



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:00 PM GMT

PDB ID : 4ZAB
Title : Structure of A. niger Fdc1 in complex with alpha-fluoro cinnamic acid
Authors : Payne, K.A.P.; Leys, D.
Deposited on : 2015-04-13
Resolution : 1.16 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

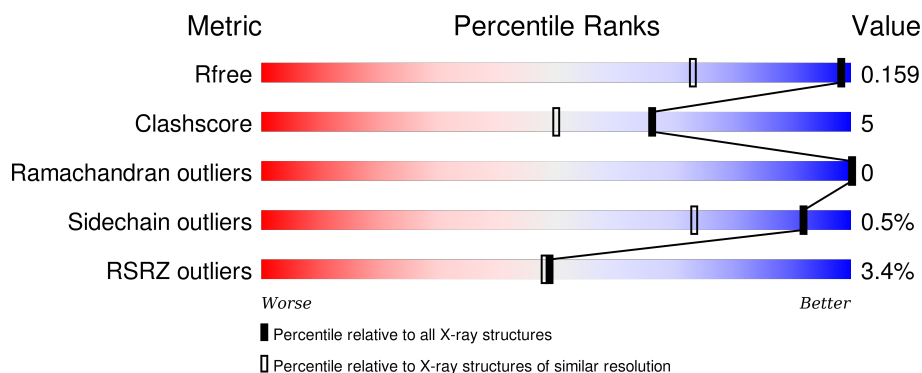
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.16 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	1036 (1.20-1.12)
Clashscore	102246	1109 (1.20-1.12)
Ramachandran outliers	100387	1058 (1.20-1.12)
Sidechain outliers	100360	1058 (1.20-1.12)
RSRZ outliers	91569	1038 (1.20-1.12)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	508	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	4LW	A	606	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4694 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.

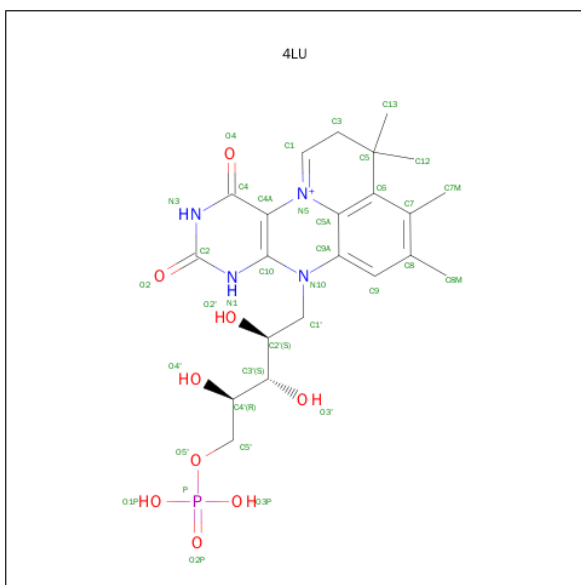
- Molecule 1 is a protein called A niger Fdc1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	500	4122	2608	711	773	30	0	37	1

There are 8 discrepancies between the modelled and reference sequences:

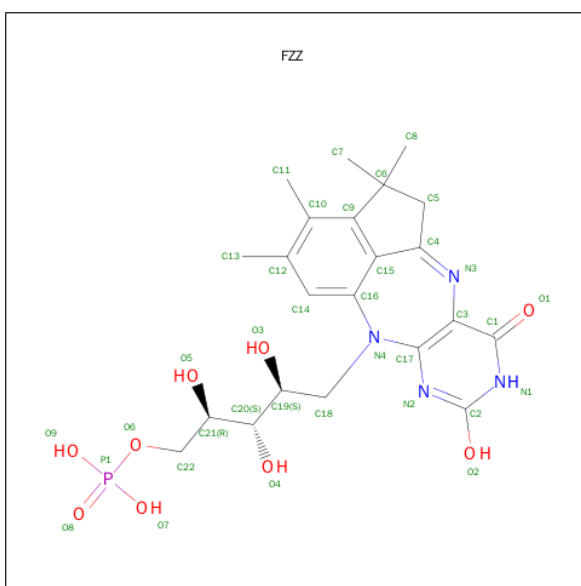
Chain	Residue	Modelled	Actual	Comment	Reference
A	501	LEU	-	expression tag	UNP A2QHE5
A	502	GLU	-	expression tag	UNP A2QHE5
A	503	HIS	-	expression tag	UNP A2QHE5
A	504	HIS	-	expression tag	UNP A2QHE5
A	505	HIS	-	expression tag	UNP A2QHE5
A	506	HIS	-	expression tag	UNP A2QHE5
A	507	HIS	-	expression tag	UNP A2QHE5
A	508	HIS	-	expression tag	UNP A2QHE5

- Molecule 2 is 1-deoxy-5-O-phosphono-1-(3,3,4,5-tetramethyl-9,11-dioxo-2,3,8,9,10,11-hexahydro-7H-quinolino[1,8-fg]pteridin-12-ium-7-yl)-D-ribitol (three-letter code: 4LU) (formula: C₂₂H₃₀N₄O₉P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	1
			36	22	4	9	1		

- Molecule 3 is 1-deoxy-5-O-phosphono-1-[(10aR)-2,2,3,4-tetramethyl-8,10-dioxo-1,2,8,9,10,10a-hexahydro-6H-indeno[1,7-e]pyrimido[4,5-b][1,4]diazepin-6-yl]-D-ribitol (three-letter code: FZZ) (formula: $C_{22}H_{29}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	1
			36	22	4	9	1		

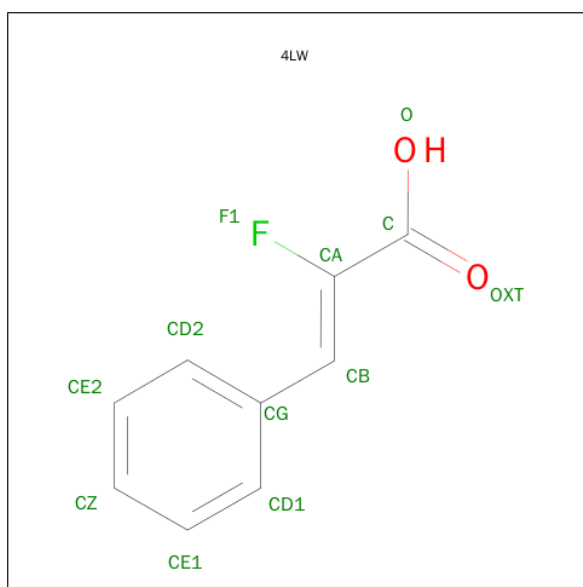
- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

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

- Molecule 6 is (2Z)-2-fluoro-3-phenylprop-2-enoic acid (three-letter code: 4LW) (formula: C₉H₇F O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	F	O	0	0
			12	9	1	2		

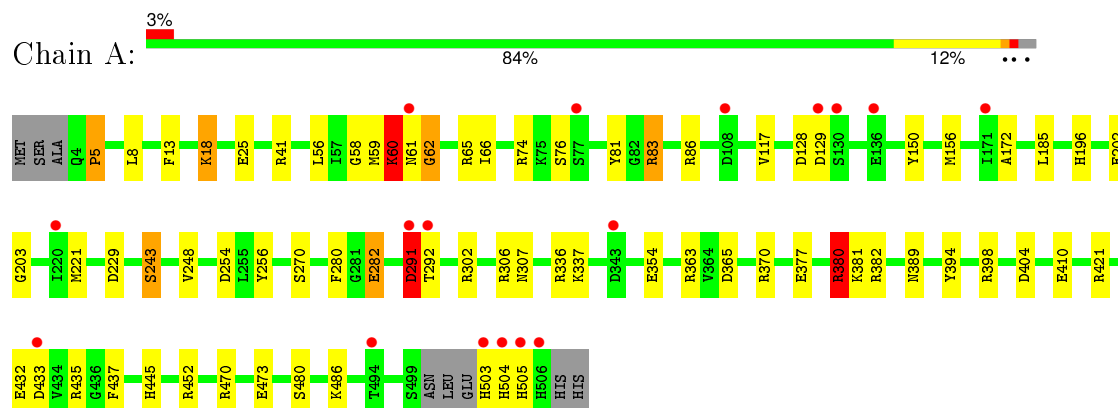
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	485	Total	O	0	0
			485	485		

3 Residue-property plots [i](#)

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: A niger Fdc1



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	95.91Å 64.20Å 87.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.28 – 1.16 32.38 – 1.16	Depositor EDS
% Data completeness (in resolution range)	100.0 (32.28-1.16) 98.7 (32.38-1.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 1.16Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.146 , 0.154 0.150 , 0.159	Depositor DCC
R_{free} test set	9183 reflections (5.23%)	DCC
Wilson B-factor (Å ²)	7.1	Xtriage
Anisotropy	1.748	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 184718 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4694	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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, 0.375 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: FZZ, K, MN, 4LW, 4LU

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	1.48	46/4223 (1.1%)	1.37	42/5737 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	65	ARG	CZ-NH1	-10.77	1.19	1.33
1	A	480[A]	SER	CB-OG	10.14	1.55	1.42
1	A	480[B]	SER	CB-OG	10.14	1.55	1.42
1	A	380	ARG	CA-CB	9.86	1.75	1.53
1	A	389	ASN	CG-ND2	-9.22	1.09	1.32
1	A	370	ARG	CZ-NH2	9.06	1.44	1.33
1	A	337	LYS	CE-NZ	-9.03	1.26	1.49
1	A	291	ASP	CB-CG	-8.17	1.34	1.51
1	A	410	GLU	CG-CD	7.57	1.63	1.51
1	A	282	GLU	CD-OE2	7.56	1.33	1.25
1	A	65	ARG	NE-CZ	7.42	1.42	1.33
1	A	202	GLU	CD-OE1	7.01	1.33	1.25
1	A	25	GLU	CD-OE2	-7.00	1.18	1.25
1	A	156	MET	CB-CG	-6.93	1.29	1.51
1	A	282	GLU	CD-OE1	6.60	1.32	1.25
1	A	202	GLU	CD-OE2	6.55	1.32	1.25
1	A	473	GLU	CA-CB	6.46	1.68	1.53
1	A	62[A]	GLY	CA-C	6.21	1.61	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	62[B]	GLY	CA-C	6.21	1.61	1.51
1	A	354	GLU	CD-OE2	6.01	1.32	1.25
1	A	410	GLU	CD-OE1	5.89	1.32	1.25
1	A	5	PRO	CB-CG	-5.87	1.20	1.50
1	A	336	ARG	CZ-NH1	-5.84	1.25	1.33
1	A	18[A]	LYS	CE-NZ	5.80	1.63	1.49
1	A	18[B]	LYS	CE-NZ	5.80	1.63	1.49
1	A	221[A]	MET	CG-SD	-5.77	1.66	1.81
1	A	221[B]	MET	CG-SD	-5.77	1.66	1.81
1	A	377[A]	GLU	CD-OE1	5.64	1.31	1.25
1	A	377[B]	GLU	CD-OE1	5.64	1.31	1.25
1	A	486	LYS	CD-CE	-5.62	1.37	1.51
1	A	370	ARG	CD-NE	-5.53	1.37	1.46
1	A	473	GLU	CD-OE1	5.52	1.31	1.25
1	A	254	ASP	CB-CG	5.49	1.63	1.51
1	A	307	ASN	C-O	-5.45	1.13	1.23
1	A	256	TYR	CZ-OH	5.42	1.47	1.37
1	A	18[A]	LYS	CD-CE	5.38	1.64	1.51
1	A	18[B]	LYS	CD-CE	5.38	1.64	1.51
1	A	370	ARG	CG-CD	5.37	1.65	1.51
1	A	243	SER	CB-OG	5.32	1.49	1.42
1	A	81	TYR	CD1-CE1	5.29	1.47	1.39
1	A	398	ARG	CZ-NH1	-5.23	1.26	1.33
1	A	410	GLU	CD-OE2	5.16	1.31	1.25
1	A	433	ASP	C-O	5.16	1.33	1.23
1	A	203	GLY	CA-C	-5.11	1.43	1.51
1	A	363	ARG	CD-NE	-5.08	1.37	1.46
1	A	432	GLU	CD-OE1	-5.00	1.20	1.25

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	ARG	NE-CZ-NH2	-14.92	112.84	120.30
1	A	404	ASP	CB-CG-OD2	13.43	130.39	118.30
1	A	398	ARG	NE-CZ-NH2	-12.13	114.24	120.30
1	A	435	ARG	NE-CZ-NH2	-11.50	114.55	120.30
1	A	65	ARG	NE-CZ-NH2	-10.82	114.89	120.30
1	A	86	ARG	NE-CZ-NH2	-9.71	115.45	120.30
1	A	363	ARG	NE-CZ-NH2	-9.55	115.53	120.30
1	A	60[A]	LYS	O-C-N	-8.83	108.58	122.70
1	A	60[B]	LYS	O-C-N	-8.83	108.58	122.70
1	A	380	ARG	NE-CZ-NH1	8.76	124.68	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	291	ASP	CB-CA-C	-8.37	93.67	110.40
1	A	380	ARG	CG-CD-NE	8.33	129.30	111.80
1	A	398	ARG	NE-CZ-NH1	8.31	124.46	120.30
1	A	337	LYS	CD-CE-NZ	8.00	130.11	111.70
1	A	156	MET	CA-CB-CG	7.72	126.43	113.30
1	A	306	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	A	377[A]	GLU	CG-CD-OE2	-7.23	103.85	118.30
1	A	377[B]	GLU	CG-CD-OE2	-7.23	103.85	118.30
1	A	65	ARG	NH1-CZ-NH2	6.80	126.88	119.40
1	A	185[A]	LEU	CB-CG-CD2	6.80	122.56	111.00
1	A	185[B]	LEU	CB-CG-CD2	6.80	122.56	111.00
1	A	302	ARG	NE-CZ-NH2	-6.78	116.91	120.30
1	A	382	ARG	NE-CZ-NH2	-6.63	116.99	120.30
1	A	282	GLU	OE1-CD-OE2	-6.51	115.48	123.30
1	A	229	ASP	CB-CG-OD1	6.45	124.11	118.30
1	A	363	ARG	NE-CZ-NH1	6.36	123.48	120.30
1	A	394	TYR	CB-CG-CD1	5.80	124.48	121.00
1	A	202	GLU	OE1-CD-OE2	5.70	130.14	123.30
1	A	382	ARG	CG-CD-NE	-5.64	99.96	111.80
1	A	421	ARG	NE-CZ-NH2	-5.63	117.48	120.30
1	A	83	ARG	CB-CA-C	5.60	121.61	110.40
1	A	365	ASP	CB-CG-OD2	-5.56	113.29	118.30
1	A	421	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	A	437	PHE	CB-CG-CD1	5.49	124.64	120.80
1	A	150	TYR	CB-CG-CD2	5.48	124.28	121.00
1	A	41	ARG	NE-CZ-NH2	5.44	123.02	120.30
1	A	13	PHE	CB-CG-CD2	-5.44	116.99	120.80
1	A	74	ARG	NH1-CZ-NH2	5.40	125.34	119.40
1	A	65	ARG	NE-CZ-NH1	-5.17	117.72	120.30
1	A	404	ASP	CB-CG-OD1	-5.13	113.69	118.30
1	A	404	ASP	CB-CA-C	-5.11	100.18	110.40
1	A	128	ASP	CB-CG-OD2	-5.10	113.71	118.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	243	SER	Mainchain
1	A	291	ASP	Sidechain
1	A	380	ARG	Sidechain
1	A	60[A]	LYS	Mainchain
1	A	83	ARG	Sidechain

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	4122	0	4000	35	12
2	A	36	0	23	7	0
3	A	36	0	20	5	0
4	A	1	0	0	0	0
5	A	2	0	0	0	0
6	A	12	0	6	0	0
7	A	485	0	0	15	14
All	All	4694	0	4049	41	14

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 5.

All (41) 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:380:ARG:CB	1:A:380:ARG:CA	1.75	1.60
1:A:18[B]:LYS:CE	1:A:18[B]:LYS:NZ	1.70	1.54
1:A:381[A]:LYS:HG3	7:A:703:HOH:O	1.53	1.06
1:A:291:ASP:O	7:A:839:HOH:O	1.82	0.96
2:A:601[A]:4LU:H14	3:A:602[A]:FZZ:H5	1.48	0.95
1:A:292:THR:HA	7:A:839:HOH:O	1.69	0.92
2:A:601[A]:4LU:H13	2:A:601[A]:4LU:H14	1.52	0.91
1:A:60[A]:LYS:O	1:A:62[A]:GLY:N	2.11	0.84
1:A:59[A]:MET:SD	7:A:1004:HOH:O	2.40	0.78
2:A:601[A]:4LU:H14	3:A:602[A]:FZZ:C7	2.12	0.77
1:A:380:ARG:CA	1:A:380:ARG:CG	2.63	0.75
1:A:380:ARG:C	1:A:380:ARG:CB	2.54	0.74
1:A:59[A]:MET:CE	7:A:1004:HOH:O	2.35	0.73
1:A:60[A]:LYS:O	1:A:61[A]:ASN:C	2.30	0.70
1:A:60[A]:LYS:C	1:A:62[A]:GLY:N	2.48	0.67
1:A:381[A]:LYS:HE3	7:A:703:HOH:O	1.95	0.64
1:A:60[A]:LYS:C	1:A:62[A]:GLY:H	1.99	0.64
1:A:196[B]:HIS:HE1	1:A:270:SER:O	1.82	0.62
1:A:381[A]:LYS:CG	7:A:703:HOH:O	2.26	0.60
1:A:381[A]:LYS:HE2	7:A:1136:HOH:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:ARG:CB	1:A:380:ARG:N	2.63	0.56
1:A:280:PHE:O	1:A:282:GLU:HG3	2.05	0.56
1:A:129[A]:ASP:O	1:A:129[A]:ASP:CG	2.42	0.56
1:A:59[A]:MET:HE3	7:A:1004:HOH:O	2.01	0.55
2:A:601[A]:4LU:C12	2:A:601[A]:4LU:H14	2.13	0.55
1:A:59[A]:MET:HG2	7:A:720:HOH:O	2.07	0.54
1:A:292:THR:CA	7:A:839:HOH:O	2.41	0.50
1:A:172:ALA:HA	3:A:602[A]:FZZ:C1	2.43	0.49
1:A:58:GLY:O	1:A:59[B]:MET:C	2.48	0.48
2:A:601[A]:4LU:H7	3:A:602[A]:FZZ:H20	1.78	0.48
1:A:196[A]:HIS:HE1	1:A:270:SER:O	1.97	0.48
1:A:196[B]:HIS:CE1	1:A:270:SER:O	2.65	0.48
1:A:445:HIS:HE1	7:A:909:HOH:O	1.99	0.46
1:A:5:PRO:HG2	1:A:8:LEU:HD12	1.96	0.46
1:A:470:ARG:NE	7:A:844:HOH:O	2.48	0.46
1:A:76[B]:SER:HB3	7:A:1034:HOH:O	2.16	0.44
1:A:381[A]:LYS:CE	7:A:703:HOH:O	2.62	0.43
1:A:56[B]:LEU:HD11	1:A:66:ILE:HG13	2.01	0.42
1:A:117:VAL:HG23	1:A:248[B]:VAL:HG22	2.02	0.41
2:A:601[A]:4LU:H17	3:A:602[A]:FZZ:H11	1.63	0.41
1:A:172:ALA:HA	2:A:601[A]:4LU:C4	2.51	0.41

All (14) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:505:HIS:CD2	7:A:800:HOH:O[3_545]	0.55	1.65
1:A:503:HIS:C	7:A:810:HOH:O[3_545]	0.58	1.62
1:A:504:HIS:N	7:A:810:HOH:O[3_545]	1.02	1.18
7:A:711:HOH:O	7:A:711:HOH:O[2_655]	1.18	1.02
1:A:505:HIS:NE2	7:A:800:HOH:O[3_545]	1.26	0.94
1:A:503:HIS:O	7:A:810:HOH:O[3_545]	1.28	0.92
1:A:503:HIS:ND1	7:A:764:HOH:O[3_545]	1.29	0.91
1:A:505:HIS:CG	7:A:800:HOH:O[3_545]	1.31	0.89
7:A:710:HOH:O	7:A:748:HOH:O[2_655]	1.61	0.59
1:A:505:HIS:CE1	7:A:800:HOH:O[3_545]	1.92	0.28
1:A:504:HIS:CA	7:A:810:HOH:O[3_545]	1.94	0.26
1:A:505:HIS:ND1	7:A:800:HOH:O[3_545]	1.95	0.25
1:A:503:HIS:CA	7:A:810:HOH:O[3_545]	2.03	0.17
1:A:503:HIS:CE1	7:A:764:HOH:O[3_545]	2.15	0.05

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	533/508 (105%)	519 (97%)	14 (3%)	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	443/430 (103%)	441 (100%)	2 (0%)	92	73

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	291	ASP
1	A	452	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	$\# Z > 2$
2	4LU	A	601[A]	3,5,4	32,39,39	3.11	15 (46%)	33,62,62	4.35	14 (42%)
3	FZZ	A	602[A]	2	31,39,39	2.77	7 (22%)	35,62,62	1.96	4 (11%)
6	4LW	A	606	-	8,12,12	2.14	3 (37%)	8,15,15	3.03	5 (62%)

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	4LU	A	601[A]	3,5,4	-	0/18/30/30	0/3/4/4
3	FZZ	A	602[A]	2	-	0/16/29/29	0/3/4/4
6	4LW	A	606	-	-	0/4/8/8	0/1/1/1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601[A]	4LU	C5-C6	-6.79	1.48	1.54
2	A	601[A]	4LU	C1'-N10	-6.35	1.41	1.48
3	A	602[A]	FZZ	C18-N4	-5.95	1.42	1.48
2	A	601[A]	4LU	C5A-C9A	-4.44	1.33	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601[A]	4LU	C10-N10	-3.88	1.34	1.39
6	A	606	4LW	CE1-CD1	-3.45	1.32	1.38
2	A	601[A]	4LU	C4-C4A	-3.30	1.34	1.41
3	A	602[A]	FZZ	C19-C20	-3.12	1.47	1.53
6	A	606	4LW	CD2-CG	-2.81	1.34	1.39
2	A	601[A]	4LU	C2'-C3'	-2.49	1.48	1.53
2	A	601[A]	4LU	C2-N3	-2.25	1.33	1.38
2	A	601[A]	4LU	C5'-C4'	2.00	1.54	1.51
3	A	602[A]	FZZ	C7-C6	2.06	1.59	1.53
3	A	602[A]	FZZ	O6-C22	2.07	1.53	1.44
2	A	601[A]	4LU	C2-N1	2.24	1.42	1.38
2	A	601[A]	4LU	C13-C5	2.50	1.61	1.53
3	A	602[A]	FZZ	C14-C12	2.69	1.45	1.37
2	A	601[A]	4LU	C6-C7	2.99	1.47	1.40
6	A	606	4LW	CE2-CD2	3.40	1.45	1.38
3	A	602[A]	FZZ	C3-C17	3.68	1.47	1.40
2	A	601[A]	4LU	C9-C8	3.77	1.48	1.37
2	A	601[A]	4LU	O2'-C2'	4.11	1.52	1.43
2	A	601[A]	4LU	C4-N3	4.96	1.42	1.33
2	A	601[A]	4LU	C9A-N10	8.10	1.50	1.38
3	A	602[A]	FZZ	C5-C4	11.54	1.60	1.50

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602[A]	FZZ	C6-C5-C4	-7.68	96.74	104.28
2	A	601[A]	4LU	C4A-C4-N3	-7.31	113.97	123.52
6	A	606	4LW	CZ-CE2-CD2	-5.50	112.56	120.20
2	A	601[A]	4LU	N3-C2-N1	-5.28	118.79	127.69
3	A	602[A]	FZZ	C3-C1-N1	-4.84	117.20	123.52
6	A	606	4LW	CE1-CD1-CG	-3.60	116.05	120.66
2	A	601[A]	4LU	C4A-N5-C5A	-2.88	115.84	120.96
2	A	601[A]	4LU	C13-C5-C6	-2.82	105.72	111.78
2	A	601[A]	4LU	O2'-C2'-C1'	-2.72	103.22	109.93
2	A	601[A]	4LU	C13-C5-C3	-2.47	104.57	108.93
2	A	601[A]	4LU	O5'-P-O2P	-2.43	100.97	107.08
2	A	601[A]	4LU	O3'-C3'-C2'	-2.30	102.76	108.73
3	A	602[A]	FZZ	O6-P1-O8	-2.29	101.32	107.08
2	A	601[A]	4LU	C9-C8-C7	2.03	121.75	119.26
6	A	606	4LW	CE2-CD2-CG	2.12	123.38	120.66
2	A	601[A]	4LU	C8-C7-C6	2.31	121.38	119.38
6	A	606	4LW	CD1-CG-CD2	2.35	121.14	117.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602[A]	FZZ	C1-C3-C17	3.52	122.19	119.94
6	A	606	4LW	CZ-CE1-CD1	3.92	125.65	120.20
2	A	601[A]	4LU	C12-C5-C6	4.30	121.01	111.78
2	A	601[A]	4LU	C4-N3-C2	8.49	122.24	115.16
2	A	601[A]	4LU	C4-C4A-C10	13.10	128.33	119.94
2	A	601[A]	4LU	C4A-C10-N10	14.30	130.90	120.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601[A]	4LU	7	0
3	A	602[A]	FZZ	5	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	500/508 (98%)	0.27	17 (3%)	49 48	7, 10, 20, 30	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	503	HIS	8.0
1	A	504	HIS	7.2
1	A	505	HIS	4.5
1	A	61[A]	ASN	4.4
1	A	506	HIS	4.0
1	A	108[A]	ASP	3.7
1	A	292	THR	3.6
1	A	433	ASP	2.8
1	A	291	ASP	2.6
1	A	220	ILE	2.5
1	A	130	SER	2.3
1	A	343	ASP	2.2
1	A	129[A]	ASP	2.2
1	A	77	SER	2.1
1	A	136	GLU	2.1
1	A	171	ILE	2.0
1	A	494	THR	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, 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	LLDF	B-factors(Å ²)	Q<0.9
6	4LW	A	606	12/12	0.90	0.15	4.14	13,18,22,23	0
2	4LU	A	601[A]	36/36	0.98	0.10	1.06	6,7,12,18	36
3	FZZ	A	602[A]	36/36	0.98	0.10	0.93	6,8,12,14	36
4	MN	A	603	1/1	1.00	0.07	-0.53	7,7,7,7	0
5	K	A	605	1/1	1.00	0.06	-1.42	7,7,7,7	0
5	K	A	604	1/1	1.00	0.07	-2.27	7,7,7,7	0

6.5 Other polymers [i](#)

There are no such residues in this entry.