



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:04 AM UTC

PDB ID : 1E3M / pdb_00001e3m
Title : The crystal structure of E. coli MutS binding to DNA with a G:T mismatch
Authors : Lamers, M.H.; Perrakis, A.; Enzlin, J.H.; Winterwerp, H.H.K.; De Wind, N.; Sixma, T.K.
Deposited on : 2000-06-19
Resolution : 2.20 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

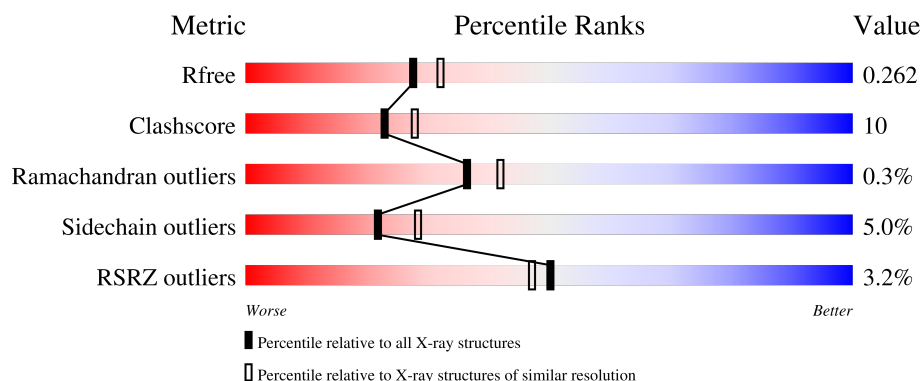
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

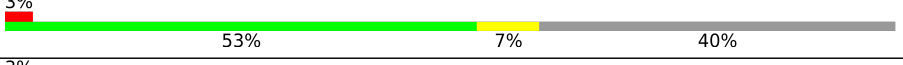
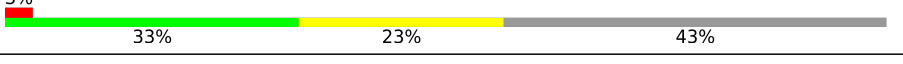
The reported resolution of this entry is 2.20 Å.

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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	800	
1	B	800	
2	E	30	
3	F	30	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13392 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 DNA MISMATCH REPAIR PROTEIN MUTS.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	788	Total	C	N	O	S	Se	0	0	0
			6207	3905	1103	1170	6	23			
1	B	754	Total	C	N	O	S	Se	0	0	0
			5964	3756	1060	1120	6	22			

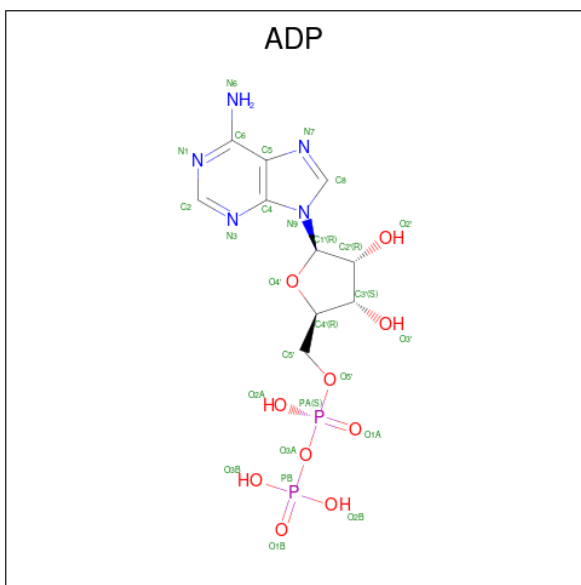
- Molecule 2 is a DNA chain called 5'-D(*AP*GP*CP*TP*GP*CP*CP*AP*GP*GP*CP*A
P*CP*CP*AP* GP*TP*GP*TP*CP*AP*GP*CP*GP*TP*CP*CP*TP*AP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	18	Total	C	N	O	P	0	0	0
			367	174	72	104	17			

- Molecule 3 is a DNA chain called 5'-D(*AP*TP*AP*GP*GP*AP*CP*GP*CP*TP*GP*A
P*CP*AP*CP* TP*GP*GP*TP*GP*CP*TP*TP*GP*GP*CP*AP*GP*CP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	17	Total	C	N	O	P	0	0	0
			347	166	62	103	16			

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Mg 2 2	0	0

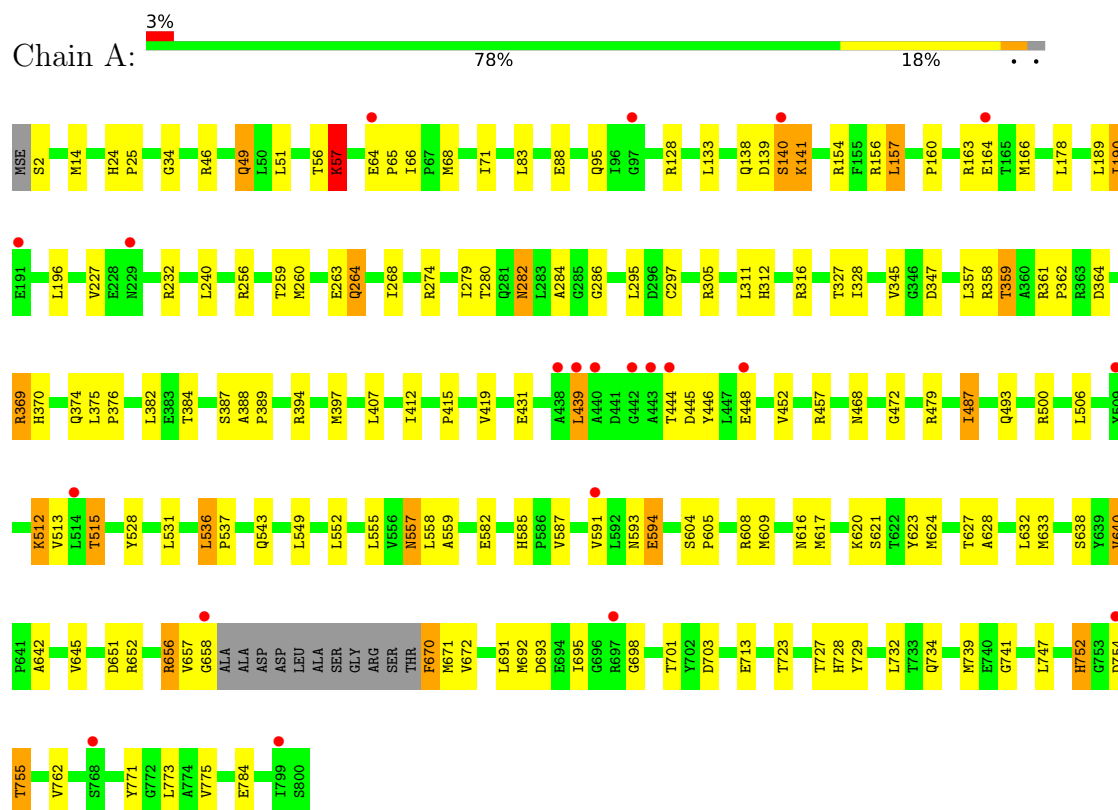
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	228	Total O 228 228	0	0
6	B	223	Total O 223 223	0	0
6	E	11	Total O 11 11	0	0
6	F	16	Total O 16 16	0	0

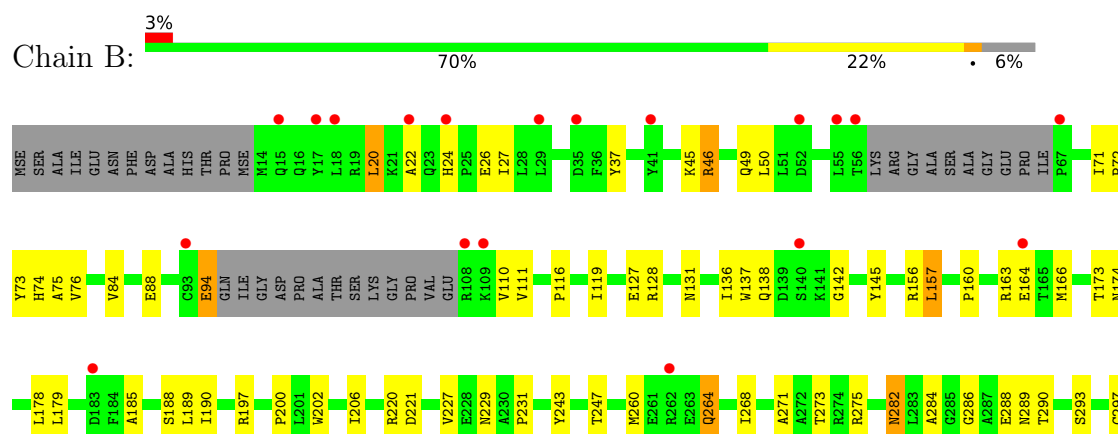
3 Residue-property plots

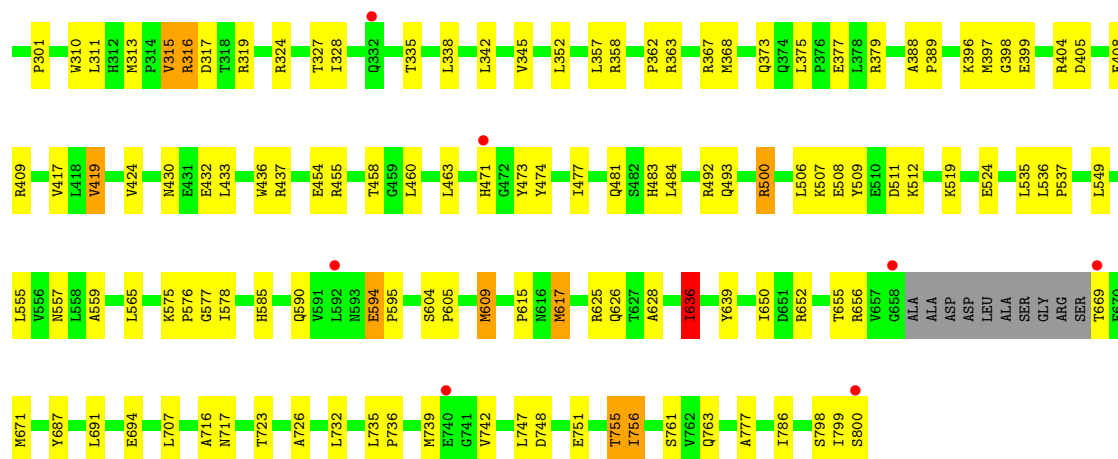
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: DNA MISMATCH REPAIR PROTEIN MUTS

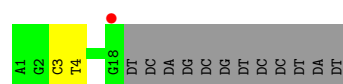


• Molecule 1: DNA MISMATCH REPAIR PROTEIN MUTS

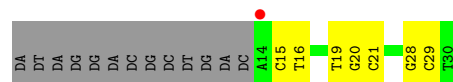
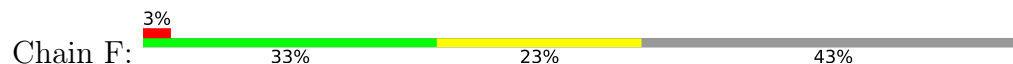




• Molecule 2: 5'-D(*AP*GP*CP*TP*GP*CP*CP*AP*GP*GP*CP*AP*CP*CP*AP* GP*TP*GP*TP*CP*AP*GP*CP*GP*TP*CP*CP*TP*AP*T)-3'



• Molecule 3: 5'-D(*AP*TP*AP*GP*GP*AP*CP*GP*CP*TP*GP*AP*CP*AP*CP* TP*GP*GP*TP*GP*CP*TP*TP*GP*GP*CP*AP*GP*CP*T)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.96Å 92.37Å 261.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.7 (20.00-2.20) 98.5 (20.00-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.228 , 0.266 0.224 , 0.262	Depositor DCC
R_{free} test set	2199 reflections (2.01%)	wwPDB-VP
Wilson B-factor (Å ²)	42.1	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 25.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.022 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13392	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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.

5 Model quality

5.1 Standard geometry

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

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	0.67	1/6290 (0.0%)	0.90	4/8475 (0.0%)
1	B	0.70	2/6041 (0.0%)	0.90	3/8136 (0.0%)
2	E	0.35	0/412	0.83	0/634
3	F	0.39	0/388	0.96	0/598
All	All	0.67	3/13131 (0.0%)	0.90	7/17843 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	609	MSE	SE-CE	-13.21	1.55	1.95
1	A	14	MSE	SE-CE	-6.38	1.76	1.95
1	B	368	MSE	SE-CE	-6.22	1.76	1.95

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	SER	N-CA-C	-7.50	102.24	111.40
1	B	206	ILE	N-CA-C	6.34	117.08	110.62
1	B	636	ILE	N-CA-C	-6.10	106.78	113.43
1	A	701	THR	N-CA-C	5.73	117.20	111.07
1	A	66	ILE	N-CA-C	5.67	113.49	107.76
1	B	142	GLY	N-CA-C	5.24	117.25	110.45
1	A	57	LYS	N-CA-C	5.00	116.67	109.07

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	6207	0	6251	122	0
1	B	5964	0	6018	139	0
2	E	367	0	202	4	0
3	F	347	0	194	4	0
4	A	27	0	12	0	0
5	A	2	0	0	0	0
6	A	228	0	0	8	0
6	B	223	0	0	12	0
6	E	11	0	0	0	0
6	F	16	0	0	0	0
All	All	13392	0	12677	266	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 10.

All (266) 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:133:LEU:HG	1:A:240:LEU:HD11	1.40	1.02
1:A:727:THR:HG21	1:A:732:LEU:HD12	1.44	0.98
1:A:141:LYS:H	1:A:141:LYS:HD2	1.31	0.93
1:A:34:GLY:H	1:A:95:GLN:HE22	1.15	0.92
1:B:84:VAL:HG13	1:B:116:PRO:HA	1.52	0.92
1:A:727:THR:HG22	1:A:729:TYR:H	1.35	0.89
1:A:282:ASN:HD22	1:A:284:ALA:H	1.26	0.84
1:B:24:HIS:CD2	1:B:110:VAL:HG21	2.13	0.84
1:B:37:TYR:CE2	1:B:76:VAL:HG21	2.14	0.83
1:B:37:TYR:HE2	1:B:76:VAL:HG21	1.44	0.82
1:B:160:PRO:HG2	1:B:166:MSE:SE	2.29	0.82
1:A:305:ARG:NH2	1:A:347:ASP:OD2	2.13	0.81
1:A:46:ARG:NH2	1:A:88:GLU:OE1	2.15	0.79
1:A:397:MSE:HE1	1:A:552:LEU:HD22	1.67	0.77
1:B:311:LEU:HD23	1:B:636:ILE:HD12	1.66	0.77
1:B:735:LEU:HD22	1:B:739:MSE:HE2	1.67	0.76
1:B:327:THR:HG21	1:B:555:LEU:HD13	1.67	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:LYS:HD3	1:A:232:ARG:HH12	1.52	0.75
1:B:264:GLN:H	1:B:264:GLN:HE21	1.32	0.75
1:B:656:ARG:O	1:B:656:ARG:HG3	1.87	0.74
1:B:669:THR:HG22	1:B:671:MSE:H	1.52	0.74
1:A:439:LEU:HD22	1:A:439:LEU:H	1.50	0.74
1:A:34:GLY:H	1:A:95:GLN:NE2	1.84	0.74
1:B:290:THR:HG23	1:B:293:SER:H	1.50	0.73
1:B:282:ASN:HD22	1:B:284:ALA:H	1.37	0.73
1:B:799:ILE:O	1:B:799:ILE:HG22	1.90	0.71
1:B:20:LEU:HD21	1:B:110:VAL:HG23	1.72	0.70
1:B:157:LEU:HD23	1:B:157:LEU:O	1.91	0.70
1:B:94:GLU:OE1	1:B:94:GLU:HA	1.90	0.69
1:B:290:THR:HG22	1:B:293:SER:HB3	1.73	0.69
1:A:752:HIS:ND1	1:A:752:HIS:N	2.40	0.69
1:B:282:ASN:HD21	1:B:286:GLY:H	1.39	0.69
1:B:575:LYS:HB3	1:B:576:PRO:HD2	1.75	0.68
1:A:754:ASP:O	1:A:755:THR:HB	1.94	0.68
1:B:288:GLU:HG2	1:B:289:ASN:ND2	2.10	0.66
1:A:616:ASN:O	1:A:617:MSE:HB2	1.95	0.66
1:B:128:ARG:NH2	6:B:2028:HOH:O	2.29	0.66
1:A:64:GLU:O	6:A:2029:HOH:O	2.14	0.65
1:B:751:GLU:OE2	6:B:2205:HOH:O	2.14	0.65
1:B:290:THR:CG2	1:B:293:SER:H	2.10	0.65
1:A:632:LEU:HD23	1:A:632:LEU:C	2.22	0.64
1:B:430:ASN:HD22	1:B:433:LEU:H	1.44	0.64
1:B:163:ARG:HH12	1:B:188:SER:HB2	1.63	0.64
1:B:282:ASN:HD22	1:B:284:ALA:N	1.95	0.63
1:A:327:THR:HG21	1:A:555:LEU:HD13	1.80	0.63
1:B:639:TYR:OH	6:B:2180:HOH:O	2.13	0.63
1:A:282:ASN:HD22	1:A:284:ALA:N	1.93	0.62
1:B:471:HIS:CE1	1:B:493:GLN:HB2	2.34	0.62
1:A:593:ASN:O	1:A:594:GLU:HB3	1.98	0.62
1:B:316:ARG:HD3	1:B:650:ILE:O	2.00	0.62
1:B:474:TYR:CE1	1:B:500:ARG:HD2	2.33	0.62
1:B:26:GLU:HG3	1:B:27:ILE:HG13	1.80	0.62
1:B:799:ILE:O	1:B:799:ILE:CG2	2.48	0.62
1:B:163:ARG:HG3	1:B:189:LEU:HD21	1.82	0.62
1:A:128:ARG:NE	6:A:2048:HOH:O	2.16	0.61
1:B:282:ASN:ND2	1:B:286:GLY:H	1.99	0.60
1:B:362:PRO:HG3	1:B:419:VAL:HG13	1.83	0.60
1:B:268:ILE:HB	1:B:652:ARG:HG2	1.82	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:ARG:HA	1:A:397:MSE:HE3	1.84	0.59
1:B:483:HIS:CE1	1:B:484:LEU:HG	2.37	0.59
1:B:328:ILE:HG23	1:B:559:ALA:HA	1.82	0.59
1:A:133:LEU:CG	1:A:240:LEU:HD11	2.25	0.59
1:A:621:SER:OG	1:A:693:ASP:OD2	2.20	0.59
1:B:157:LEU:HD23	1:B:157:LEU:C	2.26	0.59
1:B:297:CYS:H	1:B:557:ASN:HD21	1.51	0.59
1:B:363:ARG:O	1:B:367:ARG:HG3	2.04	0.58
1:B:46:ARG:NH1	1:B:88:GLU:OE2	2.37	0.58
1:A:698:GLY:HA3	1:A:703:ASP:HB3	1.86	0.58
1:A:295:LEU:HD11	1:A:633:MSE:HE3	1.85	0.58
1:B:345:VAL:HG11	1:B:549:LEU:HD13	1.86	0.58
1:B:735:LEU:HD22	1:B:739:MSE:CE	2.35	0.57
1:A:536:LEU:N	1:A:537:PRO:HD2	2.20	0.57
1:A:604:SER:HB2	1:A:605:PRO:CD	2.34	0.57
1:A:157:LEU:HD23	1:A:157:LEU:C	2.30	0.57
1:B:628:ALA:HB2	1:B:691:LEU:HD11	1.86	0.56
1:A:274:ARG:HG3	1:A:279:ILE:HD12	1.87	0.56
1:A:727:THR:HG23	6:A:2201:HOH:O	2.05	0.56
1:B:437:ARG:NH2	1:B:524:GLU:OE2	2.33	0.56
1:B:375:LEU:O	1:B:379:ARG:HG3	2.06	0.56
3:F:28:DG:H2''	3:F:29:DC:H5''	1.88	0.56
1:B:739:MSE:HE3	1:B:742:VAL:HG21	1.88	0.56
1:B:282:ASN:ND2	1:B:284:ALA:H	2.02	0.55
1:B:313:MSE:HE2	6:B:2088:HOH:O	2.06	0.55
1:A:412:ILE:HD11	1:A:415:PRO:HA	1.88	0.55
1:B:655:THR:O	1:B:655:THR:OG1	2.15	0.55
1:B:454:GLU:O	1:B:458:THR:HG23	2.07	0.54
1:A:24:HIS:N	1:A:25:PRO:HD3	2.23	0.53
1:A:609:MSE:HE2	1:A:723:THR:HG21	1.90	0.53
1:A:190:ILE:HG23	1:A:196:LEU:HD11	1.90	0.53
1:A:282:ASN:ND2	1:A:284:ALA:H	2.02	0.53
1:B:310:TRP:CB	1:B:636:ILE:HD11	2.38	0.53
1:B:73:TYR:O	1:B:76:VAL:HG23	2.09	0.53
1:B:227:VAL:HG12	1:B:260:MSE:HB2	1.90	0.53
1:B:22:ALA:HB3	6:B:2003:HOH:O	2.08	0.53
1:A:282:ASN:HD21	1:A:286:GLY:H	1.57	0.52
1:A:388:ALA:HB3	1:A:389:PRO:HD3	1.89	0.52
1:B:310:TRP:HB3	1:B:636:ILE:HD11	1.91	0.52
2:E:3:DC:H2''	2:E:4:DT:C5'	2.39	0.52
1:B:229:ASN:C	1:B:231:PRO:HD3	2.34	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:632:LEU:HD23	1:A:632:LEU:O	2.10	0.52
1:A:227:VAL:HG12	1:A:260:MSE:HB2	1.92	0.52
1:B:609:MSE:HE2	1:B:723:THR:HG21	1.92	0.52
1:B:282:ASN:HD22	1:B:282:ASN:C	2.17	0.51
1:A:670:PHE:N	1:A:670:PHE:CD2	2.79	0.51
1:A:446:TYR:C	1:A:446:TYR:CD1	2.89	0.51
1:A:692:MSE:HE3	1:A:695:ILE:HD13	1.93	0.51
1:B:271:ALA:HB1	1:B:275:ARG:HH21	1.75	0.51
1:B:273:THR:HG23	1:B:655:THR:HG23	1.93	0.51
1:B:474:TYR:HA	1:B:506:LEU:HD21	1.92	0.51
1:A:591:VAL:O	1:A:591:VAL:HG23	2.10	0.51
1:A:670:PHE:N	1:A:670:PHE:HD2	2.09	0.51
1:B:430:ASN:HD21	1:B:432:GLU:HB2	1.76	0.50
1:B:575:LYS:CB	1:B:576:PRO:HD2	2.41	0.50
3:F:15:DC:H2''	3:F:16:DT:H5'	1.93	0.50
1:B:72:PRO:HG2	1:B:75:ALA:HB3	1.93	0.50
1:A:46:ARG:NH2	1:A:88:GLU:CD	2.70	0.50
1:A:46:ARG:NE	6:A:2021:HOH:O	2.22	0.50
1:A:582:GLU:O	1:A:642:ALA:HA	2.12	0.50
1:A:394:ARG:O	1:A:397:MSE:HG2	2.10	0.50
1:A:263:GLU:HB3	1:A:268:ILE:HD11	1.93	0.50
1:B:324:ARG:O	1:B:328:ILE:HG13	2.11	0.50
1:A:628:ALA:HB2	1:A:691:LEU:HD11	1.92	0.50
1:B:24:HIS:HD2	1:B:110:VAL:HG21	1.74	0.50
2:E:3:DC:H2''	2:E:4:DT:H5''	1.93	0.50
1:A:160:PRO:HG2	1:A:166:MSE:SE	2.62	0.50
1:B:145:TYR:CG	1:B:166:MSE:HE1	2.47	0.50
1:A:56:THR:OG1	1:A:57:LYS:N	2.45	0.49
1:A:439:LEU:HB3	1:A:512:LYS:HZ1	1.76	0.49
1:A:620:LYS:HG2	1:A:747:LEU:HD13	1.94	0.49
1:A:297:CYS:H	1:A:557:ASN:ND2	2.09	0.49
1:A:357:LEU:O	1:A:358:ARG:HB2	2.10	0.49
1:A:359:THR:HG22	1:A:359:THR:O	2.12	0.49
1:A:640:VAL:HG11	1:A:645:VAL:HG21	1.94	0.49
1:B:46:ARG:O	1:B:50:LEU:HG	2.11	0.49
1:B:748:ASP:OD1	1:B:748:ASP:C	2.54	0.49
1:B:509:TYR:CD2	1:B:512:LYS:HD2	2.48	0.49
1:A:305:ARG:HH22	1:A:347:ASP:CG	2.18	0.49
1:A:623:TYR:CE1	1:A:762:VAL:HG21	2.48	0.49
1:A:370:HIS:O	1:A:374:GLN:HG2	2.12	0.49
1:A:468:ASN:O	1:A:472:GLY:N	2.36	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:ASP:OD1	1:B:319:ARG:HB2	2.11	0.48
1:B:220:ARG:HG3	1:B:221:ASP:N	2.28	0.48
1:B:594:GLU:HG3	1:B:595:PRO:HD2	1.94	0.48
1:B:138:GLN:HE22	1:B:185:ALA:HB3	1.77	0.48
1:A:156:ARG:HD3	1:A:259:THR:HB	1.96	0.48
1:A:604:SER:HB2	1:A:605:PRO:HD2	1.96	0.48
1:A:608:ARG:HG3	1:A:608:ARG:HH11	1.78	0.48
1:B:405:ASP:OD2	1:B:409:ARG:NH1	2.47	0.48
1:A:656:ARG:NH1	1:A:658:GLY:HA2	2.28	0.48
1:A:670:PHE:CG	1:A:671:MSE:N	2.82	0.48
1:A:256:ARG:HH21	1:A:543:GLN:NE2	2.12	0.48
1:B:652:ARG:NH1	6:B:2182:HOH:O	2.43	0.47
1:A:68:MSE:HE1	3:F:21:DC:H1'	1.96	0.47
1:A:154:ARG:NH2	6:A:2059:HOH:O	2.46	0.47
1:B:362:PRO:HD2	1:B:417:VAL:O	2.14	0.47
1:A:282:ASN:ND2	1:A:286:GLY:H	2.13	0.47
1:B:271:ALA:CB	1:B:275:ARG:HH21	2.27	0.47
1:B:138:GLN:NE2	1:B:185:ALA:HB3	2.29	0.47
1:B:625:ARG:NE	6:B:2177:HOH:O	2.44	0.47
1:B:164:GLU:CD	1:B:164:GLU:H	2.23	0.47
1:A:713:GLU:HG3	1:A:739:MSE:SE	2.65	0.47
1:A:512:LYS:HA	1:A:515:THR:OG1	2.14	0.46
1:A:46:ARG:NH2	6:A:2022:HOH:O	2.46	0.46
1:A:141:LYS:HD2	1:A:141:LYS:N	2.15	0.46
1:A:133:LEU:HG	1:A:240:LEU:CD1	2.29	0.46
1:B:474:TYR:CA	1:B:506:LEU:HD21	2.44	0.46
1:B:282:ASN:ND2	1:B:284:ALA:N	2.61	0.46
1:B:474:TYR:CD1	1:B:500:ARG:HD2	2.50	0.46
1:B:388:ALA:HB3	1:B:389:PRO:HD3	1.98	0.46
1:A:624:MSE:O	1:A:627:THR:HB	2.16	0.46
1:A:493:GLN:NE2	1:A:500:ARG:HH11	2.14	0.45
1:A:608:ARG:NH1	1:A:741:GLY:HA3	2.30	0.45
1:A:282:ASN:HD22	1:A:282:ASN:C	2.24	0.45
1:B:373:GLN:O	6:B:2113:HOH:O	2.21	0.45
1:B:798:SER:C	1:B:800:SER:H	2.24	0.45
1:A:369:ARG:HG3	1:A:407:LEU:HB3	1.97	0.45
1:B:243:TYR:O	1:B:247:THR:HG23	2.17	0.45
1:B:460:LEU:HD22	1:B:481:GLN:HB3	1.98	0.45
1:B:71:ILE:O	1:B:71:ILE:HG13	2.16	0.45
1:A:141:LYS:HB2	1:A:232:ARG:NH2	2.31	0.45
1:A:656:ARG:HH12	1:A:658:GLY:HA2	1.81	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:396:LYS:NZ	6:B:2119:HOH:O	2.49	0.45
1:A:651:ASP:OD1	1:A:652:ARG:N	2.50	0.45
1:A:369:ARG:HG3	1:A:407:LEU:CB	2.46	0.45
1:A:448:GLU:O	1:A:452:VAL:HG23	2.18	0.44
1:A:141:LYS:H	1:A:141:LYS:CD	2.05	0.44
1:B:111:VAL:O	1:B:111:VAL:HG22	2.16	0.44
1:B:388:ALA:N	1:B:389:PRO:CD	2.80	0.44
1:A:49:GLN:HE21	1:A:49:GLN:HB3	1.65	0.44
1:A:280:THR:OG1	1:A:312:HIS:HE1	2.01	0.44
1:A:670:PHE:CE2	1:A:672:VAL:HG23	2.52	0.44
1:B:404:ARG:O	1:B:408:GLU:HG3	2.18	0.44
1:A:2:SER:N	6:A:2001:HOH:O	2.50	0.44
1:A:51:LEU:HD21	1:A:83:LEU:HG	1.98	0.44
1:B:507:LYS:HD2	1:B:507:LYS:HA	1.66	0.44
1:A:632:LEU:C	1:A:632:LEU:CD2	2.90	0.44
1:A:773:LEU:CD2	1:B:707:LEU:HG	2.47	0.44
1:A:419:VAL:HG21	1:A:528:TYR:CE2	2.53	0.44
1:B:315:VAL:HG23	6:B:2073:HOH:O	2.18	0.44
1:B:200:PRO:HB2	1:B:202:TRP:CD1	2.52	0.43
1:B:173:THR:O	1:B:174:ASN:C	2.61	0.43
1:B:377:GLU:OE1	1:B:377:GLU:HA	2.18	0.43
1:B:536:LEU:N	1:B:537:PRO:CD	2.80	0.43
1:B:716:ALA:C	1:B:717:ASN:HD22	2.26	0.43
1:B:777:ALA:HA	1:B:786:ILE:HD11	2.00	0.43
1:A:657:VAL:O	1:A:657:VAL:HG12	2.19	0.43
1:B:310:TRP:HB3	1:B:636:ILE:CG1	2.49	0.43
1:A:457:ARG:NH2	6:A:2148:HOH:O	2.50	0.43
1:B:179:LEU:HD23	1:B:197:ARG:HB2	2.00	0.43
1:A:139:ASP:HB2	1:A:232:ARG:HH11	1.83	0.43
1:A:274:ARG:HE	1:A:312:HIS:HD2	1.64	0.43
1:A:487:ILE:HD13	1:A:487:ILE:HA	1.81	0.43
1:B:282:ASN:ND2	1:B:282:ASN:C	2.77	0.43
1:B:342:LEU:HD23	1:B:342:LEU:HA	1.93	0.43
1:B:615:PRO:HB2	1:B:617:MSE:HG2	1.99	0.43
1:A:328:ILE:HG23	1:A:559:ALA:HA	2.01	0.43
1:A:375:LEU:HB2	1:A:376:PRO:HD3	2.01	0.43
1:B:436:TRP:CH2	1:B:519:LYS:HD2	2.54	0.43
1:A:139:ASP:HB2	1:A:232:ARG:NH1	2.34	0.42
1:B:375:LEU:HD22	1:B:397:MSE:HG2	2.00	0.42
1:B:650:ILE:HA	1:B:687:TYR:O	2.19	0.42
1:A:163:ARG:HG2	1:A:189:LEU:HD21	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:TRP:CZ2	1:B:519:LYS:HD2	2.54	0.42
1:A:138:GLN:CG	1:A:139:ASP:N	2.81	0.42
1:A:728:HIS:ND1	1:A:728:HIS:N	2.67	0.42
1:B:747:LEU:HD23	1:B:761:SER:O	2.20	0.42
1:A:345:VAL:HG11	1:A:549:LEU:HD13	2.02	0.42
1:A:439:LEU:HB3	1:A:512:LYS:NZ	2.35	0.42
1:A:585:HIS:CE1	1:A:587:VAL:HB	2.55	0.42
1:B:508:GLU:O	1:B:511:ASP:HB2	2.20	0.42
1:B:20:LEU:CD2	1:B:110:VAL:HG23	2.45	0.42
1:B:131:ASN:ND2	6:B:2035:HOH:O	2.50	0.42
1:A:558:LEU:HD13	1:A:638:SER:HB2	2.02	0.42
1:B:473:TYR:O	1:B:507:LYS:NZ	2.52	0.42
1:A:361:ARG:O	1:A:364:ASP:HB2	2.20	0.41
2:E:3:DC:H2''	2:E:4:DT:H5'	2.03	0.41
1:A:493:GLN:HE22	1:A:500:ARG:HH11	1.68	0.41
1:B:358:ARG:NE	6:B:2145:HOH:O	2.50	0.41
1:A:382:LEU:HD11	1:A:552:LEU:HD21	2.01	0.41
1:A:617:MSE:SE	1:B:671:MSE:HG3	2.71	0.41
1:B:585:HIS:HA	1:B:626:GLN:HB2	2.02	0.41
1:A:362:PRO:HG3	1:A:419:VAL:HG22	2.03	0.41
1:B:290:THR:HG22	1:B:293:SER:CB	2.45	0.41
1:B:310:TRP:HB2	1:B:636:ILE:HD11	2.03	0.41
1:B:669:THR:CG2	1:B:671:MSE:HB2	2.50	0.41
2:E:3:DC:C2'	2:E:4:DT:H5''	2.50	0.41
1:A:604:SER:CB	1:A:605:PRO:CD	2.98	0.41
1:B:127:GLU:HG3	1:B:301:PRO:HB3	2.02	0.41
1:B:398:GLY:C	1:B:399:GLU:HG3	2.45	0.41
3:F:19:DT:H2'	3:F:20:DG:C8	2.56	0.41
1:A:64:GLU:HA	1:A:65:PRO:HD2	1.81	0.41
1:B:84:VAL:HG11	1:B:119:ILE:HD13	2.01	0.41
1:B:565:LEU:HD22	1:B:590:GLN:NE2	2.36	0.41
1:B:735:LEU:HB2	1:B:736:PRO:HD3	2.03	0.41
1:B:755:THR:HB	1:B:756:ILE:H	1.58	0.41
1:B:604:SER:HB2	1:B:605:PRO:HD2	2.03	0.40
1:A:359:THR:O	1:A:359:THR:CG2	2.69	0.40
1:A:771:TYR:O	1:A:775:VAL:HG23	2.21	0.40
1:A:264:GLN:O	1:A:316:ARG:HD3	2.22	0.40
1:B:37:TYR:CD2	1:B:76:VAL:HG21	2.54	0.40
1:B:455:ARG:HD2	1:B:463:LEU:O	2.22	0.40
1:B:577:GLY:C	1:B:578:ILE:HG13	2.46	0.40
1:B:655:THR:HG22	1:B:691:LEU:HD12	2.02	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:624:MSE:HE1	1:A:693:ASP:CG	2.46	0.40
1:B:72:PRO:HB2	1:B:74:HIS:ND1	2.37	0.40
1:B:694:GLU:N	1:B:726:ALA:O	2.46	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	784/800 (98%)	753 (96%)	29 (4%)	2 (0%)	36	42
1	B	746/800 (93%)	718 (96%)	25 (3%)	3 (0%)	30	34
All	All	1530/1600 (96%)	1471 (96%)	54 (4%)	5 (0%)	36	42

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	594	GLU
1	A	755	THR
1	B	46	ARG
1	B	45	LYS
1	B	617	MSE

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	656/640 (102%)	621 (95%)	35 (5%)	20	26
1	B	631/640 (99%)	602 (95%)	29 (5%)	24	32
All	All	1287/1280 (100%)	1223 (95%)	64 (5%)	22	28

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	57	LYS
1	A	71	ILE
1	A	140	SER
1	A	141	LYS
1	A	157	LEU
1	A	164	GLU
1	A	178	LEU
1	A	190	ILE
1	A	264	GLN
1	A	282	ASN
1	A	311	LEU
1	A	359	THR
1	A	369	ARG
1	A	384	THR
1	A	387	SER
1	A	431	GLU
1	A	439	LEU
1	A	444	THR
1	A	445	ASP
1	A	479	ARG
1	A	487	ILE
1	A	506	LEU
1	A	512	LYS
1	A	513	VAL
1	A	515	THR
1	A	531	LEU
1	A	536	LEU
1	A	557	ASN
1	A	640	VAL
1	A	656	ARG
1	A	670	PHE
1	A	734	GLN
1	A	752	HIS
1	A	784	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	20	LEU
1	B	49	GLN
1	B	94	GLU
1	B	136	ILE
1	B	137	TRP
1	B	156	ARG
1	B	157	LEU
1	B	178	LEU
1	B	190	ILE
1	B	264	GLN
1	B	282	ASN
1	B	315	VAL
1	B	316	ARG
1	B	335	THR
1	B	338	LEU
1	B	352	LEU
1	B	357	LEU
1	B	419	VAL
1	B	424	VAL
1	B	477	ILE
1	B	492	ARG
1	B	500	ARG
1	B	535	LEU
1	B	594	GLU
1	B	636	ILE
1	B	732	LEU
1	B	755	THR
1	B	756	ILE
1	B	763	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	86	GLN
1	A	95	GLN
1	A	242	GLN
1	A	248	GLN
1	A	281	GLN
1	A	282	ASN
1	A	312	HIS
1	A	488	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	493	GLN
1	A	543	GLN
1	A	557	ASN
1	A	590	GLN
1	A	602	ASN
1	A	606	GLN
1	A	643	GLN
1	A	683	ASN
1	A	714	ASN
1	A	734	GLN
1	B	24	HIS
1	B	131	ASN
1	B	138	GLN
1	B	211	GLN
1	B	264	GLN
1	B	282	ASN
1	B	430	ASN
1	B	488	ASN
1	B	543	GLN
1	B	557	ASN
1	B	590	GLN
1	B	717	ASN
1	B	791	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	$\# Z > 2$
4	ADP	A	1801	5	28,29,29	1.09	3 (10%)	43,45,45	1.83	10 (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
4	ADP	A	1801	5	-	0/16/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1801	ADP	C2-N1	2.47	1.38	1.33
4	A	1801	ADP	C2-N3	2.37	1.38	1.33
4	A	1801	ADP	C8-N7	2.10	1.35	1.31

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1801	ADP	N3-C2-N1	-5.85	119.73	128.58
4	A	1801	ADP	C5-C4-N3	-4.59	120.40	126.72
4	A	1801	ADP	N9-C8-N7	-3.71	108.67	113.94
4	A	1801	ADP	C2-N3-C4	3.65	120.74	111.83
4	A	1801	ADP	C5-N7-C8	3.32	108.66	103.45
4	A	1801	ADP	N3-C4-N9	2.88	132.07	127.17
4	A	1801	ADP	C4-C5-N7	-2.37	107.87	110.58
4	A	1801	ADP	C2'-C1'-N9	-2.36	107.44	113.30
4	A	1801	ADP	O2B-PB-O3A	2.31	112.39	104.64
4	A	1801	ADP	C3'-C2'-C1'	2.05	105.33	101.46

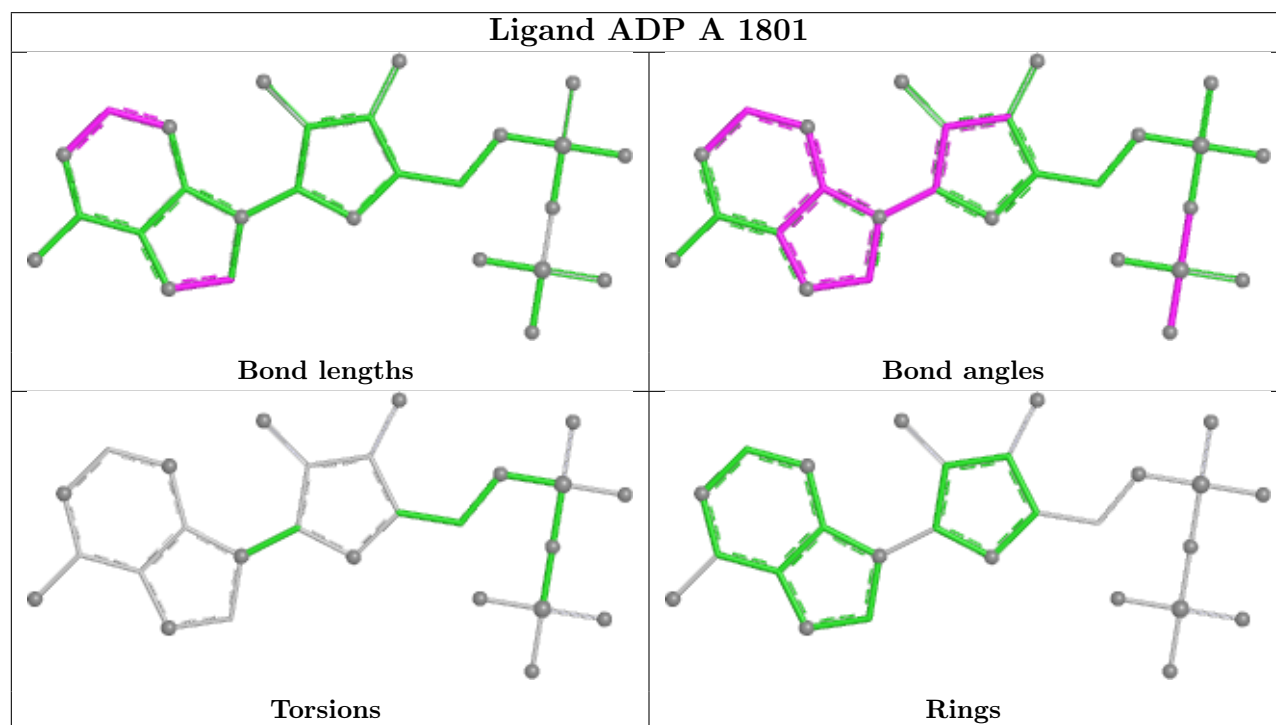
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 than 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 ⓘ

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	765/800 (95%)	0.20	21 (2%) 56 53	30, 44, 61, 82	0
1	B	732/800 (91%)	0.24	26 (3%) 46 43	26, 44, 66, 83	0
2	E	18/30 (60%)	0.08	1 (5%) 30 27	39, 50, 95, 101	0
3	F	17/30 (56%)	-0.03	1 (5%) 28 25	38, 50, 83, 92	0
All	All	1532/1660 (92%)	0.21	49 (3%) 50 47	26, 44, 65, 101	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	800	SER	4.4
1	B	41	TYR	4.0
1	A	444	THR	3.6
1	B	56	THR	3.2
1	A	439	LEU	3.2
1	B	183	ASP	3.0
1	B	55	LEU	3.0
1	A	509	TYR	2.9
1	A	754	ASP	2.9
1	A	658	GLY	2.7
1	B	669	THR	2.7
2	E	18	DG	2.7
3	F	14	DA	2.7
1	A	440	ALA	2.6
1	B	108	ARG	2.6
1	B	67	PRO	2.6
1	B	592	LEU	2.6
1	A	591	VAL	2.6
1	A	442	GLY	2.4
1	A	191	GLU	2.4
1	A	697	ARG	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	64	GLU	2.3
1	B	658	GLY	2.3
1	B	24	HIS	2.3
1	B	93	CYS	2.3
1	B	471	HIS	2.3
1	B	29	LEU	2.3
1	B	17	TYR	2.3
1	B	22	ALA	2.3
1	B	109	LYS	2.3
1	A	438	ALA	2.2
1	B	140	SER	2.2
1	A	164	GLU	2.2
1	B	740	GLU	2.2
1	B	18	LEU	2.2
1	A	768	SER	2.2
1	A	229	ASN	2.2
1	A	448	GLU	2.2
1	B	35	ASP	2.1
1	A	97	GLY	2.1
1	A	140	SER	2.1
1	B	332	GLN	2.1
1	A	443	ALA	2.0
1	B	164	GLU	2.0
1	A	514	LEU	2.0
1	B	52	ASP	2.0
1	A	799	ILE	2.0
1	B	262	ARG	2.0
1	B	15	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 oligosaccharides in this entry.

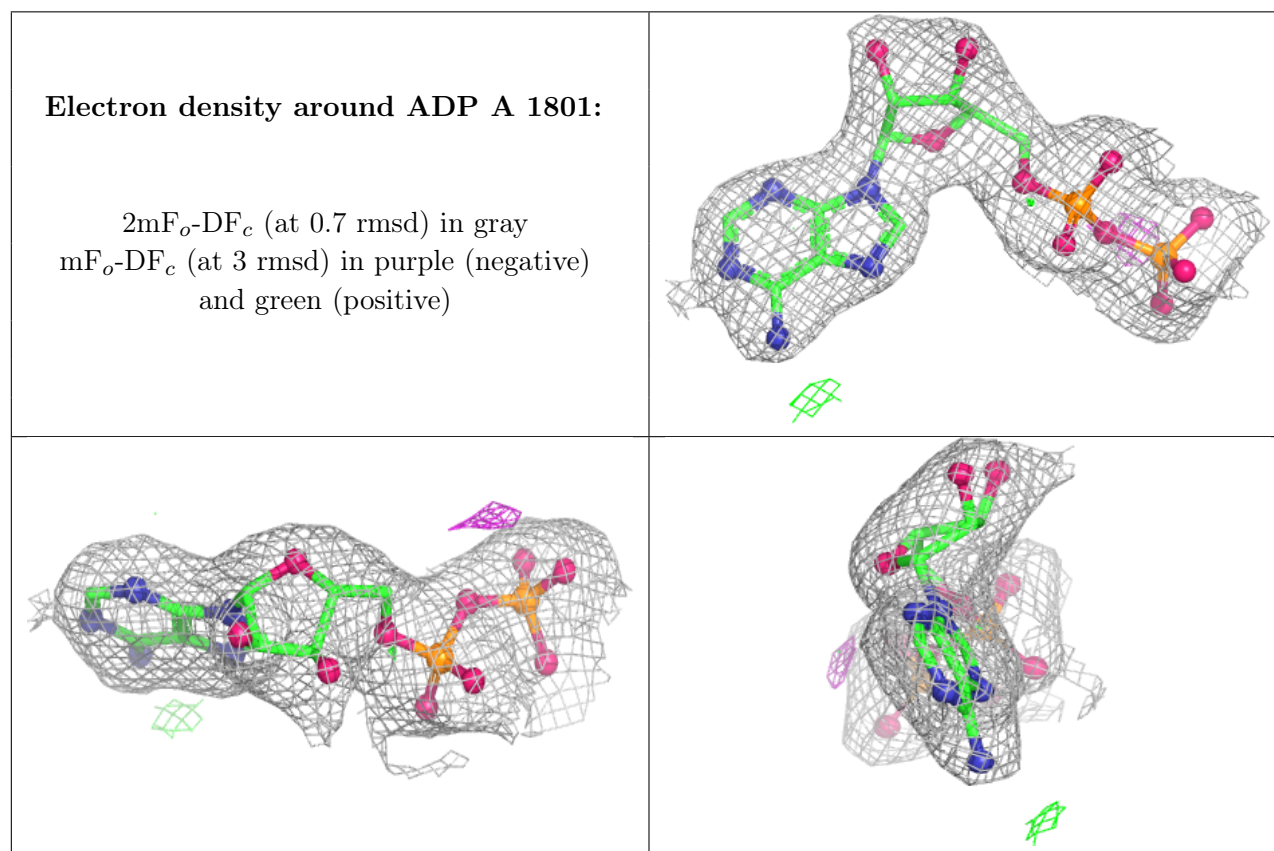
6.4 Ligands [i](#)

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,

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	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	A	1011	1/1	0.90	0.11	57,57,57,57	0
5	MG	A	1001	1/1	0.96	0.04	40,40,40,40	0
4	ADP	A	1801	27/27	0.97	0.07	40,48,49,52	0

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.