



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:13 AM UTC

PDB ID : 5A0W / pdb_00005a0w
Title : THE CRYSTAL STRUCTURE OF I-DMOI E117A IN COMPLEX WITH ITS TARGET DNA AND IN THE PRESENCE OF 2MM MN
Authors : Molina, R.; Besker, N.; Prieto, J.; Montoya, G.; D'Abramo, M.
Deposited on : 2015-04-23
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:

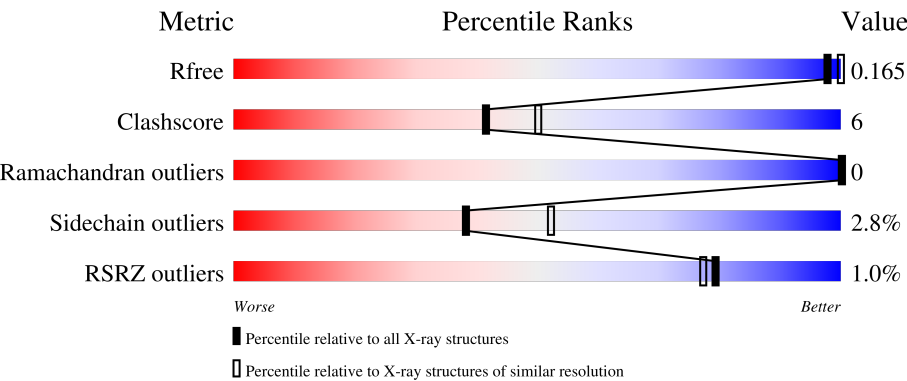
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

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	199	75%16%9%
1	D	199	% 82%13%
1	G	199	2% 80%9%11%
2	B	25	4% 72%28%
2	E	25	64%36%

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Mol	Chain	Length	Quality of chain
2	H	25	56% 40%
3	C	25	92% 8%
3	F	25	64% 36%
3	I	25	72% 28%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8087 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 HOMING ENDONUCLEASE I-DMOI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	0	6	0
			1547	998	287	259	3			
1	D	191	Total	C	N	O	S	2	3	0
			1584	1025	287	269	3			
1	G	178	Total	C	N	O	S	0	3	0
			1487	961	272	251	3			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ALA	-	expression tag	UNP P21505
A	189	ALA	-	expression tag	UNP P21505
A	190	ALA	-	expression tag	UNP P21505
A	191	ALA	-	expression tag	UNP P21505
A	192	LEU	-	expression tag	UNP P21505
A	193	GLU	-	expression tag	UNP P21505
A	194	HIS	-	expression tag	UNP P21505
A	195	HIS	-	expression tag	UNP P21505
A	196	HIS	-	expression tag	UNP P21505
A	197	HIS	-	expression tag	UNP P21505
A	198	HIS	-	expression tag	UNP P21505
A	199	HIS	-	expression tag	UNP P21505
A	117	ALA	GLU	conflict	UNP P21505
D	1	ALA	-	expression tag	UNP P21505
D	189	ALA	-	expression tag	UNP P21505
D	190	ALA	-	expression tag	UNP P21505
D	191	ALA	-	expression tag	UNP P21505
D	192	LEU	-	expression tag	UNP P21505
D	193	GLU	-	expression tag	UNP P21505
D	194	HIS	-	expression tag	UNP P21505
D	195	HIS	-	expression tag	UNP P21505
D	196	HIS	-	expression tag	UNP P21505
D	197	HIS	-	expression tag	UNP P21505

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Chain	Residue	Modelled	Actual	Comment	Reference
D	198	HIS	-	expression tag	UNP P21505
D	199	HIS	-	expression tag	UNP P21505
D	117	ALA	GLU	conflict	UNP P21505
G	1	ALA	-	expression tag	UNP P21505
G	189	ALA	-	expression tag	UNP P21505
G	190	ALA	-	expression tag	UNP P21505
G	191	ALA	-	expression tag	UNP P21505
G	192	LEU	-	expression tag	UNP P21505
G	193	GLU	-	expression tag	UNP P21505
G	194	HIS	-	expression tag	UNP P21505
G	195	HIS	-	expression tag	UNP P21505
G	196	HIS	-	expression tag	UNP P21505
G	197	HIS	-	expression tag	UNP P21505
G	198	HIS	-	expression tag	UNP P21505
G	199	HIS	-	expression tag	UNP P21505
G	117	ALA	GLU	conflict	UNP P21505

- Molecule 2 is a DNA chain called 25MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	25	Total	C	N	O	P	0	0	0
			511	242	94	151	24			
2	E	25	Total	C	N	O	P	0	0	0
			511	242	94	151	24			
2	H	25	Total	C	N	O	P	0	0	0
			511	242	94	151	24			

- Molecule 3 is a DNA chain called 25MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	25	Total	C	N	O	P	0	0	0
			508	240	99	145	24			
3	F	25	Total	C	N	O	P	0	0	0
			508	240	99	145	24			
3	I	25	Total	C	N	O	P	0	0	0
			508	240	99	145	24			

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

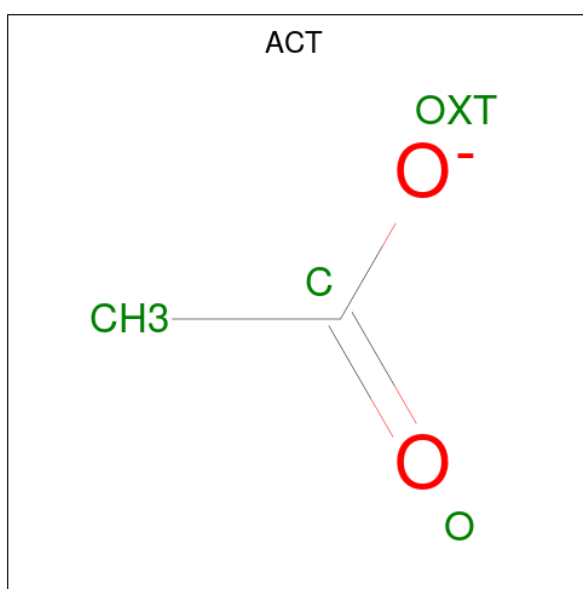
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Cl	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Cl	0	0
			1	1		
4	F	1	Total	Cl	0	0
			1	1		
4	G	2	Total	Cl	0	0
			2	2		
4	H	1	Total	Cl	0	0
			1	1		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		
5	I	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total 1	Mn 1	0	0
6	B	1	Total 1	Mn 1	0	0
6	D	1	Total 1	Mn 1	0	0
6	F	1	Total 1	Mn 1	0	0
6	G	1	Total 1	Mn 1	0	0
6	H	1	Total 1	Mn 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	108	Total 108	O 108	0	0
7	B	43	Total 43	O 43	0	0
7	C	12	Total 12	O 12	0	0
7	D	74	Total 74	O 74	0	0
7	E	21	Total 21	O 21	0	0
7	F	9	Total 9	O 9	0	0
7	G	79	Total 79	O 79	0	0
7	H	19	Total 19	O 19	0	0
7	I	14	Total 14	O 14	0	0



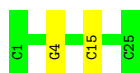
- Molecule 2: 25MER

Chain H:
56% 40% .



- Molecule 3: 25MER

Chain C:
92% 8%



- Molecule 3: 25MER

Chain F:
64% 36%



- Molecule 3: 25MER

Chain I:
72% 28%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.93Å 69.77Å 106.98Å 90.00° 120.06° 90.00°	Depositor
Resolution (Å)	29.22 – 2.20 29.22 – 2.20	Depositor EDS
% Data completeness (in resolution range)	88.5 (29.22-2.20) 96.5 (29.22-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.20Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.171 , 0.208 0.160 , 0.165	Depositor DCC
R_{free} test set	3192 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.526	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 63.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for -h-l,k,h 0.024 for l,k,-h-l 0.017 for l,-k,h 0.018 for h,-k,-h-l 0.015 for -h-l,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8087	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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: CL, ACT, MN

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	0.47	0/1590	0.71	1/2136 (0.0%)
1	D	0.47	0/1624	0.75	0/2188
1	G	0.46	0/1522	0.74	0/2044
2	B	0.28	0/572	0.98	0/882
2	E	0.29	0/572	0.91	0/882
2	H	0.29	0/572	0.95	2/882 (0.2%)
3	C	0.28	0/570	0.91	0/877
3	F	0.29	0/570	0.94	0/877
3	I	0.31	0/570	0.92	0/877
All	All	0.40	0/8162	0.83	3/11645 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	LYS	N-CA-C	-5.80	104.96	111.28
2	H	13	DA	N9-C1'-C2'	5.21	121.32	113.50
2	H	16	DT	N1-C1'-C2'	5.14	121.21	113.50

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	1547	0	1642	15	0
1	D	1584	0	1666	21	0
1	G	1487	0	1571	14	0
2	B	511	0	282	8	0
2	E	511	0	282	9	0
2	H	511	0	282	16	0
3	C	508	0	279	3	0
3	F	508	0	279	7	0
3	I	508	0	279	12	0
4	A	2	0	0	0	0
4	D	1	0	0	0	0
4	F	1	0	0	0	0
4	G	2	0	0	0	0
4	H	1	0	0	0	0
5	A	4	0	3	0	0
5	C	8	0	6	1	0
5	F	4	0	3	0	0
5	I	4	0	3	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	D	1	0	0	0	0
6	F	1	0	0	0	0
6	G	1	0	0	0	0
6	H	1	0	0	0	0
7	A	108	0	0	1	0
7	B	43	0	0	0	0
7	C	12	0	0	0	0
7	D	74	0	0	2	0
7	E	21	0	0	0	0
7	F	9	0	0	0	0
7	G	79	0	0	3	0
7	H	19	0	0	0	0
7	I	14	0	0	1	0
All	All	8087	0	6577	90	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 6.

All (90) 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)
2:H:11:DG:H22	3:I:15:DC:H42	1.07	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:11:DG:N2	3:I:15:DC:N3	2.15	0.94
2:E:11:DG:N2	3:F:15:DC:N3	2.18	0.90
2:H:13:DA:H62	3:I:13:DT:H3	1.28	0.79
2:H:11:DG:H22	3:I:15:DC:N4	1.81	0.79
3:F:2:DG:H2''	3:F:3:DC:H5''	1.65	0.77
3:I:7:DG:H2''	3:I:8:DG:H5''	1.71	0.73
1:D:19[B]:ILE:HD11	1:D:95:LEU:HD13	1.73	0.70
1:D:60[B]:ILE:HD11	1:D:87:LEU:HD22	1.72	0.70
1:D:152[B]:HIS:HD2	7:D:2066:HOH:O	1.75	0.70
2:H:11:DG:N2	3:I:15:DC:H42	1.86	0.69
1:D:155:ASP:HB3	1:D:160:VAL:HG13	1.73	0.69
2:E:12:DT:C6	2:E:12:DT:H5'	2.28	0.68
2:B:11:DG:N2	3:C:15:DC:C2	2.62	0.68
2:E:11:DG:H2''	2:E:12:DT:OP2	1.92	0.68
2:B:11:DG:N2	3:C:15:DC:N3	2.43	0.66
1:D:183:ASN:HB2	1:D:184:PRO:HD2	1.78	0.66
2:H:24:DC:H2''	2:H:25:DG:C8	2.32	0.64
1:A:42:GLN:HG2	1:A:47:LEU:HD23	1.82	0.61
2:E:12:DT:H5'	2:E:12:DT:H6	1.65	0.61
1:D:172:LYS:O	1:D:176:THR:HB	2.00	0.61
2:H:11:DG:N2	3:I:15:DC:N4	2.48	0.60
1:G:45:GLU:HB2	1:G:78:TYR:CE1	2.37	0.59
2:H:11:DG:N2	3:I:15:DC:C4	2.69	0.59
3:I:25:DC:H5''	7:I:2014:HOH:O	2.03	0.58
1:G:80:LEU:C	1:G:80:LEU:HD23	2.30	0.57
1:A:185:LEU:HD12	1:A:186:PRO:HD2	1.86	0.57
2:E:19:DC:H2'	2:E:20:DG:C8	2.41	0.56
1:D:42:GLN:HG2	1:D:47:LEU:HD23	1.90	0.54
2:B:19:DC:H2'	2:B:20:DG:C8	2.43	0.54
1:D:181:HIS:HE2	1:D:193:GLU:CD	2.15	0.54
1:D:19[B]:ILE:HD12	1:D:109:PHE:CZ	2.43	0.53
1:A:152[B]:HIS:HD2	7:A:2082:HOH:O	1.92	0.52
1:D:33:ARG:NH1	2:E:21:DG:O6	2.41	0.52
1:D:183:ASN:HB2	1:D:184:PRO:CD	2.39	0.52
3:F:19:DG:H2''	3:F:20:DC:H5'	1.92	0.52
2:E:11:DG:H1'	2:E:12:DT:H5''	1.91	0.51
1:A:5:GLU:C	1:A:7:VAL:H	2.19	0.51
2:B:18:DC:H5''	2:B:18:DC:H6	1.75	0.50
3:F:15:DC:H2'	3:F:16:DC:O4'	2.11	0.50
3:F:23:DG:H2''	3:F:24:DG:O5'	2.11	0.50
1:D:33:ARG:HD2	3:F:4:DG:O6	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:33:ARG:NH1	2:H:21:DG:O6	2.45	0.50
1:D:120:LYS:HE2	7:D:2056:HOH:O	2.11	0.50
1:G:102:ASN:O	1:G:106:GLN:HG3	2.12	0.49
1:G:45:GLU:HB2	1:G:78:TYR:CZ	2.47	0.49
3:I:15:DC:H2'	3:I:16:DC:O4'	2.14	0.48
2:H:18:DC:H2'	2:H:19:DC:C6	2.49	0.48
1:A:49:LYS:HG2	1:A:71:ILE:HD13	1.97	0.47
2:B:18:DC:H2'	2:B:19:DC:H6	1.79	0.47
1:G:45:GLU:HG2	1:G:49:LYS:HD2	1.96	0.46
1:A:68:LYS:HE2	1:A:68:LYS:HB3	1.75	0.46
2:B:18:DC:H2'	2:B:19:DC:C6	2.50	0.46
2:H:18:DC:H2''	2:H:19:DC:H5'	1.97	0.46
3:I:8:DG:H2''	3:I:9:DA:O4'	2.15	0.46
2:H:19:DC:H2'	2:H:20:DG:C8	2.51	0.46
1:D:45:GLU:HB2	1:D:78:TYR:CE1	2.52	0.45
1:A:45:GLU:HB3	1:A:78:TYR:CE1	2.51	0.45
2:B:13:DA:C2	2:B:14:DA:C4	3.05	0.45
1:D:6:ASN:O	1:D:10:ILE:HG12	2.17	0.45
1:G:22:GLY:HA3	1:G:39:VAL:O	2.17	0.45
2:E:24:DC:H2''	2:E:25:DG:C8	2.52	0.44
3:C:4:DG:H21	5:C:1027:ACT:H3	1.81	0.44
2:E:8:DC:H2''	2:E:9:DG:C8	2.53	0.44
7:G:2006:HOH:O	5:I:1026:ACT:H3	2.18	0.44
2:H:18:DC:H2'	2:H:19:DC:H6	1.83	0.44
1:A:53:ALA:HB3	1:A:54:PRO:HD3	2.00	0.44
1:G:5:GLU:N	7:G:2001:HOH:O	2.50	0.44
1:D:163:LEU:HD23	1:D:163:LEU:C	2.44	0.43
1:G:142:LEU:HD23	1:G:142:LEU:HA	1.83	0.43
1:D:10:ILE:HD13	1:D:184:PRO:CG	2.49	0.43
1:G:163:LEU:C	1:G:163:LEU:HD23	2.44	0.43
1:G:178:LEU:HD23	1:G:178:LEU:HA	1.76	0.43
2:H:21:DG:H2''	2:H:22:DC:C5'	2.49	0.43
1:A:163:LEU:C	1:A:163:LEU:HD23	2.43	0.43
1:A:93:ASN:O	1:A:97[B]:ARG:HG3	2.19	0.42
1:A:183:ASN:N	1:A:184:PRO:HD3	2.35	0.42
1:G:120:LYS:HG2	7:G:2048:HOH:O	2.19	0.42
1:D:128:TRP:CZ3	3:F:16:DC:H2'	2.54	0.42
2:H:13:DA:N6	3:I:13:DT:H3	2.07	0.42
1:D:19[B]:ILE:HD12	1:D:109:PHE:CE1	2.55	0.42
1:D:22:GLY:HA3	1:D:39:VAL:O	2.19	0.42
1:D:19[B]:ILE:HD12	1:D:109:PHE:HZ	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:170:ARG:O	1:G:174:VAL:HG23	2.20	0.41
1:A:132:LYS:HD2	1:A:161:TYR:CE1	2.55	0.41
1:A:66:LYS:O	1:A:66:LYS:HG2	2.20	0.40
1:A:155:ASP:HB3	1:A:160:VAL:HB	2.02	0.40
2:H:13:DA:C8	2:H:14:DA:N9	2.89	0.40
1:A:26:LYS:NZ	1:A:28:LYS:NZ	2.70	0.40
2:B:22:DC:H2'	1:G:97:ARG:NH2	2.37	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	186/199 (94%)	181 (97%)	5 (3%)	0	100	100
1	D	192/199 (96%)	188 (98%)	4 (2%)	0	100	100
1	G	179/199 (90%)	176 (98%)	3 (2%)	0	100	100
All	All	557/597 (93%)	545 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	170/177 (96%)	162 (95%)	8 (5%)	23	31
1	D	173/177 (98%)	169 (98%)	4 (2%)	44	59
1	G	162/177 (92%)	160 (99%)	2 (1%)	63	78
All	All	505/531 (95%)	491 (97%)	14 (3%)	38	52

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	7	VAL
1	A	27	LEU
1	A	48	ILE
1	A	66	LYS
1	A	165	ILE
1	A	167	LEU
1	A	177	ILE
1	D	48	ILE
1	D	160	VAL
1	D	170	ARG
1	D	183	ASN
1	G	5	GLU
1	G	48	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	57	GLN
1	D	42	GLN
1	G	70	GLN
1	G	156	HIS
1	G	158	HIS
1	G	181	HIS

5.3.3 RNA ⓘ

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 18 ligands modelled in this entry, 13 are monoatomic - leaving 5 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	$\# Z > 2$
5	ACT	C	1026	-	3,3,3	0.77	0	3,3,3	1.35	0
5	ACT	I	1026	-	3,3,3	0.74	0	3,3,3	1.40	0
5	ACT	F	1027	-	3,3,3	0.81	0	3,3,3	1.29	0
5	ACT	C	1027	-	3,3,3	0.82	0	3,3,3	1.37	0
5	ACT	A	1188	-	3,3,3	0.77	0	3,3,3	1.33	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	1026	ACT	1	0
5	C	1027	ACT	1	0

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	182/199 (91%)	-0.53	0 100 100	16, 33, 64, 97	9 (4%)
1	D	191/199 (95%)	-0.44	2 (1%) 79 77	20, 37, 66, 90	11 (5%)
1	G	178/199 (89%)	-0.34	4 (2%) 62 59	17, 36, 73, 133	6 (3%)
2	B	25/25 (100%)	-0.70	1 (4%) 42 39	32, 44, 65, 91	0
2	E	25/25 (100%)	-0.40	0 100 100	33, 52, 91, 98	0
2	H	25/25 (100%)	-0.73	0 100 100	35, 51, 63, 71	0
3	C	25/25 (100%)	-0.75	0 100 100	29, 50, 62, 67	0
3	F	25/25 (100%)	-0.40	0 100 100	29, 55, 93, 96	0
3	I	25/25 (100%)	-0.57	0 100 100	34, 52, 70, 73	0
All	All	701/747 (93%)	-0.47	7 (0%) 79 77	16, 39, 73, 133	26 (3%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	182	LEU	5.8
1	G	31	GLY	5.6
1	D	156	HIS	3.1
1	G	181	HIS	2.8
1	D	195	HIS	2.7
2	B	11	DG	2.5
1	G	5	GLU	2.4

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [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	ACT	C	1026	4/4	0.73	0.26	51,62,65,68	0
4	CL	H	1026	1/1	0.77	0.20	86,86,86,86	0
5	ACT	C	1027	4/4	0.78	0.27	57,60,68,73	0
5	ACT	A	1188	4/4	0.79	0.21	59,62,66,69	0
5	ACT	F	1027	4/4	0.81	0.24	73,73,75,75	0
4	CL	G	1300	1/1	0.83	0.23	90,90,90,90	0
4	CL	A	1300	1/1	0.86	0.17	86,86,86,86	0
5	ACT	I	1026	4/4	0.88	0.18	54,66,68,70	0
4	CL	F	1026	1/1	0.90	0.16	78,78,78,78	0
4	CL	G	1301	1/1	0.91	0.20	89,89,89,89	0
4	CL	A	1187	1/1	0.92	0.15	78,78,78,78	0
4	CL	D	1196	1/1	0.94	0.18	72,72,72,72	0
6	MN	D	1197	1/1	0.99	0.02	31,31,31,31	0
6	MN	F	1028	1/1	0.99	0.03	45,45,45,45	0
6	MN	G	1183	1/1	0.99	0.03	32,32,32,32	0
6	MN	H	1027	1/1	0.99	0.03	43,43,43,43	0
6	MN	A	1189	1/1	1.00	0.02	26,26,26,26	0
6	MN	B	1026	1/1	1.00	0.01	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.