



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:27 PM UTC

PDB ID : 4GOO / pdb_00004goo
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

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

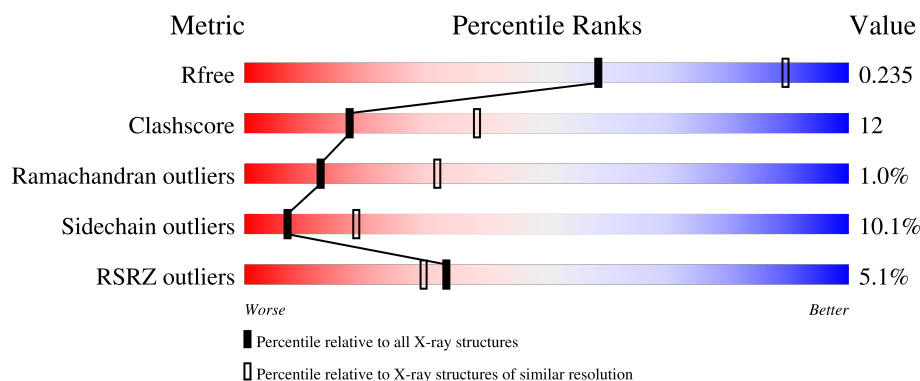
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (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 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	D	278	4% 63% 19% • • 13%
1	E	278	5% 58% 25% • • 12%
1	F	278	4% 61% 23% • 12%

2 Entry composition [i](#)

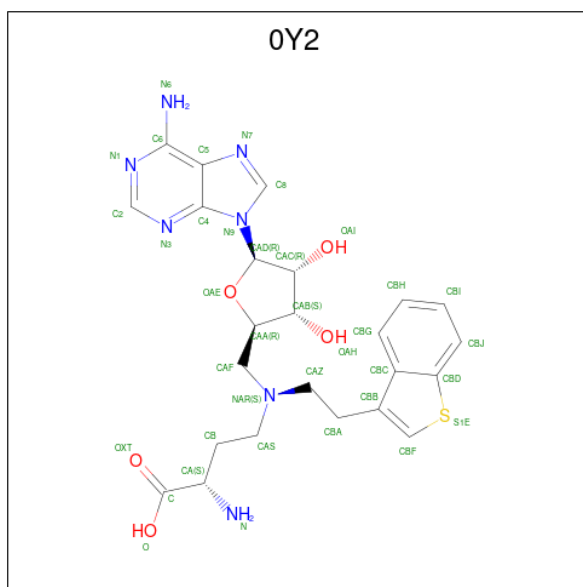
There are 3 unique types of molecules in this entry. The entry contains 6154 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (CCD ID: 0Y2) (formula: C₂₄H₂₉N₇O₅S).



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		
2	F	1	Total	C	N	O	S	0	0
			37	24	7	5	1		

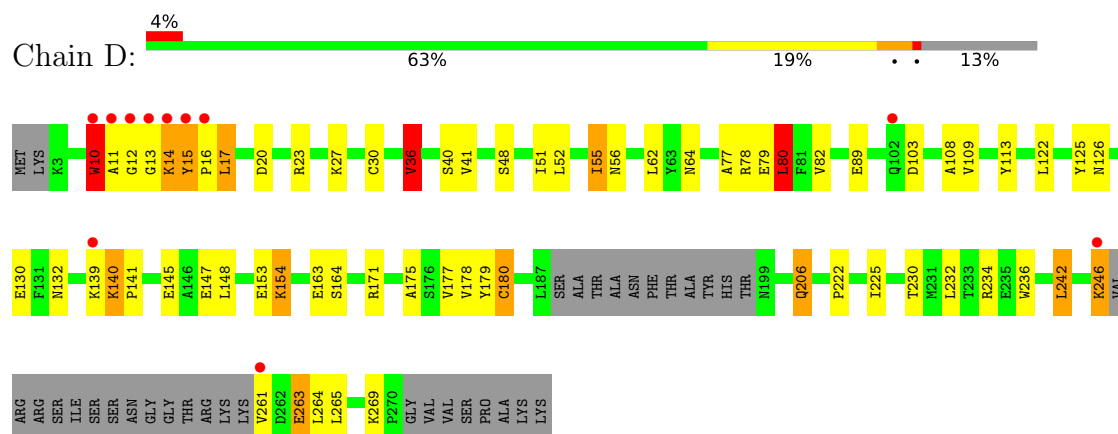
- Molecule 3 is water.

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

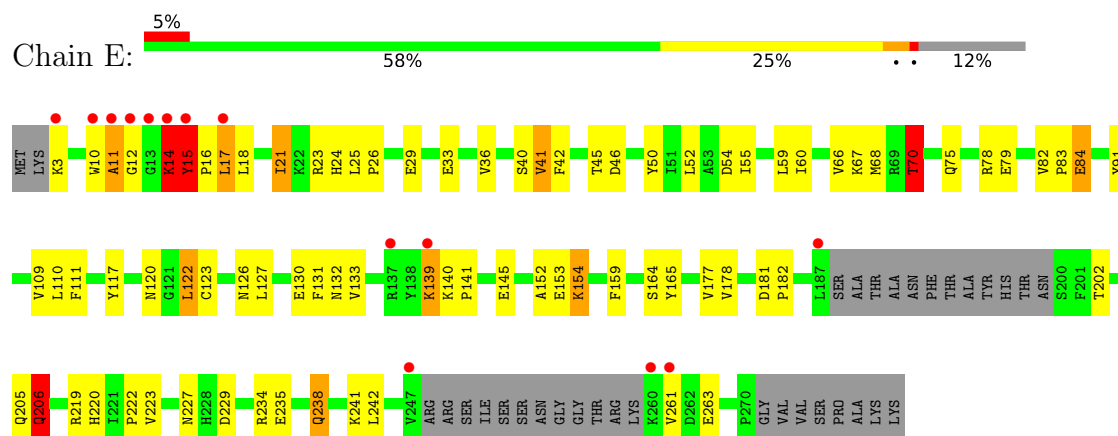
3 Residue-property plots [i](#)

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.

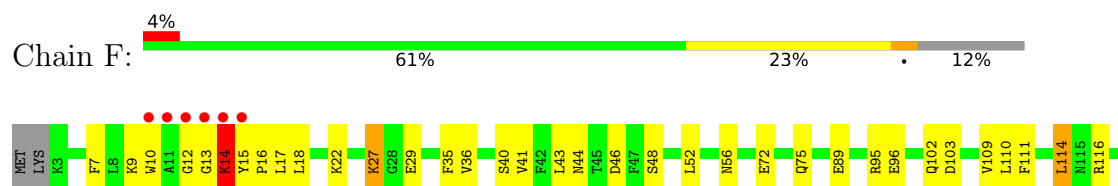
- Molecule 1: DNA adenine methylase

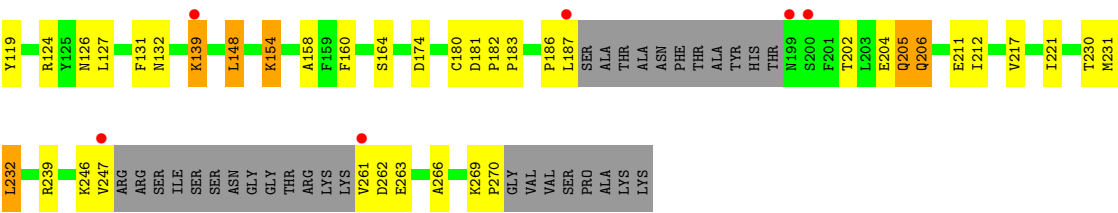


- Molecule 1: DNA adenine methylase



- Molecule 1: DNA adenine methylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.46Å 161.46Å 95.21Å 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) 99.1 (29.06-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 2.72Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.211 , 0.241 0.204 , 0.235	Depositor DCC
R_{free} test set	2009 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	64.2	Xtriage
Anisotropy	0.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6154	wwPDB-VP
Average B, all atoms (Å ²)	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.*

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: 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	1.35	8/2063 (0.4%)	1.42	17/2794 (0.6%)
1	E	1.25	3/2071 (0.1%)	1.32	13/2804 (0.5%)
1	F	1.26	5/2070 (0.2%)	1.30	10/2804 (0.4%)
All	All	1.29	16/6204 (0.3%)	1.35	40/8402 (0.5%)

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 (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	14	LYS	CG-CD	12.18	1.89	1.52
1	D	10	TRP	CA-C	6.93	1.61	1.52
1	D	178	VAL	CA-CB	6.68	1.62	1.54
1	E	152	ALA	CA-CB	-6.08	1.43	1.53
1	F	181	ASP	N-CA	-5.93	1.39	1.46
1	D	103	ASP	CA-C	5.60	1.58	1.52
1	E	159	PHE	N-CA	5.41	1.52	1.46
1	D	108	ALA	CA-C	5.38	1.59	1.52
1	F	206	GLN	CA-C	-5.31	1.45	1.52
1	D	40	SER	N-CA	-5.29	1.39	1.46
1	F	266	ALA	CA-CB	-5.25	1.46	1.53
1	D	62	LEU	N-CA	5.23	1.52	1.46
1	D	154	LYS	CA-C	5.20	1.59	1.52
1	F	116	ARG	CA-C	5.16	1.59	1.52
1	D	132	ASN	CA-C	5.11	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	182	PRO	CA-C	-5.05	1.47	1.52

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	140	LYS	CA-C-N	-18.40	102.55	120.21
1	D	140	LYS	C-N-CA	-18.40	102.55	120.21
1	E	70	THR	N-CA-C	8.90	120.60	111.07
1	D	79	GLU	CA-C-N	-8.77	109.45	122.86
1	D	79	GLU	C-N-CA	-8.77	109.45	122.86
1	D	80	LEU	CB-CA-C	7.55	123.28	111.13
1	D	36	VAL	CB-CA-C	-7.39	102.36	112.04
1	F	109	VAL	N-CA-C	-7.38	103.59	110.53
1	D	64	ASN	N-CA-C	-6.92	103.51	112.23
1	F	40	SER	N-CA-C	6.75	118.64	111.28
1	D	82	VAL	CA-C-N	-6.52	112.97	119.56
1	D	82	VAL	C-N-CA	-6.52	112.97	119.56
1	E	133	VAL	N-CA-C	6.51	116.06	108.96
1	E	132	ASN	N-CA-CB	-6.27	102.56	110.90
1	E	14	LYS	N-CA-C	6.25	118.10	109.29
1	E	181	ASP	CA-C-N	6.25	126.82	120.38
1	E	181	ASP	C-N-CA	6.25	126.82	120.38
1	D	36	VAL	N-CA-CB	6.06	118.05	110.47
1	D	10	TRP	N-CA-C	6.01	123.39	113.89
1	E	41	VAL	N-CA-C	-5.91	104.98	110.53
1	E	206	GLN	CA-CB-CG	-5.89	102.33	114.10
1	F	180	CYS	N-CA-C	5.74	118.31	109.07
1	F	46	ASP	N-CA-C	5.71	117.29	110.44
1	F	205	GLN	N-CA-C	-5.62	104.54	111.40
1	D	36	VAL	N-CA-C	5.59	116.99	110.62
1	F	12	GLY	N-CA-C	5.59	124.24	115.08
1	F	72	GLU	N-CA-C	-5.58	105.20	111.28
1	E	29	GLU	N-CA-C	5.57	119.25	112.23
1	F	182	PRO	CA-C-N	-5.35	114.27	119.78
1	F	182	PRO	C-N-CA	-5.35	114.27	119.78
1	E	33	GLU	CA-C-N	5.32	124.93	119.56
1	E	33	GLU	C-N-CA	5.32	124.93	119.56
1	E	46	ASP	N-CA-C	5.30	116.81	109.54
1	D	206	GLN	N-CA-C	-5.24	105.47	111.07
1	D	180	CYS	N-CA-C	5.22	117.03	108.52
1	D	175	ALA	N-CA-C	5.22	118.75	112.38
1	D	103	ASP	CB-CA-C	5.15	115.72	109.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	158	ALA	N-CA-C	5.10	117.56	109.50
1	E	153	GLU	N-CA-C	-5.07	105.83	111.36
1	D	153	GLU	N-CA-C	-5.02	105.46	111.03

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	54	0
1	E	2016	0	1960	55	0
1	F	2015	0	1953	38	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	151	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 (151) 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	1.12
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:11:ALA:HB2	2:E:301:OY2:H6	1.45	0.98
1:D:10:TRP:HB3	2:D:301:OY2: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:D:14:LYS:C	1:D:16:PRO:HD2	2.04	0.82
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: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:OY2:H23	2:D:301:OY2:H6	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:OY2:CBF	2.20	0.71
1:E:14:LYS:C	1:E:16:PRO:CD	2.63	0.68
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:HD23	1:D:80:LEU:O	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:OY2:S1E	2.39	0.61
1:E:10:TRP:HE1	1:E:40:SER:CB	2.13	0.61
1:F:95:ARG:HH21	1:F:132:ASN:C	2.09	0.60
1:D:10:TRP:H	1:D:11:ALA:C	2.09	0.60
1:F:29:GLU:O	1:F:48:SER:HB3	2.02	0.60
2:D:301:OY2:H6	2:D:301:OY2: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
1:D:246:LYS:O	1:D:246:LYS:HD3	2.03	0.59
1:E:10:TRP:NE1	1:E:40:SER:HG	2.00	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:H	1:F:139:LYS:CD	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:22:LYS:NZ	1:F:44:ASN:O	2.21	0.56
1:D:56:ASN:OD1	1:D:56:ASN:C	2.49	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:126:ASN:OD1	1:F:126:ASN:C	2.49	0.54
1:D:80:LEU:HD23	1:D:80:LEU:C	2.33	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:F:13:GLY:O	1:F:14:LYS:HB2	2.07	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:D:177:VAL:HG22	1:D:222:PRO:HG2	1.91	0.52
1:E:178:VAL:HB	1:E:223:VAL:HG13	1.89	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: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: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:E:120:ASN:HB2	1:E:122:LEU:HD22	1.92	0.49
1:E:206:GLN:NE2	1:E:227:ASN:HD21	2.09	0.49
1:F:246:LYS:O	1:F:247:VAL:CG2	2.53	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:230:THR:O	1:D:234:ARG:HG3	2.14	0.47
1:E:67:LYS:HG2	1:E:68:MET:HG3	1.96	0.47
1:D:10:TRP:CG	2:D:301:OY2:S1E	3.03	0.47
1:D:80:LEU:C	1:D:80:LEU:CD2	2.88	0.47
1:E:66:VAL:O	1:E:70:THR:HB	2.15	0.47
1:F:103:ASP:OD1	1:F:103:ASP:C	2.58	0.47
1:E:126:ASN:OD1	1:E:126:ASN:C	2.58	0.47
1:E:202:THR:H	1:E:205:GLN:NE2	2.12	0.46
1:E:110:LEU:O	1:E:111:PHE:C	2.59	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:F:56:ASN:OD1	1:F:56:ASN:C	2.58	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:HE2	1:F:154:LYS:HB3	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:E:164:SER:O	1:E:165:TYR:C	2.59	0.45
1:F:95:ARG:NH2	1:F:132:ASN:C	2.75	0.44
1:F:114:LEU:HG	1:F:131:PHE:CE2	2.52	0.44
1:D:140:LYS:O	1:D:141:PRO:C	2.59	0.44
1:F:35:PHE:O	1:F:36:VAL:C	2.60	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:229:ASP:CG	1:E:234:ARG:HH21	2.23	0.43
1:F:211:GLU:O	1:F:212:ILE:C	2.58	0.43
1:D:180:CYS:HB2	1:D:225:ILE:HG12	2.01	0.43
1:E:11:ALA:HB3	1:E:14:LYS:HZ2	1.83	0.43
1:D:147:GLU:H	1:D:147:GLU:CD	2.24	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HD2	1:F:27:LYS:HA	1.70	0.43
1:F:10:TRP:CD1	2:F:301:OY2:CBF	3.02	0.42
1:E:14:LYS:HB3	1:E:14:LYS:HE3	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
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:E:36:VAL:HG12	1:E:54:ASP:HB2	2.02	0.42
1:E:91:TYR:CD2	1:E:91:TYR:C	2.98	0.42
1:E:140:LYS:HA	1:E:141:PRO:HD3	1.87	0.41
1:F:230:THR:HG22	1:F:231:MET:N	2.34	0.41
1:F:10:TRP:CE3	1:F:13:GLY:O	2.73	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:D:10:TRP:N	1:D:11:ALA:C	2.79	0.41
1:F:110:LEU:O	1:F:111:PHE:C	2.65	0.40
1:D:36:VAL:CG1	1:D:52:LEU:HB3	2.51	0.40
1:D:126:ASN:OD1	1:D:126:ASN:C	2.64	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%)	30	54
1	E	238/278 (86%)	220 (92%)	14 (6%)	4 (2%)	7	19
1	F	238/278 (86%)	220 (92%)	16 (7%)	2 (1%)	16	37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	713/834 (86%)	662 (93%)	44 (6%)	7 (1%)	12	32

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	11	ALA
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%)	190 (90%)	21 (10%)	7	19
1	E	212/239 (89%)	189 (89%)	23 (11%)	6	16
1	F	212/239 (89%)	192 (91%)	20 (9%)	8	21
All	All	635/717 (89%)	571 (90%)	64 (10%)	7	18

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

Mol	Chain	Res	Type
1	D	10	TRP
1	D	14	LYS
1	D	17	LEU
1	D	27	LYS
1	D	36	VAL
1	D	48	SER
1	D	55	ILE
1	D	80	LEU
1	D	89	GLU

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Mol	Chain	Res	Type
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
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	75	GLN
1	E	79	GLU
1	E	84	GLU
1	E	122	LEU
1	E	127	LEU
1	E	130	GLU
1	E	139	LYS
1	E	154	LYS
1	E	206	GLN
1	E	241	LYS
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

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Mol	Chain	Res	Type
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	154	LYS
1	F	164	SER
1	F	174	ASP
1	F	204	GLU
1	F	206	GLN
1	F	232	LEU
1	F	239	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	24	HIS
1	D	238	GLN
1	E	205	GLN
1	E	206	GLN
1	F	4	ASN
1	F	120	ASN
1	F	205	GLN
1	F	238	GLN
1	F	243	HIS

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

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	E	301	-	40,41,41	2.18	11 (27%)	51,59,59	2.93	20 (39%)
2	0Y2	F	301	-	40,41,41	2.44	12 (30%)	51,59,59	3.09	20 (39%)
2	0Y2	D	301	-	40,41,41	2.42	13 (32%)	51,59,59	3.35	17 (33%)

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	E	301	-	-	14/22/38/38	0/5/5/5
2	0Y2	F	301	-	-	6/22/38/38	0/5/5/5
2	0Y2	D	301	-	-	11/22/38/38	0/5/5/5

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	0Y2	CBF-S1E	-8.35	1.61	1.72
2	F	301	0Y2	CBF-S1E	-7.38	1.62	1.72
2	E	301	0Y2	CBF-S1E	-6.52	1.63	1.72
2	E	301	0Y2	CBC-CBD	6.33	1.50	1.41
2	D	301	0Y2	CBF-CBB	-6.00	1.30	1.36
2	F	301	0Y2	CBF-CBB	-5.97	1.30	1.36
2	F	301	0Y2	CBC-CBD	5.07	1.48	1.41
2	D	301	0Y2	CBC-CBD	4.17	1.47	1.41
2	F	301	0Y2	C6-N6	4.15	1.44	1.34
2	D	301	0Y2	C6-N6	3.88	1.44	1.34
2	F	301	0Y2	CBA-CBB	3.75	1.57	1.50
2	E	301	0Y2	C6-N6	3.65	1.43	1.34
2	D	301	0Y2	OAI-CAC	-3.46	1.34	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	0Y2	CBA-CBB	3.43	1.56	1.50
2	F	301	0Y2	CAZ-NAR	3.34	1.55	1.47
2	E	301	0Y2	CBA-CBB	3.14	1.56	1.50
2	E	301	0Y2	OAI-CAC	-3.14	1.35	1.43
2	F	301	0Y2	OAI-CAC	-2.95	1.35	1.43
2	F	301	0Y2	CBG-CBC	-2.85	1.35	1.39
2	E	301	0Y2	CBD-S1E	2.83	1.81	1.73
2	E	301	0Y2	CBF-CBB	-2.71	1.33	1.36
2	D	301	0Y2	CBJ-CBD	-2.70	1.35	1.39
2	D	301	0Y2	C5-C4	2.68	1.43	1.39
2	E	301	0Y2	C4-N3	2.44	1.39	1.34
2	E	301	0Y2	C5-C4	2.44	1.43	1.39
2	F	301	0Y2	C5-N7	-2.37	1.34	1.39
2	D	301	0Y2	OAH-CAB	-2.33	1.37	1.43
2	D	301	0Y2	CAZ-NAR	2.31	1.52	1.47
2	E	301	0Y2	CAC-CAB	-2.31	1.47	1.53
2	D	301	0Y2	CAF-CAA	-2.29	1.44	1.51
2	F	301	0Y2	CAF-NAR	2.28	1.51	1.47
2	F	301	0Y2	OAH-CAB	-2.28	1.37	1.43
2	E	301	0Y2	C8-N7	2.22	1.35	1.31
2	D	301	0Y2	C4-N3	2.17	1.38	1.34
2	F	301	0Y2	CAC-CAB	-2.09	1.47	1.53
2	D	301	0Y2	CAC-CAB	-2.09	1.47	1.53

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	0Y2	CBD-S1E-CBF	14.58	99.07	91.16
2	D	301	0Y2	CBC-CBD-S1E	-13.77	102.06	111.46
2	F	301	0Y2	CBC-CBD-S1E	-11.81	103.40	111.46
2	E	301	0Y2	CBC-CBD-S1E	-10.17	104.52	111.46
2	F	301	0Y2	CBD-S1E-CBF	9.02	96.05	91.16
2	F	301	0Y2	CAZ-CBA-CBB	8.39	127.94	112.58
2	E	301	0Y2	CBD-S1E-CBF	7.46	95.21	91.16
2	E	301	0Y2	CBJ-CBD-S1E	6.13	138.14	127.36
2	E	301	0Y2	C5-C4-N3	-5.48	119.17	126.72
2	E	301	0Y2	CAZ-CBA-CBB	4.65	121.08	112.58
2	D	301	0Y2	CBJ-CBD-S1E	4.55	135.35	127.36
2	F	301	0Y2	C5-C4-N3	-4.44	120.60	126.72
2	D	301	0Y2	N1-C2-N3	-4.26	122.14	128.58
2	F	301	0Y2	CBJ-CBD-S1E	4.01	134.40	127.36
2	E	301	0Y2	N3-C4-N9	3.99	133.96	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	0Y2	C5-C4-N3	-3.99	121.23	126.72
2	E	301	0Y2	CBJ-CBD-CBC	-3.94	117.34	121.47
2	E	301	0Y2	N1-C2-N3	-3.88	122.72	128.58
2	F	301	0Y2	N1-C2-N3	-3.86	122.74	128.58
2	E	301	0Y2	CBG-CBC-CBB	-3.77	128.23	133.85
2	E	301	0Y2	CBA-CBB-CBC	-3.76	118.96	126.18
2	F	301	0Y2	N3-C4-N9	3.67	133.41	127.17
2	F	301	0Y2	N9-C8-N7	-3.53	108.92	113.94
2	E	301	0Y2	CAC-CAD-N9	3.39	121.72	113.30
2	F	301	0Y2	C5-N7-C8	3.37	108.75	103.45
2	F	301	0Y2	O-C-OXT	-3.36	116.47	124.08
2	D	301	0Y2	CAZ-CBA-CBB	3.35	118.71	112.58
2	E	301	0Y2	C2-N3-C4	3.31	119.92	111.83
2	D	301	0Y2	N3-C4-N9	3.20	132.61	127.17
2	E	301	0Y2	CAZ-NAR-CAF	-3.14	104.58	111.89
2	E	301	0Y2	C5-N7-C8	3.10	108.32	103.45
2	E	301	0Y2	C4-C5-N7	-3.09	107.05	110.58
2	F	301	0Y2	CAS-NAR-CAF	-3.05	104.79	111.89
2	F	301	0Y2	C4-N9-C8	2.78	108.66	105.74
2	F	301	0Y2	CBG-CBC-CBB	-2.77	129.73	133.85
2	F	301	0Y2	C2-N3-C4	2.71	118.44	111.83
2	D	301	0Y2	C2-N1-C6	2.59	122.98	118.73
2	E	301	0Y2	O-C-OXT	-2.54	118.32	124.08
2	F	301	0Y2	C4-C5-N7	-2.53	107.69	110.58
2	E	301	0Y2	CB-CAS-NAR	2.53	120.34	113.69
2	D	301	0Y2	CBA-CAZ-NAR	2.52	123.34	113.12
2	D	301	0Y2	C5-N7-C8	2.50	107.38	103.45
2	D	301	0Y2	O-C-OXT	-2.47	118.47	124.08
2	D	301	0Y2	C2-N3-C4	2.47	117.86	111.83
2	E	301	0Y2	N9-C8-N7	-2.44	110.47	113.94
2	F	301	0Y2	CB-CAS-NAR	2.42	120.06	113.69
2	D	301	0Y2	CAS-NAR-CAF	-2.41	106.27	111.89
2	F	301	0Y2	CAC-CAB-CAA	2.38	107.22	102.61
2	F	301	0Y2	CAS-NAR-CAZ	-2.38	105.77	111.44
2	D	301	0Y2	C4-C5-N7	-2.35	107.90	110.58
2	D	301	0Y2	CBG-CBC-CBB	-2.27	130.47	133.85
2	E	301	0Y2	CBG-CBC-CBD	2.22	121.29	118.77
2	D	301	0Y2	N9-C8-N7	-2.21	110.81	113.94
2	F	301	0Y2	C5-C6-N6	-2.18	117.89	123.29
2	F	301	0Y2	N6-C6-N1	2.17	123.20	118.38
2	D	301	0Y2	CBB-CBF-S1E	-2.05	111.44	113.75
2	E	301	0Y2	CAC-CAB-CAA	2.04	106.55	102.61

There are no chirality outliers.

All (31) torsion outliers are listed below:

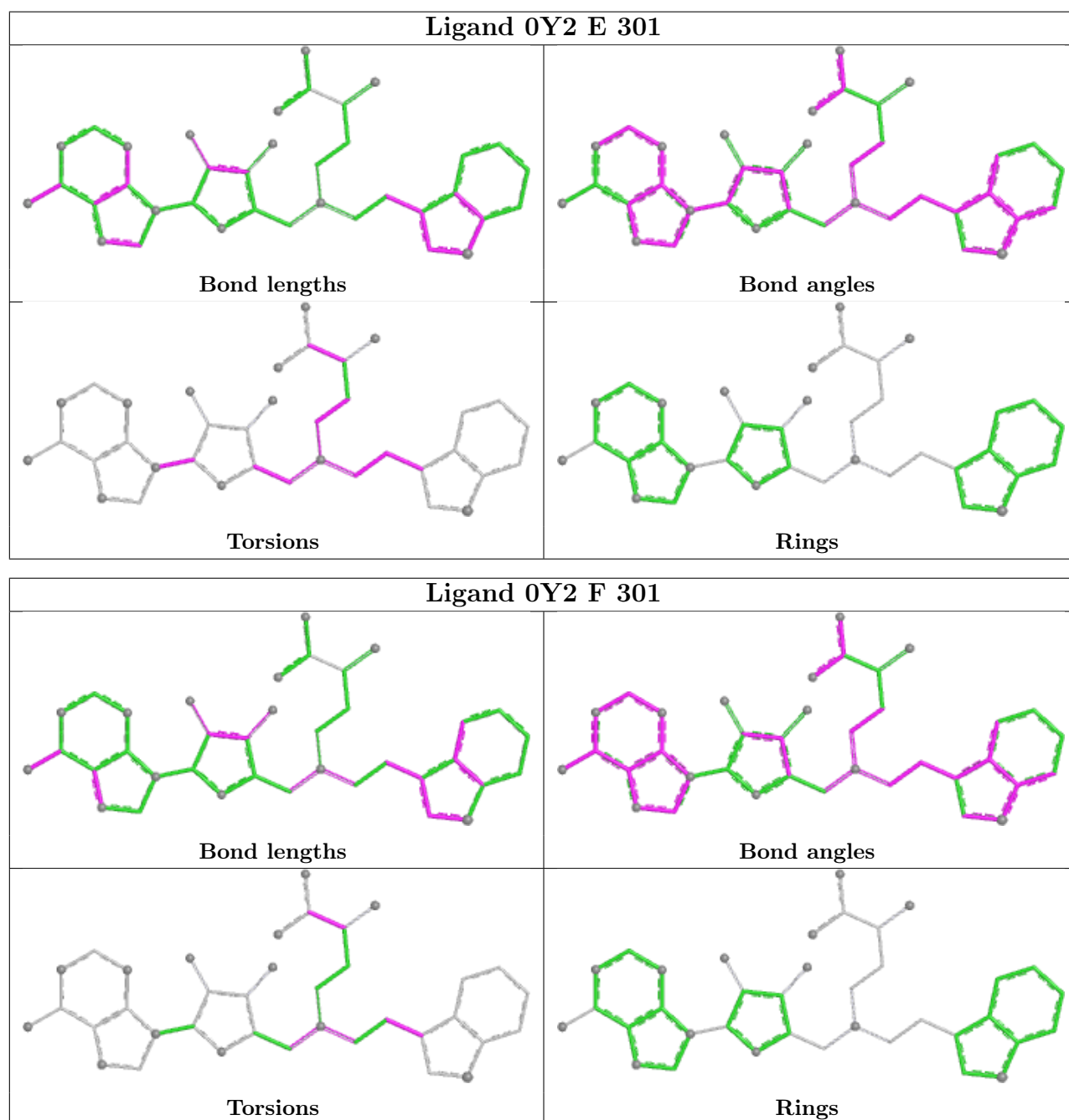
Mol	Chain	Res	Type	Atoms
2	D	301	0Y2	C-CA-CB-CAS
2	D	301	0Y2	OAE-CAA-CAF-NAR
2	D	301	0Y2	CAB-CAA-CAF-NAR
2	E	301	0Y2	NAR-CAZ-CBA-CBB
2	E	301	0Y2	NAR-CAS-CB-CA
2	E	301	0Y2	OXT-C-CA-N
2	E	301	0Y2	OAE-CAA-CAF-NAR
2	E	301	0Y2	CAB-CAA-CAF-NAR
2	F	301	0Y2	CAZ-CBA-CBB-CBC
2	F	301	0Y2	CAZ-CBA-CBB-CBF
2	F	301	0Y2	CAA-CAF-NAR-CAZ
2	D	301	0Y2	CBA-CAZ-NAR-CAF
2	E	301	0Y2	CB-CAS-NAR-CAZ
2	E	301	0Y2	O-C-CA-N
2	F	301	0Y2	O-C-CA-N
2	E	301	0Y2	CBA-CAZ-NAR-CAS
2	E	301	0Y2	CAA-CAF-NAR-CAZ
2	E	301	0Y2	CBA-CAZ-NAR-CAF
2	F	301	0Y2	CBA-CAZ-NAR-CAS
2	F	301	0Y2	OXT-C-CA-N
2	D	301	0Y2	CB-CAS-NAR-CAZ
2	D	301	0Y2	NAR-CAS-CB-CA
2	E	301	0Y2	CAC-CAD-N9-C8
2	D	301	0Y2	CBA-CAZ-NAR-CAS
2	D	301	0Y2	CB-CAS-NAR-CAF
2	D	301	0Y2	O-C-CA-N
2	E	301	0Y2	OAE-CAD-N9-C8
2	E	301	0Y2	CAZ-CBA-CBB-CBF
2	E	301	0Y2	CAZ-CBA-CBB-CBC
2	D	301	0Y2	CAC-CAD-N9-C8
2	D	301	0Y2	OXT-C-CA-N

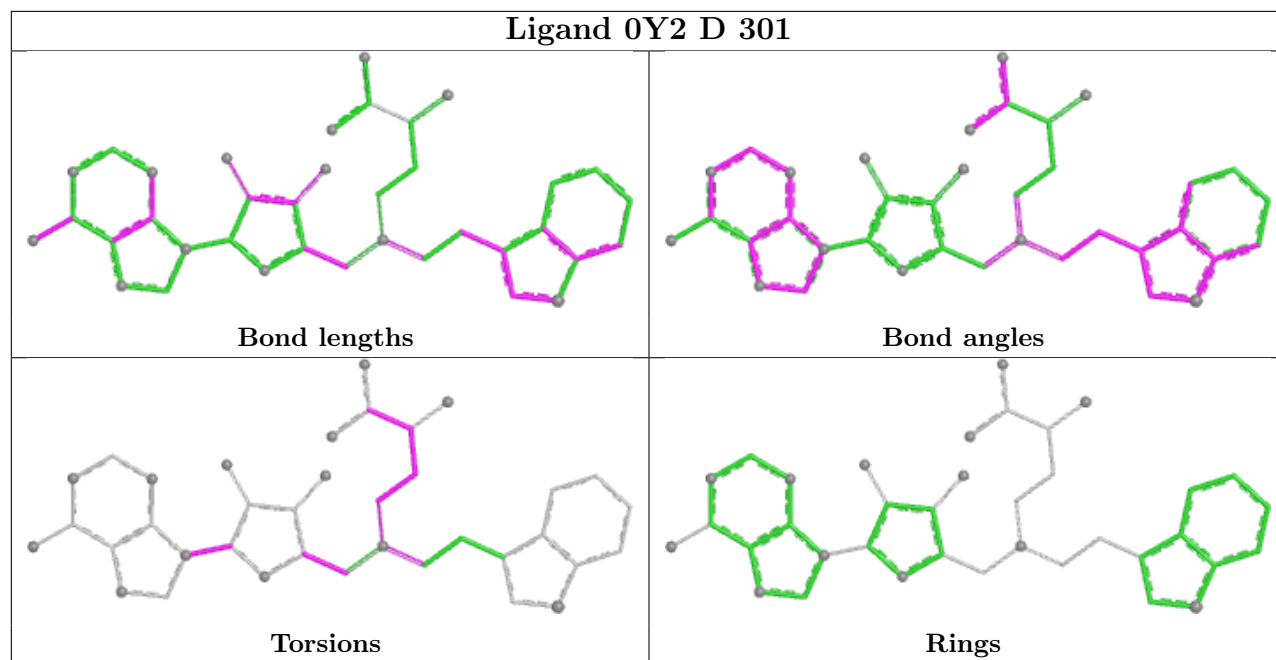
There are no ring outliers.

3 monomers are involved in 15 short contacts:

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

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	D	243/278 (87%)	-0.05	11 (4%) 38 34	35, 49, 77, 107	0
1	E	244/278 (87%)	0.16	14 (5%) 29 26	39, 56, 87, 104	0
1	F	244/278 (87%)	0.14	12 (4%) 35 31	36, 56, 79, 103	1 (0%)
All	All	731/834 (87%)	0.08	37 (5%) 33 29	35, 54, 83, 107	1 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	11	ALA	5.8
1	E	10	TRP	5.7
1	F	187	LEU	5.4
1	F	11	ALA	5.4
1	E	247	VAL	5.3
1	F	14	LYS	5.1
1	F	12	GLY	4.8
1	D	12	GLY	4.7
1	D	10	TRP	4.4
1	E	14	LYS	4.3
1	D	13	GLY	4.3
1	E	3	LYS	3.7
1	D	261	VAL	3.6
1	D	139	LYS	3.5
1	F	15	TYR	3.4
1	F	10	TRP	3.3
1	F	247	VAL	3.2
1	D	14	LYS	3.2
1	E	139	LYS	3.2
1	E	261	VAL	3.2
1	F	139	LYS	3.1
1	E	15	TYR	2.9
1	E	12	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	13	GLY	2.7
1	E	17	LEU	2.6
1	D	15	TYR	2.5
1	F	199	ASN	2.4
1	E	11	ALA	2.3
1	E	260	LYS	2.2
1	D	16	PRO	2.2
1	E	187	LEU	2.2
1	F	261	VAL	2.2
1	D	102	GLN	2.1
1	E	13	GLY	2.0
1	D	246	LYS	2.0
1	F	200	SER	2.0
1	E	137	ARG	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](#)

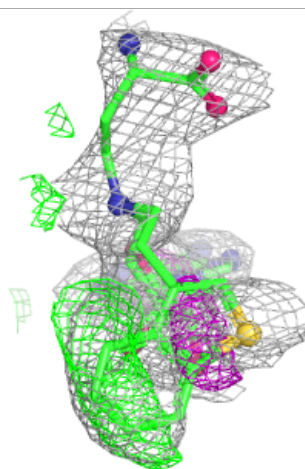
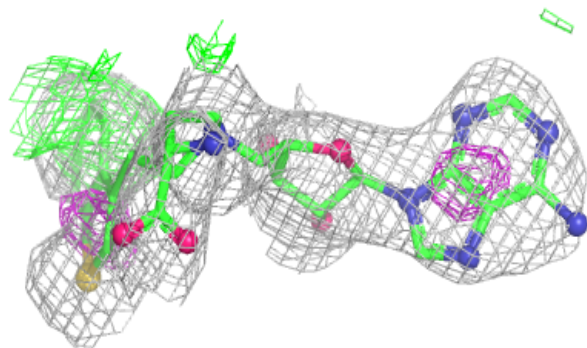
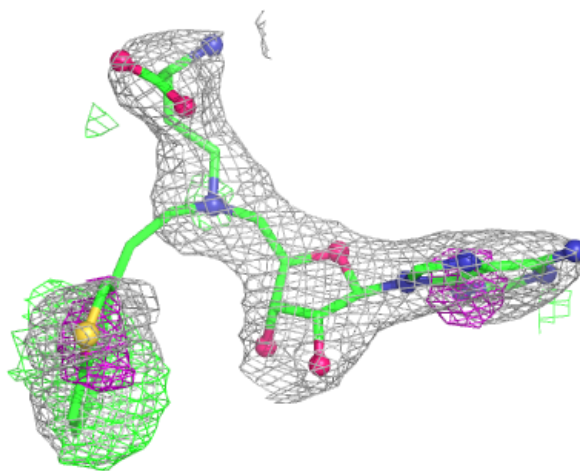
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	0Y2	E	301	37/37	0.82	0.21	77,94,107,108	0
2	0Y2	F	301	37/37	0.88	0.16	72,83,90,92	0
2	0Y2	D	301	37/37	0.89	0.16	76,84,91,92	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.

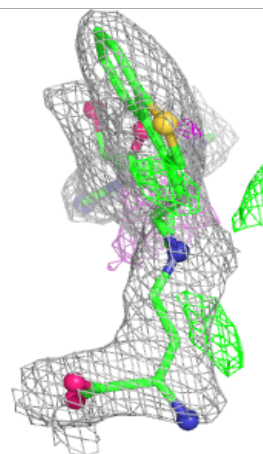
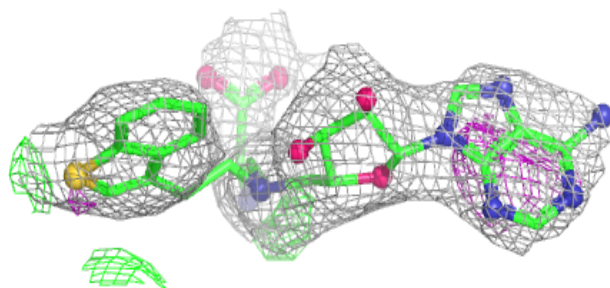
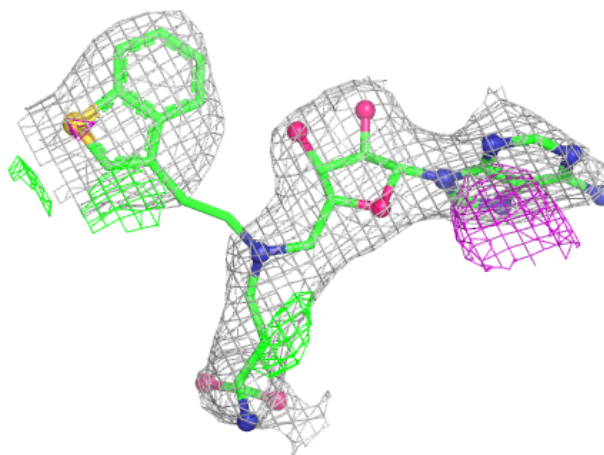
Electron density around 0Y2 E 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



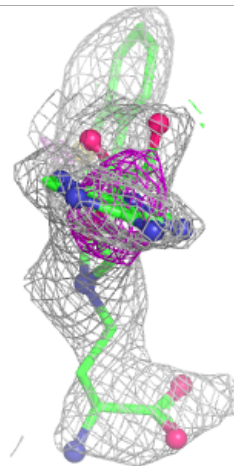
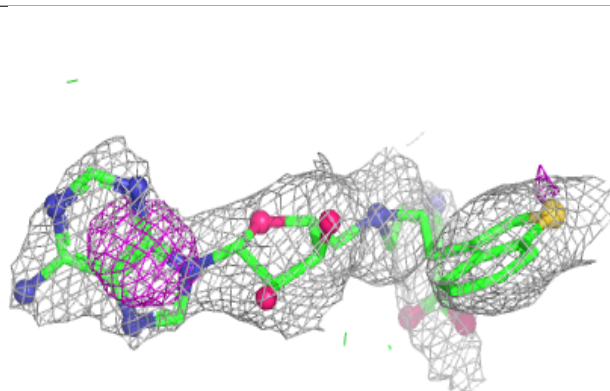
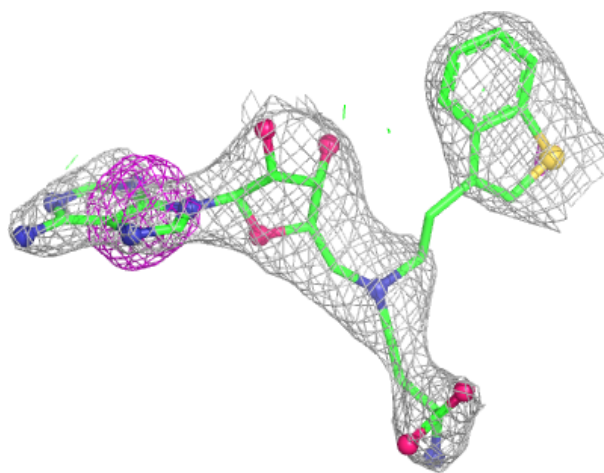
Electron density around 0Y2 F 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 0Y2 D 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.