Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:10 PM GMT

PDB ID	:	4BRE
Title	:	Legionella pneumophila NTPDase1 crystal form II (closed) in complex with transition state mimic adenosine 5'phosphovanadate
Authors	:	Zebisch, M.; Schaefer, P.; Lauble, P.; Straeter, N.
Deposited on	:	2013-06-04
Resolution	:	1.60 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

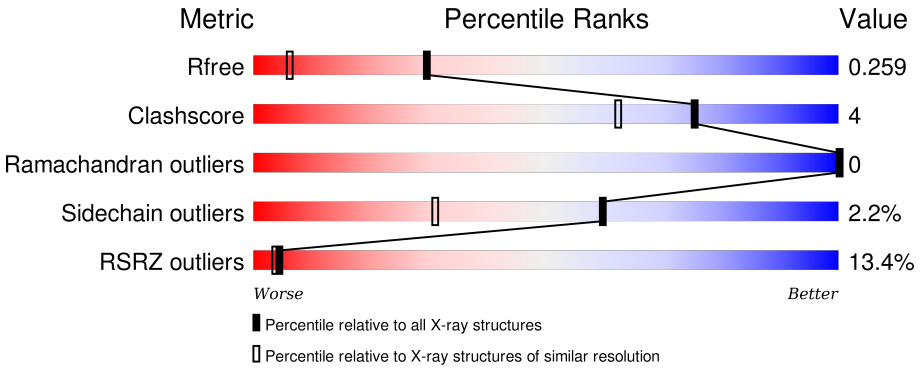
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 1.60 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}	91344	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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. 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	A	368	17% 86% 10%
1	B	368	9% 90% 6%

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
2	NA	A	1395	-	-	-	X
3	50T	B	1395[A]	-	-	-	X
3	50T	B	1395[B]	-	-	-	X
4	MG	A	1397	-	-	-	X
4	MG	B	1396	-	-	-	X
5	MES	A	1398	-	-	-	X
5	MES	B	1397	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 6544 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 ECTONUCLEOSIDE TRIPHOSPHATE DIPHOSPHOHYDROLASE I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	358	Total	C	N	O	S	0	23	0
			2984	1908	486	574	16			
1	B	358	Total	C	N	O	S	0	23	0
			2984	1904	489	575	16			

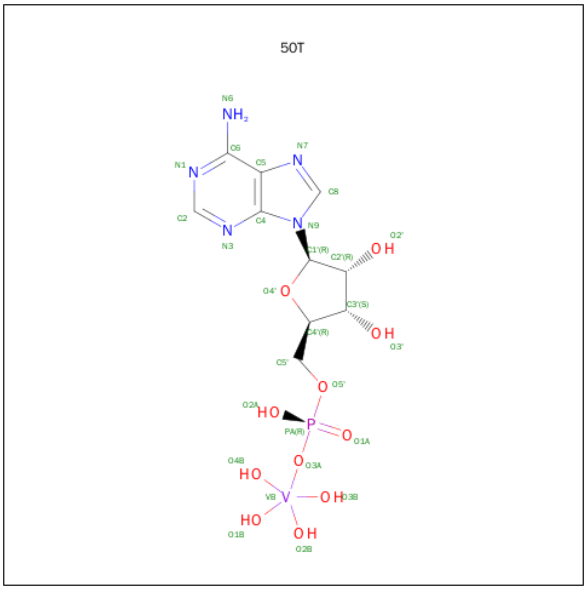
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	MET	-	EXPRESSION TAG	UNP Q5ZUA2
A	137	ASP	GLU	CONFLICT	UNP Q5ZUA2
A	149	VAL	ALA	CONFLICT	UNP Q5ZUA2
A	394	LEU	-	EXPRESSION TAG	UNP Q5ZUA2
A	395	GLU	-	EXPRESSION TAG	UNP Q5ZUA2
A	396	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
A	397	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
A	398	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
A	399	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
A	400	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
A	401	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
B	34	MET	-	EXPRESSION TAG	UNP Q5ZUA2
B	137	ASP	GLU	CONFLICT	UNP Q5ZUA2
B	149	VAL	ALA	CONFLICT	UNP Q5ZUA2
B	394	LEU	-	EXPRESSION TAG	UNP Q5ZUA2
B	395	GLU	-	EXPRESSION TAG	UNP Q5ZUA2
B	396	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
B	397	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
B	398	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
B	399	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
B	400	HIS	-	EXPRESSION TAG	UNP Q5ZUA2
B	401	HIS	-	EXPRESSION TAG	UNP Q5ZUA2

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0

- Molecule 3 is ADENOSINE-5'-PHOSPHOVANADATE (three-letter code: 50T) (formula: $C_{10}H_{17}N_5O_{11}PV$).

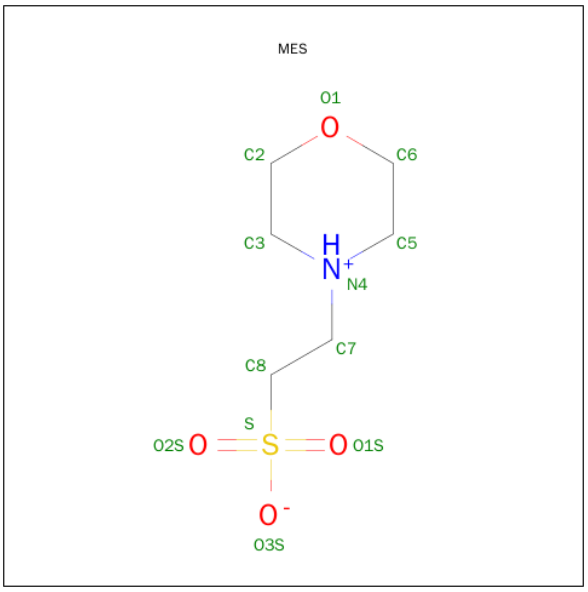


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O P V 28 10 5 11 1 1	0	0
3	B	1	Total C N O P V 47 20 10 15 1 1	0	1

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mg 1 1	0	0
4	A	1	Total Mg 1 1	0	0

- Molecule 5 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
5	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 6 is water.

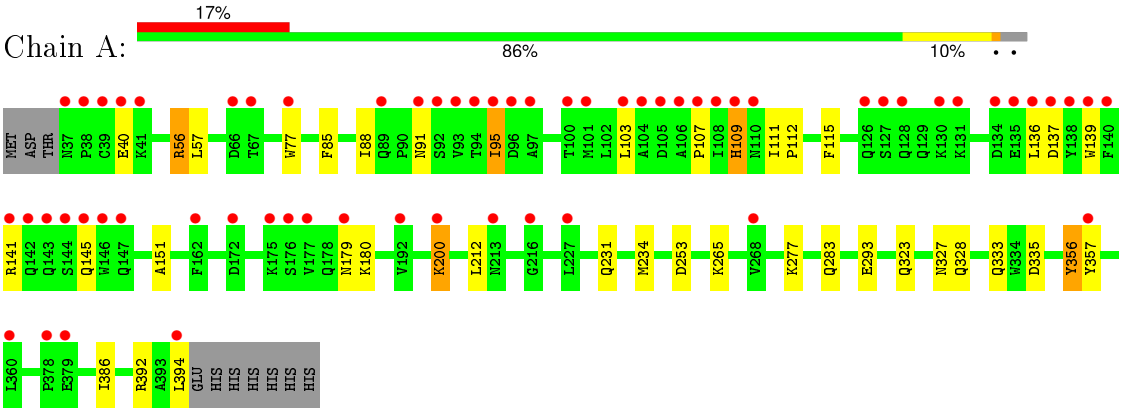
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	237	Total	O	0	3
			240	240		
6	B	232	Total	O	0	2
			234	234		

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 errors 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.

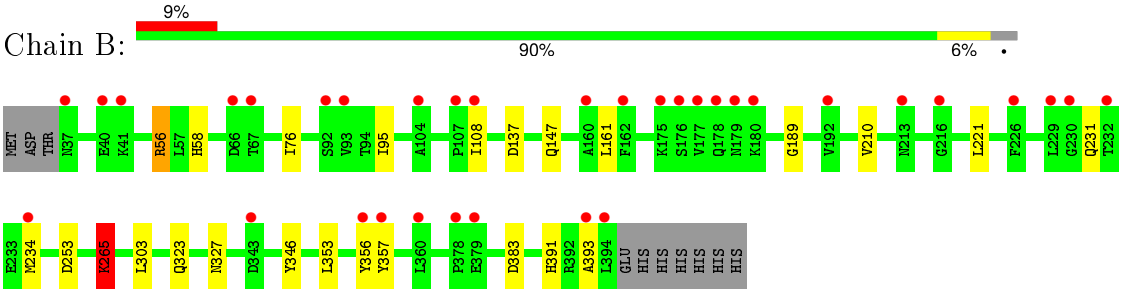
- Molecule 1: ECTONUCLEOSIDE TRIPHOSPHATE DIPHOSPHOHYDROLASE I

Chain A:



- Molecule 1: ECTONUCLEOSIDE TRIPHOSPHATE DIPHOSPHOHYDROLASE I

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	62.68Å 86.20Å 70.88Å 90.00° 105.94° 90.00°	Depositor
Resolution (Å)	29.33 – 1.60 29.33 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.33-1.60) 99.9 (29.33-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.62 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.172 , 0.208 0.226 , 0.259	Depositor DCC
R_{free} test set	1404 reflections (1.49%)	DCC
Wilson B-factor (Å^2)	18.4	Xtriage
Anisotropy	0.623	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.44 , 51.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 95350 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6544	wwPDB-VP
Average B, all atoms (Å^2)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: NA, 50T, MG, MES

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	0.90	0/3128	0.93	4/4262 (0.1%)
1	B	0.91	1/3125 (0.0%)	0.95	5/4256 (0.1%)
All	All	0.90	1/6253 (0.0%)	0.94	9/8518 (0.1%)

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 mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	189	GLY	N-CA	5.41	1.54	1.46

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	253	ASP	CB-CG-OD2	7.26	124.83	118.30
1	A	356	TYR	CB-CG-CD2	-6.48	117.11	121.00
1	A	277	LYS	CD-CE-NZ	-6.12	97.62	111.70
1	B	56	ARG	NE-CZ-NH1	-6.11	117.24	120.30
1	A	56	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	B	253	ASP	CB-CG-OD2	5.77	123.49	118.30
1	B	137	ASP	CB-CG-OD1	5.56	123.30	118.30
1	B	265	LYS	CD-CE-NZ	5.42	124.17	111.70
1	B	253	ASP	CB-CA-C	5.32	121.04	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	391[B]	HIS	Peptide

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	2984	0	2893	26	0
1	B	2984	0	2889	19	0
2	A	1	0	0	0	0
3	A	28	0	12	1	0
3	B	47	0	24	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	12	0	13	1	0
5	B	12	0	13	0	0
6	A	240	0	0	9	0
6	B	234	0	0	3	0
All	All	6544	0	5844	47	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 4.

All (47) 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:333[B]:GLN:OE1	1:A:335[B]:ASP:OD1	1.55	1.24
1:B:323[B]:GLN:OE1	1:B:327:ASN:ND2	1.73	1.19
1:A:392[B]:ARG:NH2	6:A:2238:HOH:O	1.90	1.03
1:A:283[B]:GLN:OE1	6:A:2168:HOH:O	2.02	0.76
1:B:323[B]:GLN:CD	1:B:327:ASN:HD22	1.92	0.73
1:A:293[A]:GLU:OE1	6:A:2181:HOH:O	2.05	0.73
1:A:137:ASP:HB3	1:A:141:ARG:HH12	1.57	0.68
1:A:231[B]:GLN:HG3	6:A:2091:HOH:O	1.94	0.67
1:B:231[B]:GLN:HG3	6:B:2103:HOH:O	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231[B]:GLN:CG	6:B:2103:HOH:O	2.45	0.64
1:A:109:HIS:HB3	1:A:145:GLN:HB3	1.82	0.60
1:A:231[B]:GLN:CG	6:A:2091:HOH:O	2.48	0.60
1:A:200:LYS:HG3	6:A:2096:HOH:O	2.01	0.59
1:B:231[A]:GLN:HG3	1:B:353:LEU:HD21	1.85	0.58
1:A:91:ASN:O	1:A:95:ILE:HG13	2.04	0.58
1:B:231[B]:GLN:OE1	1:B:357[B]:TYR:OH	2.23	0.56
1:B:231[B]:GLN:HG2	1:B:356:TYR:CE2	2.41	0.56
1:B:303:LEU:HD13	1:B:357[A]:TYR:HE2	1.71	0.56
3:A:1396:50T:N1	6:A:2187:HOH:O	2.34	0.54
1:A:137:ASP:HB3	1:A:141:ARG:NH1	2.23	0.54
1:B:161:LEU:HD22	1:B:210[A]:VAL:HG11	1.91	0.53
1:A:323[B]:GLN:OE1	1:A:327:ASN:ND2	2.42	0.52
1:A:77:TRP:CD2	1:A:107:PRO:HD3	2.45	0.51
5:A:1398:MES:H32	5:A:1398:MES:H81	1.52	0.48
1:A:328:GLN:HG3	6:A:2194:HOH:O	2.14	0.48
1:A:136:LEU:C	1:A:136:LEU:HD23	2.33	0.48
1:A:115:PHE:O	1:A:151:ALA:HA	2.15	0.45
1:A:231[B]:GLN:OE1	1:A:357[B]:TYR:OH	2.34	0.45
1:B:76:ILE:HD12	1:B:108:ILE:HD13	1.98	0.45
1:A:231[B]:GLN:HG2	1:A:356:TYR:CE2	2.53	0.44
1:A:111:ILE:HB	1:A:112:PRO:CD	2.47	0.44
1:B:231[B]:GLN:CD	1:B:357[B]:TYR:OH	2.56	0.44
1:B:231[B]:GLN:HG2	1:B:356:TYR:HE2	1.81	0.43
1:A:392[B]:ARG:CZ	6:A:2238:HOH:O	2.50	0.43
1:B:393:ALA:HB2	6:B:2232:HOH:O	2.19	0.43
1:A:111:ILE:HB	1:A:112:PRO:HD2	2.00	0.42
1:A:179:ASN:O	1:A:179:ASN:OD1	2.37	0.42
1:A:57:LEU:HD22	1:A:103:LEU:HD23	2.02	0.41
1:B:231[A]:GLN:HE22	1:B:346:TYR:HE2	1.68	0.41
1:B:234[B]:MET:HG2	1:B:356:TYR:CE1	2.55	0.41
1:B:210[B]:VAL:HG23	1:B:221:LEU:HB2	2.03	0.41
1:B:265:LYS:HE3	1:B:265:LYS:HB3	1.88	0.41
1:A:394:LEU:HA	1:A:394:LEU:HD23	1.81	0.41
1:A:85:PHE:O	1:A:88:ILE:HG12	2.21	0.41
1:A:103:LEU:HD12	1:A:139:TRP:CH2	2.56	0.41
1:B:231[A]:GLN:HG3	1:B:353:LEU:CD2	2.51	0.40
1:B:58:HIS:NE2	1:B:383:ASP:OD2	2.54	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	379/368 (103%)	365 (96%)	14 (4%)	0	100	100
1	B	379/368 (103%)	365 (96%)	14 (4%)	0	100	100
All	All	758/736 (103%)	730 (96%)	28 (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	341/328 (104%)	330 (97%)	11 (3%)	46	18
1	B	341/328 (104%)	336 (98%)	5 (2%)	72	50
All	All	682/656 (104%)	666 (98%)	16 (2%)	60	29

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	GLU
1	A	56	ARG
1	A	95	ILE
1	A	109	HIS
1	A	180	LYS
1	A	200	LYS
1	A	212	LEU
1	A	234[A]	MET

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Mol	Chain	Res	Type
1	A	234[B]	MET
1	A	265	LYS
1	A	386	ILE
1	B	56	ARG
1	B	95	ILE
1	B	147[A]	GLN
1	B	147[B]	GLN
1	B	265	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	HIS
1	A	179	ASN
1	A	327	ASN
1	A	328	GLN
1	B	319	GLN
1	B	327	ASN
1	B	370	ASN

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)

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\#\|Z\| > 2$
3	50T	A	1396	4	19,30,30	1.82	4 (21%)	21,49,49	2.70	4 (19%)
5	MES	A	1398	-	11,12,12	1.01	0	14,16,16	3.40	5 (35%)
3	50T	B	1395[A]	-	19,30,30	1.68	4 (21%)	21,49,49	2.66	6 (28%)
3	50T	B	1395[B]	-	19,30,30	1.66	3 (15%)	21,49,49	3.15	5 (23%)
5	MES	B	1397	-	11,12,12	0.73	0	14,16,16	3.87	4 (28%)

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	50T	A	1396	4	-	0/6/33/33	0/3/3/3
5	MES	A	1398	-	-	0/6/14/14	0/1/1/1
3	50T	B	1395[A]	-	-	0/6/33/33	0/3/3/3
3	50T	B	1395[B]	-	-	0/6/33/33	0/3/3/3
5	MES	B	1397	-	-	0/6/14/14	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1396	50T	O3'-C3'	2.10	1.48	1.43
3	B	1395[B]	50T	PA-O1A	2.13	1.54	1.48
3	B	1395[A]	50T	PA-O1A	2.13	1.54	1.48
3	B	1395[A]	50T	C2-N3	2.58	1.36	1.32
3	A	1396	50T	C2-N3	2.97	1.37	1.32
3	B	1395[B]	50T	PA-O2A	3.53	1.57	1.48
3	B	1395[A]	50T	PA-O2A	3.53	1.57	1.48
3	B	1395[A]	50T	C6-N6	4.06	1.47	1.34
3	A	1396	50T	C6-N6	4.15	1.47	1.34
3	A	1396	50T	PA-O2A	4.70	1.59	1.48
3	B	1395[B]	50T	C6-N6	5.10	1.50	1.34

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	1397	MES	O1S-S-C8	-11.83	96.81	106.91
3	B	1395[B]	50T	N3-C2-N1	-11.45	120.13	128.89
3	B	1395[A]	50T	N3-C2-N1	-9.65	121.51	128.89
3	A	1396	50T	N3-C2-N1	-9.50	121.62	128.89
3	A	1396	50T	C2'-C1'-N9	-4.80	106.96	114.29
3	B	1395[B]	50T	C4'-O4'-C1'	-4.58	104.69	109.72
3	A	1396	50T	C4'-O4'-C1'	-3.52	105.85	109.72
3	B	1395[A]	50T	C2'-C1'-N9	-3.45	109.02	114.29
3	B	1395[B]	50T	O2A-PA-O1A	-3.40	108.58	118.70
3	B	1395[A]	50T	O2A-PA-O1A	-3.40	108.58	118.70
5	B	1397	MES	O3S-S-O2S	-3.35	103.80	111.61
3	B	1395[A]	50T	C1'-N9-C4	-2.53	123.12	126.94
3	B	1395[A]	50T	O4'-C1'-N9	2.41	113.14	108.10
5	B	1397	MES	C6-O1-C2	2.43	118.08	109.89
5	A	1398	MES	C2-C3-N4	2.64	114.12	110.12
5	A	1398	MES	O3S-S-O1S	2.65	117.78	111.61
3	B	1395[B]	50T	C2-N1-C6	2.86	123.87	118.77
3	A	1396	50T	PA-O5'-C5'	2.88	130.16	120.25
3	B	1395[A]	50T	N6-C6-N1	3.10	125.85	119.20
5	A	1398	MES	O1S-S-C8	4.25	110.53	106.91
5	A	1398	MES	C5-N4-C3	4.33	118.29	108.90
3	B	1395[B]	50T	O4'-C1'-N9	4.47	117.47	108.10
5	B	1397	MES	O2S-S-C8	6.22	112.21	106.91
5	A	1398	MES	O2S-S-C8	10.06	115.49	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1396	50T	1	0
5	A	1398	MES	1	0

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^{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(Å^2)	Q<0.9
1	A	358/368 (97%)	0.92	62 (17%)	2	2	3, 8, 18, 30	0
1	B	358/368 (97%)	0.61	34 (9%)	10	9	3, 7, 17, 27	0
All	All	716/736 (97%)	0.76	96 (13%)	4	4	3, 8, 18, 30	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	138	TYR	7.4
1	A	144	SER	7.2
1	A	107	PRO	6.5
1	A	104	ALA	6.0
1	B	177	VAL	5.7
1	B	394	LEU	5.7
1	A	37	ASN	5.5
1	A	108	ILE	4.9
1	A	177	VAL	4.9
1	B	66	ASP	4.8
1	A	145	GLN	4.7
1	B	93	VAL	4.6
1	A	93	VAL	4.5
1	A	92	SER	4.4
1	A	140	PHE	4.4
1	A	66	ASP	4.3
1	B	393	ALA	4.3
1	A	96	ASP	4.2
1	A	179	ASN	4.2
1	A	105	ASP	4.2
1	A	77	TRP	4.1
1	B	179	ASN	4.1
1	B	104	ALA	3.8
1	A	136	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	394	LEU	3.8
1	A	109	HIS	3.7
1	A	41	LYS	3.7
1	A	40	GLU	3.7
1	A	143	GLN	3.7
1	A	91	ASN	3.6
1	A	141	ARG	3.6
1	B	108	ILE	3.5
1	A	216	GLY	3.5
1	A	130	LYS	3.5
1	B	176	SER	3.5
1	A	139	TRP	3.3
1	A	131	LYS	3.2
1	A	97	ALA	3.2
1	A	100	THR	3.2
1	B	107	PRO	3.1
1	B	216	GLY	3.1
1	A	135	GLU	3.1
1	B	37	ASN	3.1
1	A	142	GLN	3.0
1	B	229	LEU	3.0
1	A	38	PRO	2.9
1	A	95	ILE	2.9
1	B	343	ASP	2.9
1	A	379	GLU	2.8
1	A	67	THR	2.8
1	A	94	THR	2.8
1	A	175	LYS	2.8
1	B	180	LYS	2.8
1	B	41	LYS	2.8
1	B	175	LYS	2.8
1	A	110	ASN	2.7
1	A	137	ASP	2.7
1	A	127	SER	2.7
1	A	360	LEU	2.7
1	B	213	ASN	2.7
1	A	106	ALA	2.6
1	B	379	GLU	2.6
1	A	192	VAL	2.6
1	B	192	VAL	2.6
1	B	357[A]	TYR	2.5
1	A	126	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	39	CYS	2.5
1	A	268	VAL	2.5
1	A	103	LEU	2.5
1	B	160	ALA	2.5
1	A	357[A]	TYR	2.4
1	B	378	PRO	2.4
1	A	134	ASP	2.4
1	B	162	PHE	2.4
1	B	67	THR	2.4
1	A	101	MET	2.3
1	A	176	SER	2.3
1	B	178	GLN	2.3
1	A	213	ASN	2.3
1	A	378	PRO	2.3
1	A	146	TRP	2.3
1	A	227	LEU	2.3
1	B	92	SER	2.2
1	B	232	THR	2.2
1	B	230	GLY	2.2
1	B	356	TYR	2.2
1	B	234[A]	MET	2.2
1	A	89	GLN	2.1
1	A	200	LYS	2.1
1	B	226	PHE	2.1
1	A	147	GLN	2.0
1	B	360	LEU	2.0
1	A	172	ASP	2.0
1	B	40	GLU	2.0
1	A	162	PHE	2.0
1	A	128	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å^2)	Q<0.9
2	NA	A	1395	1/1	0.96	0.28	9.80	37,37,37,37	0
4	MG	A	1397	1/1	0.99	0.26	6.59	16,16,16,16	0
4	MG	B	1396	1/1	0.99	0.23	3.97	15,15,15,15	0
5	MES	B	1397	12/12	0.94	0.23	3.79	25,32,38,40	12
5	MES	A	1398	12/12	0.89	0.22	2.56	24,34,41,43	12
3	50T	B	1395[A]	28/28	0.95	0.19	2.09	14,19,21,24	19
3	50T	B	1395[B]	28/28	0.95	0.19	2.09	14,17,18,19	19
3	50T	A	1396	28/28	0.95	0.17	1.48	14,23,34,38	0

6.5 Other polymers (i)

There are no such residues in this entry.