



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:56 PM GMT

PDB ID : 4GOO  
Title : Crystal Structure of E. coli DNA Adenine Methyltransferase in Complex with Benzothiophene Aza-SAM  
Authors : Harmer, J.E.; Roach, P.L.  
Deposited on : 2012-08-20  
Resolution : 2.70 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](mailto: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.

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

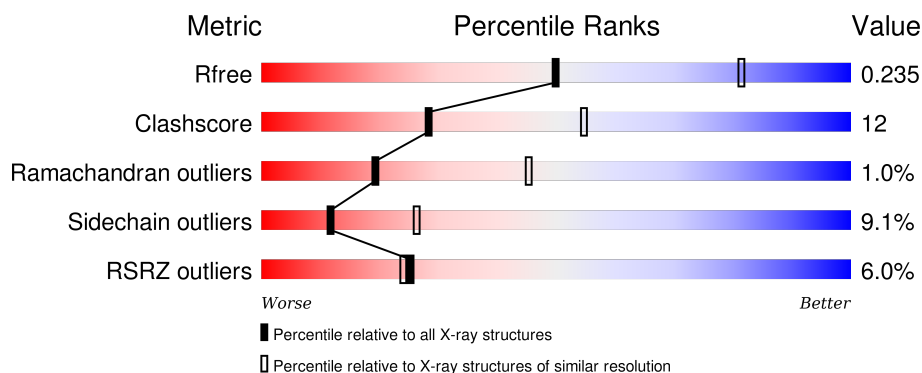
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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. 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	D	278	<div> <div>4%</div> <div>67% 17% • 13%</div> </div>
1	E	278	<div> <div>4%</div> <div>62% 22% • • 12%</div> </div>
1	F	278	<div> <div>8%</div> <div>66% 18% • • 12%</div> </div>

## 2 Entry composition [i](#)

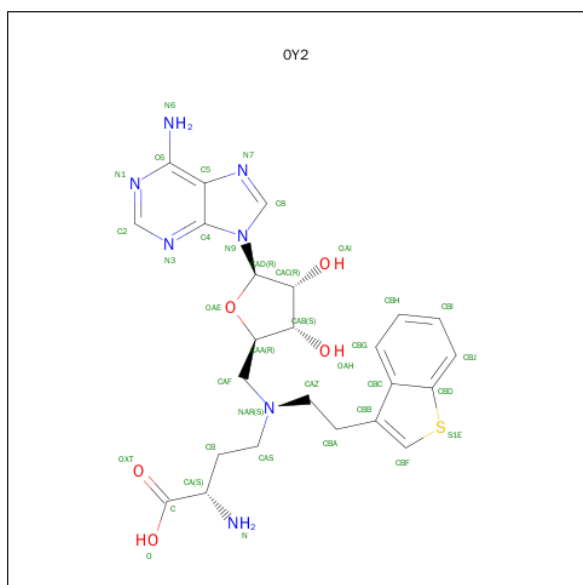
There are 3 unique types of molecules in this entry. The entry contains 6154 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 adenine methylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	243	Total	C	N	O	S	0	0	0
			2008	1300	338	362	8			
1	E	244	Total	C	N	O	S	0	0	0
			2016	1307	339	362	8			
1	F	244	Total	C	N	O	S	3	0	0
			2015	1305	339	363	8			

- Molecule 2 is 5'-{[(3S)-3-AMINO-3-CARBOXYPROPYL][2-(1-BENZOTHIOPHEN-3-YL)ETHYL]AMINO}-5'-DEOXYADENOSINE (three-letter code: 0Y2) (formula: C<sub>24</sub>H<sub>29</sub>N<sub>7</sub>O<sub>5</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	D	1	Total	C	N	O	S	0	0
			37	24	7	5	1		
2	E	1	Total	C	N	O	S	0	0
			37	24	7	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	F	1	Total	C	N	O	S	0	0
			37	24	7	5	1		

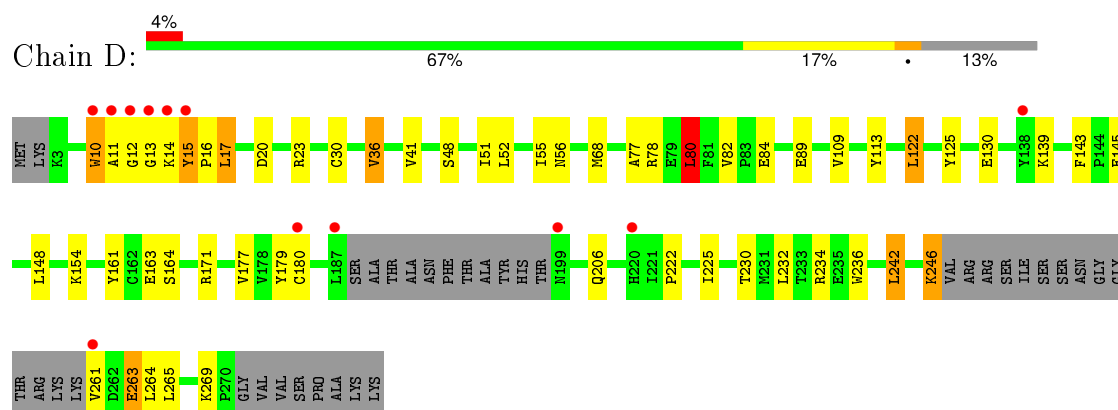
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	4	Total	O	0	0
			4	4		

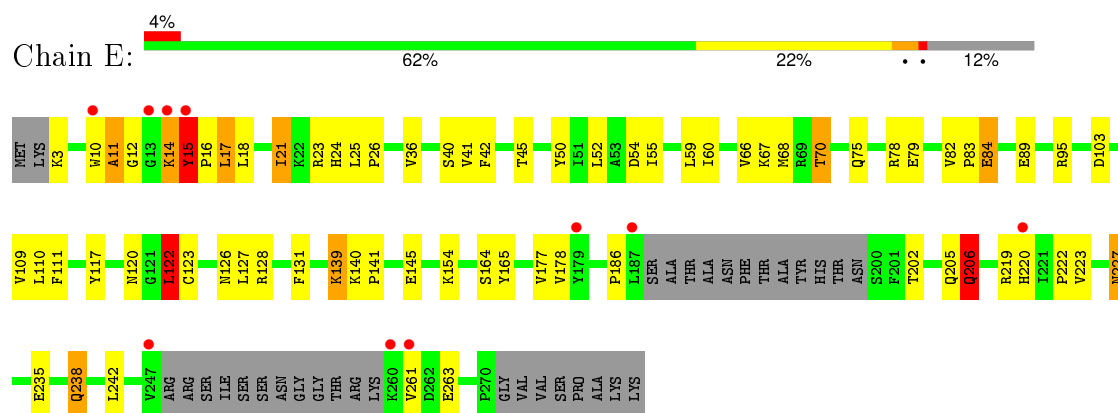
### 3 Residue-property plots

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.

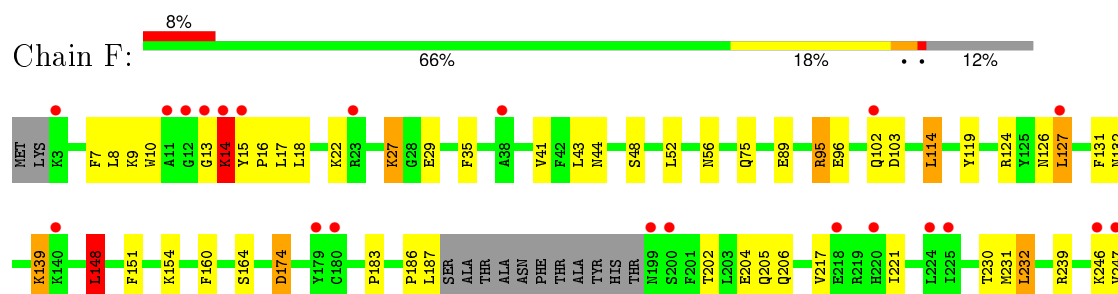
- Molecule 1: DNA adenine methylase



- Molecule 1: DNA adenine methylase



- Molecule 1: DNA adenine methylase



ARG	ARG	ILE	SER	ASN	GLY	THR	ARG	LYS	V261	D262	E263	K269	P270	GLY	VAL	VAL	SER	PRO	ALA	LYS	LYS
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	161.46 Å   161.46 Å   95.21 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	29.06 – 2.70 29.06 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.06-2.70) 99.2 (29.06-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.10 (at 2.72 Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.211 , 0.241 0.204 , 0.235	Depositor DCC
$R_{free}$ test set	2009 reflections (5.40%)	DCC
Wilson B-factor (Å <sup>2</sup> )	64.2	Xtriage
Anisotropy	0.226	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 54.9	EDS
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 39360 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6154	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 0Y2

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	D	1.19	4/2063 (0.2%)	1.04	3/2794 (0.1%)
1	E	1.08	2/2071 (0.1%)	0.99	6/2804 (0.2%)
1	F	1.09	4/2070 (0.2%)	1.01	5/2804 (0.2%)
All	All	1.12	10/6204 (0.2%)	1.01	14/8402 (0.2%)

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	E	0	1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	14	LYS	CG-CD	10.74	1.89	1.52
1	F	35	PHE	CE2-CZ	6.79	1.50	1.37
1	E	89	GLU	CG-CD	5.59	1.60	1.51
1	D	161	TYR	CE1-CZ	5.52	1.45	1.38
1	D	143	PHE	CE2-CZ	5.49	1.47	1.37
1	D	84	GLU	CG-CD	5.38	1.60	1.51
1	D	82	VAL	CB-CG1	5.11	1.63	1.52
1	E	75	GLN	CG-CD	5.05	1.62	1.51
1	F	151	PHE	CE1-CZ	5.04	1.47	1.37
1	F	174	ASP	CB-CG	5.01	1.62	1.51

All (14) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	128	ARG	NE-CZ-NH2	-7.32	116.64	120.30
1	F	127	LEU	CA-CB-CG	7.28	132.05	115.30
1	D	80	LEU	CB-CA-C	6.88	123.28	110.20
1	E	122	LEU	CA-CB-CG	6.76	130.86	115.30
1	E	95	ARG	NE-CZ-NH1	6.57	123.59	120.30
1	F	174	ASP	CB-CG-OD1	6.47	124.13	118.30
1	E	103	ASP	CB-CG-OD1	6.27	123.94	118.30
1	E	128	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	F	8	LEU	CB-CG-CD1	-5.82	101.11	111.00
1	F	95	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	D	122	LEU	CA-CB-CG	5.52	128.00	115.30
1	D	68	MET	CG-SD-CE	5.06	108.30	100.20
1	E	206	GLN	CA-CB-CG	-5.03	102.33	113.40
1	F	148	LEU	CB-CG-CD1	5.02	119.54	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	131	PHE	Peptide

## 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	D	2008	0	1944	50	0
1	E	2016	0	1960	52	0
1	F	2015	0	1953	34	0
2	D	37	0	27	8	0
2	E	37	0	27	2	0
2	F	37	0	27	5	0
3	D	4	0	0	0	0
All	All	6154	0	5938	140	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 12.

All (140) 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:D:10:TRP:HB2	2:D:301:0Y2:S1E	1.53	1.46
1:D:10:TRP:CB	2:D:301:0Y2:S1E	2.14	1.33
1:E:14:LYS:C	1:E:16:PRO:HD3	1.75	1.07
1:D:14:LYS:HB3	1:D:16:PRO:CD	1.85	1.05
1:D:14:LYS:HB3	1:D:16:PRO:HD2	1.39	1.01
1:F:230:THR:HG22	1:F:232:LEU:H	1.26	0.98
1:E:11:ALA:HB2	2:E:301:0Y2:H6	1.45	0.98
1:D:10:TRP:HB3	2:D:301:0Y2:S1E	2.00	0.97
1:F:246:LYS:O	1:F:247:VAL:HG23	1.72	0.88
1:D:16:PRO:O	1:D:17:LEU:HB2	1.74	0.87
1:D:13:GLY:N	1:D:14:LYS:HA	1.93	0.83
1:F:139:LYS:HD3	1:F:139:LYS:H	1.44	0.81
1:F:17:LEU:CD1	1:F:263:GLU:HG2	2.11	0.81
1:D:14:LYS:HB3	1:D:16:PRO:HD3	1.62	0.80
1:E:11:ALA:HB3	1:E:14:LYS:NZ	1.98	0.78
1:D:242:LEU:HG	1:D:264:LEU:HD11	1.66	0.77
1:E:10:TRP:HE1	1:E:40:SER:HB3	1.50	0.77
1:D:14:LYS:C	1:D:16:PRO:HD2	2.04	0.77
1:E:14:LYS:O	1:E:16:PRO:HD3	1.84	0.77
1:D:14:LYS:CB	1:D:16:PRO:HD2	2.16	0.74
1:D:13:GLY:H	1:D:14:LYS:HG3	1.52	0.73
2:D:301:0Y2:H6	2:D:301:0Y2:H23	1.69	0.73
1:E:82:VAL:HB	1:E:83:PRO:HD2	1.71	0.72
1:D:10:TRP:HB2	2:D:301:0Y2:CBF	2.20	0.71
1:E:235:GLU:O	1:E:238:GLN:HG3	1.94	0.68
1:F:230:THR:HG22	1:F:232:LEU:N	2.06	0.68
1:D:80:LEU:HD21	1:D:113:TYR:HB2	1.76	0.67
1:D:78:ARG:NH1	1:D:145:GLU:OE1	2.30	0.65
1:F:17:LEU:HD11	1:F:263:GLU:HG2	1.77	0.65
1:D:206:GLN:HG2	1:D:236:TRP:CH2	2.34	0.63
1:E:11:ALA:HB3	1:E:14:LYS:HZ3	1.61	0.63
1:F:18:LEU:O	1:F:22:LYS:HG3	1.99	0.63
1:D:80:LEU:O	1:D:80:LEU:HD23	1.99	0.63
1:D:10:TRP:O	1:D:10:TRP:CE3	2.53	0.62
1:D:10:TRP:O	1:D:10:TRP:HE3	1.82	0.62
1:E:12:GLY:H	1:E:14:LYS:HZ2	1.48	0.62
1:F:10:TRP:HB2	2:F:301:0Y2:S1E	2.39	0.61
1:E:10:TRP:HE1	1:E:40:SER:CB	2.13	0.61
1:F:29:GLU:O	1:F:48:SER:HB3	2.02	0.60
2:D:301:0Y2:H6	2:D:301:0Y2:CAB	2.31	0.60
1:D:13:GLY:N	1:D:14:LYS:CA	2.64	0.60
1:E:235:GLU:O	1:E:238:GLN:CG	2.50	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:246:LYS:HD3	1:D:246:LYS:O	2.03	0.59
1:E:177:VAL:HG22	1:E:222:PRO:HG2	1.83	0.58
1:D:15:TYR:N	1:D:16:PRO:HD2	2.18	0.58
2:F:301:OY2:H8	2:F:301:OY2:CAB	2.34	0.58
1:E:14:LYS:C	1:E:16:PRO:CD	2.63	0.57
1:D:17:LEU:HD11	1:D:263:GLU:HG2	1.86	0.57
1:F:15:TYR:HB2	1:F:16:PRO:HD3	1.86	0.57
1:F:139:LYS:CD	1:F:139:LYS:H	2.12	0.57
1:E:12:GLY:H	1:E:14:LYS:NZ	2.03	0.56
1:E:206:GLN:HE21	1:E:227:ASN:HD21	1.53	0.56
1:F:95:ARG:HH21	1:F:132:ASN:C	2.09	0.56
1:D:10:TRP:H	1:D:11:ALA:C	2.09	0.56
1:F:202:THR:H	1:F:205:GLN:HE21	1.54	0.56
1:D:13:GLY:H	1:D:14:LYS:CG	2.19	0.55
1:F:7:PHE:CG	1:F:148:LEU:HD13	2.42	0.55
1:E:45:THR:HG23	1:E:50:TYR:OH	2.07	0.55
1:D:206:GLN:CG	1:D:236:TRP:CZ3	2.90	0.55
1:F:13:GLY:O	1:F:14:LYS:HB2	2.07	0.54
1:D:206:GLN:HG3	1:D:236:TRP:CZ3	2.42	0.53
1:E:42:PHE:CD1	1:E:52:LEU:HD11	2.44	0.53
1:D:55:ILE:HG13	2:D:301:OY2:C4	2.38	0.53
1:D:163:GLU:OE2	1:D:171:ARG:NH2	2.36	0.53
1:D:15:TYR:N	1:D:16:PRO:CD	2.72	0.53
2:F:301:OY2:H1	2:F:301:OY2:OAH	2.09	0.53
1:E:178:VAL:HB	1:E:223:VAL:HG13	1.89	0.52
1:D:177:VAL:HG22	1:D:222:PRO:HG2	1.91	0.52
1:E:17:LEU:HD21	1:E:263:GLU:HG2	1.91	0.52
1:E:139:LYS:HG2	1:E:140:LYS:H	1.74	0.52
1:D:125:TYR:HA	1:D:130:GLU:O	2.09	0.52
1:E:78:ARG:NH1	1:E:145:GLU:OE2	2.43	0.52
1:E:10:TRP:CE3	1:E:14:LYS:HD3	2.45	0.52
1:F:22:LYS:NZ	1:F:44:ASN:O	2.21	0.51
1:E:202:THR:H	1:E:205:GLN:HE21	1.58	0.51
1:D:20:ASP:HA	1:D:23:ARG:HG3	1.91	0.51
1:E:219:ARG:O	1:E:220:HIS:HB2	2.10	0.51
1:D:56:ASN:OD1	1:D:56:ASN:C	2.49	0.51
1:E:24:HIS:O	1:E:26:PRO:HD3	2.10	0.50
1:E:139:LYS:H	1:E:139:LYS:HD2	1.77	0.50
1:F:217:VAL:HA	1:F:221:ILE:O	2.11	0.50
1:D:10:TRP:H	1:D:12:GLY:N	2.08	0.50
1:F:126:ASN:C	1:F:126:ASN:OD1	2.49	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:206:GLN:NE2	1:E:227:ASN:HD21	2.09	0.49
1:E:120:ASN:HB2	1:E:122:LEU:HD22	1.92	0.49
1:F:246:LYS:O	1:F:247:VAL:CG2	2.53	0.49
1:D:80:LEU:C	1:D:80:LEU:HD23	2.33	0.49
1:E:41:VAL:O	1:E:45:THR:HG22	2.12	0.49
1:E:15:TYR:N	1:E:16:PRO:CD	2.76	0.49
1:F:186:PRO:O	1:F:187:LEU:HG	2.12	0.49
1:D:13:GLY:N	1:D:14:LYS:HG3	2.25	0.48
1:D:77:ALA:HA	1:D:109:VAL:HG13	1.96	0.48
1:D:14:LYS:CA	1:D:16:PRO:HD2	2.43	0.48
1:E:15:TYR:N	1:E:16:PRO:HD3	2.24	0.48
1:D:206:GLN:HG2	1:D:236:TRP:CZ3	2.48	0.48
1:D:13:GLY:H	1:D:14:LYS:CA	2.26	0.48
1:F:202:THR:H	1:F:205:GLN:NE2	2.11	0.48
1:E:67:LYS:HG2	1:E:68:MET:HG3	1.96	0.47
1:D:230:THR:O	1:D:234:ARG:HG3	2.14	0.47
1:D:10:TRP:CG	2:D:301:OY2:S1E	3.03	0.47
1:E:66:VAL:O	1:E:70:THR:HB	2.15	0.47
1:E:202:THR:H	1:E:205:GLN:NE2	2.12	0.46
1:D:30:CYS:SG	1:D:51:ILE:HD12	2.55	0.46
1:E:42:PHE:O	1:E:154:LYS:HE3	2.15	0.46
1:F:52:LEU:O	1:F:160:PHE:HA	2.16	0.46
1:E:117:TYR:CZ	1:E:141:PRO:HG2	2.51	0.45
1:F:13:GLY:O	1:F:14:LYS:CB	2.64	0.45
1:F:154:LYS:HB3	1:F:154:LYS:HE2	1.85	0.45
1:F:269:LYS:HG3	1:F:270:PRO:HD2	1.98	0.45
1:E:10:TRP:HZ2	1:E:40:SER:H	1.65	0.45
1:F:114:LEU:HG	1:F:131:PHE:CE2	2.52	0.44
1:E:123:CYS:HB2	2:E:301:OY2:H4	2.00	0.44
1:E:14:LYS:O	1:E:15:TYR:HD2	2.01	0.43
1:E:17:LEU:O	1:E:21:ILE:HG13	2.18	0.43
1:E:11:ALA:HB3	1:E:14:LYS:HZ2	1.83	0.43
1:D:180:CYS:HB2	1:D:225:ILE:HG12	2.01	0.43
1:D:12:GLY:N	1:D:13:GLY:CA	2.82	0.43
1:E:83:PRO:HG2	1:E:84:GLU:OE1	2.19	0.43
1:E:139:LYS:H	1:E:139:LYS:CD	2.32	0.43
1:F:27:LYS:HA	1:F:27:LYS:HD2	1.70	0.43
1:F:10:TRP:CD1	2:F:301:OY2:CBF	3.02	0.42
1:E:14:LYS:HE3	1:E:14:LYS:HB3	1.49	0.42
1:F:261:VAL:HG22	1:F:262:ASP:H	1.84	0.42
1:D:41:VAL:HG21	1:D:179:TYR:CE2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:82:VAL:HB	1:E:83:PRO:CD	2.44	0.42
1:E:139:LYS:HD2	1:E:139:LYS:N	2.35	0.42
1:F:103:ASP:OD1	1:F:103:ASP:C	2.58	0.42
1:D:80:LEU:C	1:D:80:LEU:CD2	2.88	0.42
1:E:36:VAL:HG12	1:E:54:ASP:HB2	2.02	0.42
1:E:126:ASN:C	1:E:126:ASN:OD1	2.58	0.42
1:E:110:LEU:O	1:E:111:PHE:C	2.59	0.41
1:F:230:THR:HG22	1:F:231:MET:N	2.34	0.41
1:E:140:LYS:HA	1:E:141:PRO:HD3	1.87	0.41
1:F:10:TRP:CE3	1:F:13:GLY:O	2.73	0.41
1:F:56:ASN:OD1	1:F:56:ASN:C	2.58	0.41
1:E:59:LEU:HA	1:E:59:LEU:HD12	1.83	0.41
1:D:13:GLY:H	1:D:14:LYS:CB	2.34	0.41
1:F:183:PRO:HA	2:F:301:OY2:OXT	2.21	0.41
1:E:164:SER:O	1:E:165:TYR:C	2.59	0.41
1:D:36:VAL:CG1	1:D:52:LEU:HB3	2.51	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	D	237/278 (85%)	222 (94%)	14 (6%)	1 (0%)	39	69
1	E	238/278 (86%)	220 (92%)	14 (6%)	4 (2%)	11	29
1	F	238/278 (86%)	220 (92%)	16 (7%)	2 (1%)	24	51
All	All	713/834 (86%)	662 (93%)	44 (6%)	7 (1%)	19	45

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	11	ALA

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Mol	Chain	Res	Type
1	E	15	TYR
1	F	14	LYS
1	F	119	TYR
1	E	238	GLN
1	D	15	TYR
1	E	109	VAL

### 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	D	211/239 (88%)	193 (92%)	18 (8%)	13	30
1	E	212/239 (89%)	191 (90%)	21 (10%)	10	22
1	F	212/239 (89%)	193 (91%)	19 (9%)	12	27
All	All	635/717 (89%)	577 (91%)	58 (9%)	12	26

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	10	TRP
1	D	17	LEU
1	D	36	VAL
1	D	48	SER
1	D	80	LEU
1	D	89	GLU
1	D	122	LEU
1	D	139	LYS
1	D	148	LEU
1	D	154	LYS
1	D	164	SER
1	D	232	LEU
1	D	242	LEU
1	D	246	LYS
1	D	261	VAL
1	D	263	GLU

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Mol	Chain	Res	Type
1	D	265	LEU
1	D	269	LYS
1	E	3	LYS
1	E	14	LYS
1	E	15	TYR
1	E	17	LEU
1	E	18	LEU
1	E	21	ILE
1	E	23	ARG
1	E	25	LEU
1	E	55	ILE
1	E	60	ILE
1	E	70	THR
1	E	79	GLU
1	E	84	GLU
1	E	122	LEU
1	E	127	LEU
1	E	139	LYS
1	E	186	PRO
1	E	206	GLN
1	E	227	ASN
1	E	242	LEU
1	E	261	VAL
1	F	9	LYS
1	F	27	LYS
1	F	41	VAL
1	F	43	LEU
1	F	75	GLN
1	F	89	GLU
1	F	96	GLU
1	F	102	GLN
1	F	114	LEU
1	F	124	ARG
1	F	127	LEU
1	F	139	LYS
1	F	148	LEU
1	F	164	SER
1	F	174	ASP
1	F	204	GLU
1	F	206	GLN
1	F	232	LEU
1	F	239	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	205	GLN
1	E	206	GLN
1	F	4	ASN
1	F	205	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](#)

3 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	0Y2	D	301	-	30,41,41	3.09	11 (36%)	27,59,59	2.20	4 (14%)
2	0Y2	E	301	-	30,41,41	2.18	10 (33%)	27,59,59	2.45	7 (25%)
2	0Y2	F	301	-	30,41,41	3.00	13 (43%)	27,59,59	2.59	8 (29%)

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	0Y2	D	301	-	-	0/14/38/38	0/5/5/5
2	0Y2	E	301	-	-	0/14/38/38	0/5/5/5
2	0Y2	F	301	-	-	0/14/38/38	0/5/5/5

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	0Y2	CBF-CBB	-11.16	1.30	1.37
2	F	301	0Y2	CBF-CBB	-11.11	1.30	1.37
2	D	301	0Y2	CBF-S1E	-7.76	1.61	1.70
2	F	301	0Y2	CBF-S1E	-6.71	1.62	1.70
2	E	301	0Y2	CBF-S1E	-5.78	1.63	1.70
2	E	301	0Y2	CBF-CBB	-5.70	1.33	1.37
2	D	301	0Y2	CBJ-CBD	-4.14	1.35	1.40
2	D	301	0Y2	OAI-CAC	-3.58	1.34	1.43
2	E	301	0Y2	OAI-CAC	-3.25	1.35	1.43
2	F	301	0Y2	CBG-CBC	-3.25	1.35	1.42
2	F	301	0Y2	OAI-CAC	-3.06	1.35	1.43
2	F	301	0Y2	CBJ-CBD	-2.90	1.36	1.40
2	D	301	0Y2	CAF-CAA	-2.80	1.44	1.51
2	D	301	0Y2	CBG-CBC	-2.58	1.37	1.42
2	E	301	0Y2	CBG-CBC	-2.57	1.37	1.42
2	D	301	0Y2	OAH-CAB	-2.41	1.37	1.43
2	F	301	0Y2	OAH-CAB	-2.36	1.37	1.43
2	E	301	0Y2	CAC-CAB	-2.31	1.47	1.53
2	E	301	0Y2	CAF-CAA	-2.12	1.46	1.51
2	F	301	0Y2	CAC-CAB	-2.09	1.47	1.53
2	D	301	0Y2	CAC-CAB	-2.09	1.47	1.53
2	E	301	0Y2	OAH-CAB	-2.02	1.38	1.43
2	D	301	0Y2	CBA-CBB	2.02	1.56	1.51
2	F	301	0Y2	CBH-CBG	2.16	1.41	1.36
2	D	301	0Y2	CAZ-NAR	2.24	1.52	1.47
2	F	301	0Y2	CBA-CBB	2.26	1.57	1.51
2	F	301	0Y2	CAF-NAR	2.28	1.51	1.47
2	E	301	0Y2	C4-N3	2.29	1.39	1.35
2	E	301	0Y2	C2-N3	2.58	1.36	1.32
2	F	301	0Y2	C2-N3	2.64	1.36	1.32
2	E	301	0Y2	C6-N6	2.78	1.43	1.34
2	D	301	0Y2	C6-N6	2.97	1.44	1.34
2	F	301	0Y2	C6-N6	3.20	1.44	1.34
2	F	301	0Y2	CAZ-NAR	3.23	1.55	1.47

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	0Y2	N3-C2-N1	-8.83	122.14	128.89
2	E	301	0Y2	N3-C2-N1	-8.07	122.72	128.89
2	F	301	0Y2	N3-C2-N1	-8.04	122.74	128.89
2	F	301	0Y2	CAA-OAE-CAD	-3.99	105.34	109.72
2	E	301	0Y2	CAZ-NAR-CAF	-3.13	104.58	112.25
2	F	301	0Y2	CAS-NAR-CAF	-3.04	104.79	112.25
2	E	301	0Y2	CAA-OAE-CAD	-2.83	106.61	109.72
2	E	301	0Y2	C4-C5-N7	-2.65	107.05	109.48
2	D	301	0Y2	CAS-NAR-CAF	-2.44	106.27	112.25
2	F	301	0Y2	CAS-NAR-CAZ	-2.37	105.77	111.45
2	F	301	0Y2	CBB-CBF-S1E	2.21	114.62	112.26
2	F	301	0Y2	CAC-CAB-CAA	2.24	107.22	102.61
2	D	301	0Y2	C2-N1-C6	2.36	122.98	118.77
2	D	301	0Y2	CAZ-CBA-CBB	2.58	118.71	112.86
2	F	301	0Y2	CB-CAS-NAR	2.92	120.06	114.53
2	E	301	0Y2	CB-CAS-NAR	3.06	120.34	114.53
2	E	301	0Y2	CAZ-CBA-CBB	3.63	121.08	112.86
2	E	301	0Y2	CAC-CAD-N9	4.86	121.72	114.29
2	F	301	0Y2	CAZ-CBA-CBB	6.65	127.94	112.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	301	0Y2	8	0
2	E	301	0Y2	2	0
2	F	301	0Y2	5	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

### 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	D	243/278 (87%)	0.03	12 (4%) 33 32	35, 49, 77, 107	0
1	E	244/278 (87%)	0.07	10 (4%) 41 41	39, 56, 87, 104	0
1	F	244/278 (87%)	0.26	22 (9%) 12 9	36, 56, 80, 103	1 (0%)
All	All	731/834 (87%)	0.12	44 (6%) 25 24	35, 54, 84, 107	1 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	12	GLY	6.9
1	F	12	GLY	6.5
1	F	199	ASN	5.7
1	D	13	GLY	5.2
1	F	11	ALA	4.9
1	D	14	LYS	4.4
1	F	179	TYR	4.3
1	E	260	LYS	4.1
1	E	10	TRP	4.0
1	D	199	ASN	3.7
1	D	11	ALA	3.6
1	E	247	VAL	3.6
1	D	10	TRP	3.6
1	E	220	HIS	3.5
1	E	13	GLY	3.1
1	F	200	SER	3.1
1	E	14	LYS	3.1
1	D	261	VAL	3.1
1	F	127	LEU	3.0
1	D	187	LEU	3.0
1	F	180	CYS	2.8
1	F	13	GLY	2.8
1	E	15	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	220	HIS	2.6
1	D	15	TYR	2.6
1	F	38	ALA	2.5
1	F	3	LYS	2.5
1	F	140	LYS	2.5
1	D	180	CYS	2.3
1	F	247	VAL	2.3
1	E	187	LEU	2.2
1	F	14	LYS	2.2
1	F	261	VAL	2.2
1	D	220	HIS	2.2
1	E	179	TYR	2.2
1	F	102	GLN	2.1
1	F	224	LEU	2.1
1	F	15	TYR	2.1
1	D	138	TYR	2.0
1	F	246	LYS	2.0
1	F	225	ILE	2.0
1	F	23	ARG	2.0
1	F	218	GLU	2.0
1	E	261	VAL	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	0Y2	D	301	37/37	0.77	0.28	1.42	76,84,91,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	0Y2	E	301	37/37	0.68	0.31	1.26	77,94,107,108	0
2	0Y2	F	301	37/37	0.84	0.24	0.52	72,83,90,92	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.