



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 04:32 PM GMT

PDB ID : 4DPH
Title : Quadruple mutant (N51I+C59R+S108N+I164L)Plasmodium falciparum dihydrofolate reductase-thymidylatesynthase (PfDHFR-TS) complexed with P65 and NADPH
Authors : Yuthavong, Y.; Vilaivan, T.; Kamchonwongpaisan, S.; Charman, S.A.; McLennan, D.N.; White, K.L.; Vivas, L.; Bongard, E.; Chitnumsub, P.; Tarnchompoo, B.; Thongphanchang, C.; Taweechai, S.; Vanichtanakul, J.; Arwon, U.; Fantauzzi, P.; Yuvaniyama, J.; Charman, W.N.; Matthews, D.
Deposited on : 2012-02-13
Resolution : 2.38 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

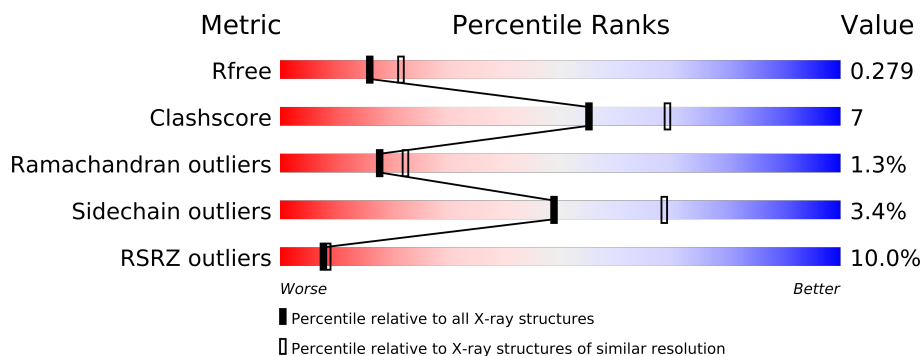
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.38 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	2963 (2.40-2.36)
Clashscore	79885	3668 (2.40-2.36)
Ramachandran outliers	78287	3600 (2.40-2.36)
Sidechain outliers	78261	3602 (2.40-2.36)
RSRZ outliers	66119	2966 (2.40-2.36)

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. 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	608	
1	B	608	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	BME	A	704	-	X
5	BME	A	705	-	X
5	BME	A	706	-	X

2 Entry composition i

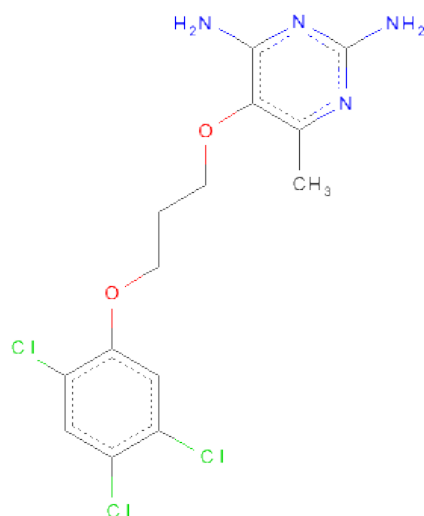
There are 6 unique types of molecules in this entry. The entry contains 9473 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Bifunctional dihydrofolate reductase-thymidylatesynthase.

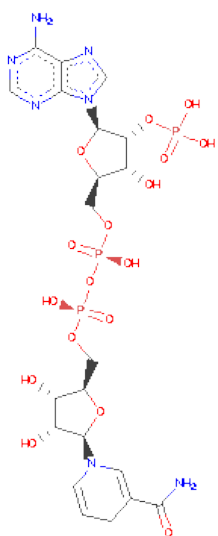
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	538	Total	C	N	O	S	0	0	0
			4472	2890	740	817	25			
1	B	529	Total	C	N	O	S	0	0	0
			4395	2845	725	800	25			

- Molecule 2 is 2,4-DIAMINO-6-METHYL-5-[3-(2,4,5-TRICHLOROPHENOXY)PROPYLOXY]PYRIMIDINE (three-letter code: P65) (formula: C₁₄H₁₅Cl₃N₄O₂).



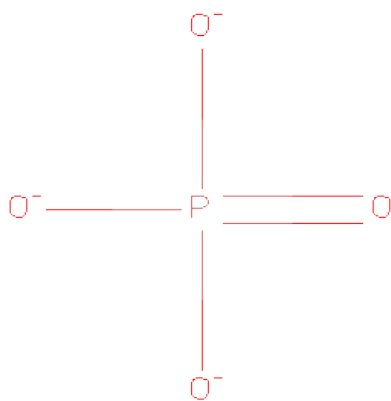
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			23	14	3	4	2		
2	B	1	Total	C	Cl	N	O	0	0
			23	14	3	4	2		

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



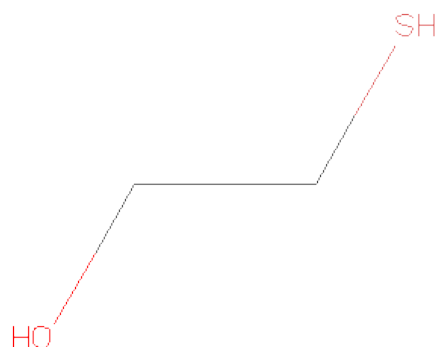
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

- Molecule 5 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C_2H_6OS).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	229	Total	O	0	0
			229	229		
6	B	209	Total	O	0	0
			209	209		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.95Å 156.45Å 164.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.66 – 2.38 51.79 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.1 (54.66-2.38) 99.1 (51.79-2.38)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.219 , 0.286 0.214 , 0.279	Depositor DCC
R_{free} test set	3072 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	45.5	Xtriage
Anisotropy	0.376	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.6	EDS
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	1 of 60475 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9473	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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: NDP, PO4, P65, BME

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	0.71	0/4576	0.74	4/6178 (0.1%)
1	B	0.67	1/4497 (0.0%)	0.71	1/6071 (0.0%)
All	All	0.69	1/9073 (0.0%)	0.73	5/12249 (0.0%)

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	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	494	CYS	CB-SG	-5.35	1.73	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	470	ARG	NE-CZ-NH2	-5.67	117.46	120.30
1	A	487	LEU	CA-CB-CG	5.42	127.78	115.30
1	A	361	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	B	383	LEU	CA-CB-CG	5.21	127.29	115.30
1	A	59	ARG	NE-CZ-NH1	5.11	122.85	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	165	GLY	Peptide

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	4472	0	4426	49	0
1	B	4395	0	4346	73	0
2	A	23	0	15	1	0
2	B	23	0	15	0	0
3	A	48	0	26	2	0
3	B	48	0	26	5	0
4	A	5	0	0	1	0
4	B	5	0	0	0	0
5	A	12	0	18	0	0
5	B	4	0	6	0	0
6	A	229	0	0	6	0
6	B	209	0	0	12	0
All	All	9473	0	8878	120	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 7.

All (120) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:425:ASP:HA	1:B:444:MET:CE	1.76	1.13
1:B:425:ASP:HA	1:B:444:MET:HE3	1.21	1.11
1:B:392:ASN:HA	1:B:444:MET:HE2	1.40	1.03
1:B:376:LEU:HD22	1:B:379:ILE:HD11	1.43	0.98
1:B:547:LEU:HG	6:B:957:HOH:O	1.61	0.97
1:B:584:GLU:HG3	6:B:803:HOH:O	1.66	0.94
1:A:461:ASN:HB3	6:A:973:HOH:O	1.70	0.90
1:B:425:ASP:CA	1:B:444:MET:HE3	2.03	0.89
1:B:392:ASN:HA	1:B:444:MET:CE	2.07	0.85
1:B:307:ILE:HG12	6:B:996:HOH:O	1.77	0.83
1:B:524:SER:HB2	6:B:915:HOH:O	1.80	0.82

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:169:VAL:HG23	3:B:702:NDP:O1A	1.81	0.79
1:B:425:ASP:HA	1:B:444:MET:HE1	1.67	0.74
1:B:171:GLN:HE21	1:B:175:GLU:HG3	1.53	0.73
1:B:210:VAL:HB	6:B:948:HOH:O	1.90	0.72
1:A:12:TYR:HD1	1:A:181:LYS:HB2	1.54	0.71
1:A:512:CYS:SG	1:A:547:LEU:HD22	2.31	0.70
1:B:512:CYS:SG	1:B:547:LEU:HD22	2.33	0.69
1:A:4:GLN:HB2	6:A:847:HOH:O	1.94	0.68
1:A:59:ARG:O	1:A:63:THR:HG23	1.94	0.67
1:B:7:ASP:HA	1:B:180:LYS:HE3	1.76	0.66
1:B:165:GLY:HA3	1:B:170:TYR:CZ	2.32	0.64
1:A:493:LEU:HD12	1:A:493:LEU:C	2.18	0.64
1:A:109:TRP:CZ2	1:A:117:LYS:HD2	2.34	0.63
1:A:312:PHE:CE1	1:A:564:LEU:HD23	2.34	0.63
1:B:210:VAL:O	1:B:210:VAL:HG23	1.97	0.62
1:B:425:ASP:CA	1:B:444:MET:CE	2.65	0.62
1:B:424:ASN:O	1:B:444:MET:HE1	2.00	0.61
1:A:74:LYS:C	6:A:959:HOH:O	2.39	0.60
1:B:43:LYS:N	1:B:194:ASP:OD2	2.29	0.60
1:B:167:SER:HB3	3:B:702:NDP:O2N	2.02	0.59
1:B:392:ASN:CA	1:B:444:MET:HE2	2.26	0.58
1:B:132:LYS:HD2	1:B:133:LYS:HD2	1.85	0.57
1:A:100:ASN:OD1	1:A:159:TYR:HB3	2.03	0.57
1:A:166:GLY:HA3	3:A:702:NDP:O1A	2.05	0.57
1:B:45:VAL:HG12	1:B:46:LEU:H	1.71	0.56
1:B:336:MET:HE1	1:B:560:LEU:HB2	1.87	0.56
1:A:12:TYR:CD1	1:A:181:LYS:HB2	2.39	0.56
1:A:572:PRO:HB3	1:A:596:TYR:HA	1.87	0.56
1:B:373:LYS:HG3	1:B:598:HIS:CE1	2.41	0.55
1:B:493:LEU:HD12	1:B:493:LEU:C	2.26	0.55
1:B:407:ASN:ND2	6:B:991:HOH:O	2.39	0.55
1:B:425:ASP:CB	1:B:444:MET:HE3	2.38	0.54
1:B:171:GLN:NE2	1:B:175:GLU:HG3	2.22	0.53
1:A:312:PHE:HE1	1:A:564:LEU:HD23	1.74	0.53
1:B:109:TRP:CZ2	1:B:117:LYS:HD2	2.45	0.51
1:A:499:PHE:CE1	1:B:340:ASN:HB3	2.44	0.51
1:B:482:LEU:HD22	1:B:488:PRO:HB3	1.93	0.51
1:A:350:VAL:HG12	1:A:553:TYR:CD1	2.46	0.51
1:B:307:ILE:HG23	6:B:958:HOH:O	2.11	0.51
1:B:307:ILE:CG2	1:B:312:PHE:HE2	2.23	0.51
1:B:41:GLY:HA2	1:B:47:PRO:HD3	1.94	0.50
1:A:70:TYR:OH	1:A:157:ASN:OD1	2.28	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:428:PRO:HG2	1:A:481:ASP:HB3	1.93	0.50
1:B:65:VAL:HG22	1:B:159:TYR:CB	2.41	0.50
1:A:214:TYR:O	1:A:220:THR:HA	2.11	0.50
1:B:122:ARG:O	1:B:124:ASN:ND2	2.42	0.49
1:B:201:ASN:HD21	1:B:203:ASN:HB2	1.77	0.49
1:B:5:VAL:N	6:B:967:HOH:O	2.46	0.49
1:A:27:LYS:HG2	1:A:27:LYS:O	2.13	0.49
1:A:514:LEU:HD21	1:A:550:ALA:HB1	1.95	0.49
1:B:376:LEU:HD22	1:B:379:ILE:CD1	2.31	0.49
1:A:581:LYS:HA	1:A:581:LYS:HE3	1.94	0.49
1:B:506:ILE:HG12	1:B:544:ILE:HB	1.94	0.48
1:A:51:ILE:HD13	1:A:187:ILE:HD12	1.94	0.48
1:A:376:LEU:HD12	1:A:593:ILE:HD11	1.95	0.48
1:B:12:TYR:HE1	1:B:180:LYS:HD2	1.78	0.48
1:B:59:ARG:O	1:B:63:THR:HG23	2.12	0.48
1:A:55:MET:HE1	2:A:701:P65:CL1	2.51	0.48
1:A:311:ASP:OD2	1:A:561:LYS:NZ	2.42	0.48
1:B:152:LEU:O	1:B:156:LEU:HD12	2.13	0.48
1:A:166:GLY:HA3	3:A:702:NDP:PA	2.54	0.47
1:B:336:MET:CE	1:B:560:LEU:HB2	2.43	0.47
1:B:210:VAL:HG12	1:B:224:ILE:HG22	1.95	0.47
1:B:108:ASN:ND2	3:B:702:NDP:H5N	2.30	0.46
1:B:344:ASP:O	1:B:346:THR:N	2.48	0.46
1:B:553:TYR:O	1:B:556:HIS:HB2	2.15	0.46
1:A:512:CYS:SG	1:A:547:LEU:CD2	3.04	0.46
1:B:411:GLU:HB2	6:B:869:HOH:O	2.15	0.45
1:A:105:GLY:CA	1:A:169:VAL:HG21	2.46	0.45
1:B:582:ASN:HB3	6:B:803:HOH:O	2.15	0.45
1:B:171:GLN:HG3	1:B:175:GLU:OE2	2.16	0.45
1:B:43:LYS:HE3	1:B:43:LYS:HB2	1.58	0.45
1:A:354:PHE:CE2	1:B:506:ILE:HG13	2.51	0.45
1:A:355:GLY:HA2	1:A:547:LEU:O	2.16	0.45
1:B:214:TYR:O	1:B:220:THR:HA	2.17	0.45
1:B:114:LYS:HD2	1:B:115:LYS:N	2.33	0.44
1:A:318:LEU:HD11	1:A:326:TYR:OH	2.16	0.44
1:A:35:TYR:CZ	1:A:38:ARG:HD2	2.53	0.44
1:B:106:ARG:HE	3:B:702:NDP:P2B	2.40	0.44
1:B:312:PHE:CE1	1:B:564:LEU:HD23	2.52	0.44
1:B:109:TRP:CE2	1:B:117:LYS:HD2	2.53	0.44
1:B:165:GLY:HA3	1:B:170:TYR:CE2	2.53	0.43
1:A:105:GLY:N	1:A:169:VAL:HG21	2.33	0.43
1:A:29:ASN:OD1	1:A:373:LYS:NZ	2.51	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:312:PHE:HB2	1:A:316:ASN:ND2	2.34	0.43
1:B:344:ASP:O	1:B:347:GLY:N	2.52	0.43
1:B:210:VAL:O	1:B:210:VAL:CG2	2.67	0.43
1:A:392:ASN:OD1	1:A:394:ASN:HB2	2.19	0.43
1:A:312:PHE:HB2	1:A:316:ASN:HD22	1.84	0.43
1:B:470:ARG:NE	6:B:974:HOH:O	2.52	0.42
1:A:109:TRP:CH2	1:A:117:LYS:HD2	2.54	0.42
1:B:58:PHE:HE1	1:B:164:LEU:HD22	1.84	0.42
1:B:314:ILE:HB	1:B:565:ASN:HB3	2.02	0.42
1:A:77:ARG:N	6:A:959:HOH:O	2.53	0.41
1:A:70:TYR:HA	1:A:73:LEU:HD12	2.01	0.41
1:A:353:LYS:HE2	1:A:353:LYS:HB3	1.65	0.41
1:B:62:THR:HA	1:B:162:PHE:CE2	2.56	0.41
1:A:485:MET:SD	1:A:489:PRO:HD3	2.61	0.41
1:A:209:SER:OG	1:A:320:TYR:HD2	2.04	0.41
1:A:77:ARG:HG3	6:A:995:HOH:O	2.21	0.41
1:A:455:GLN:HB3	1:A:474:LEU:HD12	2.02	0.41
1:A:566:ARG:HD3	1:A:599:HIS:CG	2.56	0.41
1:A:493:LEU:CD1	1:A:493:LEU:C	2.89	0.40
1:B:157:ASN:HB3	6:B:904:HOH:O	2.21	0.40
1:B:133:LYS:HD2	1:B:133:LYS:H	1.87	0.40
1:A:286:GLU:HG2	1:B:319:LYS:HD3	2.02	0.40
1:B:108:ASN:HD21	3:B:702:NDP:H5N	1.87	0.40
4:A:703:PO4:O1	6:A:999:HOH:O	2.22	0.40
1:B:284:ASP:OD1	1:B:284:ASP:N	2.47	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	530/608 (87%)	491 (93%)	31 (6%)	8 (2%)	15	18
1	B	518/608 (85%)	474 (92%)	38 (7%)	6 (1%)	19	24
All	All	1048/1216 (86%)	965 (92%)	69 (7%)	14 (1%)	18	22

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	49	LYS
1	B	345	ARG
1	A	139	ASP
1	A	49	LYS
1	A	284	ASP
1	A	430	TYR
1	A	22	SER
1	A	26	GLY
1	A	310	ASN
1	B	119	LEU
1	A	309	PRO
1	B	309	PRO
1	B	123	ILE
1	B	429	ILE

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	501/570 (88%)	484 (97%)	17 (3%)	49	68
1	B	492/570 (86%)	475 (96%)	17 (4%)	48	67
All	All	993/1140 (87%)	959 (97%)	34 (3%)	49	68

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	VAL
1	A	49	LYS
1	A	97	LYS
1	A	132	LYS
1	A	144	ASN
1	A	341	LYS
1	A	345	ARG
1	A	348	VAL
1	A	353	LYS
1	A	376	LEU

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Mol	Chain	Res	Type
1	A	449	GLU
1	A	487	LEU
1	A	581	LYS
1	A	592	THR
1	A	593	ILE
1	A	605	ASP
1	A	606	MET
1	B	43	LYS
1	B	49	LYS
1	B	52	SER
1	B	104	MET
1	B	107	THR
1	B	114	LYS
1	B	132	LYS
1	B	133	LYS
1	B	155	LYS
1	B	196	PHE
1	B	208	ILE
1	B	231	ASN
1	B	284	ASP
1	B	344	ASP
1	B	457	LYS
1	B	564	LEU
1	B	605	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	144	ASN
1	A	294	ASN
1	A	316	ASN
1	A	394	ASN
1	A	424	ASN
1	A	554	ASN
1	B	99	GLN
1	B	171	GLN
1	B	201	ASN
1	B	203	ASN
1	B	305	ASN
1	B	316	ASN
1	B	394	ASN

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Mol	Chain	Res	Type
1	B	424	ASN

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 ⓘ

10 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	P65	A	701	-	24,24,24	3.04	5 (20%)	33,33,33	2.03	10 (30%)
3	NDP	A	702	-	52,52,52	1.43	6 (11%)	80,80,80	1.61	11 (13%)
4	PO4	A	703	-	4,4,4	0.30	0	6,6,6	0.34	0
5	BME	A	704	-	3,3,3	0.41	0	2,2,2	1.36	0
5	BME	A	705	-	3,3,3	0.53	0	2,2,2	1.07	0
5	BME	A	706	-	3,3,3	0.37	0	2,2,2	1.46	0
2	P65	B	701	-	24,24,24	3.24	4 (16%)	33,33,33	1.52	4 (12%)
3	NDP	B	702	-	52,52,52	1.37	5 (9%)	80,80,80	1.51	7 (8%)
4	PO4	B	703	-	4,4,4	0.25	0	6,6,6	0.32	0
5	BME	B	704	-	3,3,3	0.47	0	2,2,2	2.04	1 (50%)

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	P65	A	701	-	-	0/8/8/8	0/2/2/2
3	NDP	A	702	-	-	0/35/77/77	0/3/5/5
4	PO4	A	703	-	-	0/0/0/0	0/0/0/0
5	BME	A	704	-	-	0/1/1/1	0/0/0/0
5	BME	A	705	-	-	0/1/1/1	0/0/0/0
5	BME	A	706	-	-	0/1/1/1	0/0/0/0
2	P65	B	701	-	-	0/8/8/8	0/2/2/2
3	NDP	B	702	-	-	0/35/77/77	0/3/5/5
4	PO4	B	703	-	-	0/0/0/0	0/0/0/0
5	BME	B	704	-	-	0/1/1/1	0/0/0/0

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	P65	C15-CL2	-9.33	1.49	1.73
2	B	701	P65	C13-CL1	-9.17	1.50	1.73
2	A	701	P65	C15-CL2	-8.58	1.51	1.73
2	B	701	P65	C16-CL3	-8.32	1.52	1.73
2	A	701	P65	C16-CL3	-7.83	1.53	1.73
2	A	701	P65	C13-CL1	-7.50	1.54	1.73
3	A	702	NDP	O7N-C7N	6.44	1.41	1.24
3	B	702	NDP	O7N-C7N	6.37	1.41	1.24
2	A	701	P65	C6-C5	-3.60	1.38	1.41
3	A	702	NDP	C6N-C5N	3.37	1.40	1.33
3	B	702	NDP	C2A-N3A	3.17	1.38	1.32
3	A	702	NDP	C2A-N3A	3.12	1.38	1.32
3	A	702	NDP	C2N-C3N	3.11	1.40	1.34
3	B	702	NDP	C6N-C5N	2.98	1.39	1.33
3	B	702	NDP	C2N-C3N	2.65	1.39	1.34
3	B	702	NDP	C2A-N1A	2.56	1.39	1.33
3	A	702	NDP	C2A-N1A	2.43	1.38	1.33
2	A	701	P65	C1-N2	2.39	1.37	1.34
2	B	701	P65	C6-C5	-2.21	1.39	1.41
3	A	702	NDP	C7N-C3N	2.19	1.52	1.47

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	702	NDP	N3A-C2A-N1A	-10.27	120.12	128.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	NDP	N3A-C2A-N1A	-8.71	121.43	128.71
2	A	701	P65	C3-N2-C1	5.03	120.60	116.57
2	B	701	P65	C3-N2-C1	4.89	120.48	116.57
3	A	702	NDP	O4B-C1B-N9A	4.71	112.82	108.44
2	A	701	P65	C3-N4-C5	4.17	121.69	116.96
2	A	701	P65	C17-C16-CL3	3.89	124.67	118.55
3	A	702	NDP	C4A-C5A-N7A	-3.53	106.50	109.52
3	B	702	NDP	N3A-C4A-N9A	3.51	131.78	125.43
2	B	701	P65	C3-N4-C5	3.38	120.80	116.96
2	A	701	P65	N4-C3-N2	-3.31	119.90	125.55
2	A	701	P65	C14-C15-CL2	3.30	123.74	118.55
3	A	702	NDP	N3A-C4A-N9A	3.01	130.87	125.43
2	A	701	P65	NH1-C3-N2	2.97	122.32	117.19
5	B	704	BME	C1-C2-S2	-2.88	108.07	113.03
2	A	701	P65	CM1-C1-N2	2.83	120.97	116.42
3	B	702	NDP	O4B-C1B-N9A	2.72	110.97	108.44
3	A	702	NDP	O2B-C2B-C1B	-2.72	100.28	110.36
2	B	701	P65	N4-C3-N2	-2.71	120.92	125.55
3	A	702	NDP	N6A-C6A-N1A	-2.63	114.20	119.36
3	A	702	NDP	C1D-N1N-C2N	-2.42	116.91	121.02
2	B	701	P65	C10-O11-C12	2.38	123.50	117.65
2	A	701	P65	C16-C15-CL2	-2.26	114.34	120.90
3	B	702	NDP	O2A-PA-O3	2.25	115.81	105.14
2	A	701	P65	C14-C13-CL1	2.19	122.00	118.55
2	A	701	P65	C14-C13-C12	-2.19	117.64	120.94
3	A	702	NDP	O4D-C1D-N1N	2.16	112.62	108.05
3	A	702	NDP	N7A-C8A-N9A	-2.15	108.29	114.36
3	B	702	NDP	C5A-C4A-N3A	-2.12	121.08	125.70
3	B	702	NDP	PN-O3-PA	-2.10	125.51	131.68
3	A	702	NDP	C5A-C4A-N3A	-2.07	121.19	125.70
3	A	702	NDP	O2N-PN-O1N	2.06	123.74	112.21
3	B	702	NDP	C2A-N3A-C4A	2.05	119.84	114.01

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, 95th 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	538/608 (88%)	0.47	40 (7%) 14 15	26, 39, 81, 98	0
1	B	529/608 (87%)	0.67	68 (12%) 4 4	24, 43, 96, 116	0
All	All	1067/1216 (87%)	0.57	108 (10%) 8 8	24, 41, 91, 116	0

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	85	THR	9.4
1	B	136	PHE	8.1
1	B	138	GLU	7.7
1	A	31	VAL	7.3
1	B	232	LYS	7.3
1	A	26	GLY	6.9
1	A	25	GLU	6.7
1	A	24	ASN	6.5
1	B	607	ALA	6.3
1	B	231	ASN	6.2
1	B	345	ARG	6.2
1	A	608	ALA	5.7
1	B	346	THR	5.5
1	A	231	ASN	5.0
1	B	151	VAL	4.7
1	B	75	TYR	4.7
1	A	230	ASN	4.5
1	B	118	PRO	4.5
1	A	606	MET	4.4
1	B	29	ASN	4.3
1	A	29	ASN	4.3
1	A	283	ASP	4.2
1	B	80	TYR	4.1
1	A	607	ALA	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	608	ALA	4.1
1	A	345	ARG	4.1
1	B	310	ASN	4.1
1	A	35	TYR	4.0
1	A	310	ASN	4.0
1	B	116	PHE	3.8
1	B	307	ILE	3.8
1	A	200	ILE	3.7
1	B	230	ASN	3.7
1	B	62	THR	3.7
1	B	146	VAL	3.7
1	B	135	ASP	3.6
1	A	28	LYS	3.6
1	B	306	SER	3.6
1	B	9	PHE	3.6
1	B	103	VAL	3.5
1	B	11	ILE	3.4
1	B	130	THR	3.4
1	B	204	GLU	3.4
1	A	27	LYS	3.3
1	A	346	THR	3.2
1	B	309	PRO	3.1
1	A	306	SER	3.1
1	B	45	VAL	3.0
1	B	119	LEU	3.0
1	B	152	LEU	3.0
1	B	140	VAL	2.9
1	B	347	GLY	2.9
1	A	22	SER	2.9
1	B	110	GLU	2.9
1	B	8	VAL	2.9
1	B	7	ASP	2.9
1	A	3	GLU	2.8
1	B	178	LEU	2.8
1	B	58	PHE	2.8
1	A	307	ILE	2.8
1	B	109	TRP	2.8
1	B	131	LEU	2.8
1	B	126	ILE	2.8
1	A	97	LYS	2.8
1	B	157	ASN	2.7
1	B	150	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	23	LYS	2.7
1	A	205	TYR	2.7
1	A	313	GLN	2.7
1	B	134	GLU	2.7
1	B	137	ASP	2.7
1	B	102	VAL	2.7
1	B	173	PHE	2.6
1	A	4	GLN	2.6
1	B	22	SER	2.5
1	B	202	GLU	2.5
1	B	164	LEU	2.5
1	B	38	ARG	2.5
1	A	75	TYR	2.5
1	B	141	TYR	2.4
1	B	285	GLU	2.4
1	A	202	GLU	2.4
1	A	203	ASN	2.4
1	A	8	VAL	2.4
1	B	5	VAL	2.4
1	A	375	PHE	2.4
1	B	81	LEU	2.3
1	B	605	ASP	2.3
1	A	206	GLN	2.3
1	B	46	LEU	2.3
1	B	76	LYS	2.3
1	B	179	ILE	2.2
1	B	107	THR	2.2
1	A	33	ASN	2.2
1	A	604	MET	2.2
1	B	104	MET	2.2
1	B	6	CYS	2.2
1	A	308	HIS	2.1
1	A	34	ASN	2.1
1	B	111	SER	2.1
1	B	174	LEU	2.1
1	A	96	LYS	2.1
1	B	205	TYR	2.1
1	B	229	THR	2.1
1	A	6	CYS	2.0
1	B	35	TYR	2.0
1	B	59	ARG	2.0
1	B	97	LYS	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, 95th 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 ²)	Q<0.9
5	BME	A	706	4/4	0.33	10.54	77,77,77,78	0
5	BME	A	705	4/4	0.30	9.55	89,90,90,91	0
5	BME	A	704	4/4	0.19	7.35	90,92,93,94	0
5	BME	B	704	4/4	0.15	1.33	65,67,68,71	0
2	P65	A	701	23/23	0.13	0.01	26,31,53,58	0
3	NDP	A	702	48/48	0.12	-0.31	43,50,53,54	0
3	NDP	B	702	48/48	0.19	-0.32	83,88,102,104	0
2	P65	B	701	23/23	0.17	-0.48	59,66,77,79	0
4	PO4	B	703	5/5	0.12	-0.58	48,48,50,51	0
4	PO4	A	703	5/5	0.13	-1.44	61,62,62,62	0

6.5 Other polymers ⓘ

There are no such residues in this entry.