



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:06 AM UTC

PDB ID : 5SIN / pdb\_00005sin  
Title : CRYSTAL STRUCTURE OF HUMAN PHOSPHODIESTERASE 10 IN COMPLEX WITH c14c(c(nn1c2ncccc2)C)cc(NC(=O)c3ccccc3)s4, micromolar IC<sub>50</sub>=0.104457  
Authors : Joseph, C.; Benz, J.; Flohr, A.; Peters, J.; Rudolph, M.G.  
Deposited on : 2022-02-01  
Resolution : 2.40 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](mailto: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>.

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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

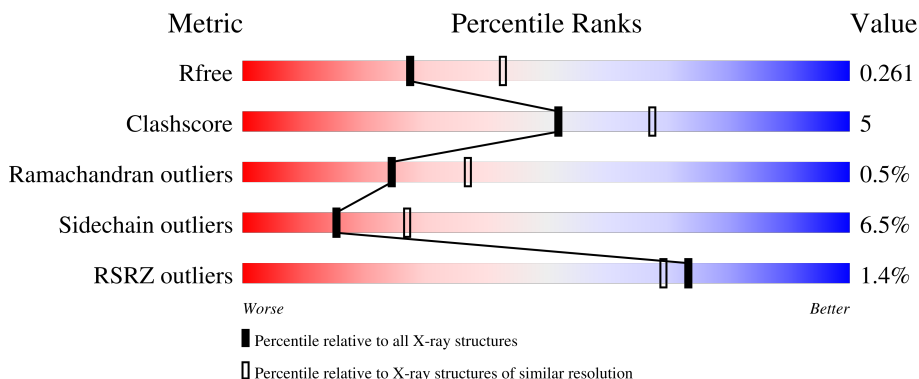
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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  $\geq 3$ , 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	343	
1	B	343	
1	C	343	
1	D	343	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10545 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 cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	313	Total 2547	C 1628	N 432	O 463	S 24	0	1	0
1	B	315	Total 2551	C 1630	N 434	O 463	S 24	0	0	0
1	C	313	Total 2549	C 1629	N 435	O 461	S 24	0	1	0
1	D	314	Total 2542	C 1625	N 434	O 459	S 24	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	447	GLY	-	expression tag	UNP Q9Y233
A	448	SER	-	expression tag	UNP Q9Y233
B	447	GLY	-	expression tag	UNP Q9Y233
B	448	SER	-	expression tag	UNP Q9Y233
C	447	GLY	-	expression tag	UNP Q9Y233
C	448	SER	-	expression tag	UNP Q9Y233
D	447	GLY	-	expression tag	UNP Q9Y233
D	448	SER	-	expression tag	UNP Q9Y233

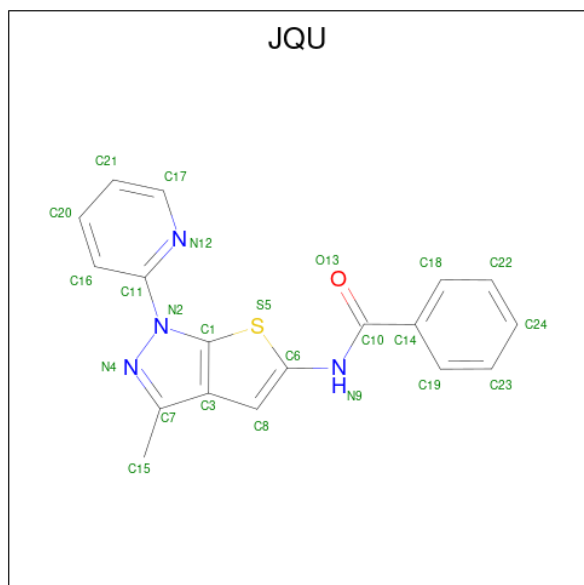
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0
2	C	1	Total 1	Zn 1	0	0
2	D	1	Total 1	Zn 1	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0

- Molecule 4 is N-[3-methyl-1-(pyridin-2-yl)-1H-thieno[2,3-c]pyrazol-5-yl]benzamide (CCD ID: JQU) (formula: C<sub>18</sub>H<sub>14</sub>N<sub>4</sub>OS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O S 24 18 4 1 1	0	0
4	B	1	Total C N O S 24 18 4 1 1	0	0
4	C	1	Total C N O S 24 18 4 1 1	0	0
4	D	1	Total C N O S 24 18 4 1 1	0	0

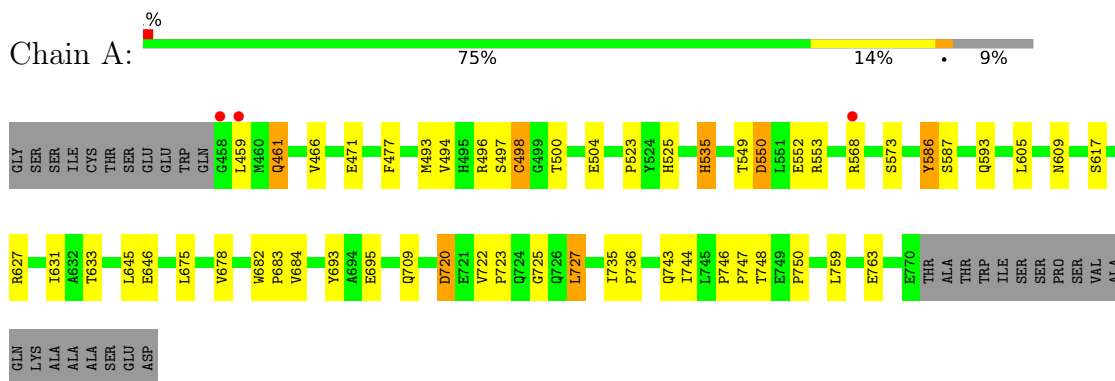
- Molecule 5 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	A	70	Total 70	O 70	0	0
5	B	82	Total 82	O 82	0	0
5	C	65	Total 65	O 65	0	0
5	D	35	Total 35	O 35	0	0

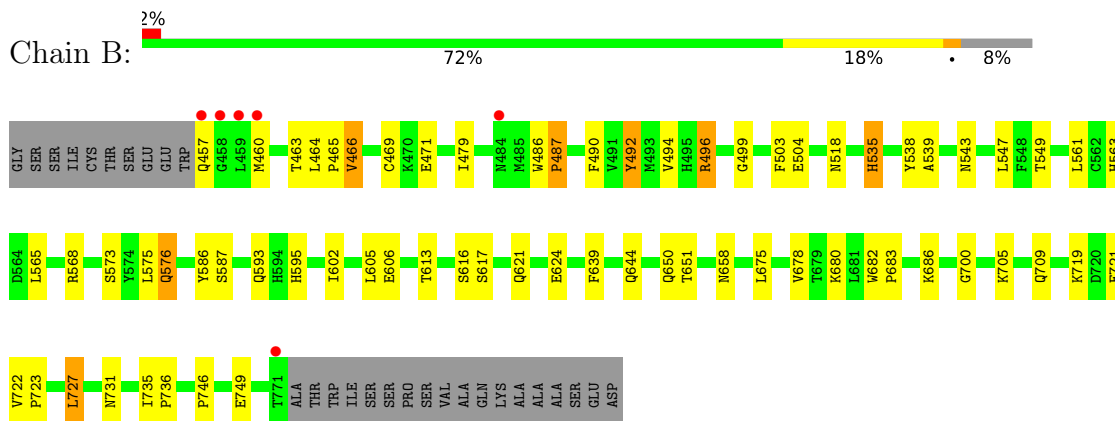
### 3 Residue-property plots [i](#)

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

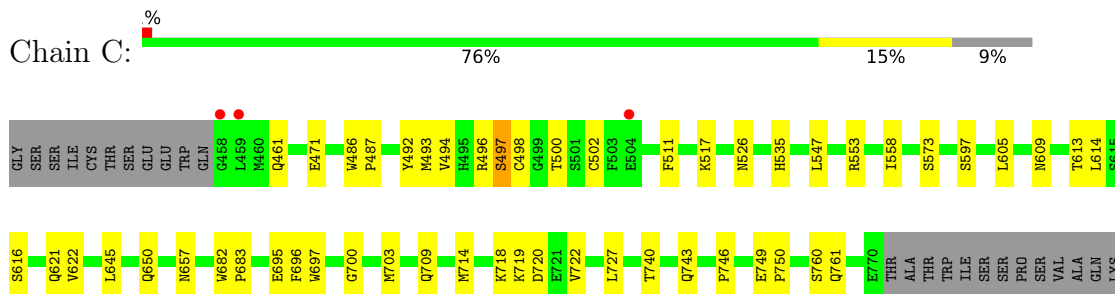
- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



ALA  
ALA  
ALA  
SER  
GLU  
ASP

- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A

Chain D:  2% 72% 19% 8%

GLY SER SER ILE CYS THR SER GLU GLU W456 Q457 G458 L459 M460 V466 E471 I479 N484 M485 W486 Y492 M493 R496 S497 C498 G499 T500 E504 C509 K517 V533 A534 H535 T546 L547 R553 C560 S573 Y574 L575 Q576 T588 E592

H595 F596 S597 S601 L605 S617 E618 Y619 V622 L623 E624 I625 R626 R627 I631 E646 E649 Y649 O650 T651 C673 D674 L675 V678 W682 K686 W697 G700 I708 Q709 M714 D715 R716 K719 D720 E721 V722 P723 L727 Y730 A734

L745 F746 P747 I748 E749 P750 L759 W762 I766 R767 G768 E769 GLU THR ALA THR TRP ILE SER SER PRO VAL ALA GLN LYS ALA ALA ALA SER GLU ASP

## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	135.88Å 135.88Å 234.97Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.64 – 2.40 43.64 – 2.40	Depositor EDS
% Data completeness (in resolution range)	90.1 (43.64-2.40) 90.1 (43.64-2.40)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.20 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.193 , 0.262 0.197 , 0.261	Depositor DCC
$R_{free}$ test set	2945 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.7	Xtrriage
Anisotropy	0.110	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 29.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.028 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10545	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: JQU, ZN, CME, MG

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	1.07	0/2601	1.46	7/3519 (0.2%)
1	B	1.13	2/2602 (0.1%)	1.52	13/3521 (0.4%)
1	C	1.08	1/2603 (0.0%)	1.51	8/3521 (0.2%)
1	D	1.11	0/2593	1.52	6/3509 (0.2%)
All	All	1.10	3/10399 (0.0%)	1.50	34/14070 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	645	LEU	C-O	-7.02	1.16	1.24
1	B	563	HIS	CE1-NE2	6.06	1.38	1.32
1	B	487	PRO	C-O	-5.58	1.17	1.24

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	746	PRO	N-CA-C	8.19	120.69	110.70
1	C	746	PRO	N-CA-C	6.99	119.22	110.70
1	C	502	CYS	N-CA-CB	6.72	120.70	110.28
1	B	535	HIS	CA-CB-CG	-6.35	107.45	113.80
1	B	576	GLN	CB-CA-C	6.09	120.45	110.88
1	A	609	ASN	N-CA-CB	6.05	119.05	110.46
1	A	466	VAL	CA-C-O	-5.90	115.27	121.41
1	D	588	THR	CA-CB-OG1	-5.86	100.81	109.60
1	D	597	SER	CA-C-N	5.71	127.94	120.28
1	D	597	SER	C-N-CA	5.71	127.94	120.28
1	C	535	HIS	CA-CB-CG	-5.49	108.31	113.80
1	C	718	LYS	CA-C-N	5.49	127.57	120.44
1	C	718	LYS	C-N-CA	5.49	127.57	120.44
1	B	499	GLY	CA-C-O	-5.47	118.67	122.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	492	TYR	N-CA-C	-5.44	105.35	111.28
1	B	466	VAL	CA-C-O	-5.42	115.77	121.41
1	A	759	LEU	N-CA-C	-5.32	105.48	111.28
1	C	502	CYS	CB-CA-C	-5.31	100.47	110.67
1	B	680	LYS	CA-C-O	-5.30	115.69	121.89
1	A	633	THR	N-CA-C	-5.28	106.62	113.16
1	D	535	HIS	CA-CB-CG	-5.27	108.53	113.80
1	B	639	PHE	CA-C-N	5.25	125.77	120.00
1	B	639	PHE	C-N-CA	5.25	125.77	120.00
1	C	609	ASN	CA-CB-CG	-5.23	107.37	112.60
1	D	746	PRO	N-CA-C	5.20	117.04	110.70
1	D	574	TYR	CA-C-O	-5.17	115.39	120.82
1	A	535	HIS	CA-CB-CG	-5.13	108.67	113.80
1	B	721	GLU	CA-C-N	5.08	124.39	120.33
1	B	721	GLU	C-N-CA	5.08	124.39	120.33
1	B	700	GLY	CA-C-N	5.01	127.31	120.54
1	B	700	GLY	C-N-CA	5.01	127.31	120.54
1	A	609	ASN	CA-CB-CG	-5.01	107.59	112.60
1	C	657	ASN	N-CA-C	-5.01	107.30	113.41
1	A	693	TYR	N-CA-C	-5.01	106.01	111.82

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	2547	0	2519	28	0
1	B	2551	0	2515	26	0
1	C	2549	0	2524	19	0
1	D	2542	0	2513	30	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	24	0	0	0	0
4	B	24	0	0	1	0
4	C	24	0	0	0	0
4	D	24	0	0	1	0
5	A	70	0	0	1	1
5	B	82	0	0	2	0
5	C	65	0	0	1	1
5	D	35	0	0	1	0
All	All	10545	0	10071	104	1

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 (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:C:493:MET:O	1:C:497:SER:HB2	1.76	0.84
1:B:651:THR:OG1	5:B:901:HOH:O	2.00	0.78
1:A:494:VAL:O	1:A:498:CYS:HB2	1.85	0.77
1:D:730:TYR:HA	1:D:734:ALA:HB3	1.71	0.70
1:A:493:MET:O	1:A:497:SER:HB2	1.93	0.68
1:C:749:GLU:HB3	1:C:750:PRO:HD3	1.80	0.64
1:D:727:LEU:HD12	1:D:766:ILE:CD1	2.27	0.64
1:A:725:GLY:O	5:A:901:HOH:O	2.16	0.63
1:A:461:GLN:NE2	1:A:461:GLN:HA	2.14	0.62
1:B:727:LEU:CD2	1:B:731:ASN:HD21	2.14	0.61
1:D:509:CME:HB2	5:D:933:HOH:O	2.01	0.59
1:A:586:TYR:HB3	1:A:593:GLN:NE2	2.17	0.59
1:D:492:TYR:CZ	1:D:496:ARG:HD2	2.39	0.57
1:A:722:VAL:HB	1:A:723:PRO:HD3	1.88	0.55
1:C:486:TRP:N	1:C:487:PRO:CD	2.70	0.55
1:C:682:TRP:CG	1:C:761:GLN:HE21	2.25	0.55
1:A:735:ILE:HB	1:A:736:PRO:HD3	1.89	0.54
4:B:803:JQU:O13	4:B:803:JQU:S5	2.65	0.53
4:D:803:JQU:O13	4:D:803:JQU:S5	2.67	0.53
1:B:457:GLN:HA	1:B:460:MET:HE3	1.90	0.53
1:B:651:THR:CB	5:B:901:HOH:O	2.55	0.53
1:B:675:LEU:O	1:B:678:VAL:HG22	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:727:LEU:CD2	1:B:731:ASN:ND2	2.71	0.52
1:C:682:TRP:N	1:C:683:PRO:CD	2.72	0.52
1:D:745:LEU:C	1:D:747:PRO:HD2	2.35	0.52
1:A:627:ARG:O	1:A:631:ILE:HG12	2.10	0.52
1:D:749:GLU:N	1:D:750:PRO:CD	2.73	0.51
1:A:586:TYR:HB3	1:A:593:GLN:HE21	1.75	0.51
1:A:461:GLN:CD	1:A:500:THR:HG21	2.37	0.50
1:A:645:LEU:HD22	1:A:744:ILE:CD1	2.42	0.50
1:A:461:GLN:NE2	1:A:500:THR:HG21	2.27	0.50
1:B:539:ALA:O	1:B:543:ASN:ND2	2.44	0.49
1:C:461:GLN:NE2	1:C:500:THR:HG21	2.28	0.49
1:D:619:TYR:O	1:D:622:VAL:HG12	2.13	0.49
1:B:565:LEU:O	1:B:595:HIS:ND1	2.43	0.48
1:D:727:LEU:HD12	1:D:766:ILE:HD12	1.94	0.48
1:C:740:THR:HA	1:C:743:GLN:NE2	2.28	0.48
1:D:700:GLY:HA3	1:D:714:MET:O	2.13	0.48
1:D:697:TRP:HB3	1:D:716:ARG:HB3	1.95	0.48
1:A:746:PRO:HG2	1:A:747:PRO:HD3	1.95	0.48
1:D:719:LYS:O	1:D:722:VAL:HG23	2.14	0.48
1:B:490:PHE:CZ	1:B:561:LEU:HD22	2.49	0.47
1:D:479:ILE:HD12	1:D:486:TRP:CD1	2.49	0.47
1:D:533:VAL:HG12	1:D:560:CYS:HB3	1.95	0.47
1:A:675:LEU:O	1:A:678:VAL:HG22	2.15	0.47
1:D:682:TRP:CZ2	1:D:686:LYS:HD3	2.50	0.47
1:A:523:PRO:HD2	1:A:695:GLU:HG2	1.97	0.46
1:D:745:LEU:C	1:D:747:PRO:CD	2.88	0.46
1:A:684:VAL:HG22	1:B:705:LYS:HG2	1.98	0.46
1:C:511:PHE:CE2	1:C:558:ILE:HD12	2.51	0.46
1:C:700:GLY:O	1:C:703:MET:HB2	2.15	0.46
1:A:682:TRP:HB3	1:A:683:PRO:HD3	1.98	0.45
1:B:735:ILE:HB	1:B:736:PRO:HD3	1.97	0.45
1:B:492:TYR:CZ	1:B:496:ARG:HD2	2.51	0.45
1:B:479:ILE:HD12	1:B:486:TRP:CD1	2.51	0.45
1:C:494:VAL:O	1:C:498:CYS:HB3	2.16	0.45
1:D:484:ASN:N	1:D:484:ASN:HD22	2.15	0.45
1:D:546:THR:OG1	1:D:547:LEU:HD12	2.17	0.45
1:D:675:LEU:O	1:D:678:VAL:HG22	2.17	0.44
1:A:497:SER:O	1:A:553:ARG:HD2	2.17	0.44
1:B:535:HIS:O	1:B:538:TYR:HB3	2.17	0.44
1:D:592:GLU:HA	1:D:595:HIS:HD2	1.81	0.44
1:D:722:VAL:HB	1:D:723:PRO:HD3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:645:LEU:HD22	1:A:744:ILE:HD13	1.99	0.44
1:C:492:TYR:CE2	1:C:496:ARG:HD2	2.52	0.44
1:B:486:TRP:N	1:B:487:PRO:CD	2.81	0.43
1:C:650:GLN:HA	1:C:650:GLN:OE1	2.18	0.43
1:A:727:LEU:HD21	1:A:763:GLU:HG3	1.99	0.43
1:A:461:GLN:HE22	1:A:500:THR:CG2	2.32	0.43
1:A:748:THR:C	1:A:750:PRO:HD2	2.44	0.43
1:B:465:PRO:O	1:B:466:VAL:C	2.62	0.43
1:D:533:VAL:HG22	1:D:673:CYS:O	2.19	0.43
1:D:627:ARG:O	1:D:631:ILE:HG12	2.19	0.42
1:C:553:ARG:NH1	5:C:903:HOH:O	2.51	0.42
1:A:550:ASP:OD2	1:A:550:ASP:N	2.52	0.42
1:A:646[A]:GLU:OE1	1:A:743:GLN:NE2	2.50	0.42
1:A:525:HIS:ND1	1:A:695:GLU:OE1	2.38	0.42
1:D:493:MET:O	1:D:497:SER:CB	2.68	0.42
1:D:624:GLU:OE2	1:D:627:ARG:NH1	2.47	0.42
1:A:461:GLN:OE1	1:A:500:THR:HG21	2.20	0.42
1:B:494:VAL:HG11	1:B:503:PHE:CD1	2.54	0.42
1:B:463:THR:HG22	1:B:464:LEU:N	2.35	0.42
1:D:493:MET:SD	1:D:535:HIS:HA	2.59	0.42
1:B:658:ASN:OD1	1:B:658:ASN:C	2.62	0.42
1:B:644:GLN:HE21	1:B:644:GLN:HA	1.84	0.42
1:D:749:GLU:N	1:D:750:PRO:HD2	2.35	0.42
1:B:722:VAL:HB	1:B:723:PRO:HD3	2.02	0.42
1:C:695:GLU:O	1:C:696:PHE:C	2.62	0.42
1:B:518:ASN:HD22	1:B:602:ILE:HD11	1.85	0.41
1:C:526:ASN:OD1	1:C:526:ASN:C	2.62	0.41
1:D:498:CYS:HB3	1:D:553:ARG:HB3	2.01	0.41
1:B:494:VAL:HG11	1:B:503:PHE:HD1	1.86	0.41
1:B:682:TRP:HB3	1:B:683:PRO:HD3	2.01	0.41
1:C:492:TYR:CZ	1:C:496:ARG:HD2	2.56	0.41
1:A:549:THR:HG23	1:A:552:GLU:OE2	2.21	0.41
1:B:644:GLN:HA	1:B:644:GLN:NE2	2.36	0.41
1:D:618:GLU:O	1:D:619:TYR:C	2.64	0.41
1:C:697:TRP:CZ2	1:C:719:LYS:HG3	2.56	0.41
1:B:586:TYR:HB3	1:B:593:GLN:NE2	2.36	0.41
1:C:700:GLY:HA3	1:C:714:MET:O	2.21	0.41
1:C:558:ILE:HA	1:C:558:ILE:HD13	1.81	0.40
1:D:622:VAL:HA	1:D:625:ILE:HD12	2.02	0.40
1:D:759:LEU:HD12	1:D:759:LEU:O	2.22	0.40
1:A:477:PHE:HB3	1:A:535:HIS:CE1	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:970:HOH:O	5:C:961:HOH:O[5_545]	2.13	0.07

## 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	311/343 (91%)	298 (96%)	12 (4%)	1 (0%)	36	50
1	B	312/343 (91%)	301 (96%)	10 (3%)	1 (0%)	36	50
1	C	311/343 (91%)	296 (95%)	14 (4%)	1 (0%)	36	50
1	D	311/343 (91%)	287 (92%)	21 (7%)	3 (1%)	12	20
All	All	1245/1372 (91%)	1182 (95%)	57 (5%)	6 (0%)	24	37

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	720	ASP
1	C	720	ASP
1	B	617	SER
1	D	768	GLY
1	D	466	VAL
1	D	746	PRO

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	282/305 (92%)	266 (94%)	16 (6%)	18	33
1	B	281/305 (92%)	258 (92%)	23 (8%)	10	18
1	C	282/305 (92%)	266 (94%)	16 (6%)	18	33
1	D	280/305 (92%)	262 (94%)	18 (6%)	16	28
All	All	1125/1220 (92%)	1052 (94%)	73 (6%)	15	27

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	459	LEU
1	A	461	GLN
1	A	471	GLU
1	A	496	ARG
1	A	498	CYS
1	A	504	GLU
1	A	550	ASP
1	A	568	ARG
1	A	573	SER
1	A	586	TYR
1	A	587	SER
1	A	605	LEU
1	A	617	SER
1	A	709	GLN
1	A	720	ASP
1	A	727	LEU
1	B	469	CYS
1	B	471	GLU
1	B	496	ARG
1	B	504	GLU
1	B	547	LEU
1	B	549	THR
1	B	568	ARG
1	B	573	SER
1	B	575	LEU
1	B	576	GLN
1	B	587	SER
1	B	605	LEU
1	B	606	GLU
1	B	613	THR
1	B	616	SER
1	B	621	GLN
1	B	624	GLU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	B	650	GLN
1	B	686	LYS
1	B	709	GLN
1	B	719	LYS
1	B	727	LEU
1	B	749	GLU
1	C	471	GLU
1	C	497	SER
1	C	517	LYS
1	C	547	LEU
1	C	573	SER
1	C	597	SER
1	C	605	LEU
1	C	613	THR
1	C	614	LEU
1	C	616	SER
1	C	621	GLN
1	C	622	VAL
1	C	709	GLN
1	C	722	VAL
1	C	727	LEU
1	C	760	SER
1	D	457	GLN
1	D	459	LEU
1	D	471	GLU
1	D	500	THR
1	D	504	GLU
1	D	517	LYS
1	D	573	SER
1	D	575	LEU
1	D	576	GLN
1	D	601	SER
1	D	605	LEU
1	D	617	SER
1	D	646	GLU
1	D	651	THR
1	D	708	ILE
1	D	709	GLN
1	D	720	ASP
1	D	727	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	476	HIS
1	A	484	ASN
1	A	593	GLN
1	A	598	GLN
1	A	604	GLN
1	A	644	GLN
1	A	650	GLN
1	A	709	GLN
1	A	724	GLN
1	A	726	GLN
1	A	761	GLN
1	B	476	HIS
1	B	518	ASN
1	B	593	GLN
1	B	604	GLN
1	B	644	GLN
1	B	709	GLN
1	B	726	GLN
1	B	731	ASN
1	C	476	HIS
1	C	542	GLN
1	C	576	GLN
1	C	598	GLN
1	C	604	GLN
1	C	726	GLN
1	C	743	GLN
1	D	457	GLN
1	D	476	HIS
1	D	484	ASN
1	D	576	GLN
1	D	593	GLN
1	D	604	GLN
1	D	609	ASN
1	D	621	GLN
1	D	743	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](#)

4 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	CME	A	509	1	8,9,10	0.58	0	6,9,11	1.20	0
1	CME	D	509	1	8,9,10	0.51	0	6,9,11	0.92	0
1	CME	B	509	1	8,9,10	0.52	0	6,9,11	0.85	0
1	CME	C	509	1	8,9,10	0.27	0	6,9,11	0.81	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
1	CME	A	509	1	-	0/5/8/10	-
1	CME	D	509	1	-	2/5/8/10	-
1	CME	B	509	1	-	1/5/8/10	-
1	CME	C	509	1	-	4/5/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	509	CME	SD-CE-CZ-OH
1	C	509	CME	CE-SD-SG-CB
1	C	509	CME	SD-CE-CZ-OH
1	D	509	CME	SD-CE-CZ-OH
1	C	509	CME	CZ-CE-SD-SG
1	D	509	CME	CZ-CE-SD-SG
1	C	509	CME	N-CA-CB-SG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	509	CME	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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
4	JQU	C	803	-	27,27,27	1.39	5 (18%)	32,38,38	2.61	14 (43%)
4	JQU	A	803	-	27,27,27	1.51	5 (18%)	32,38,38	3.27	14 (43%)
4	JQU	D	803	-	27,27,27	1.30	2 (7%)	32,38,38	2.69	13 (40%)
4	JQU	B	803	-	27,27,27	1.49	5 (18%)	32,38,38	2.43	12 (37%)

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
4	JQU	C	803	-	-	0/12/12/12	0/4/4/4
4	JQU	A	803	-	-	1/12/12/12	0/4/4/4
4	JQU	D	803	-	-	0/12/12/12	0/4/4/4
4	JQU	B	803	-	-	0/12/12/12	0/4/4/4

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	803	JQU	C6-S5	3.92	1.81	1.75
4	A	803	JQU	C6-S5	3.69	1.81	1.75
4	B	803	JQU	C6-S5	3.23	1.80	1.75
4	B	803	JQU	C8-C6	-2.94	1.32	1.37
4	A	803	JQU	C1-S5	2.76	1.79	1.74
4	C	803	JQU	C11-N2	-2.73	1.37	1.41
4	A	803	JQU	C8-C6	-2.62	1.33	1.37
4	C	803	JQU	C15-C7	2.59	1.54	1.49
4	C	803	JQU	C6-S5	2.59	1.79	1.75
4	B	803	JQU	C18-C14	2.58	1.43	1.39
4	B	803	JQU	C10-N9	-2.45	1.33	1.37
4	B	803	JQU	C1-S5	2.43	1.78	1.74
4	D	803	JQU	C8-C6	-2.42	1.33	1.37
4	A	803	JQU	C15-C7	2.37	1.54	1.49
4	C	803	JQU	C8-C3	2.28	1.48	1.42
4	A	803	JQU	C8-C3	2.09	1.47	1.42
4	C	803	JQU	C21-C20	2.08	1.42	1.38

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	803	JQU	C11-N2-N4	10.27	132.84	117.84
4	D	803	JQU	C11-N2-N4	7.82	129.25	117.84
4	C	803	JQU	C11-N2-N4	7.09	128.19	117.84
4	A	803	JQU	C11-N2-C1	-6.74	118.48	130.40
4	A	803	JQU	C15-C7-N4	5.95	130.94	119.39
4	B	803	JQU	O13-C10-N9	-5.78	113.83	122.27
4	B	803	JQU	C11-N2-N4	5.72	126.19	117.84
4	D	803	JQU	C10-N9-C6	-5.00	119.06	125.45
4	D	803	JQU	C15-C7-N4	4.67	128.45	119.39
4	C	803	JQU	C11-N2-C1	-4.58	122.29	130.40
4	A	803	JQU	C15-C7-C3	-4.56	120.81	127.84
4	B	803	JQU	C14-C10-N9	4.51	122.75	116.25
4	A	803	JQU	C19-C14-C18	4.48	124.26	118.57
4	A	803	JQU	C17-N12-C11	4.43	123.24	116.81
4	C	803	JQU	C17-N12-C11	4.38	123.17	116.81
4	A	803	JQU	C23-C19-C14	-4.31	116.12	120.36
4	A	803	JQU	N12-C11-N2	4.26	120.64	115.78
4	D	803	JQU	C7-N4-N2	4.20	111.09	105.75
4	B	803	JQU	C15-C7-N4	4.12	127.38	119.39
4	C	803	JQU	C21-C17-N12	-3.99	117.11	123.42
4	C	803	JQU	C15-C7-N4	3.94	127.04	119.39
4	D	803	JQU	C11-N2-C1	-3.81	123.65	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	803	JQU	C1-S5-C6	3.74	91.90	89.76
4	C	803	JQU	C14-C10-N9	3.74	121.64	116.25
4	B	803	JQU	C11-N2-C1	-3.37	124.43	130.40
4	D	803	JQU	C19-C14-C18	3.33	122.80	118.57
4	D	803	JQU	C3-C7-N4	-3.24	107.20	111.73
4	B	803	JQU	C10-N9-C6	-3.11	121.48	125.45
4	D	803	JQU	C1-N2-N4	-3.10	106.87	112.83
4	A	803	JQU	C21-C17-N12	-3.10	118.51	123.42
4	D	803	JQU	C23-C19-C14	-3.10	117.31	120.36
4	C	803	JQU	C20-C16-C11	-2.87	113.67	117.61
4	B	803	JQU	N12-C11-N2	2.77	118.95	115.78
4	D	803	JQU	C14-C10-N9	2.75	120.21	116.25
4	C	803	JQU	C23-C19-C14	-2.68	117.73	120.36
4	B	803	JQU	C15-C7-C3	-2.67	123.72	127.84
4	C	803	JQU	C3-C7-N4	-2.67	108.01	111.73
4	A	803	JQU	C10-N9-C6	-2.60	122.13	125.45
4	C	803	JQU	C10-N9-C6	-2.52	122.23	125.45
4	A	803	JQU	C3-C7-N4	-2.49	108.25	111.73
4	C	803	JQU	C19-C14-C18	2.48	121.72	118.57
4	A	803	JQU	C22-C18-C14	-2.39	118.02	120.36
4	C	803	JQU	O13-C10-C14	-2.28	116.39	120.90
4	D	803	JQU	C15-C7-C3	-2.26	124.35	127.84
4	D	803	JQU	N12-C11-N2	2.25	118.35	115.78
4	A	803	JQU	C1-N2-N4	-2.24	108.53	112.83
4	B	803	JQU	C23-C19-C14	-2.24	118.17	120.36
4	B	803	JQU	C24-C22-C18	-2.08	117.67	120.24
4	C	803	JQU	N12-C11-N2	2.07	118.15	115.78
4	B	803	JQU	C3-C7-N4	-2.03	108.90	111.73
4	B	803	JQU	C21-C20-C16	-2.01	117.75	120.24
4	A	803	JQU	C7-N4-N2	2.01	108.31	105.75
4	D	803	JQU	C3-C1-N2	2.01	108.53	105.54

There are no chirality outliers.

All (1) torsion outliers are listed below:

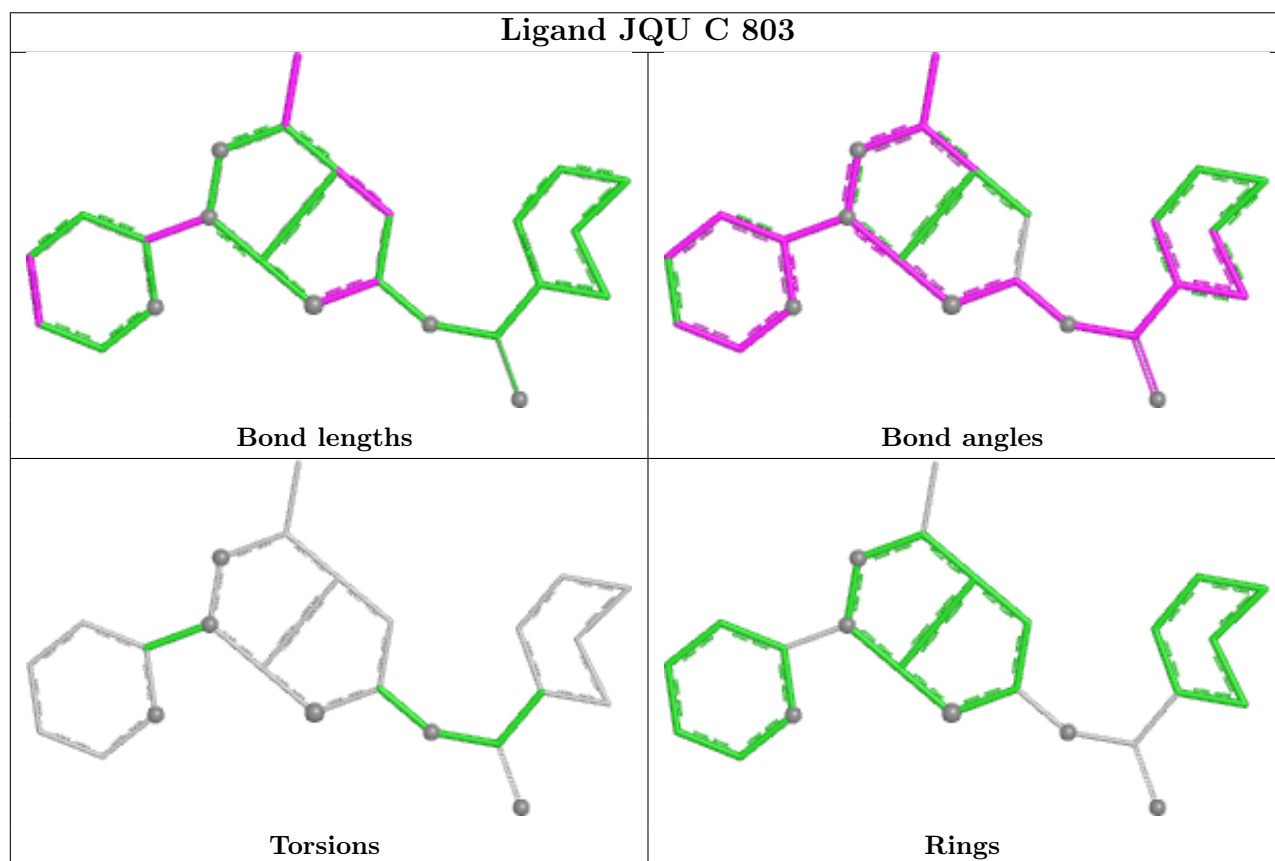
Mol	Chain	Res	Type	Atoms
4	A	803	JQU	S5-C6-N9-C10

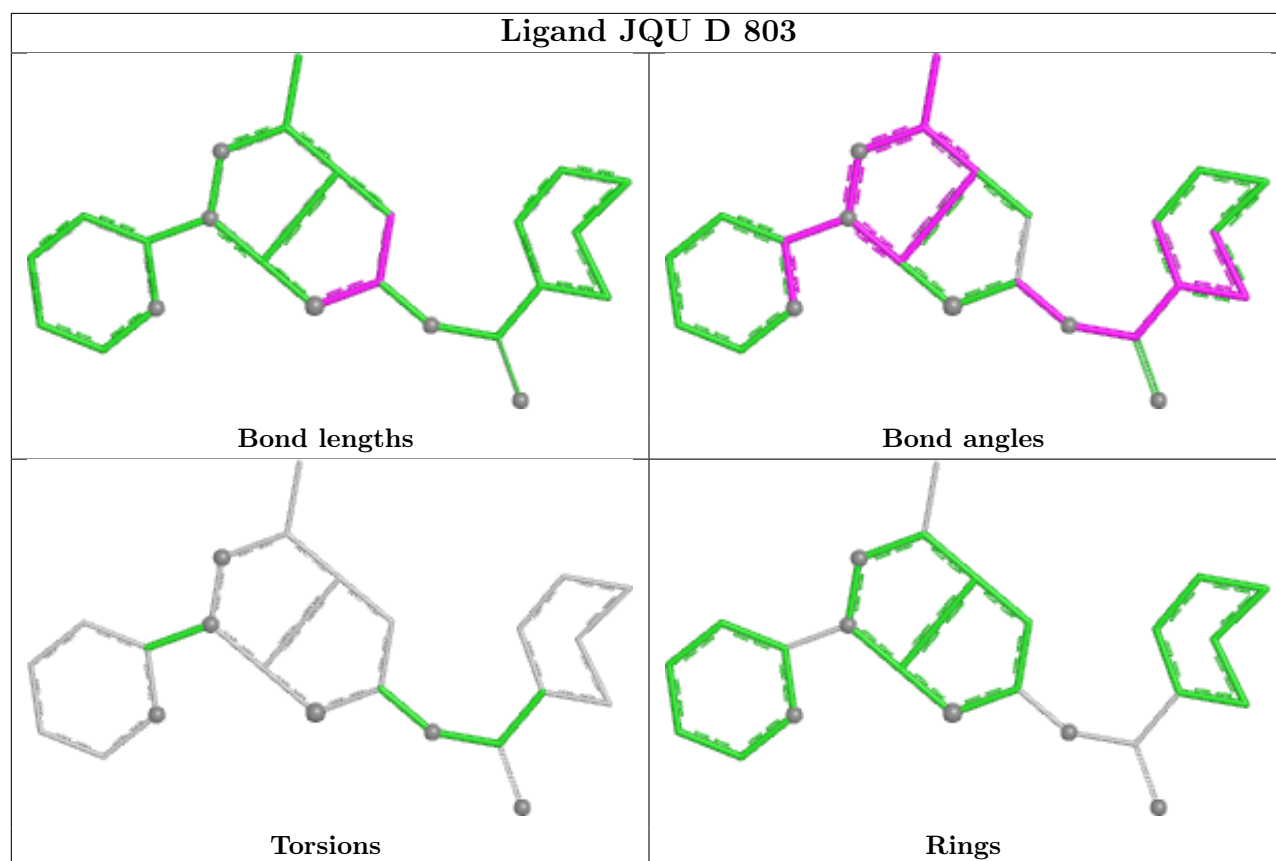
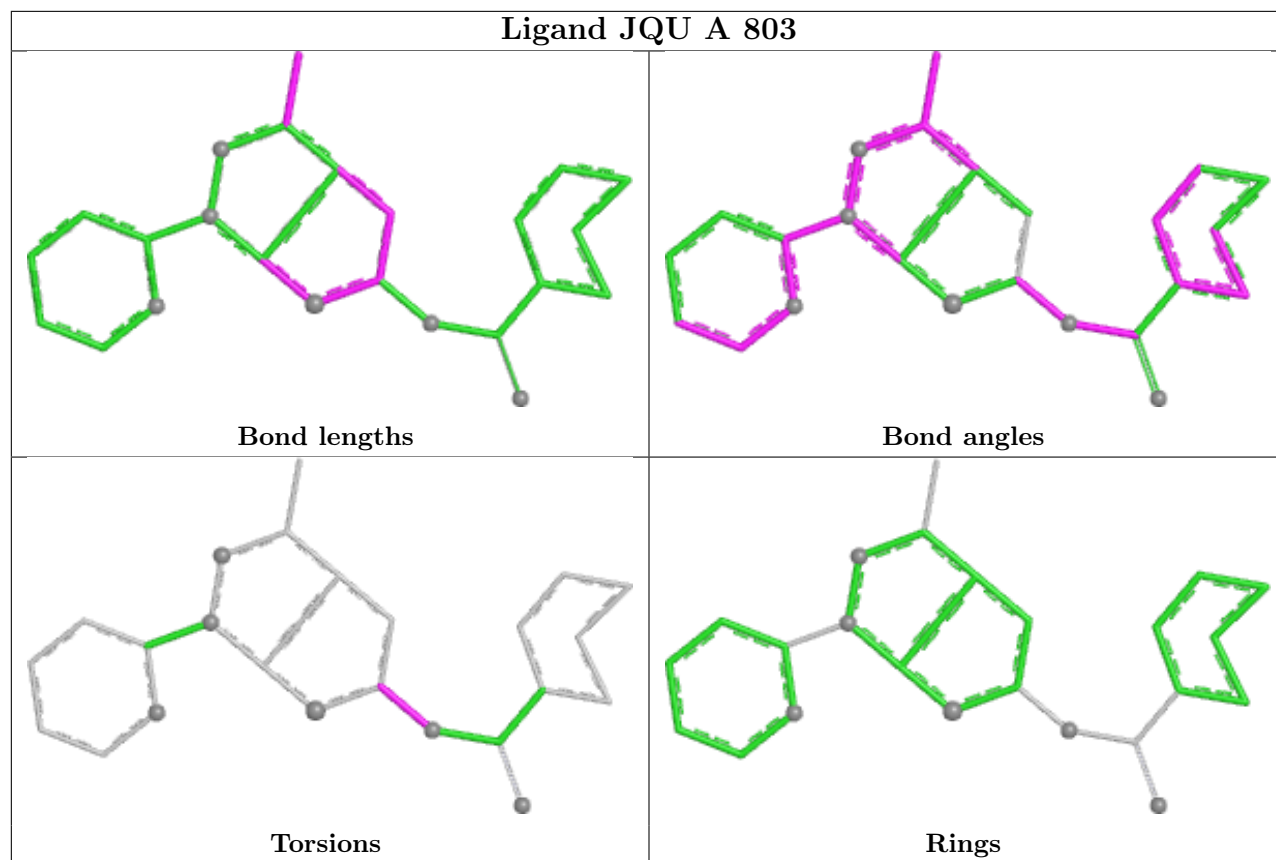
There are no ring outliers.

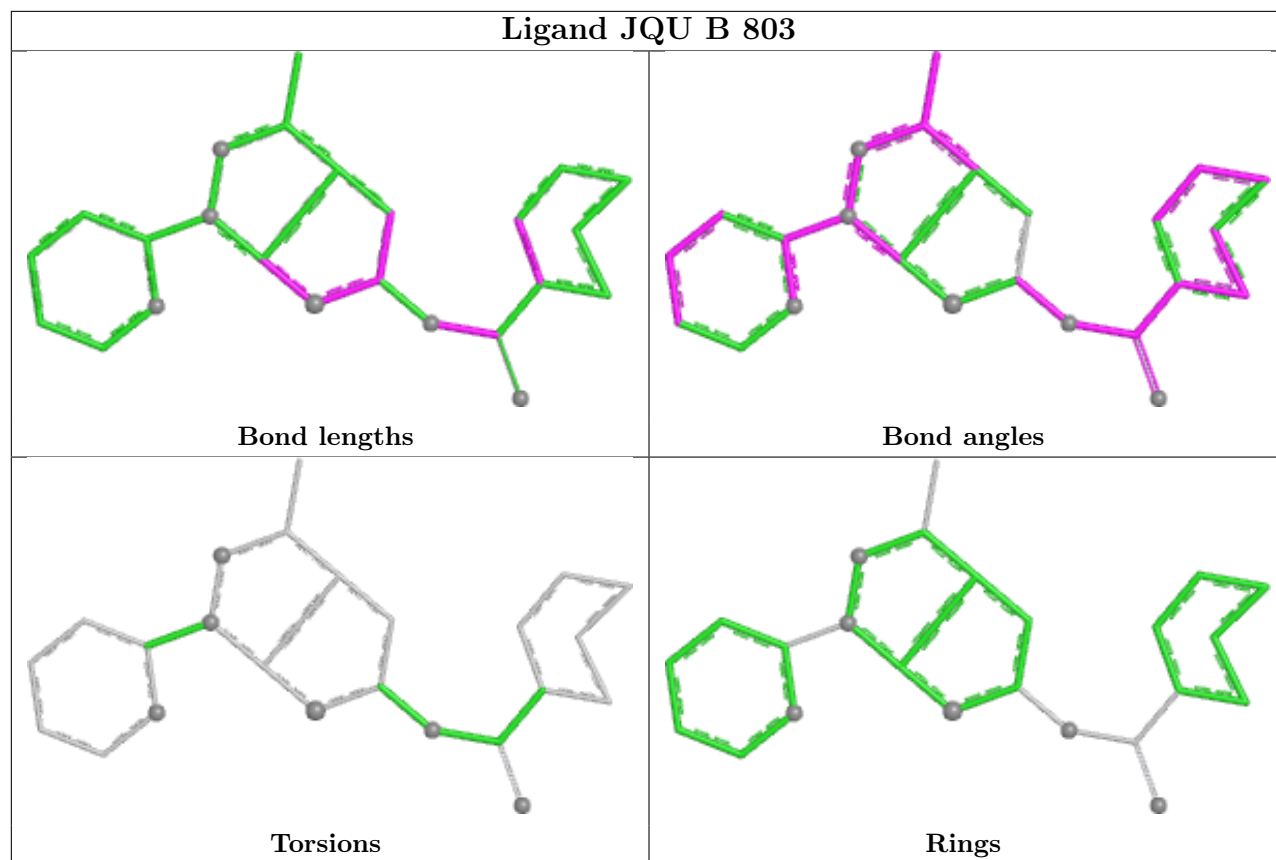
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	803	JQU	1	0
4	B	803	JQU	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	312/343 (90%)	-0.01	3 (0%) 79 76	28, 46, 75, 116	1 (0%)
1	B	314/343 (91%)	-0.03	6 (1%) 66 62	25, 44, 78, 109	0
1	C	312/343 (90%)	-0.10	3 (0%) 79 76	25, 44, 72, 100	1 (0%)
1	D	313/343 (91%)	0.26	6 (1%) 66 62	36, 58, 86, 123	0
All	All	1251/1372 (91%)	0.03	18 (1%) 73 69	25, 48, 79, 123	2 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	456	TRP	5.0
1	C	459	LEU	3.7
1	B	459	LEU	3.4
1	B	460	MET	3.1
1	B	771	THR	3.0
1	B	457	GLN	3.0
1	B	484	ASN	2.9
1	A	458	GLY	2.9
1	A	568	ARG	2.7
1	C	458	GLY	2.6
1	A	459	LEU	2.5
1	C	504	GLU	2.2
1	B	458	GLY	2.2
1	D	460	MET	2.2
1	D	649	TYR	2.2
1	D	459	LEU	2.1
1	D	762	TRP	2.1
1	D	457	GLN	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	CME	B	509	10/11	0.81	0.18	42,59,96,108	0
1	CME	D	509	10/11	0.82	0.17	54,71,107,113	0
1	CME	C	509	10/11	0.86	0.15	44,56,95,97	0
1	CME	A	509	10/11	0.86	0.16	43,61,106,109	0

## 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands [i](#)

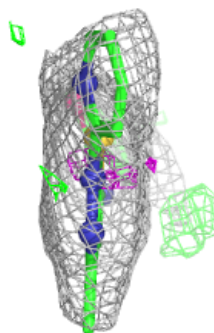
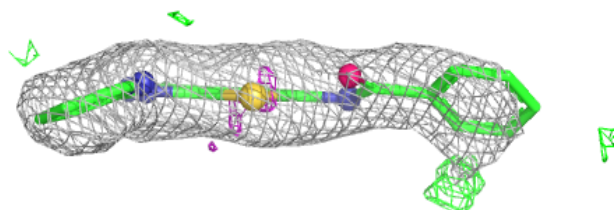
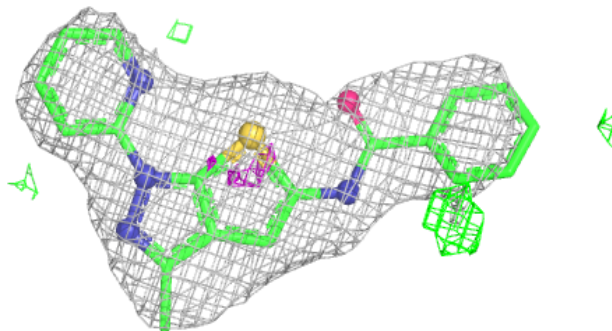
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	JQU	D	803	24/24	0.90	0.14	49,63,83,91	0
4	JQU	A	803	24/24	0.92	0.11	40,49,75,81	0
4	JQU	C	803	24/24	0.93	0.10	44,50,72,74	0
4	JQU	B	803	24/24	0.95	0.09	37,49,64,69	0
2	ZN	C	801	1/1	0.99	0.02	34,34,34,34	0
3	MG	B	802	1/1	1.00	0.01	23,23,23,23	0
3	MG	C	802	1/1	1.00	0.01	23,23,23,23	0
3	MG	D	802	1/1	1.00	0.03	41,41,41,41	0
2	ZN	B	801	1/1	1.00	0.01	33,33,33,33	0
2	ZN	A	801	1/1	1.00	0.01	39,39,39,39	0
2	ZN	D	801	1/1	1.00	0.01	47,47,47,47	0
3	MG	A	802	1/1	1.00	0.04	24,24,24,24	0

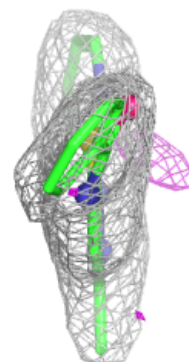
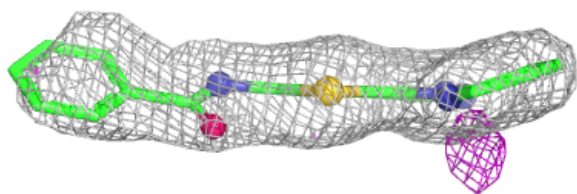
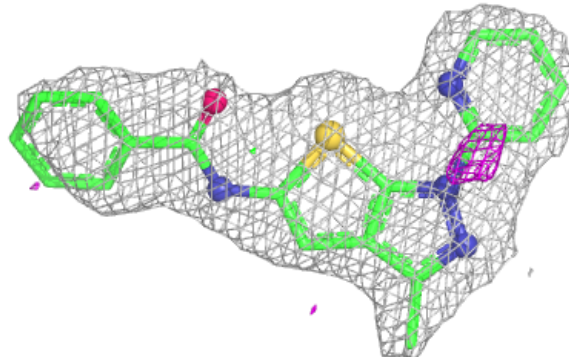
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around JQU D 803:**

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

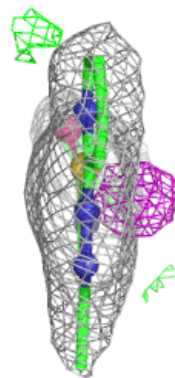
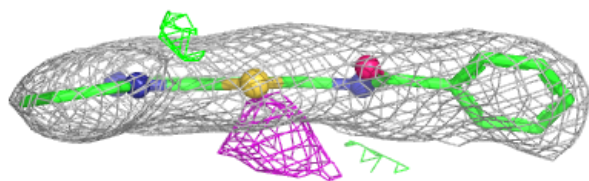
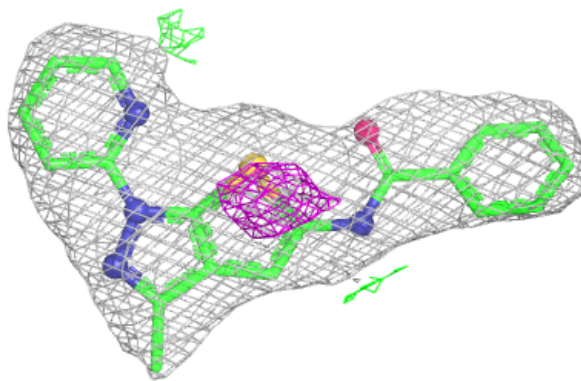
**Electron density around JQU A 803:**

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

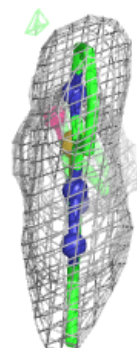
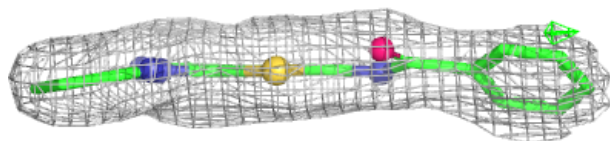
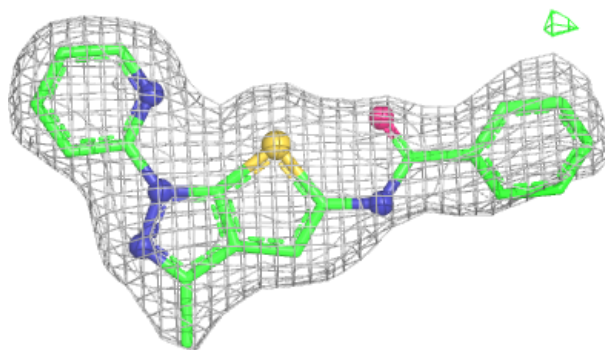


**Electron density around JQU C 803:**

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

**Electron density around JQU B 803:**

$2mF_o-DF_c$  (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.