



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 10, 2022 – 12:03 PM EST

PDB ID : 7MDJ
Title : The structure of KcsA in complex with a synthetic Fab
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Deposited on : 2021-04-05
Resolution : 2.75 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
Xtriage (Phenix)	:	1.13
EDS	:	2.26
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)	:	2.26

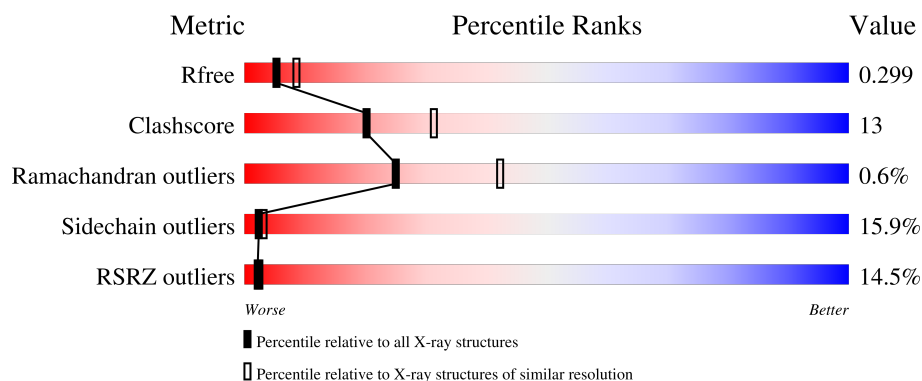
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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	229	
2	B	215	
3	C	124	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4005 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 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1599	1010	269	315	5			

- Molecule 2 is a protein called Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	196	Total	C	N	O	S	0	0	0
			1516	950	255	306	5			

- Molecule 3 is a protein called pH-gated potassium channel KcsA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	102	Total	C	N	O	S	0	0	0
			759	498	129	130	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	2	ALA	PRO	conflict	UNP P0A334
C	90	CYS	LEU	conflict	UNP P0A334

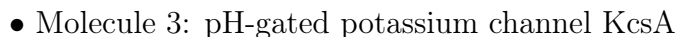
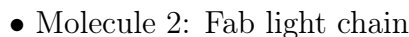
- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

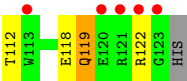
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	4	Total	K	0	0
			4	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	59	Total 59	O 59	0	0
5	B	42	Total 42	O 42	0	0
5	C	26	Total 26	O 26	0	0

- Molecule 1: Fab heavy chain





4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	138.41Å 138.41Å 71.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.01 – 2.75 97.87 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.01-2.75) 99.7 (97.87-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.201 , 0.265 0.218 , 0.299	Depositor DCC
R_{free} test set	833 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	103.5	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 91.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4005	wwPDB-VP
Average B, all atoms (Å ²)	131.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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: K

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	0.71	0/1638	0.87	0/2231
2	B	0.75	0/1548	0.96	0/2099
3	C	0.76	0/778	0.96	0/1070
All	All	0.73	0/3964	0.92	0/5400

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

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	1599	0	1546	33	0
2	B	1516	0	1477	44	0
3	C	759	0	768	29	0
4	C	4	0	0	1	0
5	A	59	0	0	3	0
5	B	42	0	0	11	0
5	C	26	0	0	9	0
All	All	4005	0	3791	96	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 13.

All (96) 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)
3:C:122:ARG:CD	5:C:301:HOH:O	1.91	1.15
3:C:122:ARG:NE	5:C:301:HOH:O	1.84	1.10
3:C:22:SER:N	3:C:25:HIS:CD2	2.22	1.07
1:A:55:ILE:HD11	3:C:49:LEU:HD13	1.65	0.78
2:B:140:PHE:CZ	5:B:325:HOH:O	2.36	0.78
3:C:122:ARG:HD3	5:C:301:HOH:O	1.66	0.75
2:B:34:ILE:HG21	2:B:72:PHE:CD1	2.24	0.72
1:A:5:VAL:HA	1:A:29:GLY:HA3	1.75	0.69
3:C:122:ARG:CZ	5:C:301:HOH:O	2.29	0.67
3:C:25:HIS:CD2	3:C:112:THR:HB	2.31	0.66
3:C:119:GLN:NE2	5:C:303:HOH:O	2.30	0.65
2:B:63:PHE:HB3	5:B:333:HOH:O	1.96	0.64
1:A:197:LEU:HD23	1:A:198:GLY:HA2	1.80	0.63
1:A:22:ARG:HB2	1:A:85:GLN:HE22	1.67	0.59
2:B:2:ASP:N	5:B:302:HOH:O	2.35	0.59
1:A:67:VAL:HG13	1:A:71:PHE:HB2	1.85	0.58
1:A:162:TRP:CH2	1:A:204:CYS:HB3	2.38	0.58
1:A:41:ARG:NE	1:A:51:ILE:HD11	2.19	0.58
1:A:20:SER:HA	1:A:86:MET:O	2.03	0.58
1:A:219:VAL:HA	5:A:317:HOH:O	2.04	0.57
2:B:114:PRO:HA	2:B:138:ASN:O	2.04	0.57
2:B:81:PRO:O	2:B:107:ILE:HD11	2.03	0.57
2:B:13:SER:HA	2:B:106:GLU:O	2.05	0.57
2:B:143:ARG:NH2	2:B:164:VAL:HG21	2.20	0.56
2:B:50:LYS:HG2	2:B:54:GLU:HB2	1.88	0.56
2:B:94:ARG:HD3	3:C:58:GLN:HG3	1.88	0.56
2:B:19:ARG:HG3	5:B:303:HOH:O	2.04	0.56
2:B:144:GLU:OE2	2:B:146:LYS:NZ	2.40	0.55
1:A:36:TRP:CZ3	3:C:62:TYR:OH	2.57	0.55
2:B:91:GLN:HE22	2:B:94:ARG:H	1.53	0.55
3:C:25:HIS:HE1	3:C:26:TRP:CE2	2.25	0.55
2:B:94:ARG:HD2	3:C:56:GLY:O	2.07	0.54
2:B:150:LYS:HG2	2:B:155:LEU:HA	1.90	0.54
2:B:106:GLU:OE1	2:B:174:TYR:OH	2.26	0.53
1:A:35:ASP:OD2	1:A:101:ARG:NH1	2.41	0.53
2:B:48:LEU:HB3	5:B:333:HOH:O	2.07	0.53
2:B:38:GLN:NE2	5:B:301:HOH:O	2.29	0.52
1:A:131:PRO:HD3	1:A:217:LYS:HD2	1.92	0.52
2:B:158:GLY:C	2:B:160:SER:N	2.63	0.52
1:A:162:TRP:CZ3	1:A:204:CYS:HB3	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ILE:CD1	3:C:49:LEU:HD13	2.39	0.52
3:C:64:ARG:HD2	5:C:313:HOH:O	2.10	0.52
1:A:149:LEU:HD13	1:A:151:LYS:HD2	1.92	0.51
3:C:22:SER:N	3:C:25:HIS:NE2	2.57	0.51
1:A:172:HIS:CD2	2:B:175:SER:HG	2.28	0.51
1:A:36:TRP:HZ3	3:C:62:TYR:CZ	2.28	0.51
2:B:112:ALA:O	2:B:140:PHE:CD1	2.64	0.51
3:C:44:SER:OG	3:C:66:LEU:HA	2.11	0.50
3:C:82:TYR:HB2	3:C:83:PRO:HD2	1.92	0.50
1:A:189:VAL:HG11	2:B:136:LEU:HD11	1.94	0.50
1:A:90:ARG:O	1:A:119:VAL:HG11	2.12	0.49
3:C:62:TYR:N	3:C:63:PRO:CD	2.76	0.49
1:A:186:LEU:HB2	5:A:316:HOH:O	2.12	0.49
2:B:38:GLN:HG3	2:B:87:TYR:CE2	2.47	0.49
2:B:152:ASP:OD2	2:B:190:HIS:HB3	2.13	0.49
2:B:140:PHE:CE1	5:B:325:HOH:O	2.61	0.49
2:B:94:ARG:CD	3:C:56:GLY:O	2.62	0.48
2:B:119:PHE:HB2	2:B:134:VAL:HB	1.96	0.48
2:B:133:VAL:HG12	2:B:149:TRP:CZ3	2.49	0.47
2:B:12:LEU:HD21	5:B:320:HOH:O	2.14	0.47
3:C:78:TYR:N	5:C:304:HOH:O	2.47	0.47
3:C:62:TYR:HB2	3:C:63:PRO:HD3	1.96	0.47
1:A:63:TYR:HE1	1:A:73:ILE:HG13	1.80	0.46
1:A:94:THR:HG23	1:A:118:THR:HA	1.97	0.46
2:B:126:LEU:HD11	2:B:187:TYR:HE2	1.81	0.46
1:A:209:LYS:HD3	1:A:209:LYS:HA	1.49	0.46
2:B:67:ARG:HB3	2:B:72:PHE:HA	1.97	0.46
3:C:66:LEU:HD12	3:C:69:SER:OG	2.17	0.45
2:B:140:PHE:HE2	2:B:144:GLU:O	1.99	0.45
3:C:118:GLU:HB3	5:C:314:HOH:O	2.16	0.45
1:A:38:HIS:HA	1:A:52:GLY:O	2.17	0.45
2:B:67:ARG:HD2	2:B:67:ARG:HA	1.63	0.45
1:A:214:LYS:HE3	1:A:214:LYS:HB2	1.53	0.45
3:C:64:ARG:HB2	5:C:306:HOH:O	2.16	0.45
1:A:211:SER:O	1:A:212:ASN:C	2.55	0.44
2:B:81:PRO:HA	2:B:107:ILE:HD13	2.00	0.44
2:B:158:GLY:HA2	5:B:305:HOH:O	2.17	0.44
1:A:42:GLN:OE1	1:A:46:LYS:O	2.36	0.44
2:B:30:ILE:HD11	2:B:72:PHE:CE2	2.53	0.44
2:B:78:SER:N	5:B:303:HOH:O	2.46	0.44
1:A:197:LEU:HA	1:A:198:GLY:HA2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:44:SER:HA	3:C:65:ALA:HB1	1.99	0.44
2:B:33:ASP:HB3	2:B:92:SER:OG	2.18	0.43
2:B:67:ARG:HE	2:B:68:SER:H	1.65	0.43
1:A:172:HIS:CD2	2:B:175:SER:OG	2.71	0.43
2:B:81:PRO:O	2:B:107:ILE:CD1	2.66	0.43
1:A:125:LYS:HD2	1:A:125:LYS:HA	1.44	0.42
1:A:186:LEU:HD12	1:A:186:LEU:H	1.83	0.42
5:B:306:HOH:O	3:C:55:PRO:HB3	2.18	0.42
2:B:133:VAL:CG1	2:B:149:TRP:CZ3	3.03	0.41
2:B:141:TYR:CD1	2:B:142:PRO:HA	2.55	0.41
1:A:152:ASP:OD1	5:A:301:HOH:O	2.21	0.41
3:C:75:THR:O	3:C:75:THR:OG1	2.38	0.41
4:C:201:K:K	4:C:203:K:K	2.25	0.41
1:A:104:GLY:HA3	3:C:62:TYR:CE1	2.56	0.41
2:B:164:VAL:HG22	2:B:176:LEU:HD12	2.03	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	208/229 (91%)	186 (89%)	21 (10%)	1 (0%)	29	47
2	B	192/215 (89%)	171 (89%)	20 (10%)	1 (0%)	29	47
3	C	100/124 (81%)	94 (94%)	5 (5%)	1 (1%)	15	27
All	All	500/568 (88%)	451 (90%)	46 (9%)	3 (1%)	25	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	77	GLY
2	B	78	SER

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Mol	Chain	Res	Type
1	A	134	PRO

5.3.2 Protein sidechains ⓘ

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	175/190 (92%)	138 (79%)	37 (21%)	1	1
2	B	174/191 (91%)	148 (85%)	26 (15%)	3	4
3	C	72/91 (79%)	68 (94%)	4 (6%)	21	36
All	All	421/472 (89%)	354 (84%)	67 (16%)	2	3

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	15	VAL
1	A	16	GLN
1	A	20	SER
1	A	24	SER
1	A	30	TYR
1	A	31	THR
1	A	32	PHE
1	A	33	THR
1	A	51	ILE
1	A	55	ILE
1	A	57	SER
1	A	58	TYR
1	A	65	ASP
1	A	67	VAL
1	A	74	SER
1	A	77	THR
1	A	78	SER
1	A	87	ASN
1	A	101	ARG
1	A	105	ASP

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Mol	Chain	Res	Type
1	A	116	LEU
1	A	118	THR
1	A	120	SER
1	A	121	SER
1	A	125	LYS
1	A	159	THR
1	A	168	THR
1	A	186	LEU
1	A	189	VAL
1	A	192	VAL
1	A	194	SER
1	A	199	THR
1	A	200	GLN
1	A	209	LYS
1	A	214	LYS
1	A	218	LYS
2	B	6	THR
2	B	8	SER
2	B	16	VAL
2	B	19	ARG
2	B	23	THR
2	B	43	LYS
2	B	50	LYS
2	B	67	ARG
2	B	68	SER
2	B	74	LEU
2	B	77	SER
2	B	90	GLN
2	B	91	GLN
2	B	104	LYS
2	B	110	THR
2	B	115	SER
2	B	143	ARG
2	B	147	VAL
2	B	148	GLN
2	B	150	LYS
2	B	169	SER
2	B	170	LYS
2	B	184	LYS
2	B	186	ASP
2	B	190	HIS
2	B	192	VAL

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Mol	Chain	Res	Type
3	C	25	HIS
3	C	71	GLU
3	C	105	LEU
3	C	119	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	42	GLN
1	A	85	GLN
1	A	87	ASN
1	A	172	HIS
1	A	179	GLN
1	A	200	GLN
2	B	38	GLN
2	B	39	GLN
2	B	90	GLN
2	B	91	GLN
2	B	139	ASN
2	B	148	GLN
3	C	25	HIS
3	C	119	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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

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	214/229 (93%)	0.75	32 (14%) 2 2	88, 141, 191, 217	0
2	B	196/215 (91%)	0.69	28 (14%) 2 2	78, 124, 167, 205	0
3	C	102/124 (82%)	1.05	14 (13%) 3 3	73, 103, 203, 236	0
All	All	512/568 (90%)	0.79	74 (14%) 2 2	73, 129, 187, 236	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	121	ARG	10.2
2	B	193	TYR	10.1
3	C	123	GLY	8.6
3	C	23	ALA	7.8
3	C	22	SER	7.7
1	A	68	LYS	7.7
1	A	63	TYR	6.1
1	A	65	ASP	5.5
2	B	133	VAL	5.4
2	B	192	VAL	5.4
1	A	32	PHE	5.3
1	A	103	ARG	5.1
1	A	30	TYR	5.1
2	B	4	GLN	5.0
1	A	61	ALA	4.8
2	B	151	VAL	4.8
3	C	120	GLU	4.7
2	B	120	PRO	4.5
1	A	66	SER	4.4
2	B	131	ALA	4.4
2	B	101	GLN	4.3
1	A	67	VAL	4.1
1	A	69	GLY	4.0

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Mol	Chain	Res	Type	RSRZ
2	B	132	SER	4.0
2	B	68	SER	4.0
3	C	122	ARG	3.9
2	B	127	LYS	3.8
1	A	97	TYR	3.7
1	A	105	ASP	3.7
1	A	39	TRP	3.6
3	C	49	LEU	3.4
3	C	87	TRP	3.3
3	C	24	LEU	3.3
1	A	107	TYR	3.2
2	B	152	ASP	3.2
1	A	78	SER	3.2
2	B	121	PRO	3.2
1	A	31	THR	3.1
2	B	197	VAL	3.1
1	A	83	TYR	3.0
1	A	71	PHE	3.0
1	A	41	ARG	3.0
1	A	51	ILE	3.0
2	B	161	GLN	3.0
3	C	27	ARG	2.9
3	C	113	TRP	2.9
2	B	187	TYR	2.8
1	A	73	ILE	2.7
3	C	57	ALA	2.7
2	B	134	VAL	2.7
1	A	77	THR	2.7
2	B	56	ILE	2.6
3	C	25	HIS	2.5
2	B	3	ILE	2.5
3	C	86	LEU	2.5
2	B	191	LYS	2.4
2	B	88	TYR	2.4
2	B	94	ARG	2.4
2	B	117	PHE	2.4
2	B	19	ARG	2.3
1	A	60	ARG	2.3
2	B	153	ASN	2.3
1	A	58	TYR	2.3
1	A	133	ALA	2.2
1	A	48	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	171	VAL	2.2
1	A	37	ILE	2.1
1	A	84	LEU	2.1
1	A	36	TRP	2.1
2	B	50	LYS	2.1
2	B	54	GLU	2.1
1	A	190	VAL	2.0
2	B	194	ALA	2.0
1	A	35	ASP	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](#)

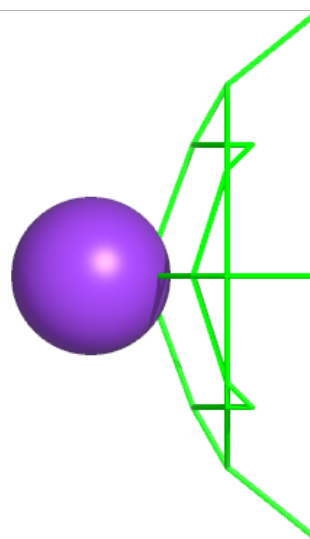
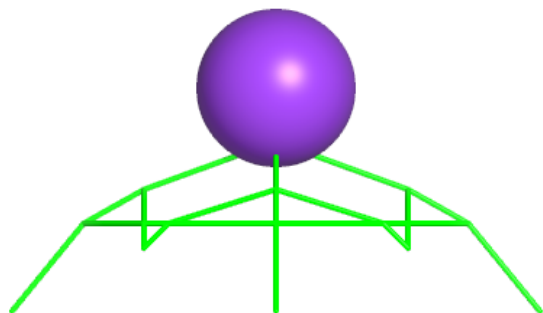
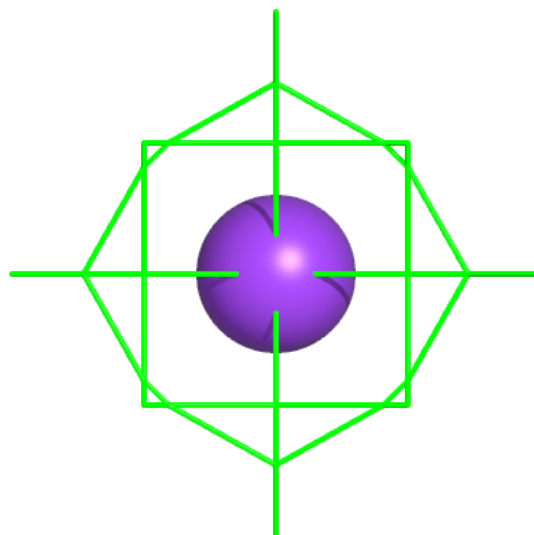
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	K	C	202	1/1	0.47	0.29	143,143,143,143	1
4	K	C	203	1/1	0.94	0.25	137,137,137,137	1
4	K	C	201	1/1	0.97	0.32	109,109,109,109	1
4	K	C	204	1/1	1.00	0.39	127,127,127,127	1

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.

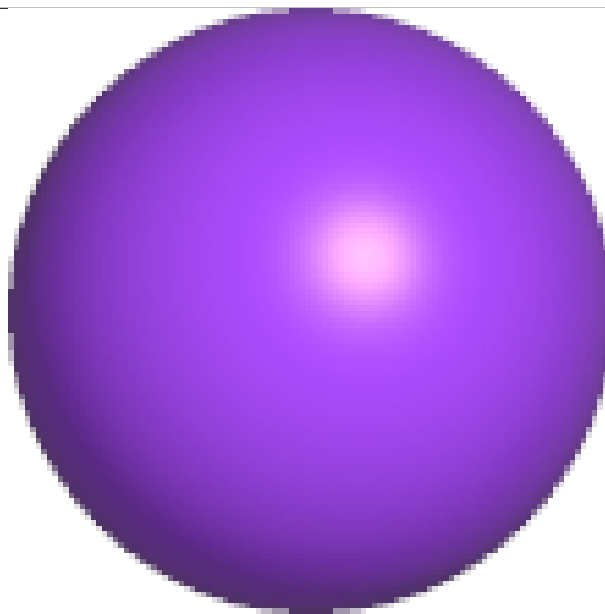
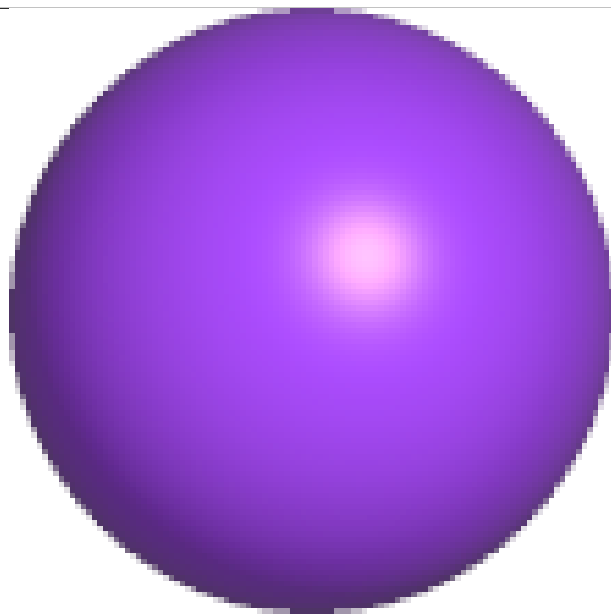
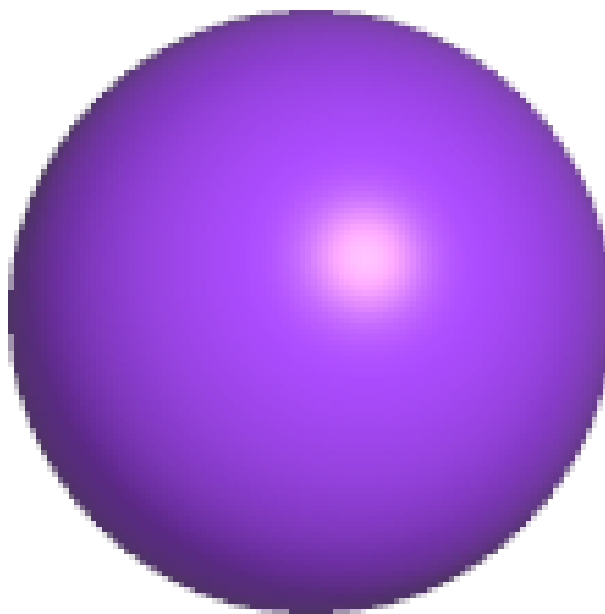
Electron density around K C 202:

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



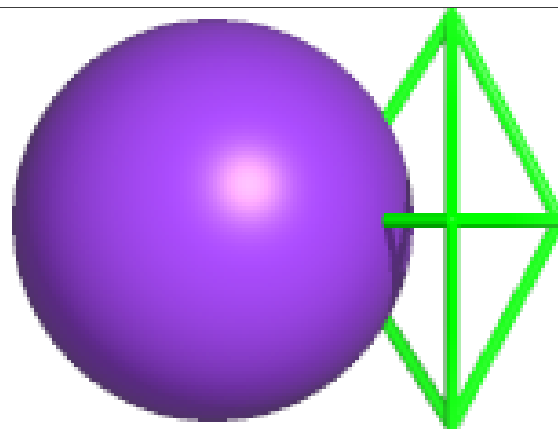
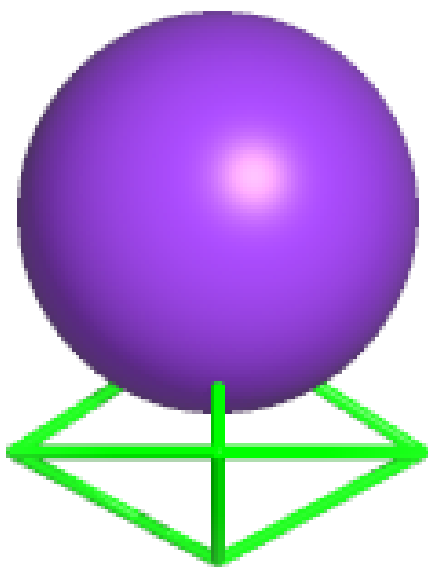
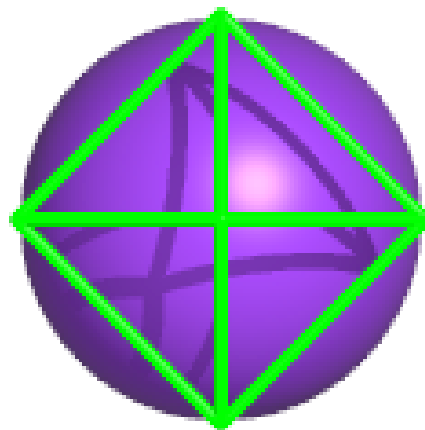
Electron density around K C 203:

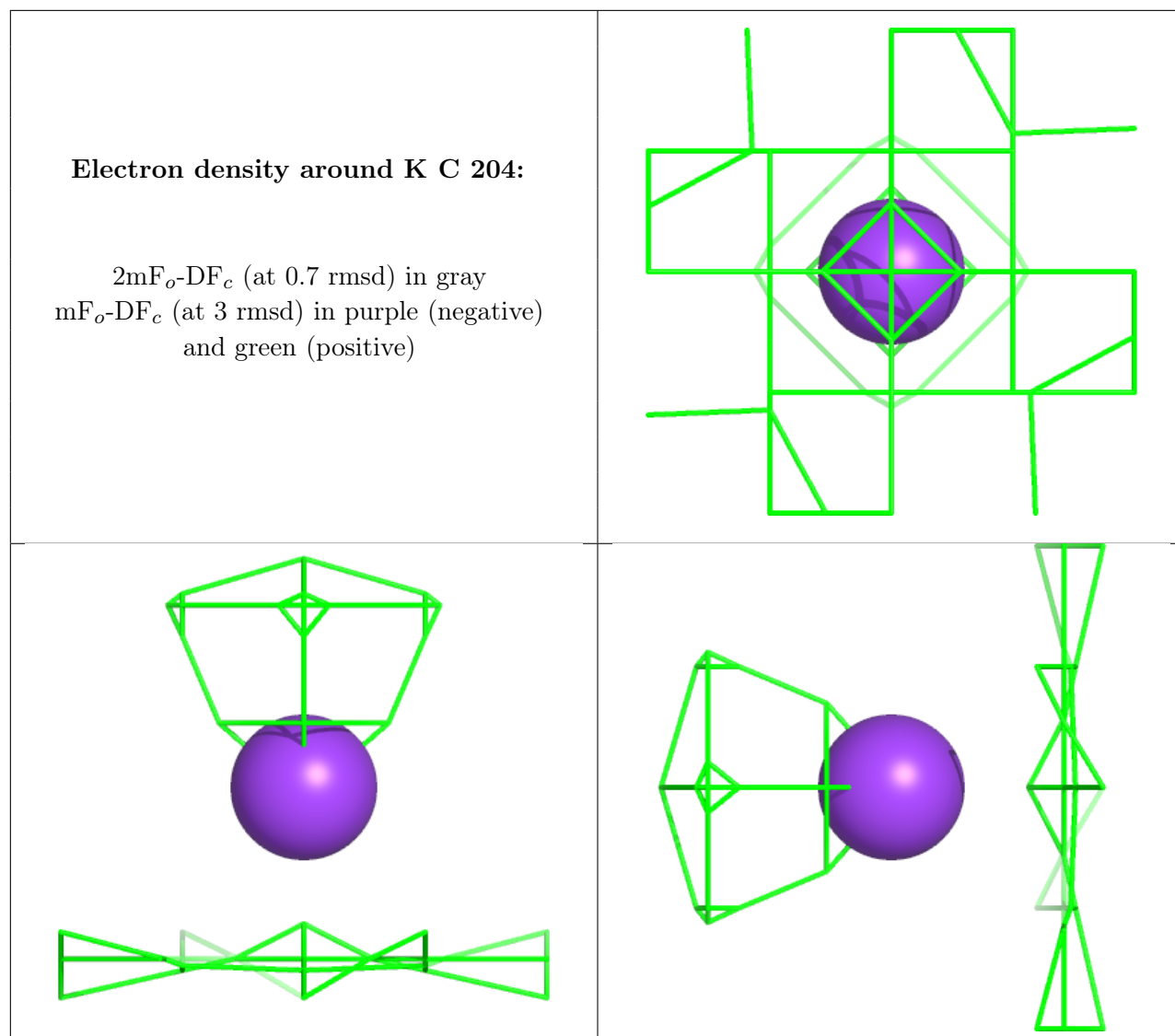
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around K C 201:

$2mF_o-DF_c$ (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.