



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:42 PM GMT

PDB ID : 1E2M  
Title : HPT + HMTT  
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Deposited on : 2000-05-23  
Resolution : 2.20 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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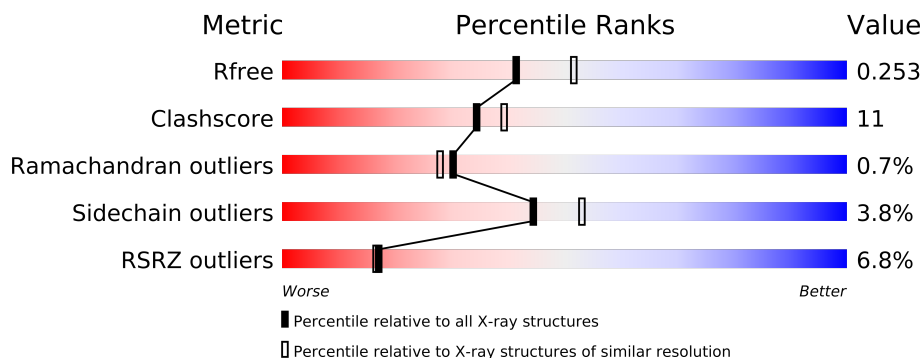
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.20 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))
$R_{free}$	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	331	
1	B	331	

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 4965 atoms, of which 0 are hydrogen and 0 are deuterium.

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 THYMIDINE KINASE.

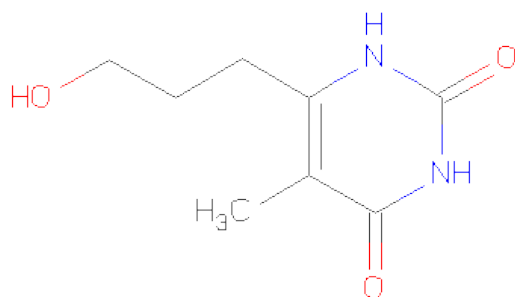
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2341	1490	410	425	16			
1	B	310	Total	C	N	O	S	0	0	0
			2359	1502	410	431	16			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 3 is 6-HYDROXYPROPYLTHYMINE (three-letter code: HPT) (formula: C<sub>8</sub>H<sub>12</sub>N<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	8	2	3		
3	B	1	Total	C	N	O	0	0
			13	8	2	3		

- Molecule 4 is water.

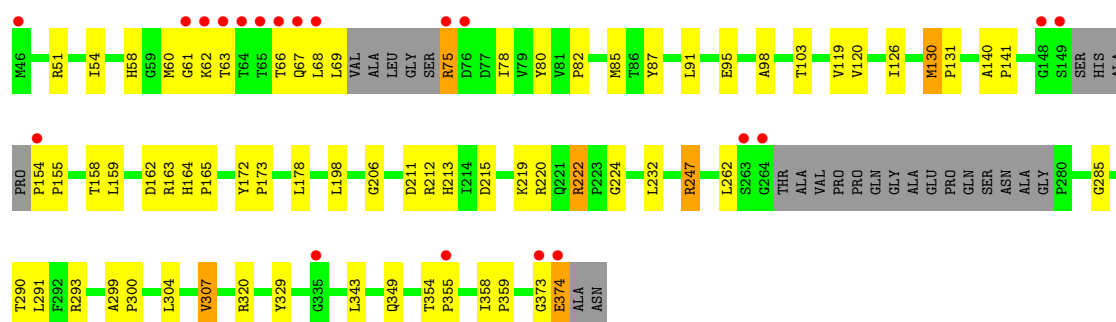
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	118	Total	O	0	0
			118	118		
4	B	111	Total	O	0	0
			111	111		

### 3 Residue-property plots

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.

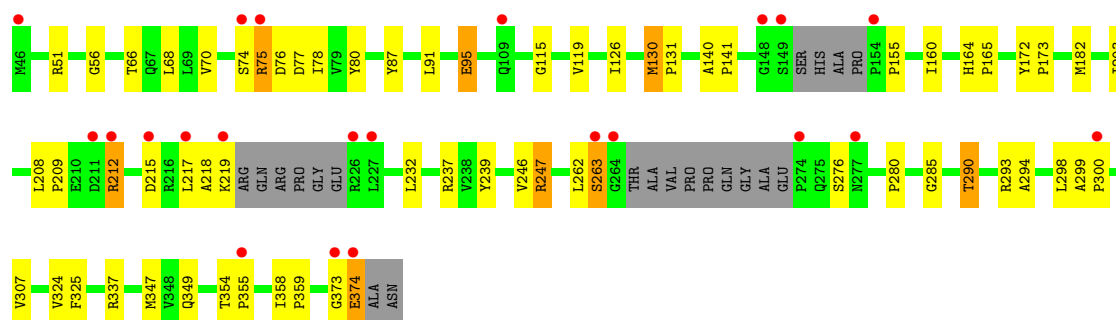
#### • Molecule 1: THYMIDINE KINASE

Chain A: 



#### • Molecule 1: THYMIDINE KINASE

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	113.70Å 117.20Å 108.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 24.89 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (20.00-2.20) 98.6 (24.89-2.20)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.67 (at 2.19Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.217 , 0.262 0.208 , 0.253	Depositor DCC
$R_{free}$ test set	1854 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.5	Xtriage
Anisotropy	0.614	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 26.3	EDS
Estimated twinning fraction	0.034 for -k,-h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	3 of 36530 reflections (0.008%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4965	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, HPT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.41	0/2394	0.90	7/3261 (0.2%)
1	B	0.41	0/2412	0.85	4/3288 (0.1%)
All	All	0.41	0/4806	0.88	11/6549 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	320	ARG	NE-CZ-NH2	-10.10	115.25	120.30
1	A	320	ARG	NE-CZ-NH1	9.90	125.25	120.30
1	A	320	ARG	CD-NE-CZ	7.47	134.05	123.60
1	A	51	ARG	CD-NE-CZ	7.32	133.85	123.60
1	B	51	ARG	CD-NE-CZ	7.17	133.63	123.60
1	B	237	ARG	CD-NE-CZ	6.07	132.10	123.60
1	A	222	ARG	CD-NE-CZ	5.67	131.54	123.60
1	A	51	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	A	307	VAL	CB-CA-C	-5.26	101.40	111.40
1	B	95	GLU	CA-C-N	5.16	128.56	117.20
1	B	237	ARG	NE-CZ-NH1	5.05	122.82	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2341	0	2357	53	0
1	B	2359	0	2374	51	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	13	0	12	0	0
3	B	13	0	12	0	0
4	A	118	0	0	8	0
4	B	111	0	0	2	0
All	All	4965	0	4755	102	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (102) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:119:VAL:HG12	1:B:119:VAL:HG12	1.56	0.87
1:A:130:MET:HB3	1:A:131:PRO:HD3	1.58	0.86
1:A:358:ILE:HB	1:A:359:PRO:HD3	1.61	0.82
1:B:130:MET:HB3	1:B:131:PRO:HD3	1.62	0.81
1:A:126:ILE:HD11	1:B:126:ILE:HD11	1.64	0.78
1:B:358:ILE:HB	1:B:359:PRO:HD3	1.67	0.77
1:A:60:MET:HE3	1:A:329:TYR:CD2	2.21	0.76
1:B:75:ARG:HE	1:B:75:ARG:H	1.34	0.75
1:A:60:MET:HE1	1:A:206:GLY:HA3	1.71	0.73
1:A:61:GLY:HA2	1:A:220:ARG:NH1	2.05	0.72
1:A:60:MET:CE	1:A:206:GLY:HA3	2.21	0.71
1:B:276:SER:HB2	1:B:324:VAL:HB	1.73	0.70
1:A:140:ALA:HB3	1:A:141:PRO:HD3	1.73	0.69
1:A:54:ILE:HD12	1:A:66:THR:HG22	1.75	0.68
1:B:285:GLY:HA2	1:B:290:THR:HG23	1.75	0.68
1:A:354:THR:HB	1:A:355:PRO:HD2	1.74	0.67
1:B:262:LEU:O	1:B:263:SER:HB2	1.95	0.67
1:A:154:PRO:HB2	1:A:155:PRO:HD2	1.77	0.66
1:B:75:ARG:NE	1:B:75:ARG:H	1.93	0.66
1:B:217:LEU:HD22	1:B:232:LEU:HD13	1.78	0.66
1:B:66:THR:HG23	1:B:160:ILE:HG21	1.78	0.65
1:B:66:THR:HG21	1:B:80:TYR:CE1	2.32	0.65
1:A:211:ASP:HB2	4:A:2072:HOH:O	1.97	0.65
1:A:62:LYS:HE2	4:A:2004:HOH:O	1.97	0.64
1:B:354:THR:HB	1:B:355:PRO:HD2	1.78	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:285:GLY:HA2	1:A:290:THR:HG23	1.80	0.63
1:A:285:GLY:HA2	1:A:290:THR:CG2	2.29	0.62
1:B:246:VAL:HG21	1:B:324:VAL:HG21	1.81	0.61
1:A:154:PRO:HB2	1:A:155:PRO:CD	2.31	0.59
1:A:54:ILE:CD1	1:A:66:THR:HG22	2.32	0.59
1:A:262:LEU:HD22	1:A:290:THR:HG22	1.85	0.58
1:A:78:ILE:HG12	1:A:158:THR:HB	1.85	0.58
1:B:373:GLY:O	1:B:374:GLU:HB2	2.05	0.56
1:A:130:MET:HB3	1:A:131:PRO:CD	2.35	0.56
1:B:70:VAL:HA	1:B:78:ILE:HD13	1.88	0.56
1:B:247:ARG:HH21	1:B:280:PRO:HD2	1.70	0.56
1:B:212:ARG:HD3	4:B:2057:HOH:O	2.05	0.56
1:B:212:ARG:HA	1:B:212:ARG:HE	1.70	0.55
1:B:285:GLY:HA2	1:B:290:THR:CG2	2.35	0.55
1:B:164:HIS:CG	1:B:165:PRO:HD2	2.41	0.55
1:A:62:LYS:HE3	4:A:2008:HOH:O	2.07	0.55
1:B:212:ARG:HH21	1:B:215:ASP:HB3	1.73	0.54
1:A:62:LYS:NZ	1:A:163:ARG:NH2	2.56	0.54
1:B:262:LEU:HD22	1:B:290:THR:HG23	1.91	0.53
1:A:164:HIS:CG	1:A:165:PRO:HD2	2.44	0.53
1:A:87:TYR:HA	1:A:91:LEU:HB2	1.90	0.53
1:A:62:LYS:HG3	4:A:2008:HOH:O	2.08	0.52
1:A:75:ARG:HB2	4:A:2014:HOH:O	2.09	0.52
1:A:60:MET:HE3	1:A:329:TYR:CE2	2.43	0.52
1:A:62:LYS:O	1:A:66:THR:HG23	2.10	0.52
1:B:70:VAL:HG21	1:B:80:TYR:HB2	1.92	0.51
1:B:325:PHE:CE1	1:B:347:MET:HG2	2.45	0.51
1:B:218:ALA:O	1:B:219:LYS:HB2	2.11	0.51
1:A:373:GLY:O	1:A:374:GLU:HB2	2.11	0.51
1:B:262:LEU:HD22	1:B:290:THR:CG2	2.42	0.50
1:B:87:TYR:HA	1:B:91:LEU:HB2	1.94	0.50
1:A:60:MET:HE2	4:A:2004:HOH:O	2.12	0.50
1:B:246:VAL:HG11	1:B:324:VAL:HG23	1.95	0.49
1:B:115:GLY:O	1:B:119:VAL:HG23	2.13	0.48
1:A:85:MET:HE3	1:A:222:ARG:HD3	1.95	0.48
1:A:60:MET:CE	4:A:2004:HOH:O	2.61	0.48
1:B:208:LEU:HD12	1:B:209:PRO:HD2	1.95	0.48
1:A:358:ILE:HB	1:A:359:PRO:CD	2.38	0.48
1:A:61:GLY:HA2	1:A:220:ARG:HH11	1.79	0.48
1:A:85:MET:CE	1:A:222:ARG:HD3	2.44	0.47
1:A:62:LYS:HZ2	1:A:163:ARG:NH2	2.12	0.47
1:B:74:SER:HB3	1:B:77:ASP:OD2	2.13	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:140:ALA:HA	1:A:198:LEU:HD11	1.96	0.47
1:B:140:ALA:HB3	1:B:141:PRO:CD	2.44	0.47
1:A:67:GLN:O	1:A:69:LEU:N	2.48	0.46
1:B:172:TYR:N	1:B:173:PRO:CD	2.79	0.46
1:A:299:ALA:HB1	1:A:300:PRO:HD2	1.98	0.46
1:A:58:HIS:HB3	1:A:163:ARG:NH2	2.32	0.45
1:B:74:SER:HA	1:B:75:ARG:HH21	1.81	0.45
1:A:247:ARG:HD3	4:A:2086:HOH:O	2.17	0.44
1:A:178:LEU:HD21	1:A:291:LEU:HD22	1.98	0.44
1:A:63:THR:O	1:A:67:GLN:HG3	2.18	0.44
1:B:74:SER:HB3	1:B:77:ASP:CG	2.38	0.43
1:B:75:ARG:HE	1:B:75:ARG:N	2.08	0.43
1:B:140:ALA:HB3	1:B:141:PRO:HD3	1.99	0.43
1:A:155:PRO:HB3	1:A:349:GLN:NE2	2.33	0.43
1:A:103:THR:OG1	1:A:120:VAL:HG11	2.18	0.43
1:B:294:ALA:O	1:B:298:LEU:HD13	2.18	0.43
1:A:293:ARG:HB3	1:A:293:ARG:NH1	2.34	0.43
1:A:80:TYR:O	1:A:82:PRO:HD3	2.18	0.43
1:B:155:PRO:HB3	1:B:349:GLN:NE2	2.34	0.43
1:B:293:ARG:NH1	1:B:293:ARG:HB3	2.35	0.42
1:B:203:ILE:CG2	1:B:324:VAL:HG22	2.49	0.42
1:B:247:ARG:NH1	1:B:247:ARG:HG2	2.34	0.42
1:A:172:TYR:N	1:A:173:PRO:CD	2.83	0.41
1:A:213:HIS:CE1	1:A:232:LEU:HD11	2.56	0.41
1:B:182:MET:HG3	4:B:2030:HOH:O	2.20	0.41
1:A:98:ALA:HB2	1:A:224:GLY:HA3	2.03	0.41
1:B:87:TYR:OH	1:B:130:MET:HG3	2.21	0.41
1:B:130:MET:HB3	1:B:131:PRO:CD	2.41	0.41
1:A:60:MET:HE1	1:A:206:GLY:CA	2.45	0.41
1:B:262:LEU:O	1:B:263:SER:CB	2.66	0.41
1:B:299:ALA:HB1	1:B:300:PRO:CD	2.51	0.41
1:B:56:GLY:HA2	1:B:239:TYR:CD1	2.56	0.40
1:A:162:ASP:O	1:A:163:ARG:HB2	2.21	0.40
1:B:218:ALA:O	1:B:219:LYS:CB	2.68	0.40
1:B:68:LEU:HB3	1:B:337:ARG:HD3	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	297/331 (90%)	284 (96%)	11 (4%)	2 (1%)	30	28
1	B	302/331 (91%)	291 (96%)	9 (3%)	2 (1%)	30	28
All	All	599/662 (90%)	575 (96%)	20 (3%)	4 (1%)	30	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	68	LEU
1	B	263	SER
1	A	95	GLU
1	B	95	GLU

### 5.3.2 Protein sidechains ⓘ

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	247/264 (94%)	236 (96%)	11 (4%)	38	44
1	B	249/264 (94%)	241 (97%)	8 (3%)	51	62
All	All	496/528 (94%)	477 (96%)	19 (4%)	44	53

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	ARG
1	A	130	MET
1	A	159	LEU
1	A	212	ARG
1	A	215	ASP

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Mol	Chain	Res	Type
1	A	219	LYS
1	A	247	ARG
1	A	304	LEU
1	A	307	VAL
1	A	343	LEU
1	A	374	GLU
1	B	75	ARG
1	B	76	ASP
1	B	130	MET
1	B	212	ARG
1	B	247	ARG
1	B	290	THR
1	B	307	VAL
1	B	374	GLU

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

Mol	Chain	Res	Type
1	A	349	GLN
1	B	349	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	A	400	-	4,4,4	1.00	0	6,6,6	0.13	0
3	HPT	A	500	-	13,13,13	1.75	4 (30%)	13,17,17	1.75	4 (30%)
2	SO4	B	400	-	4,4,4	1.02	0	6,6,6	0.24	0
3	HPT	B	500	-	13,13,13	1.76	5 (38%)	13,17,17	1.67	3 (23%)

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	SO4	A	400	-	-	0/0/0/0	0/0/0/0
3	HPT	A	500	-	-	0/4/4/4	0/1/1/1
2	SO4	B	400	-	-	0/0/0/0	0/0/0/0
3	HPT	B	500	-	-	0/4/4/4	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	HPT	C2-N2	3.97	1.43	1.37
3	B	500	HPT	C2-N2	3.64	1.43	1.37
3	A	500	HPT	C1-N2	2.75	1.42	1.37
3	B	500	HPT	C5-N1	2.53	1.41	1.36
3	B	500	HPT	C2-C3	2.39	1.48	1.42
3	B	500	HPT	C1-N2	2.37	1.41	1.37
3	A	500	HPT	C5-N1	2.35	1.40	1.36
3	A	500	HPT	C1-N1	2.31	1.41	1.37
3	B	500	HPT	C1-N1	2.23	1.41	1.37

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	HPT	C2-C3-C5	4.46	121.18	116.51
3	B	500	HPT	C2-C3-C5	3.95	120.64	116.51
3	B	500	HPT	C4-C3-C2	-2.98	115.46	121.20
3	A	500	HPT	C2-N2-C1	-2.35	120.60	125.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	HPT	C12-C11-C5	2.24	118.67	112.91
3	B	500	HPT	C2-N2-C1	-2.11	121.08	125.36
3	A	500	HPT	C4-C3-C2	-2.07	117.23	121.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	305/331 (92%)	0.10	20 (6%) 18 17	12, 24, 57, 91	0
1	B	310/331 (93%)	0.02	22 (7%) 16 15	11, 24, 55, 73	0
All	All	615/662 (92%)	0.06	42 (6%) 17 16	11, 24, 57, 91	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	GLY	8.2
1	A	75	ARG	7.9
1	A	65	THR	7.3
1	B	154	PRO	5.9
1	A	149	SER	5.8
1	B	264	GLY	5.3
1	B	149	SER	5.2
1	A	264	GLY	5.1
1	B	227	LEU	4.9
1	B	374	GLU	4.8
1	B	263	SER	4.4
1	A	62	LYS	4.4
1	A	76	ASP	4.2
1	A	373	GLY	4.0
1	A	61	GLY	4.0
1	B	148	GLY	3.9
1	B	217	LEU	3.7
1	A	68	LEU	3.7
1	A	335	GLY	3.6
1	A	374	GLU	3.4
1	B	74	SER	3.3
1	B	373	GLY	3.2
1	A	63	THR	3.2
1	A	154	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	300	PRO	2.9
1	B	211	ASP	2.7
1	B	75	ARG	2.7
1	B	46	MET	2.7
1	A	66	THR	2.6
1	B	274	PRO	2.6
1	B	226	ARG	2.5
1	B	277	ASN	2.4
1	A	64	THR	2.4
1	B	219	LYS	2.3
1	A	46	MET	2.3
1	B	109	GLN	2.3
1	B	355	PRO	2.2
1	A	67	GLN	2.1
1	A	355	PRO	2.1
1	A	263	SER	2.1
1	B	215	ASP	2.1
1	B	212	ARG	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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<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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	HPT	A	500	13/13	0.15	1.66	11,18,29,33	0
3	HPT	B	500	13/13	0.12	1.37	18,21,30,33	0
2	SO4	B	400	5/5	0.12	0.19	26,27,29,34	0
2	SO4	A	400	5/5	0.16	-0.53	42,42,44,46	0



## 6.5 Other polymers ⓘ

There are no such residues in this entry.