

# Full wwPDB X-ray Structure Validation Report (i)

Dec 8, 2023 - 05:24 am GMT

PDB ID	:	2JEV
Title	:	Crystal structure of human spermine, spermidine acetyltransferase in complex
		with a bisubstrate analog (N1-acetylspermine-S-CoA).
Authors	:	Hegde, S.S.; Chandler, J.; Vetting, M.W.; Yu, M.; Blanchard, J.S.
Deposited on	:	2007-01-23
Resolution	:	2.30  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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

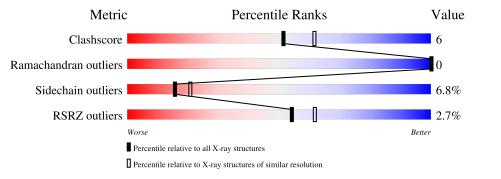
MolProbity	:	4.02b-467
5		1.8.4, CSD as541be (2020)
Xtriage (Phenix)		
EDS		2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

# 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 2.30 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	$\begin{array}{l} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 <=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	А	174	80%	13%					
1	В	174	3% 82%	14%	•••				



# 2 Entry composition (i)

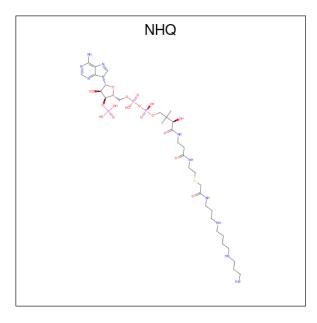
There are 3 unique types of molecules in this entry. The entry contains 2981 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	169	Total	С	Ν	0	S	0	0	0
	A	109	1392	905	223	254	10	0	0	0
1	р	160	Total	С	Ν	0	S	0	0	0
	I B	169	1392	905	223	254	10	0		0

• Molecule 1 is a protein called DIAMINE ACETYLTRANSFERASE 1.

• Molecule 2 is (3R)-27-AMINO-3-HYDROXY-2,2-DIMETHYL-4,8,14-TRIOXO-12-THI A-5,9,15,19,24-PENTAAZAHEPTACOS-1-YL [(2S,3R,4S,5S)-5-(6-AMINO-9H-PURIN -9-YL)-4-HYDROXY-3-(PHOSPHONOOXY)TETRAHYDROFURAN-2-YL]METHYL DIHYDROGEN DIPHOSPHATE (three-letter code: NHQ) (formula: C<sub>33</sub>H<sub>62</sub>N<sub>11</sub>O<sub>17</sub>P<sub>3</sub>S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf			
9	2 1	1	Total	С	Ν	Ο	Р	$\mathbf{S}$	0	0	
	1	65	33	11	17	3	1	0	0		
9	o D	В	1	Total	С	Ν	Ο	Р	$\mathbf{S}$	0	0
	1	65	33	11	17	3	1	0	U		

• Molecule 3 is water.

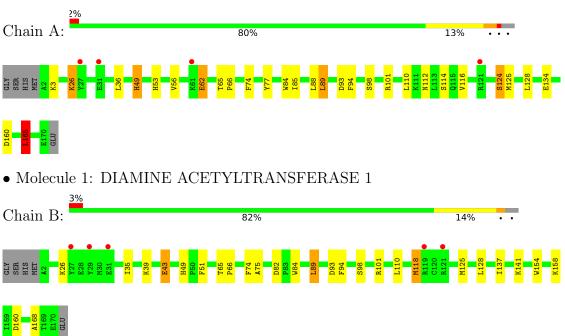


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	28	TotalO2828	0	0
3	В	39	Total O 39 39	0	0



# 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: DIAMINE ACETYLTRANSFERASE 1



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	$\begin{array}{rrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrr$	Depositor
Resolution (Å)	$\begin{array}{rrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrr$	Depositor EDS
% Data completeness (in resolution range)	98.3 (47.95-2.30) 98.4 (21.09-2.30)	Depositor EDS
R <sub>merge</sub>	0.04 (Not available)	Depositor Depositor
$\frac{\mathbf{R}_{sym}}{< I/\sigma(I) > 1}$	3.33 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
$R, R_{free}$	$ \begin{array}{rcl} 0.224 & , & 0.275 \\ 0.224 & , & (Not available) \end{array} $	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor $(Å^2)$	45.3	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.39, $31.9$	EDS
L-test for $twinning^2$	$< L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	0.487 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2981	wwPDB-VP
Average B, all atoms $(Å^2)$	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.40% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: NHQ

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	А	0.55	0/1431	0.62	1/1931~(0.1%)	
1	В	0.54	0/1431	0.64	0/1931	
All	All	0.55	0/2862	0.63	1/3862~(0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	165	LEU	CA-CB-CG	5.32	127.53	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	1392	0	1348	16	0
1	В	1392	0	1348	19	0
2	А	65	0	60	5	0
2	В	65	0	60	3	0
3	А	28	0	0	2	0
3	В	39	0	0	0	0
All	All	2981	0	2816	32	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 6.

All $(32)$ close contacts	within the same	e asymmetric unit	are listed below,	sorted by their clash
magnitude.				

Atom-1	Atom-2	Interatomic	Clash
		distance (Å)	overlap (Å)
2:A:1171:NHQ:HEM1	3:A:2021:HOH:O	1.79	0.82
1:B:39:LYS:O	1:B:43:GLU:HG3	1.82	0.80
1:A:165:LEU:HD13	1:B:118:MET:SD	2.32	0.69
1:A:49:HIS:CE1	1:B:49:HIS:HE1	2.11	0.68
1:A:65:THR:HB	1:A:66:PRO:HD2	1.76	0.66
1:B:74:PHE:O	1:B:94:PHE:HB2	1.97	0.65
1:B:137:ILE:O	1:B:141:LYS:HG2	2.01	0.61
1:A:128:LEU:O	2:A:1171:NHQ:H4M2	2.01	0.61
1:A:74:PHE:O	1:A:94:PHE:HB2	2.02	0.59
1:A:26:LYS:HE3	3:A:2006:HOH:O	2.02	0.58
2:A:1171:NHQ:HCM1	1:B:154:TRP:CZ2	2.40	0.56
1:B:101:ARG:HG2	2:B:1171:NHQ:O10	2.09	0.53
1:B:75:ALA:HA	1:B:94:PHE:HB3	1.91	0.51
1:A:114:SER:HB3	1:B:168:ALA:HB2	1.94	0.50
1:A:85:ILE:HG12	1:A:85:ILE:O	2.12	0.49
1:B:39:LYS:O	1:B:43:GLU:CG	2.59	0.49
1:B:98:SER:HA	1:B:101:ARG:HD3	1.95	0.48
1:A:84:TRP:CD2	2:B:1171:NHQ:H9M2	2.49	0.48
1:B:128:LEU:O	2:B:1171:NHQ:H4M2	2.13	0.48
2:A:1171:NHQ:H8M1	1:B:84:TRP:CD2	2.50	0.47
1:B:65:THR:HB	1:B:66:PRO:HD2	1.97	0.46
1:B:158:LYS:HE2	1:B:160:ASP:HB3	1.97	0.46
1:A:112:ASN:O	1:A:116:VAL:HG23	2.17	0.44
2:A:1171:NHQ:HEM2	1:B:82:ASP:OD1	2.18	0.44
1:A:53:HIS:HB2	1:A:77:TYR:CZ	2.54	0.43
1:A:98:SER:HA	1:A:101:ARG:NE	2.34	0.43
1:A:88:LEU:HD22	1:A:124:SER:HB2	1.99	0.42
1:A:62:GLU:H	1:A:62:GLU:HG3	1.49	0.42
1:A:49:HIS:CE1	1:B:49:HIS:CE1	3.01	0.41
1:B:89:LEU:O	1:B:125:MET:HA	2.21	0.41
1:B:49:HIS:HD2	1:B:51:PHE:CE2	2.39	0.41
1:A:89:LEU:O	1:A:125:MET:HA	2.22	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	Percent	iles
1	А	167/174~(96%)	161 (96%)	6 (4%)	0	100 1	100
1	В	167/174~(96%)	162 (97%)	5(3%)	0	100 1	100
All	All	334/348~(96%)	323~(97%)	11 (3%)	0	100 1	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	А	148/152~(97%)	135~(91%)	13 (9%)	10 12		
1	В	148/152~(97%)	141 (95%)	7~(5%)	26 37		
All	All	296/304~(97%)	276~(93%)	20 (7%)	16 21		

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	3	LYS
1	А	26	LYS
1	А	36	LEU
1	А	49	HIS
1	А	56	VAL
1	А	62	GLU
1	А	89	LEU
1	А	93	ASP

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Mol	Chain	Res	Type					
1	А	110	LEU					
1	А	124	SER					
1	А	134	GLU					
1	А	160	ASP					
1	А	165	LEU					
1	В	26	LYS					
1	В	35	ILE					
1	В	43	GLU					
1	В	89	LEU					
1	В	93	ASP					
1	В	110	LEU					
1	В	118	MET					

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Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	49	HIS
1	В	49	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 monosaccharides in this entry.

### 5.6 Ligand geometry (i)

2 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



1	it is the root-mean-square of an Z scores of the bond lengths (of angles).										
	Mol Type Chain Re		Dec	т.1	Bo	Bond lengths			Bond angles		
	MOI	туре	Unam	nes	Link	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
	2	NHQ	А	1171	-	59,67,67	1.61	5 (8%)	69,93,93	1.08	4 (5%)
ſ	2	NHQ	В	1171	-	59,67,67	1.60	4 (6%)	69,93,93	1.17	7 (10%)

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. BMSZ is the root-mean-square of all Z scores of the bond lengths (or angles)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NHQ	А	1171	-	-	19/62/82/82	0/3/3/3
2	NHQ	В	1171	-	-	18/62/82/82	0/3/3/3

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	1171	NHQ	CP-C1M	-9.65	1.40	1.51
2	В	1171	NHQ	CP-C1M	-9.60	1.40	1.51
2	А	1171	NHQ	O4X-C1X	4.49	1.47	1.41
2	В	1171	NHQ	O4X-C1X	3.86	1.46	1.41
2	В	1171	NHQ	P3X-O3X	2.89	1.64	1.59
2	В	1171	NHQ	P3X-O7A	2.32	1.63	1.54
2	А	1171	NHQ	P3X-O3X	2.06	1.63	1.59
2	А	1171	NHQ	P3X-08A	2.05	1.62	1.54
2	А	1171	NHQ	P3X-07A	2.01	1.62	1.54

All (9) bond length outliers are listed below:

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	В	1171	NHQ	N3A-C2A-N1A	-4.70	121.33	128.68
2	А	1171	NHQ	N3A-C2A-N1A	-4.40	121.81	128.68
2	В	1171	NHQ	C3-N4-C5	2.85	128.12	122.84
2	В	1171	NHQ	C2-C3-N4	-2.69	106.77	112.42
2	В	1171	NHQ	C5X-C4X-C3X	-2.37	106.54	114.40
2	А	1171	NHQ	C2-C3-N4	-2.32	107.55	112.42
2	В	1171	NHQ	O6A-C12-C11	2.17	114.03	110.55
2	А	1171	NHQ	O6A-C12-C11	2.10	113.92	110.55
2	В	1171	NHQ	P2A-O3A-P1A	-2.07	125.72	132.83

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$		
2	В	1171	NHQ	O1M-C1M-CP	-2.06	118.02	121.58		
2	А	1171	NHQ	P2A-O3A-P1A	-2.01	125.93	132.83		

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There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	А	1171	NHQ	CP-C1M-N50-C5M
2	А	1171	NHQ	C1M-CP-S1-C2
2	А	1171	NHQ	C5X-O5X-P1A-O2A
2	А	1171	NHQ	C5X-O5X-P1A-O1A
2	В	1171	NHQ	CP-C1M-N50-C5M
2	В	1171	NHQ	C1M-CP-S1-C2
2	В	1171	NHQ	C3X-C4X-C5X-O5X
2	А	1171	NHQ	N51-C3M-C4M-C5M
2	А	1171	NHQ	C6-C5-N4-C3
2	В	1171	NHQ	C6-C5-N4-C3
2	В	1171	NHQ	N51-C3M-C4M-C5M
2	А	1171	NHQ	N51-CBM-CCM-CDM
2	В	1171	NHQ	O4X-C4X-C5X-O5X
2	А	1171	NHQ	O1M-C1M-N50-C5M
2	А	1171	NHQ	O5-C5-N4-C3
2	В	1171	NHQ	O1M-C1M-N50-C5M
2	В	1171	NHQ	O5-C5-N4-C3
2	А	1171	NHQ	C3M-C4M-C5M-N50
2	В	1171	NHQ	C3M-C4M-C5M-N50
2	В	1171	NHQ	N51-CBM-CCM-CDM
2	А	1171	NHQ	O4X-C4X-C5X-O5X
2	А	1171	NHQ	C3X-C4X-C5X-O5X
2	В	1171	NHQ	C8M-C9M-N52-CEM
2	А	1171	NHQ	N53-C7M-C8M-C9M
2	В	1171	NHQ	N53-C7M-C8M-C9M
2	А	1171	NHQ	CBM-CCM-CDM-CEM
2	В	1171	NHQ	CBM-CCM-CDM-CEM
2	В	1171	NHQ	C4M-C3M-N51-CBM
2	А	1171	NHQ	C5X-O5X-P1A-O3A
2	В	1171	NHQ	C3X-O3X-P3X-O7A
2	В	1171	NHQ	CCM-CDM-CEM-N52
2	В	1171	NHQ	C2X-C3X-O3X-P3X
2	А	1171	NHQ	C4M-C3M-N51-CBM
2	В	1171	NHQ	C4X-C3X-O3X-P3X
2	А	1171	NHQ	O10-C10-C9-O9

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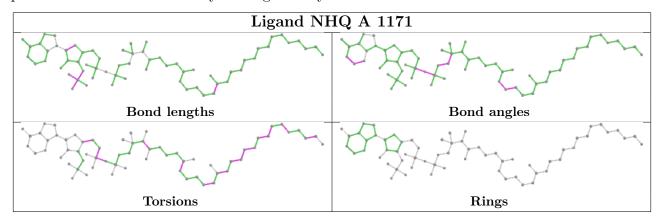
Mol	Chain	$\mathbf{Res}$	Type	Atoms
2	А	1171	NHQ	CDM-CEM-N52-C9M
2	А	1171	NHQ	P2A-O3A-P1A-O2A

There are no ring outliers.

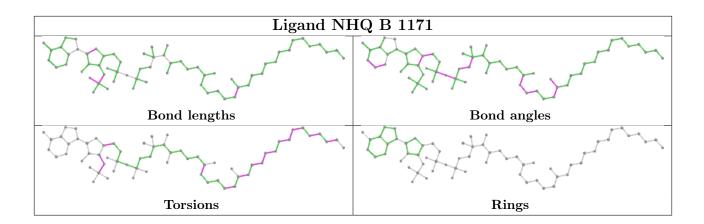
2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	1171	NHQ	5	0
2	В	1171	NHQ	3	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 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	$\langle RSRZ \rangle$	# RSRZ > 2	$OWAB(Å^2)$	Q < 0.9
1	А	169/174~(97%)	0.27	4 (2%) 59 66	28, 44, 62, 66	0
1	В	169/174~(97%)	0.33	5 (2%) 50 57	27, 43, 62, 66	0
All	All	338/348~(97%)	0.30	9 (2%) 54 62	27, 44, 62, 66	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	29	TYR	2.7
1	В	121	ARG	2.7
1	А	121	ARG	2.5
1	А	27	TYR	2.4
1	В	27	TYR	2.4
1	А	61	LYS	2.3
1	В	119	ARG	2.3
1	А	31	GLU	2.3
1	В	31	GLU	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 monosaccharides in this entry.

### 6.4 Ligands (i)

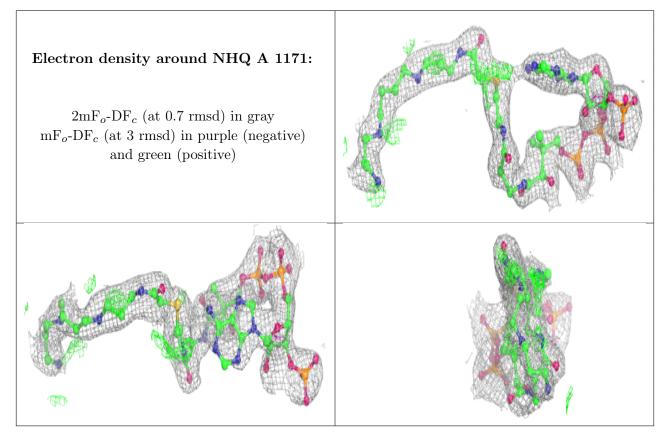
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



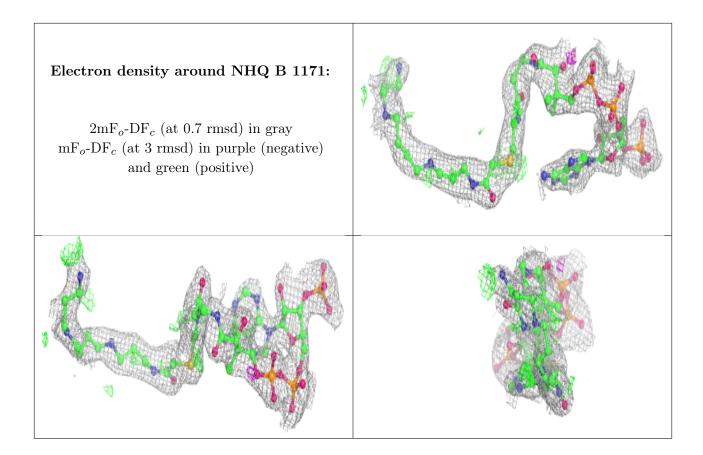
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B$ -factors( $Å^2$ )	Q<0.9
2	NHQ	А	1171	65/65	0.88	0.19	$55,\!68,\!78,\!79$	0
2	NHQ	В	1171	65/65	0.88	0.19	55,66,79,80	0

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.

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.

