



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:33 AM GMT

PDB ID : 2AOV
Title : Histamine Methyltransferase Complexed with the Antifolate Drug Metoprine
Authors : Horton, J.R.; Sawada, K.; Nishibori, M.; Cheng, X.
Deposited on : 2005-08-14
Resolution : 2.48 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

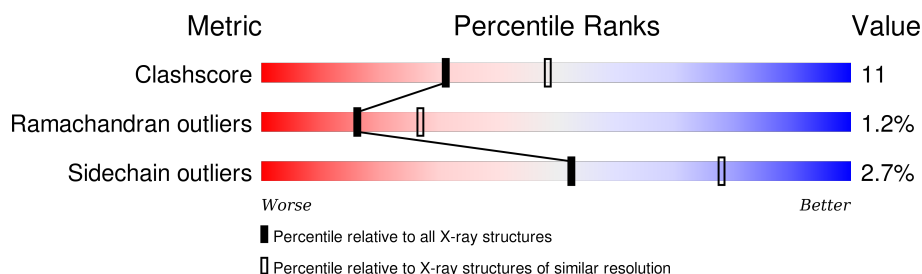
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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))
Clashscore	102246	5050 (2.50-2.46)
Ramachandran outliers	100387	4961 (2.50-2.46)
Sidechain outliers	100360	4963 (2.50-2.46)

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 $\leq 5\%$. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	292	
1	B	292	

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
3	C2M	A	401	-	X	-	-
3	C2M	B	400	-	X	-	-

2 Entry composition [i](#)

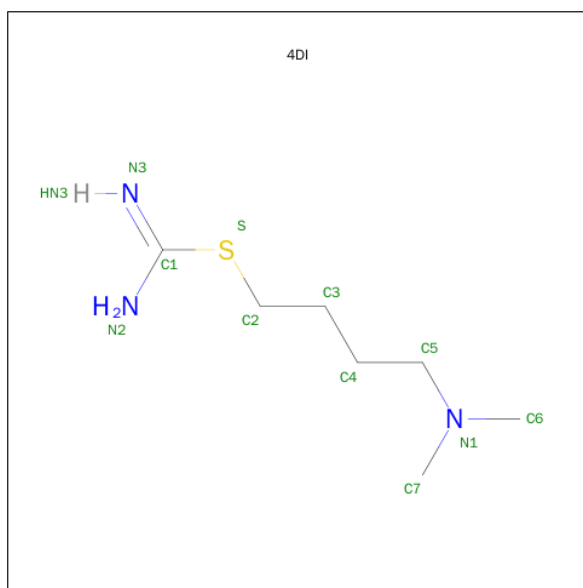
There are 4 unique types of molecules in this entry. The entry contains 4676 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Histamine N-methyltransferase.

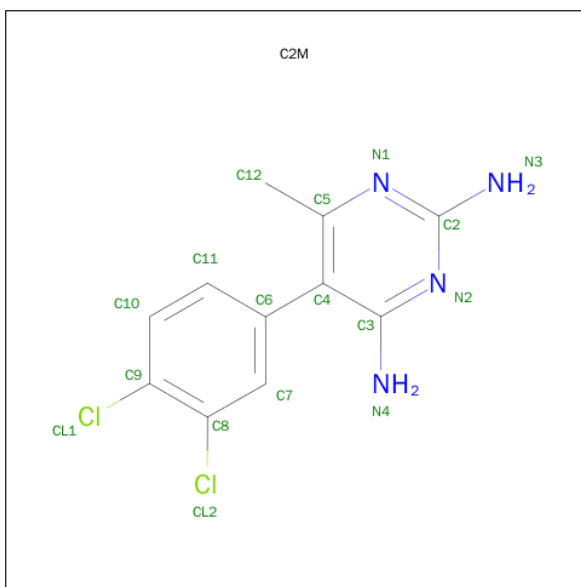
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	4	0
			2254	1435	370	433	16			
1	B	288	Total	C	N	O	S	0	2	0
			2275	1454	369	437	15			

- Molecule 2 is 4-(DIMETHYLAMINO)BUTYL IMIDOTHIOCARBAMATE (three-letter code: 4DI) (formula: C₇H₁₇N₃S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	S	0	0
			11	7	3	1		
2	A	1	Total	C	N	S	0	0
			11	7	3	1		

- Molecule 3 is 5-(3,4-DICHLOROPHENYL)-6-METHYLPYRIMIDINE-2,4-DIAMINE (three-letter code: C2M) (formula: C₁₁H₁₀Cl₂N₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	Cl	N	0	0
			17	11	2	4		
3	A	1	Total	C	Cl	N	0	0
			17	11	2	4		

- Molecule 4 is water.

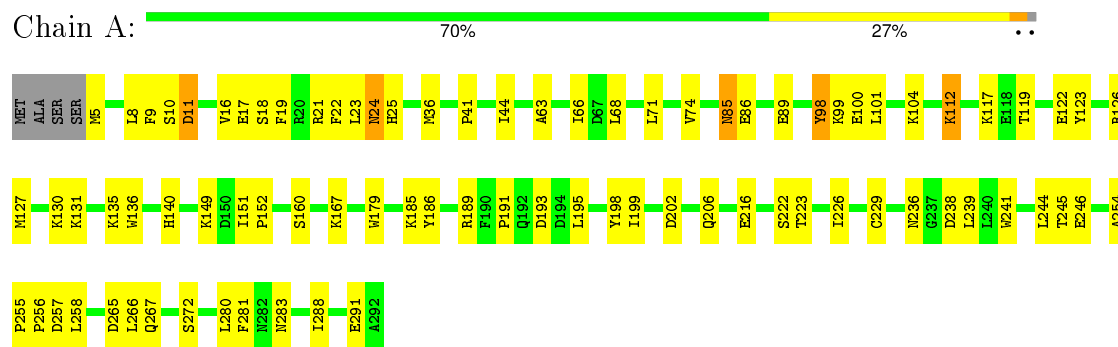
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	47	Total	O	0	0
			47	47		
4	B	44	Total	O	0	0
			44	44		

3 Residue-property plots [i](#)

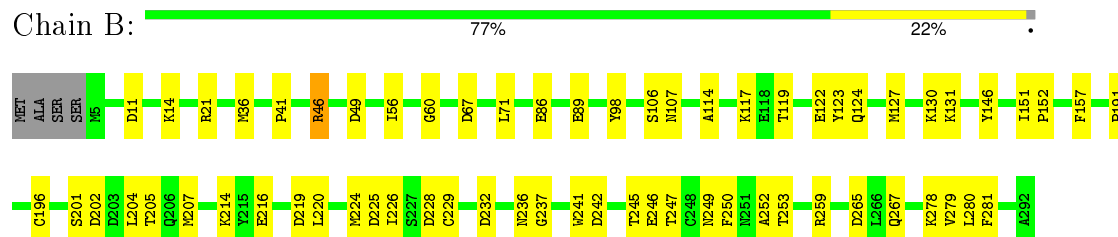
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

• Molecule 1: Histamine N-methyltransferase



• Molecule 1: Histamine N-methyltransferase



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α , β , γ	132.89Å 132.89Å 65.05Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.00 – 2.48	Depositor
% Data completeness (in resolution range)	(Not available) (28.00-2.48)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.211 , 0.254	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4676	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 4DI, C2M

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	0.39	0/2299	0.58	0/3108
1	B	0.36	0/2324	0.55	0/3140
All	All	0.37	0/4623	0.56	0/6248

There are no bond length outliers.

There are no bond angle outliers.

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	2254	0	2151	55	0
1	B	2275	0	2186	43	0
2	A	11	0	17	0	0
2	B	11	0	17	1	0
3	A	17	0	10	0	0
3	B	17	0	10	0	0
4	A	47	0	0	5	0
4	B	44	0	0	0	0
All	All	4676	0	4391	97	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 11.

All (97) 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:86:GLU:HG3	1:A:136:TRP:HH2	1.39	0.87
1:B:86:GLU:HG2	1:B:114:ALA:HB3	1.59	0.84
1:A:86:GLU:HG3	1:A:136:TRP:CH2	2.18	0.79
1:A:8:LEU:C	1:A:10:SER:H	1.86	0.76
1:B:249:ASN:O	1:B:253:THR:HG22	1.85	0.76
1:A:272:SER:OG	1:A:281:PHE:HA	1.87	0.74
1:B:225:ASP:OD2	1:B:278:LYS:HD2	1.87	0.74
1:B:11:ASP:HB3	1:B:14:LYS:HB2	1.73	0.71
1:A:236:ASN:HA	1:A:239:LEU:HD12	1.74	0.69
1:A:11:ASP:HA	4:A:425:HOH:O	1.94	0.67
1:A:185:LYS:HD3	1:A:186:TYR:CE1	2.30	0.66
1:A:10:SER:O	1:A:11:ASP:CB	2.44	0.66
1:B:249:ASN:HB3	1:B:252:ALA:HB3	1.78	0.65
1:A:63:ALA:O	1:A:98:TYR:HA	1.97	0.64
1:B:226:ILE:HG13	1:B:279:VAL:HG13	1.81	0.63
1:A:19:PHE:O	1:A:22:PHE:HB3	1.97	0.63
1:A:185:LYS:HD3	1:A:186:TYR:HE1	1.66	0.60
1:A:21:ARG:O	1:A:25:HIS:HB2	2.01	0.60
1:B:151:ILE:HB	1:B:152:PRO:HD3	1.85	0.59
1:A:112:LYS:HB3	1:A:112:LYS:NZ	2.18	0.58
1:A:8:LEU:C	1:A:10:SER:N	2.57	0.58
1:B:89:GLU:O	1:B:117:LYS:HA	2.04	0.57
1:A:100:GLU:O	1:A:104:LYS:HG3	2.04	0.57
1:B:21:ARG:HD2	1:B:236:ASN:ND2	2.20	0.57
1:B:86:GLU:CG	1:B:114:ALA:HB3	2.32	0.56
1:A:245:THR:O	1:A:246:GLU:HB3	2.05	0.56
1:A:226:ILE:HB	1:A:229:CYS:SG	2.45	0.56
1:B:46:ARG:HG2	1:B:49:ASP:OD1	2.05	0.56
1:A:8:LEU:O	1:A:10:SER:N	2.39	0.55
1:A:24:ASN:OD1	1:A:25:HIS:ND1	2.37	0.55
1:A:167:LYS:HE2	1:A:291:GLU:OE1	2.07	0.55
1:A:123:TYR:O	1:A:127:MET:HG2	2.07	0.55
1:A:265:ASP:C	1:A:267:GLN:H	2.11	0.55
1:A:246:GLU:O	1:A:246:GLU:HG3	2.07	0.54
1:A:151:ILE:HB	1:A:152:PRO:HD3	1.89	0.54
1:B:219:ASP:O	1:B:220:LEU:HD23	2.09	0.53
1:A:179:TRP:CH2	1:A:244:LEU:HD23	2.43	0.53
1:A:89:GLU:O	1:A:117:LYS:HA	2.08	0.53
1:A:288:ILE:HD12	1:A:288:ILE:N	2.24	0.53
1:B:60:GLY:O	2:B:300:4DI:H52	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:ASP:C	1:B:267:GLN:H	2.13	0.52
1:A:44:ILE:HD12	1:A:74:VAL:HA	1.91	0.51
1:B:11:ASP:HB3	1:B:14:LYS:CB	2.41	0.50
1:A:68:LEU:HD12	1:A:101:LEU:HD21	1.93	0.50
1:A:68:LEU:HD11	1:A:101:LEU:HG	1.94	0.50
1:A:122:GLU:HG2	4:A:446:HOH:O	2.12	0.49
1:B:123:TYR:O	1:B:127:MET:HG2	2.12	0.49
1:A:254:ALA:HB1	1:A:258:LEU:HD23	1.96	0.48
1:A:5:MET:HG2	1:A:191:PRO:HG3	1.95	0.48
1:A:223:THR:HB	1:A:280:LEU:HB3	1.95	0.48
1:B:226:ILE:HG13	1:B:279:VAL:CG1	2.43	0.48
1:A:198:TYR:O	1:A:199:ILE:HG13	2.14	0.48
1:B:124:GLN:HG3	1:B:157:PHE:CD2	2.49	0.47
1:B:119:THR:OG1	1:B:122:GLU:HG3	2.15	0.47
1:B:46:ARG:CG	1:B:49:ASP:OD1	2.62	0.47
1:A:238:ASP:O	1:A:241:TRP:HB2	2.14	0.47
1:A:140:HIS:HE1	4:A:414:HOH:O	1.97	0.47
1:A:99:LYS:NZ	1:B:202:ASP:OD2	2.47	0.47
1:A:16:VAL:HG23	1:A:17:GLU:N	2.30	0.47
1:A:222:SER:HB3	1:A:283:ASN:O	2.15	0.47
1:B:279:VAL:HG13	1:B:279:VAL:O	2.15	0.47
1:A:122:GLU:O	1:A:126[B]:ARG:HB3	2.14	0.47
1:B:106:SER:O	1:B:107:ASN:HB2	2.16	0.46
1:B:204:LEU:HD23	1:B:207:MET:HE3	1.98	0.46
1:A:206:GLN:HB2	4:A:418:HOH:O	2.17	0.45
1:B:250:PHE:C	1:B:252:ALA:H	2.19	0.45
1:A:8:LEU:HB3	1:A:246:GLU:OE1	2.17	0.45
1:A:236:ASN:O	1:A:239:LEU:HB2	2.17	0.45
1:B:259:ARG:NH1	1:B:259:ARG:HB3	2.32	0.44
1:B:201:SER:O	1:B:205:THR:HG23	2.17	0.44
1:B:124:GLN:HG3	1:B:157:PHE:CG	2.54	0.43
1:A:71:LEU:HD21	1:A:85:ASN:OD1	2.19	0.43
1:B:280:LEU:N	1:B:280:LEU:HD22	2.33	0.43
1:B:130:LYS:HB3	1:B:131:LYS:H	1.70	0.43
1:B:228:ASP:O	1:B:237:GLY:HA3	2.19	0.43
1:A:135:LYS:HE2	1:A:160:SER:O	2.18	0.43
1:B:191:PRO:CG	1:B:247:THR:HG22	2.48	0.43
1:B:245:THR:OG1	1:B:247:THR:HG23	2.19	0.42
1:B:242:ASP:O	1:B:246:GLU:N	2.52	0.42
1:A:119:THR:OG1	1:A:122:GLU:HG3	2.19	0.42
1:A:23:LEU:C	1:A:25:HIS:H	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:MET:HB3	1:B:281:PHE:HB3	2.00	0.42
1:B:226:ILE:O	1:B:229:CYS:HB2	2.20	0.42
1:B:36:MET:O	1:B:41:PRO:HD3	2.19	0.42
1:A:36:MET:O	1:A:41:PRO:HD3	2.19	0.41
1:B:146:TYR:O	1:B:196:CYS:HA	2.20	0.41
1:A:140:HIS:HD2	4:A:402:HOH:O	2.02	0.41
1:A:189[A]:ARG:CZ	1:A:258:LEU:HD13	2.50	0.41
1:B:265:ASP:C	1:B:267:GLN:N	2.74	0.41
1:B:237:GLY:O	1:B:241:TRP:HD1	2.02	0.41
1:A:255:PRO:HA	1:A:256:PRO:HD3	1.98	0.41
1:B:124:GLN:HG3	1:B:157:PHE:CE2	2.55	0.41
1:A:66:ILE:HA	1:A:66:ILE:HD13	1.91	0.41
1:A:85:ASN:C	1:A:85:ASN:HD22	2.24	0.41
1:A:193:ASP:C	1:A:195:LEU:H	2.24	0.41
1:B:56:ILE:HD12	1:B:71:LEU:CD2	2.51	0.40
1:B:191:PRO:HD2	1:B:247:THR:HG22	2.03	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	290/292 (99%)	266 (92%)	18 (6%)	6 (2%)	9	13
1	B	288/292 (99%)	267 (93%)	20 (7%)	1 (0%)	46	66
All	All	578/584 (99%)	533 (92%)	38 (7%)	7 (1%)	16	27

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	PHE
1	A	130	LYS

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Mol	Chain	Res	Type
1	A	11	ASP
1	A	131	LYS
1	B	232	ASP
1	A	24	ASN
1	A	266	LEU

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	238/260 (92%)	230 (97%)	8 (3%)	44	70
1	B	245/260 (94%)	240 (98%)	5 (2%)	63	85
All	All	483/520 (93%)	470 (97%)	13 (3%)	52	78

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	SER
1	A	85	ASN
1	A	98	TYR
1	A	112	LYS
1	A	149	LYS
1	A	202	ASP
1	A	216	GLU
1	A	257	ASP
1	B	46	ARG
1	B	67	ASP
1	B	98	TYR
1	B	214	LYS
1	B	216	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	ASN
1	A	140	HIS
1	A	251	ASN
1	B	24	ASN
1	B	25	HIS
1	B	75	GLN
1	B	140	HIS
1	B	165	ASN
1	B	206	GLN
1	B	267	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](#)

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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	4DI	A	301	-	9,10,10	2.20	2 (22%)	10,11,11	4.91	7 (70%)
3	C2M	A	401	-	17,18,18	3.95	13 (76%)	23,26,26	2.63	7 (30%)
2	4DI	B	300	-	9,10,10	2.21	2 (22%)	10,11,11	2.71	4 (40%)
3	C2M	B	400	-	17,18,18	3.84	14 (82%)	23,26,26	2.31	7 (30%)

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	4DI	A	301	-	-	0/8/8/8	0/0/0/0
3	C2M	A	401	-	-	0/4/4/4	0/2/2/2
2	4DI	B	300	-	-	0/8/8/8	0/0/0/0
3	C2M	B	400	-	-	0/4/4/4	0/2/2/2

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	4DI	C1-N2	-4.16	1.26	1.34
2	B	300	4DI	C1-N2	-3.85	1.27	1.34
3	B	400	C2M	C9-CL1	-2.37	1.67	1.73
3	A	401	C2M	C8-CL2	-2.20	1.68	1.73
3	A	401	C2M	C2-N1	2.29	1.39	1.35
3	A	401	C2M	C3-N2	2.48	1.38	1.35
3	B	400	C2M	C10-C9	2.65	1.44	1.39
3	B	400	C2M	C2-N1	3.00	1.40	1.35
3	B	400	C2M	C2-N2	3.22	1.41	1.35
3	A	401	C2M	C3-N4	3.59	1.43	1.34
3	B	400	C2M	C3-N2	3.61	1.40	1.35
3	B	400	C2M	C7-C6	3.64	1.46	1.39
3	B	400	C2M	C7-C8	3.95	1.45	1.38
3	B	400	C2M	C4-C3	3.96	1.48	1.43
3	B	400	C2M	C3-N4	4.07	1.44	1.34
3	A	401	C2M	C10-C9	4.13	1.48	1.39
3	A	401	C2M	C5-N1	4.23	1.40	1.34
3	A	401	C2M	C4-C3	4.42	1.49	1.43
3	B	400	C2M	C11-C10	4.67	1.47	1.38
3	B	400	C2M	C2-N3	4.69	1.43	1.34
3	A	401	C2M	C7-C8	4.85	1.46	1.38
3	A	401	C2M	C7-C6	4.88	1.48	1.39
2	A	301	4DI	C1-N3	4.89	1.46	1.28
3	A	401	C2M	C11-C6	4.96	1.50	1.39
3	B	400	C2M	C9-C8	5.21	1.52	1.39
2	B	300	4DI	C1-N3	5.25	1.47	1.28
3	B	400	C2M	C11-C6	5.55	1.51	1.39
3	A	401	C2M	C11-C10	5.63	1.48	1.38
3	A	401	C2M	C9-C8	5.80	1.53	1.39
3	A	401	C2M	C2-N3	6.18	1.46	1.34
3	B	400	C2M	C5-N1	6.19	1.43	1.34

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	C2M	N2-C2-N1	-5.27	116.89	125.53
2	B	300	4DI	C4-C3-C2	-5.08	92.19	113.18
2	A	301	4DI	C6-N1-C5	-4.52	92.79	110.79
3	B	400	C2M	N2-C2-N1	-4.17	118.70	125.53
2	A	301	4DI	C3-C2-S	-3.91	93.09	112.98
3	B	400	C2M	C7-C8-CL2	-3.65	112.87	118.50
3	A	401	C2M	C12-C5-N1	-3.36	111.02	116.34
3	A	401	C2M	C7-C8-CL2	-2.42	114.77	118.50
3	B	400	C2M	C10-C9-CL1	-2.41	113.41	118.39
3	A	401	C2M	N3-C2-N2	2.11	120.69	117.20
2	B	300	4DI	N2-C1-N3	2.65	128.94	120.62
3	B	400	C2M	C9-C8-CL2	2.75	127.73	120.87
3	A	401	C2M	N3-C2-N1	3.15	122.41	117.20
2	A	301	4DI	C3-C4-C5	3.18	128.24	113.27
2	A	301	4DI	S-C1-N2	3.54	130.11	115.49
3	B	400	C2M	C2-N1-C5	3.73	120.17	117.01
2	B	300	4DI	C2-S-C1	4.18	109.43	100.77
3	B	400	C2M	N3-C2-N1	4.32	124.36	117.20
2	B	300	4DI	C6-N1-C5	4.49	128.64	110.79
2	A	301	4DI	C4-C5-N1	4.96	126.70	113.92
3	B	400	C2M	C2-N2-C3	5.38	123.15	117.04
3	A	401	C2M	C2-N2-C3	5.40	123.18	117.04
2	A	301	4DI	C7-N1-C5	5.82	133.94	110.79
3	A	401	C2M	C2-N1-C5	7.34	123.22	117.01
2	A	301	4DI	C2-S-C1	10.91	123.36	100.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	300	4DI	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.