

Full wwPDB X-ray Structure Validation Report (i)

May 15, 2020 – 11:12 am BST

PDB ID : 1SQA

Title : Substituted 2-Naphthamidine Inhibitors of Urokinase

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Deposited on : 2004-03-18

Resolution : 2.00 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 (i) symbol.

The following versions of software and data (see references (1)) 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)

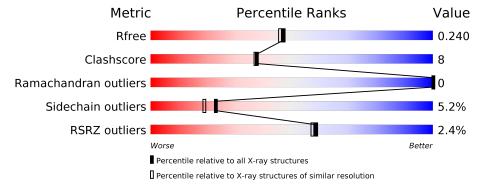
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 2.00 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar resolution} \\ (\#{\rm Entries, resolution range(\AA)}) \end{array}$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 <=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		
			2%		
1	A	245	72%	24%	• •



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2041 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.

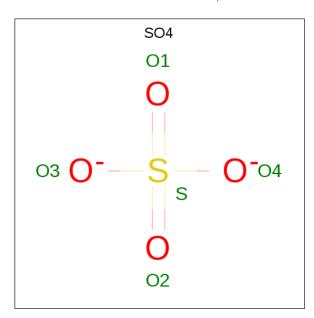
• Molecule 1 is a protein called Urokinase-type plasminogen activator.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	245	Total	С	N	О	S	0	0	0
1	A	Z40	1934	1219	338	361	16	0	0	U

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference	
Α	156	GLN	ASN	CONFLICT	UNP P00749	

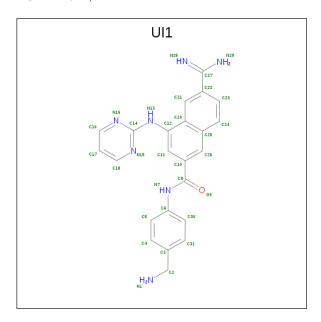
• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0



• Molecule 3 is 6-[(Z)-AMINO(IMINO)METHYL]-N-[4-(AMINOMETHYL)PHENYL] -4-(PYRIMIDIN-2-YLAMINO)-2-NAPHTHAMIDE (three-letter code: UI1) (formula: $C_{23}H_{21}N_7O$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	С	N	О	0	0
	11	_	31	23	7	1		

• Molecule 4 is water.

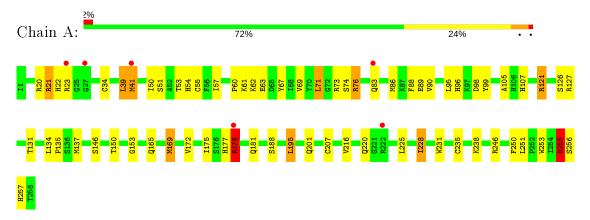
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	61	Total O 61 61	0	0



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 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: Urokinase-type plasminogen activator





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	55.16Å 53.00Å 82.30Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 - 2.00	Depositor
Resolution (A)	22.94 - 1.92	EDS
% Data completeness	67.4 (10.00-2.00)	Depositor
(in resolution range)	88.9 (22.94-1.92)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.61 (at 1.92Å)	Xtriage
Refinement program	X-PLOR 98.0	Depositor
R, R_{free}	0.210 , 0.266	Depositor
10,~10 free	0.237 , 0.240	DCC
R_{free} test set	1702 reflections (10.03%)	wwPDB-VP
Wilson B-factor (Å ²)	11.6	Xtriage
Anisotropy	0.148	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.41 , 47.9	EDS
L-test for twinning ²	$< L >=0.44, < L^2>=0.28$	Xtriage
Estimated twinning fraction	0.069 for k,h,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	2041	wwPDB-VP
Average B, all atoms (Å ²)	7.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.88% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $< L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: UI1, 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	Bo	nd lengths	Bo	ond angles
MIOI	RMSZ		# Z > 5	$\begin{array}{ c c c c c c c c c c c c c c c c c c c$	
1	A	1.36	$11/1983 \; (0.6\%)$	1.63	$36/2685 \ (1.3\%)$

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	A	51	SER	CA-CB	10.22	1.68	1.52
1	A	126	SER	CB-OG	-6.45	1.33	1.42
1	A	146	SER	CA-CB	-5.83	1.44	1.52
1	A	74	SER	CA-CB	5.79	1.61	1.52
1	A	169	MET	CA-CB	5.62	1.66	1.53
1	A	63	GLU	CD-OE1	5.59	1.31	1.25
1	A	146	SER	CB-OG	-5.59	1.34	1.42
1	A	50	ILE	CA-CB	5.35	1.67	1.54
1	A	228	ILE	CA-CB	5.27	1.67	1.54
1	A	172	VAL	CA-CB	5.18	1.65	1.54
1	A	41	MET	CB-CG	-5.03	1.35	1.51

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	A	20	ARG	NE-CZ-NH1	-13.93	113.33	120.30
1	A	20	ARG	NE-CZ-NH2	13.25	126.93	120.30



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Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
1	A	169	MET	CA-CB-CG	12.79	135.05	113.30
1	A	21	ARG	NE-CZ-NH2	-12.36	114.12	120.30
1	A	255	ARG	NE-CZ-NH2	-12.02	114.29	120.30
1	A	21	ARG	NE-CZ-NH1	10.40	125.50	120.30
1	A	73	ARG	NE-CZ-NH2	-10.37	115.11	120.30
1	A	41	MET	CA-CB-CG	-10.35	95.70	113.30
1	A	69	VAL	CG1-CB-CG2	-9.50	95.69	110.90
1	A	71	LEU	CB-CG-CD2	-8.15	97.14	111.00
1	A	255	ARG	NE-CZ-NH1	8.14	124.37	120.30
1	A	127	ARG	NE-CZ-NH2	-7.81	116.39	120.30
1	A	73	ARG	NE-CZ-NH1	7.60	124.10	120.30
1	A	175	ILE	O-C-N	-7.51	110.68	122.70
1	A	216	VAL	CG1-CB-CG2	-7.25	99.29	110.90
1	A	178	ARG	NE-CZ-NH2	-6.77	116.91	120.30
1	A	71	LEU	CB-CG-CD1	6.76	122.49	111.00
1	A	74	SER	O-C-N	-6.26	112.68	122.70
1	A	53	THR	O-C-N	-6.22	112.75	122.70
1	A	220	GLN	CA-CB-CG	-5.99	100.22	113.40
1	A	67	TYR	O-C-N	5.77	131.94	122.70
1	A	57	ILE	CG1-CB-CG2	-5.72	98.81	111.40
1	A	195	LEU	CB-CG-CD2	5.49	120.33	111.00
1	A	95	LEU	O-C-N	-5.47	113.94	122.70
1	A	175	ILE	CA-C-N	5.45	129.19	117.20
1	A	23	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	A	39	LEU	C-N-CA	5.23	134.78	121.70
1	A	238	LYS	C-N-CA	5.20	134.71	121.70
1	A	86	MET	CA-CB-CG	5.17	122.09	113.30
1	A	188	SER	O-C-N	-5.16	114.45	122.70
1	A	127	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	A	89	GLU	O-C-N	-5.10	114.54	122.70
1	A	201	GLN	O-C-N	-5.10	114.54	122.70
1	A	181	GLN	O-C-N	-5.09	114.55	122.70
1	A	175	ILE	CG1-CB-CG2	5.06	122.54	111.40
1	A	121	ARG	NE-CZ-NH1	5.00	122.80	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	153	GLY	Mainchain
1	A	250	PHE	Mainchain
1	A	255	ARG	Sidechain



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	1934	0	1880	27	0
2	A	15	0	0	0	0
3	A	31	0	20	0	0
4	A	61	0	0	3	0
All	All	2041	0	1900	27	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 8.

All (27) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \textbf{Interatomic} \\ \textbf{distance} \ (\text{\r{A}}) \end{array}$	Clash overlap (Å)
1:A:96:HIS:HD2	1:A:98:ASP:H	1.26	0.82
1:A:76:ARG:HG2	1:A:76:ARG:HH11	1.57	0.68
1:A:137:MET:HG2	1:A:251:LEU:HD12	1.80	0.64
1:A:255:ARG:HH11	1:A:255:ARG:HG2	1.61	0.64
1:A:76:ARG:NH1	1:A:76:ARG:HG2	2.18	0.59
1:A:41:MET:HE2	1:A:225:LEU:HD22	1.83	0.59
1:A:62:LYS:HD2	1:A:90:VAL:O	2.04	0.58
1:A:195:LEU:HB3	4:A:1029:HOH:O	2.05	0.55
1:A:253:TRP:O	1:A:257:HIS:HD2	1.92	0.53
1:A:195:LEU:HG	1:A:246:ARG:HB2	1.90	0.52
1:A:71:LEU:HD12	1:A:88:PHE:CE1	2.49	0.48
1:A:39:LEU:HG	1:A:131:THR:HG22	1.96	0.48
1:A:34:CYS:CB	1:A:55:CYS:SG	3.01	0.47
1:A:41:MET:HE3	1:A:134:LEU:HD21	1.97	0.47
1:A:105:ALA:HA	1:A:231:TRP:CE2	2.51	0.46
1:A:99:TYR:HA	1:A:107:HIS:O	2.16	0.45
1:A:96:HIS:CD2	1:A:98:ASP:H	2.18	0.45
1:A:41:MET:HE2	1:A:225:LEU:CD2	2.47	0.44
1:A:177:HIS:HD2	4:A:1010:HOH:O	2.01	0.43
1:A:76:ARG:HH11	1:A:165:GLN:HE21	1.68	0.42
1:A:135:PRO:O	1:A:251:LEU:HD11	2.20	0.41
1:A:178:ARG:HD3	1:A:178:ARG:HA	1.87	0.41
1:A:225:LEU:HD21	1:A:228:ILE:HD11	2.02	0.41



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Atom-1	Atom-2	$egin{aligned} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{aligned}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:121:ARG:HH11	1:A:121:ARG:HD2	1.77	0.41
1:A:207:CYS:HA	1:A:235:CYS:SG	2.61	0.40
1:A:76:ARG:HH11	1:A:165:GLN:NE2	2.19	0.40
1:A:22:HIS:HD2	4:A:1043:HOH:O	2.04	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	Percer	$_{ m itiles}$
1	A	243/245 (99%)	234 (96%)	9 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	213/214 (100%)	202 (95%)	11 (5%)	23 19		

All (11) residues with a non-rotameric sidechain are listed below:

\mathbf{Mol}	Chain	Res	Type
1	A	21	ARG
1	A	54	HIS



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Mol	Chain	Res	Type
1	A	60	PRO
1	A	61	LYS
1	A	76	ARG
1	A	83	GLN
1	A	150	THR
1	A	169	MET
1	A	178	ARG
1	A	255	ARG
1	A	256	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	22	HIS
1	A	96	HIS
1	A	107	HIS
1	A	156	GLN
1	A	165	GLN
1	A	177	HIS
1	A	257	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

4 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	Trino	Chain	Res	Link	Во	ond leng	ths	В	ond ang	gles
MIGI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	$\mid \# Z > 2$
2	SO4	A	301	-	4,4,4	1.05	0	6,6,6	1.02	0
2	SO4	A	302	-	4,4,4	0.75	0	6,6,6	2.23	2 (33%)
3	UI1	A	1001	-	34,34,34	1.45	7 (20%)	44,47,47	2.12	8 (18%)
2	SO4	A	303	-	4,4,4	0.66	0	6,6,6	0.41	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	UI1	A	1001	-	-	8/18/18/18	0/4/4/4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	${ m Observed}(m \AA)$	$\operatorname{Ideal}(\text{\AA})$
3	A	1001	UI1	C26-C10	3.59	1.43	1.37
3	A	1001	UI1	C14-N13	2.88	1.42	1.36
3	A	1001	UI1	C31-C30	2.66	1.43	1.38
3	A	1001	UI1	C24-C23	2.62	1.42	1.36
3	A	1001	UI1	C5-C6	2.47	1.43	1.39
3	A	1001	UI1	C21-C22	2.38	1.41	1.37
3	A	1001	UI1	C12-N13	2.19	1.44	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
3	A	1001	UI1	C18-N19-C14	6.11	120.87	115.45
3	A	1001	UI1	C21-C20-C12	-6.10	119.25	123.91
3	A	1001	UI1	C16-N15-C14	6.03	120.80	115.45
3	A	1001	UI1	N19-C14-N15	-5.31	120.83	126.43
2	A	302	SO4	O4-S-O3	-3.95	92.19	109.06
2	A	302	SO4	O4-S-O2	3.25	126.26	109.31
3	A	1001	UI1	C17-C16-N15	-2.29	119.69	123.43



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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
3	A	1001	UI1	C17-C18-N19	-2.23	119.78	123.43
3	A	1001	UI1	C6-N7-C8	-2.05	121.25	126.58
3	A	1001	UI1	C5-C6-N7	-2.03	113.57	120.40

There are no chirality outliers.

All (8) torsion outliers are listed below:

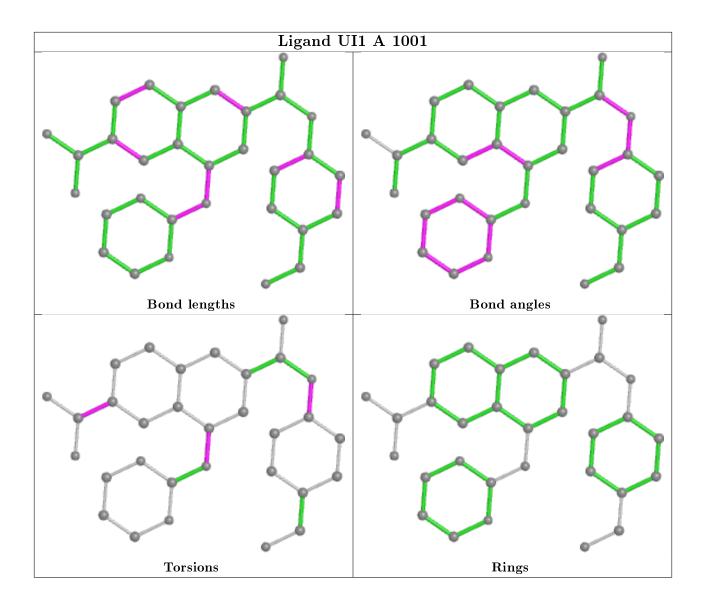
Mol	Chain	Res	Type	Atoms
3	A	1001	UI1	C5-C6-N7-C8
3	A	1001	UI1	C30-C6-N7-C8
3	A	1001	UI1	C21-C22-C27-N29
3	A	1001	UI1	C23-C22-C27-N29
3	A	1001	UI1	C20-C12-N13-C14
3	A	1001	UI1	C11-C12-N13-C14
3	A	1001	UI1	C21-C22-C27-N28
3	A	1001	UI1	C23-C22-C27-N28

There are no ring outliers.

No monomer is involved in short contacts.

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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ>2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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		$OWAB(A^2)$	Q < 0.9	
1	A	$245/245 \ (100\%)$	0.19	6 (2%)	59	57	2, 5, 18, 33	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	23	ARG	3.5
1	A	27	GLY	3.0
1	A	83	GLN	2.7
1	A	41	MET	2.5
1	A	222	ARG	2.2
1	A	178	ARG	2.2

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 carbohydrates 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^{th} 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.

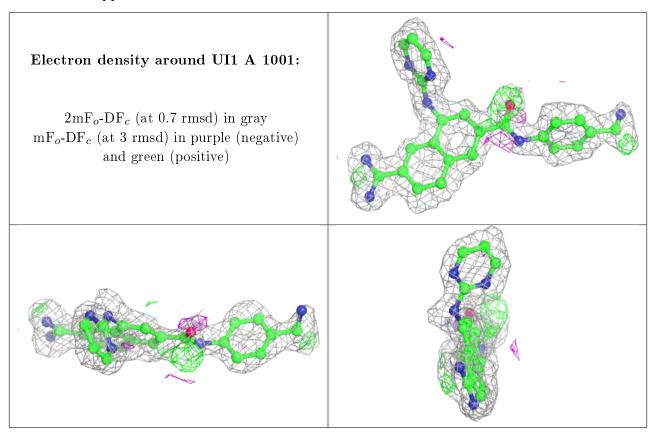
Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q < 0.9
3	UI1	A	1001	31/31	0.80	0.19	2,9,15,19	0



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Mol	Type	Chain	${f Res}$	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q < 0.9
2	SO4	A	302	5/5	0.83	0.26	21,21,26,29	0
2	SO4	A	303	5/5	0.95	0.12	35,36,37,38	0
2	SO4	A	301	5/5	0.97	0.09	8,11,13,13	0

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



6.5 Other polymers (i)

There are no such residues in this entry.

