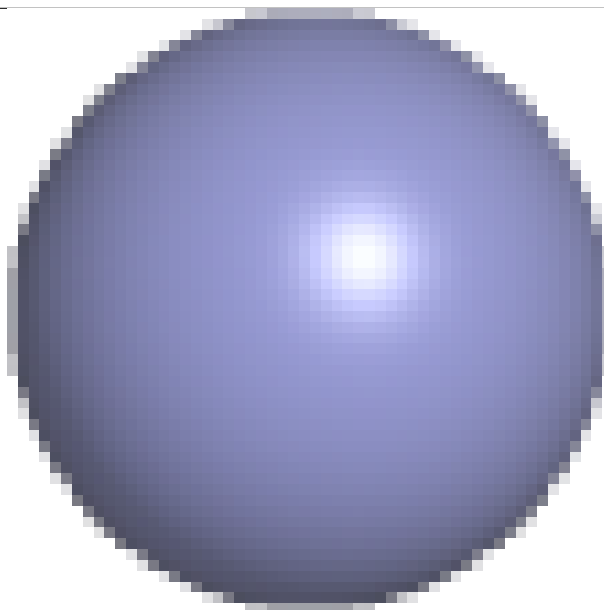
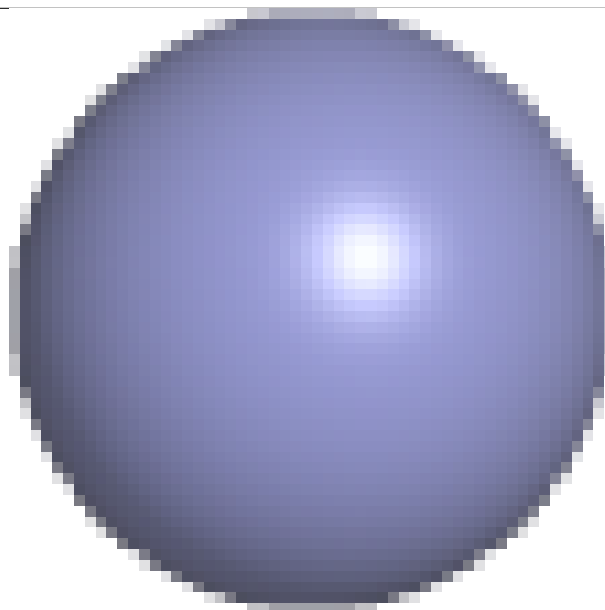
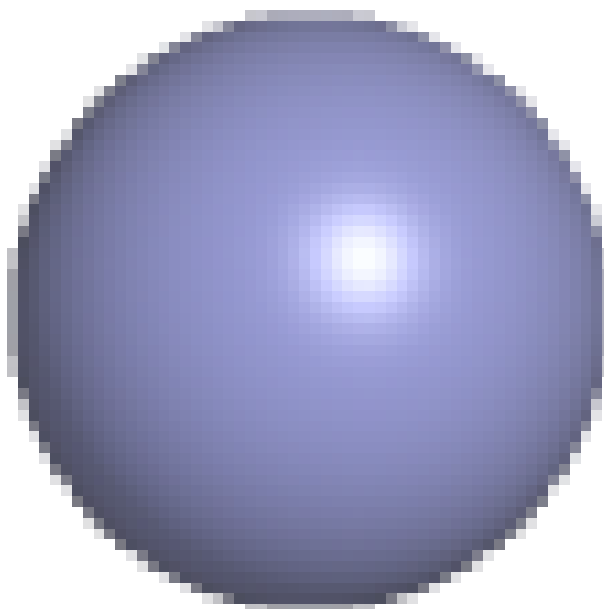
Electron density around ZN A 1001:

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



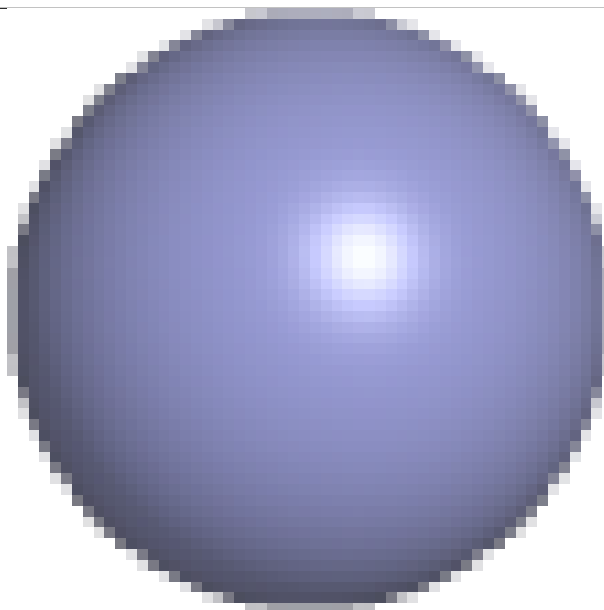
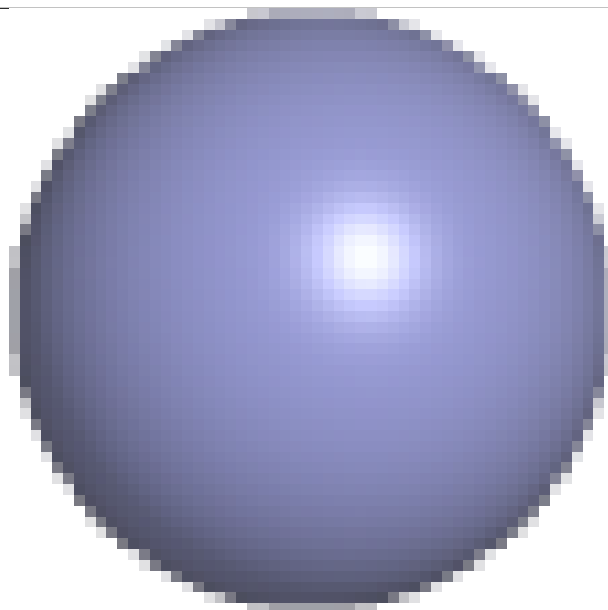
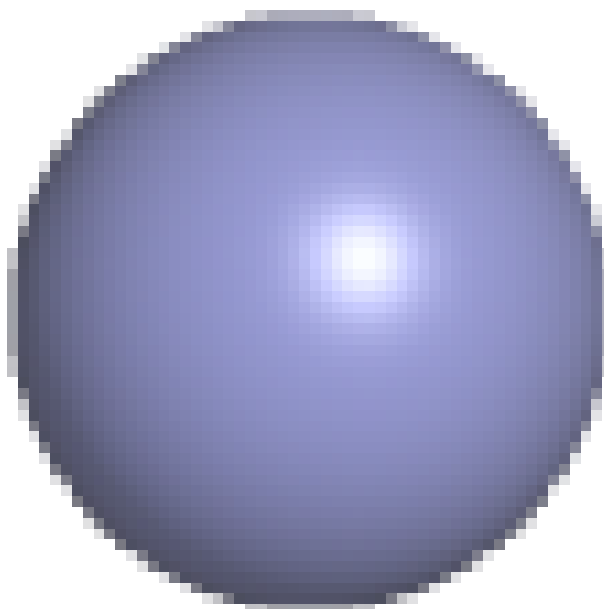
Electron density around ZN A 1005:

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



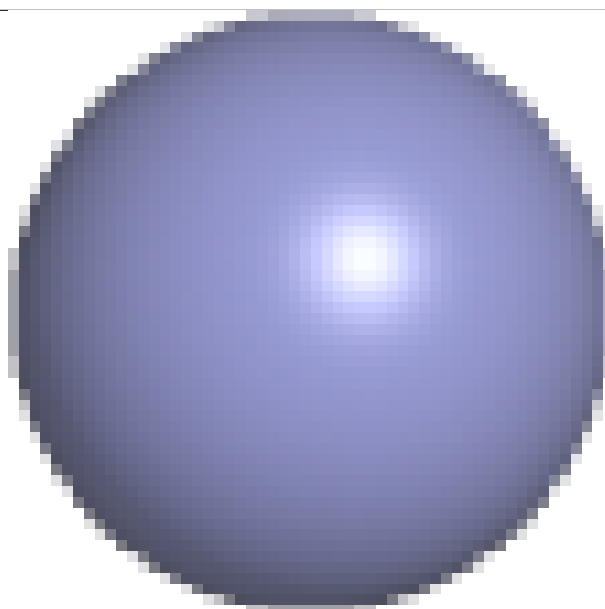
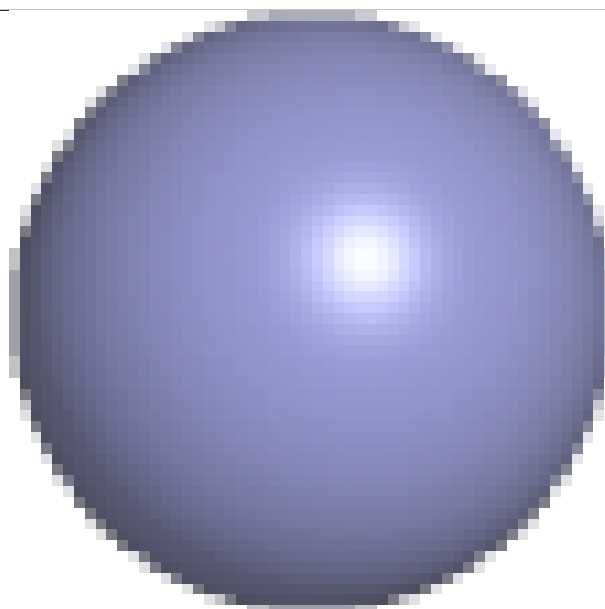
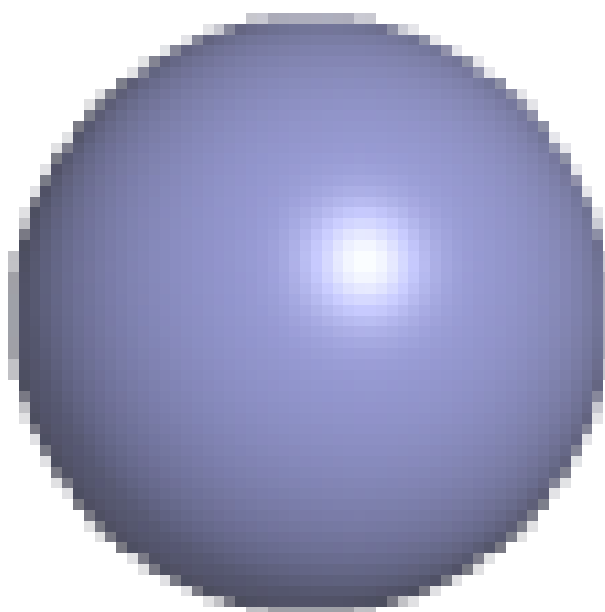
Electron density around ZN A 1004:

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



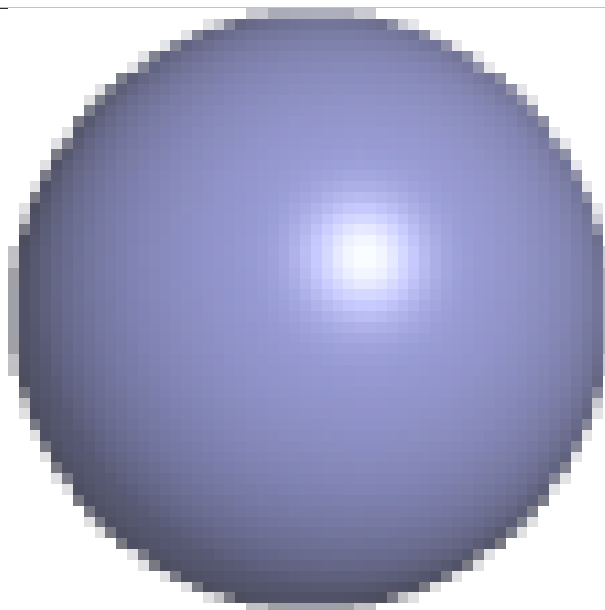
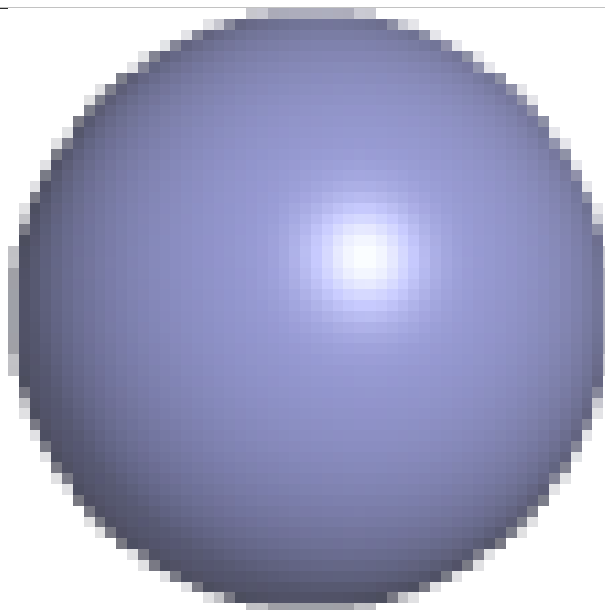
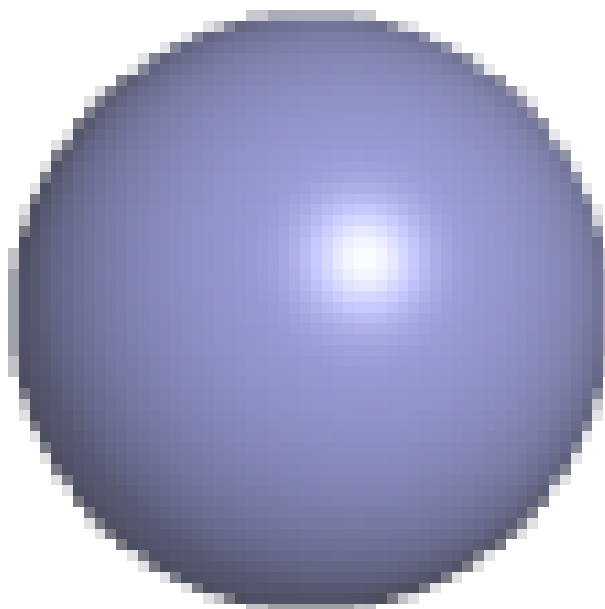
Electron density around ZN A 1003:

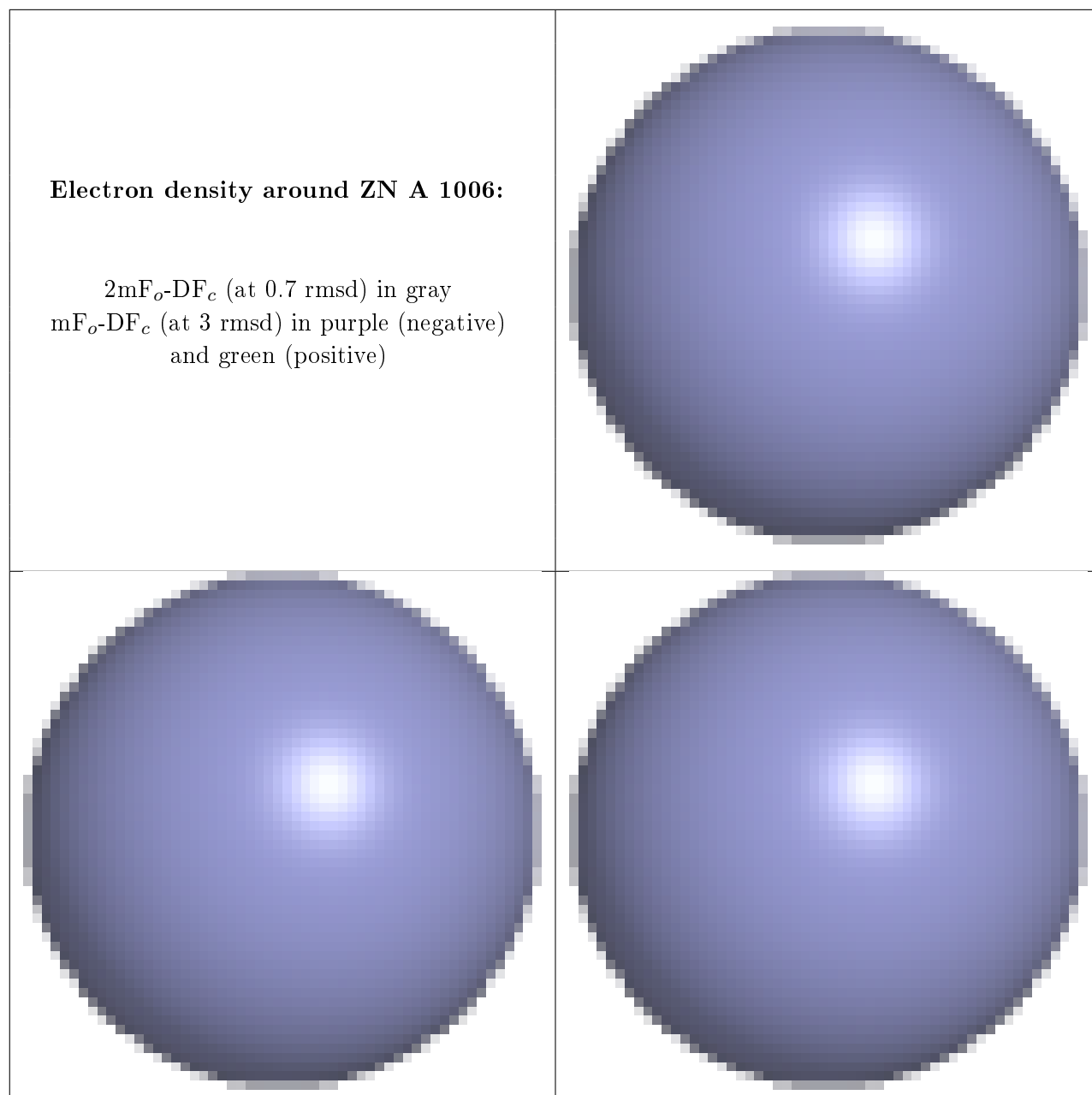
2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN A 1002:

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](#)

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