



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:23 PM UTC

PDB ID : 1K54 / pdb_00001k54
Title : OXA-10 class D beta-lactamase partially acylated with reacted 6beta-(1-hydroxy-1-methylethyl) penicillanic acid
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Deposited on : 2001-10-10
Resolution : 1.70 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

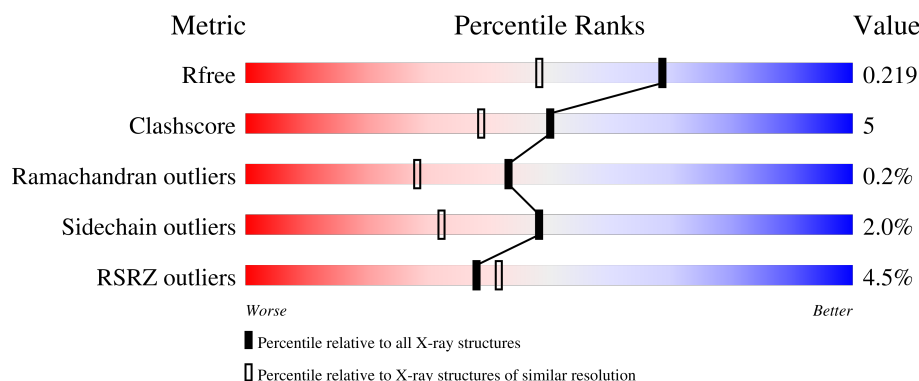
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	246	6% 87% 12% .
1	B	246	2% 91% 8%
2	C	246	5% 84% 13% ..
2	D	246	5% 90% 9% ..

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
4	HOQ	A	967	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8633 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 Beta lactamase OXA-10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	244	Total	C	N	O	S	0	4	0
			1929	1234	324	364	7			
1	B	245	Total	C	N	O	S	0	5	0
			1945	1244	323	371	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	70	KCX	LYS	modified residue	UNP P14489
B	70	KCX	LYS	modified residue	UNP P14489

- Molecule 2 is a protein called Beta lactamase OXA-10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	240	Total	C	N	O	S	0	2	0
			1877	1204	312	355	6			
2	D	244	Total	C	N	O	S	0	4	0
			1913	1222	316	369	6			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



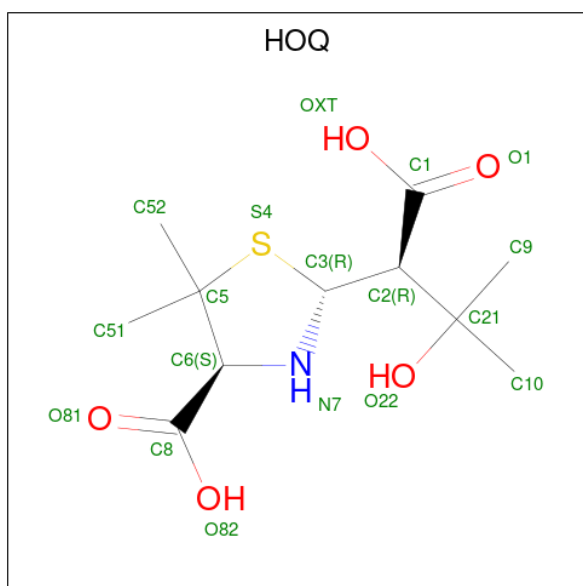
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	1
			10	8	2		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is (1R)-2-(1-CARBOXY-2-HYDROXY-2-METHYL-PROPYL)-5,5-DIMETHYL-L-THIAZOLIDINE-4-CARBOXYLIC ACID (CCD ID: HOQ) (formula: C₁₁H₁₉NO₅S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0
			17	11	1	4	1	0
4	B	1	Total	C	N	O	S	0
			17	11	1	4	1	0

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).

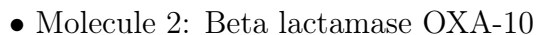


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	153	Total	O	0	0
			153	153		
6	B	231	Total	O	0	0
			231	231		
6	C	186	Total	O	0	0
			186	186		
6	D	251	Total	O	0	0
			251	251		

- Molecule 1: Beta lactamase OXA-10



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.65Å 82.48Å 101.96Å 90.00° 95.38° 90.00°	Depositor
Resolution (Å)	44.77 – 1.70 44.77 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (44.77-1.70) 99.0 (44.77-1.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 1.70Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.177 , 0.210 0.188 , 0.219	Depositor DCC
R_{free} test set	1794 reflections (1.50%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.4	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	0.95	EDS
Total number of atoms	8633	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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: HOQ, EDO, KCX, SO4

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	0.41	0/1973	0.77	2/2666 (0.1%)
1	B	0.44	0/1993	0.80	2/2693 (0.1%)
2	C	0.44	0/1924	0.76	0/2604
2	D	0.47	0/1969	0.76	0/2662
All	All	0.44	0/7859	0.77	4/10625 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	66	ALA	CA-C-N	6.69	134.31	121.54
1	B	66	ALA	C-N-CA	6.69	134.31	121.54
1	A	66	ALA	CA-C-N	6.45	133.87	121.54
1	A	66	ALA	C-N-CA	6.45	133.87	121.54

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	1929	0	1902	31	0
1	B	1945	0	1919	17	0
2	C	1877	0	1845	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1913	0	1868	19	0
3	A	20	0	0	1	0
3	B	25	0	0	1	0
3	C	25	0	0	0	0
3	D	20	0	0	0	0
4	A	17	0	17	8	0
4	B	17	0	17	1	0
5	A	8	0	12	1	0
5	B	8	0	12	1	0
5	D	8	0	12	2	0
6	A	153	0	0	6	0
6	B	231	0	0	5	0
6	C	186	0	0	9	0
6	D	251	0	0	3	0
All	All	8633	0	7604	82	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 5.

All (82) 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:155:LEU:HD22	4:A:967:HOQ:H91	1.16	1.16
1:A:155:LEU:HD22	4:A:967:HOQ:C9	1.91	0.99
5:A:1001:EDO:H22	6:A:1110:HOH:O	1.68	0.94
1:B:182:LYS:HD3	2:D:183[A]:GLU:OE2	1.76	0.86
2:D:199:GLU:HA	5:D:1005:EDO:H11	1.59	0.85
1:A:206:THR:OG1	3:A:905:SO4:O2	2.02	0.76
2:C:134:LYS:HB2	6:C:1093:HOH:O	1.84	0.76
2:C:244:GLU:CG	6:C:1036:HOH:O	2.36	0.74
2:C:255:THR:O	2:C:259:GLU:HG3	1.93	0.69
1:A:152:LYS:HD3	6:A:1108:HOH:O	1.95	0.66
1:B:206:THR:OG1	3:B:908:SO4:O1	2.13	0.66
2:D:110:GLY:O	2:D:114:VAL:HG22	1.95	0.66
2:C:110:GLY:O	2:C:114:VAL:HG22	1.97	0.64
1:A:229:GLU:OE1	2:C:104:ARG:NH1	2.31	0.63
1:B:199:GLU:HA	5:B:1003:EDO:H22	1.80	0.62
1:A:203:HIS:HD2	6:A:1023:HOH:O	1.83	0.61
6:A:1108:HOH:O	2:D:214:GLU:HG3	2.00	0.61
1:B:203:HIS:HD2	6:B:1014:HOH:O	1.84	0.59
1:A:208:PHE:O	4:A:967:HOQ:H3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:LYS:HE3	6:B:1233:HOH:O	2.03	0.58
1:B:182:LYS:HD3	2:D:183[A]:GLU:CD	2.29	0.58
1:B:99[B]:MET:SD	4:B:967:HOQ:H511	2.45	0.57
2:C:203:HIS:HD2	6:C:913:HOH:O	1.89	0.56
1:B:152:LYS:HE3	6:B:1143:HOH:O	2.05	0.56
2:D:93:ASP:OD1	2:D:94:GLY:N	2.36	0.55
2:C:243:ASN:HD22	2:C:243:ASN:C	2.15	0.55
1:A:149:GLY:HA3	2:D:214:GLU:OE2	2.08	0.54
1:A:99[B]:MET:SD	4:A:967:HOQ:H511	2.47	0.54
1:A:149:GLY:CA	2:D:214:GLU:OE2	2.55	0.54
1:A:155:LEU:CD2	4:A:967:HOQ:C9	2.78	0.53
2:D:203:HIS:HD2	6:D:1044:HOH:O	1.90	0.53
2:D:45:LYS:HE2	2:D:231[B]:GLU:OE2	2.10	0.52
1:A:99[B]:MET:HE2	6:C:1087:HOH:O	2.10	0.51
1:B:44:CYS:SG	1:B:167:VAL:HG11	2.49	0.51
1:B:100:LYS:CE	6:B:1232:HOH:O	2.58	0.51
1:B:100:LYS:HE3	6:B:1232:HOH:O	2.10	0.51
1:B:190:GLU:OE1	2:D:203:HIS:HE1	1.92	0.51
1:A:176:ASN:ND2	1:A:182:LYS:HE3	2.25	0.51
1:A:253:ILE:HB	1:A:254:PRO:HD3	1.93	0.50
2:D:229:GLU:O	5:D:1005:EDO:H12	2.12	0.49
1:A:35:GLU:HB2	1:A:37:VAL:HG23	1.93	0.49
1:A:155:LEU:CD2	4:A:967:HOQ:O22	2.60	0.49
1:A:203:HIS:HE1	2:C:190:GLU:OE1	1.97	0.47
2:C:142:GLY:H	2:C:165:ASN:ND2	2.12	0.47
1:A:152:LYS:HD3	6:D:1227:HOH:O	2.13	0.47
2:C:252:SER:O	2:C:256:LYS:HG3	2.15	0.47
2:C:256:LYS:HG3	6:C:991:HOH:O	2.14	0.47
2:C:129:GLU:HG3	2:C:146:ILE:HD11	1.96	0.47
1:A:22:ILE:O	1:A:22:ILE:HG23	2.14	0.46
2:C:259:GLU:CD	6:C:1089:HOH:O	2.59	0.46
2:C:216:ASN:HB3	6:C:1081:HOH:O	2.14	0.46
1:A:154:TRP:NE1	1:A:155:LEU:HD21	2.30	0.46
2:C:44:CYS:SG	2:C:167:VAL:HG11	2.56	0.46
2:C:253:ILE:HB	2:C:254:PRO:HD3	1.98	0.46
1:A:99[B]:MET:CE	6:C:1087:HOH:O	2.62	0.46
1:B:199:GLU:OE2	2:D:104:ARG:NH1	2.49	0.45
1:A:67:SER:H	4:A:967:HOQ:C1	2.29	0.45
2:D:142:GLY:H	2:D:165:ASN:ND2	2.15	0.45
1:A:190:GLU:OE1	2:C:203:HIS:HE1	2.00	0.45
1:A:45:LYS:O	1:A:46:SER:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:HIS:HE1	2:D:190:GLU:OE1	1.98	0.44
1:A:199:GLU:OE1	2:C:104:ARG:NH2	2.50	0.44
2:D:142:GLY:H	2:D:165:ASN:HD21	1.66	0.43
1:B:176:ASN:ND2	1:B:182:LYS:HE3	2.33	0.43
1:A:109[B]:ARG:NH2	6:A:1110:HOH:O	2.42	0.43
2:C:117:VAL:HB	2:C:118:PRO:HD3	2.00	0.43
2:C:237:PHE:CE1	2:C:250:ARG:HA	2.53	0.43
2:C:256:LYS:CG	6:C:991:HOH:O	2.67	0.43
2:D:247:LEU:N	2:D:248:PRO:CD	2.82	0.43
2:C:255:THR:HG23	2:C:264:ILE:CD1	2.49	0.42
2:D:44:CYS:SG	2:D:167:VAL:HG11	2.59	0.42
2:C:243:ASN:HD22	2:C:245:SER:H	1.68	0.42
1:B:100:LYS:HE3	1:B:100:LYS:HB2	1.83	0.42
2:C:247:LEU:N	2:C:248:PRO:CD	2.82	0.42
1:A:220:ALA:O	1:A:238:ASN:HA	2.20	0.42
2:C:243:ASN:HB3	2:C:246:LYS:HG3	2.02	0.42
1:A:199:GLU:OE2	2:C:104:ARG:CZ	2.68	0.41
2:D:87:HIS:HE1	6:D:1148:HOH:O	2.03	0.41
1:A:22:ILE:N	6:A:1155:HOH:O	2.54	0.41
1:B:110:GLY:O	1:B:114:VAL:HG22	2.21	0.41
1:A:155:LEU:HD22	4:A:967:HOQ:C21	2.49	0.41
1:A:110:GLY:O	1:A:114:VAL:HG22	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	245/246 (100%)	238 (97%)	6 (2%)	1 (0%)	30	16
1	B	247/246 (100%)	240 (97%)	6 (2%)	1 (0%)	30	16
2	C	238/246 (97%)	232 (98%)	6 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	244/246 (99%)	237 (97%)	7 (3%)	0	100	100
All	All	974/984 (99%)	947 (97%)	25 (3%)	2 (0%)	43	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	SER
1	B	67	SER

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	207/209 (99%)	205 (99%)	2 (1%)	68	58
1	B	211/209 (101%)	207 (98%)	4 (2%)	50	34
2	C	201/210 (96%)	196 (98%)	5 (2%)	42	24
2	D	206/210 (98%)	200 (97%)	6 (3%)	37	20
All	All	825/838 (98%)	808 (98%)	17 (2%)	48	30

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	166	GLN
1	B	73	ASN
1	B	100	LYS
1	B	155	LEU
1	B	166	GLN
2	C	33	SER
2	C	73	ASN
2	C	166	GLN
2	C	243	ASN
2	C	264	ILE
2	D	73	ASN

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Mol	Chain	Res	Type
2	D	166	GLN
2	D	214	GLU
2	D	231[A]	GLU
2	D	231[B]	GLU
2	D	244	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	73	ASN
1	A	158	GLN
1	A	166	GLN
1	A	203	HIS
1	B	38	ASN
1	B	73	ASN
1	B	166	GLN
1	B	203	HIS
2	C	38	ASN
2	C	73	ASN
2	C	165	ASN
2	C	166	GLN
2	C	203	HIS
2	C	243	ASN
2	D	73	ASN
2	D	121	GLN
2	D	165	ASN
2	D	166	GLN
2	D	203	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	$\# Z > 2$
1	KCX	B	70	1	10,11,12	0.79	0	6,12,14	1.24	1 (16%)
1	KCX	A	70	1	10,11,12	0.93	0	6,12,14	1.24	1 (16%)

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
1	KCX	B	70	1	-	0/9/10/12	-
1	KCX	A	70	1	-	0/9/10/12	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	KCX	OQ1-CX-NZ	-2.95	120.44	124.92
1	B	70	KCX	OQ1-CX-NZ	-2.85	120.59	124.92

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

26 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
5	EDO	A	1002	-	3,3,3	0.26	0	2,2,2	0.31	0
3	SO4	B	909	-	4,4,4	0.25	0	6,6,6	0.17	0
3	SO4	C	901[B]	-	4,4,4	0.25	0	6,6,6	0.12	0
3	SO4	C	904	-	4,4,4	0.24	0	6,6,6	0.13	0
4	HOQ	A	967	1	13,17,18	1.13	1 (7%)	16,27,29	1.02	1 (6%)
3	SO4	D	917	-	4,4,4	0.23	0	6,6,6	0.07	0
3	SO4	D	910	-	4,4,4	0.24	0	6,6,6	0.07	0
5	EDO	A	1001	-	3,3,3	0.26	0	2,2,2	0.25	0
3	SO4	D	916	-	4,4,4	0.22	0	6,6,6	0.11	0
5	EDO	B	1004	-	3,3,3	0.29	0	2,2,2	0.19	0
3	SO4	A	914	-	4,4,4	0.24	0	6,6,6	0.17	0
3	SO4	B	913	-	4,4,4	0.23	0	6,6,6	0.11	0
3	SO4	C	907	-	4,4,4	0.25	0	6,6,6	0.13	0
5	EDO	D	1005	-	3,3,3	0.28	0	2,2,2	0.42	0
3	SO4	B	912	-	4,4,4	0.26	0	6,6,6	0.06	0
3	SO4	C	902	-	4,4,4	0.25	0	6,6,6	0.07	0
3	SO4	C	901[A]	-	4,4,4	0.25	0	6,6,6	0.06	0
3	SO4	A	903	-	4,4,4	0.24	0	6,6,6	0.10	0
3	SO4	A	905	-	4,4,4	0.22	0	6,6,6	0.15	0
3	SO4	A	906	-	4,4,4	0.23	0	6,6,6	0.16	0
3	SO4	B	915	-	4,4,4	0.23	0	6,6,6	0.06	0
5	EDO	D	1006	-	3,3,3	0.24	0	2,2,2	0.24	0
5	EDO	B	1003	-	3,3,3	0.35	0	2,2,2	0.30	0
3	SO4	B	908	-	4,4,4	0.23	0	6,6,6	0.14	0
4	HOQ	B	967	1	13,17,18	1.07	1 (7%)	16,27,29	0.62	0
3	SO4	D	911	-	4,4,4	0.19	0	6,6,6	0.18	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
5	EDO	D	1006	-	-	1/1/1/1	-
5	EDO	B	1003	-	-	1/1/1/1	-
5	EDO	A	1001	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	B	1004	-	-	1/1/1/1	-
5	EDO	A	1002	-	-	0/1/1/1	-
5	EDO	D	1005	-	-	1/1/1/1	-
4	HOQ	B	967	1	-	6/11/31/33	0/1/1/1
4	HOQ	A	967	1	-	5/11/31/33	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	967	HOQ	C5-S4	-2.18	1.81	1.85
4	A	967	HOQ	C5-S4	-2.16	1.81	1.85

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	967	HOQ	C9-C21-C2	-2.35	107.95	111.82

There are no chirality outliers.

All (15) torsion outliers are listed below:

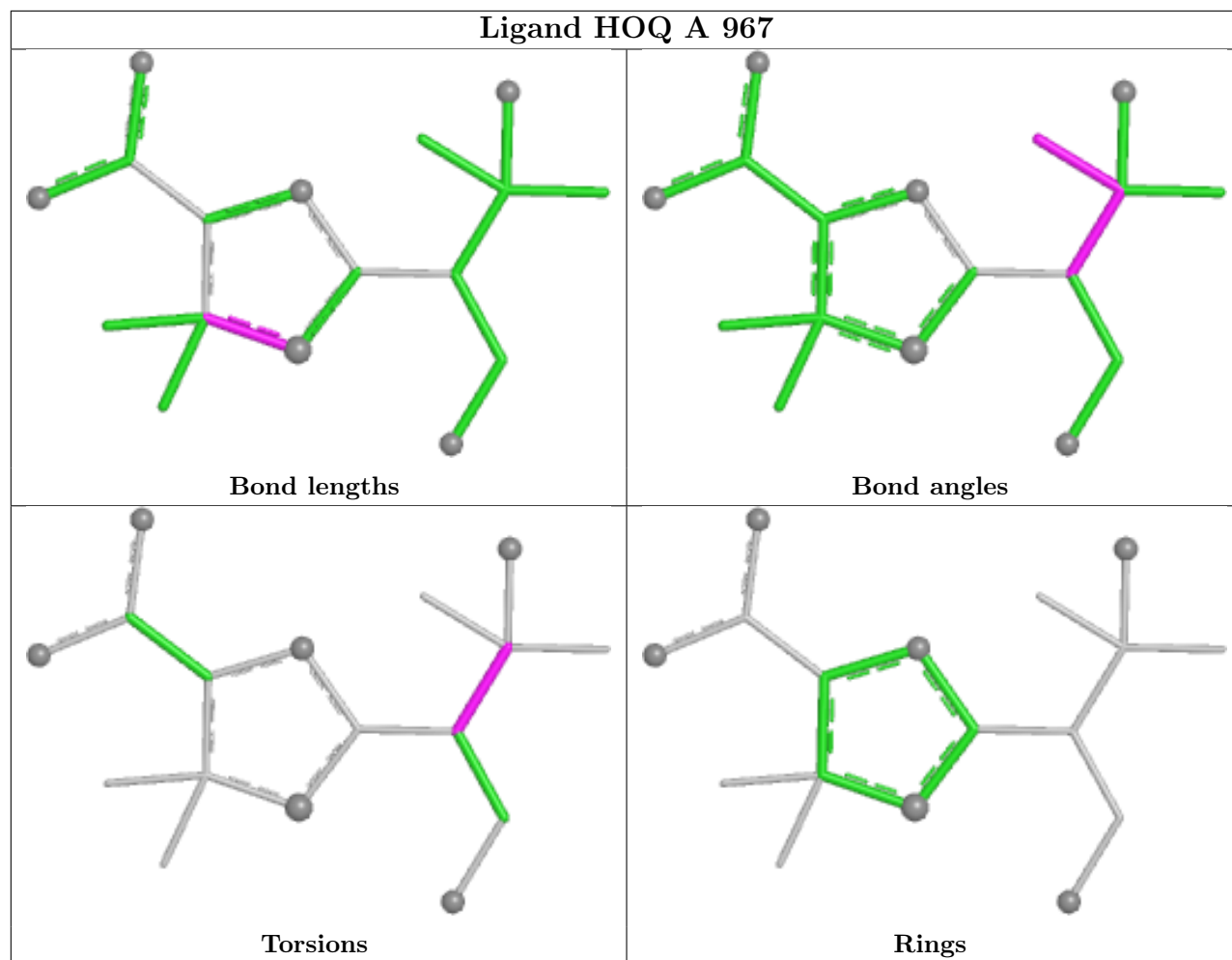
Mol	Chain	Res	Type	Atoms
4	A	967	HOQ	C1-C2-C21-O22
4	A	967	HOQ	C1-C2-C21-C9
4	A	967	HOQ	C1-C2-C21-C10
4	B	967	HOQ	C1-C2-C21-O22
4	B	967	HOQ	C1-C2-C21-C9
4	B	967	HOQ	C1-C2-C21-C10
5	D	1005	EDO	O1-C1-C2-O2
4	B	967	HOQ	N7-C6-C8-O81
4	B	967	HOQ	N7-C6-C8-O82
4	A	967	HOQ	C3-C2-C21-C10
5	D	1006	EDO	O1-C1-C2-O2
5	B	1003	EDO	O1-C1-C2-O2
5	B	1004	EDO	O1-C1-C2-O2
4	A	967	HOQ	C3-C2-C21-C9
4	B	967	HOQ	C3-C2-C21-C10

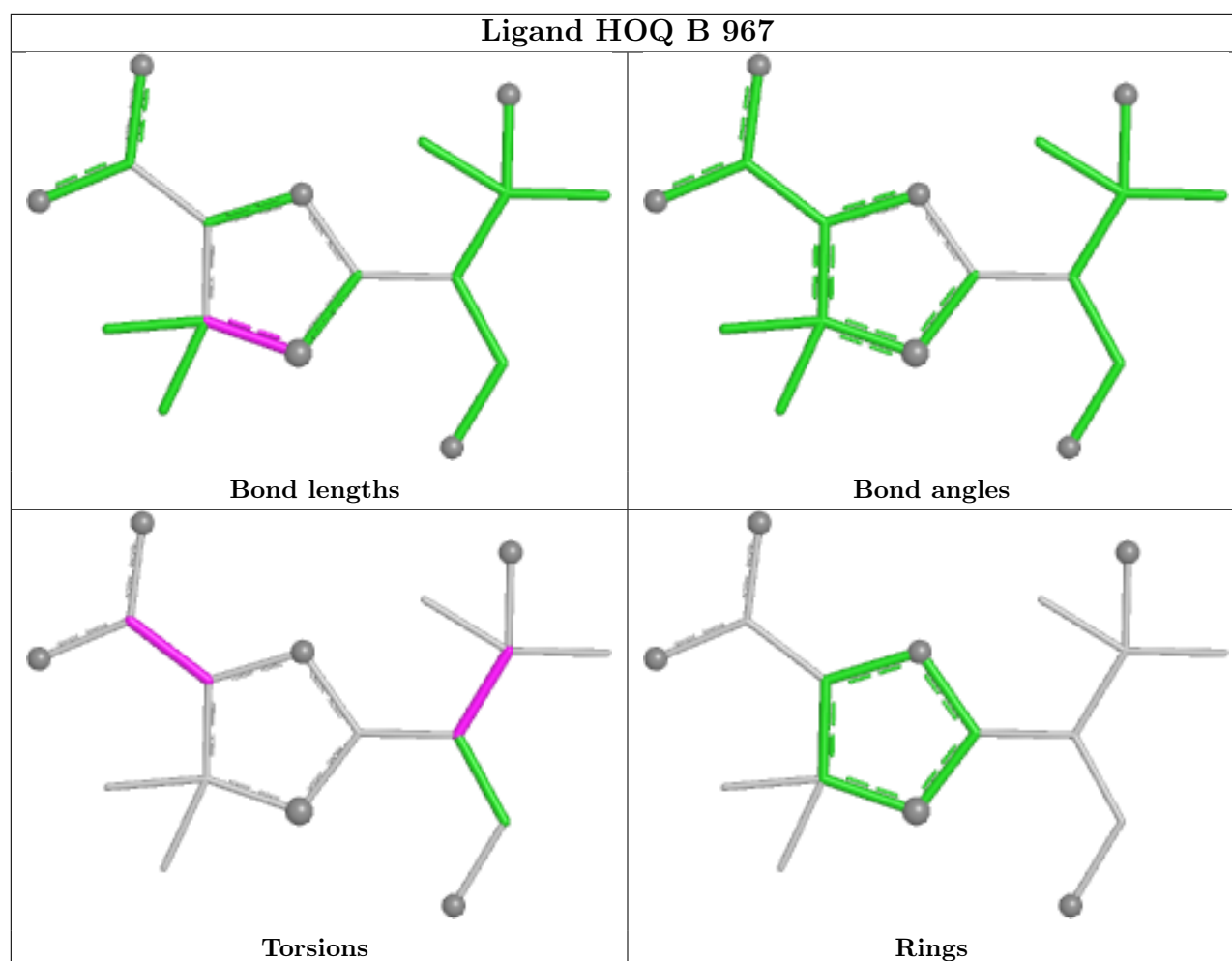
There are no ring outliers.

7 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	967	HOQ	8	0
5	A	1001	EDO	1	0
5	D	1005	EDO	2	0
3	A	905	SO4	1	0
5	B	1003	EDO	1	0
3	B	908	SO4	1	0
4	B	967	HOQ	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

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	243/246 (98%)	0.39	14 (5%) 29 32	8, 21, 34, 39	4 (1%)
1	B	244/246 (99%)	-0.01	5 (2%) 65 69	10, 16, 23, 29	5 (2%)
2	C	240/246 (97%)	0.24	12 (5%) 34 38	10, 19, 31, 40	2 (0%)
2	D	244/246 (99%)	0.01	13 (5%) 32 35	8, 14, 29, 40	4 (1%)
All	All	971/984 (98%)	0.16	44 (4%) 38 42	8, 17, 31, 40	15 (1%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	22	ILE	6.5
2	D	94	GLY	4.4
2	D	266	GLY	4.3
1	B	21	SER	3.9
2	D	21	SER	3.7
1	A	23	THR	3.7
1	A	155	LEU	3.6
2	D	98	ALA	3.4
1	A	51	CYS	3.1
2	C	102	TRP	3.0
1	A	265	GLY	2.9
2	D	214	GLU	2.9
2	C	92	TRP	2.8
1	A	216	ASN	2.8
1	A	44	CYS	2.8
2	C	98	ALA	2.7
2	C	34	ALA	2.7
1	A	49	LYS	2.7
1	A	50	SER	2.6
2	D	99	MET	2.6
2	D	97	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
2	D	93	ASP	2.5
2	C	99	MET	2.5
1	B	265	GLY	2.5
2	D	265	GLY	2.5
1	A	32	PHE	2.4
1	A	33	SER	2.4
2	C	36	ALA	2.4
2	D	50	SER	2.4
2	C	265	GLY	2.3
2	D	92	TRP	2.3
1	B	44	CYS	2.3
2	C	104	ARG	2.3
2	C	44	CYS	2.2
1	A	215[A]	SER	2.2
1	A	36	ALA	2.2
2	D	264	ILE	2.2
2	C	134	LYS	2.1
1	B	147	SER	2.1
2	C	93	ASP	2.1
1	A	35	GLU	2.1
2	C	50	SER	2.1
2	D	244	GLU	2.0
1	B	50	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	KCX	A	70	12/13	0.95	0.07	13,14,17,18	0
1	KCX	B	70	12/13	0.97	0.05	11,12,16,16	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

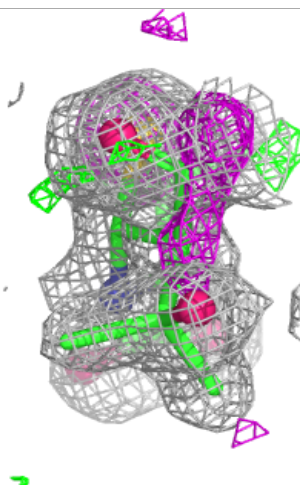
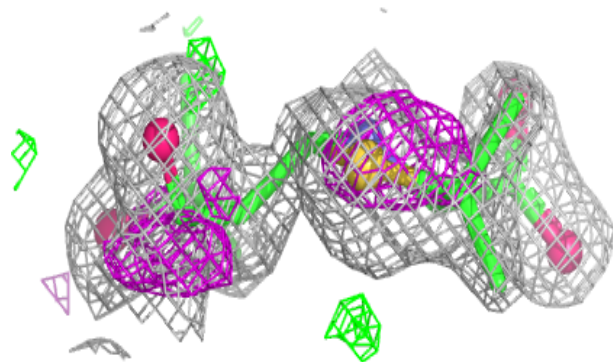
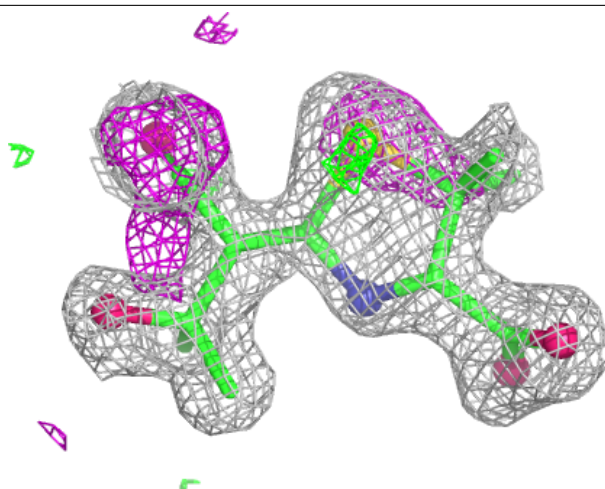
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

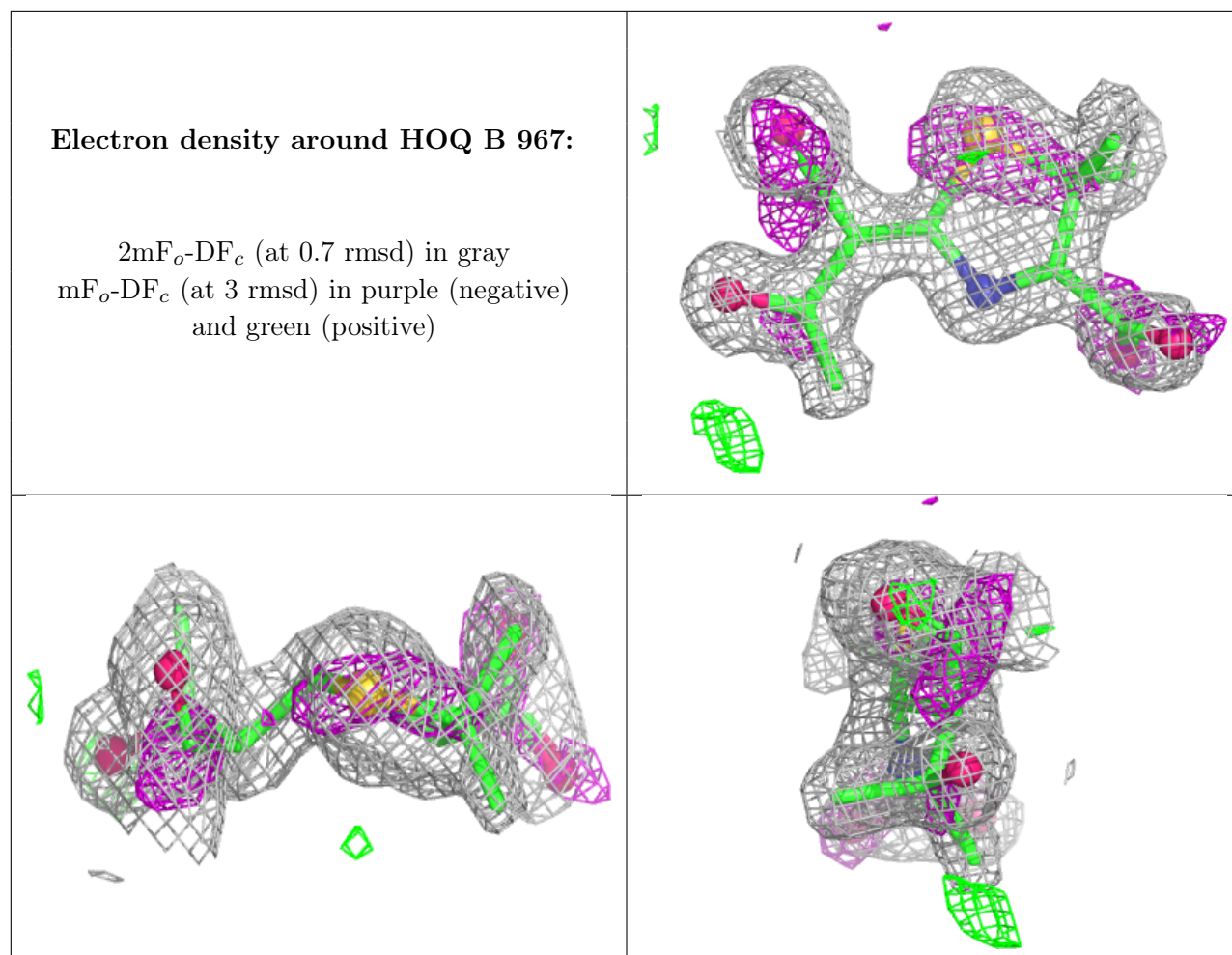
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	C	904	5/5	0.54	0.18	68,68,68,68	0
3	SO4	C	907	5/5	0.65	0.17	58,58,58,58	0
5	EDO	D	1005	4/4	0.68	0.22	25,26,26,29	0
3	SO4	A	906	5/5	0.74	0.17	37,38,38,38	0
5	EDO	B	1003	4/4	0.77	0.19	24,24,25,28	0
3	SO4	D	916	5/5	0.78	0.14	61,61,61,61	0
3	SO4	D	910	5/5	0.80	0.14	62,62,63,63	0
5	EDO	B	1004	4/4	0.81	0.18	27,27,27,28	0
3	SO4	B	915	5/5	0.82	0.10	64,65,65,65	0
3	SO4	A	914	5/5	0.82	0.14	56,56,56,56	0
3	SO4	C	901[A]	5/5	0.83	0.18	38,38,38,38	5
3	SO4	C	901[B]	5/5	0.83	0.18	63,63,63,63	5
3	SO4	A	905	5/5	0.83	0.14	44,45,45,45	0
3	SO4	B	912	5/5	0.85	0.13	44,44,45,45	0
3	SO4	B	908	5/5	0.86	0.13	37,38,38,39	0
3	SO4	D	917	5/5	0.87	0.23	59,59,60,60	0
4	HOQ	A	967	17/18	0.88	0.13	20,29,33,34	0
4	HOQ	B	967	17/18	0.89	0.12	16,26,32,32	0
3	SO4	B	913	5/5	0.89	0.12	53,53,54,54	0
3	SO4	B	909	5/5	0.90	0.18	33,34,34,34	0
3	SO4	A	903	5/5	0.90	0.10	37,37,37,37	0
5	EDO	A	1002	4/4	0.92	0.09	28,28,28,28	0
3	SO4	C	902	5/5	0.93	0.15	28,28,28,29	5
5	EDO	A	1001	4/4	0.93	0.11	20,21,23,23	0
5	EDO	D	1006	4/4	0.93	0.10	26,26,27,27	0
3	SO4	D	911	5/5	0.97	0.11	24,24,25,25	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 HOQ A 967:

$2mF_o-DF_c$ (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.