



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:42 am GMT

PDB ID : 3CLZ
Title : The set and ring associated (SRA) domain of UHRF1 bound to methylated DNA
Authors : Walker, J.R.; Avvakumov, G.V.; Xue, S.; Dong, A.; Li, Y.; Bountra, C.; Weigelt, J.; Arrowsmith, C.H.; Edwards, A.M.; Bochkarev, A.; Dhe-Paganon, S.; Structural Genomics Consortium (SGC)
Deposited on : 2008-03-20
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

<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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

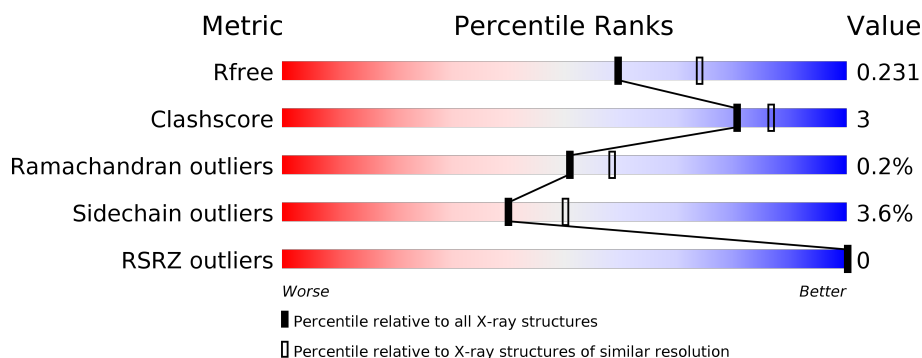
1 Overall quality at a glance i

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.






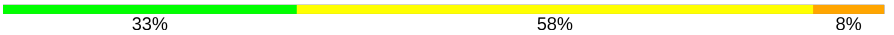
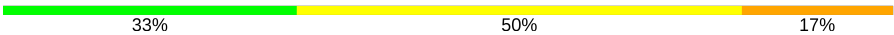
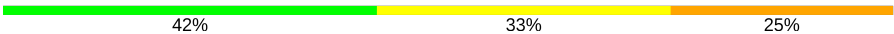
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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 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	212	<div> <div style="width: 86%;"></div> <div style="width: 9%;"></div> <div style="width: 5%;"></div> <div style="width: 0%;"></div> </div> <div>86% 9% .</div>
1	B	212	<div> <div style="width: 90%;"></div> <div style="width: 8%;"></div> <div style="width: 2%;"></div> <div style="width: 0%;"></div> </div> <div>90% 8% .</div>
1	C	212	<div> <div style="width: 88%;"></div> <div style="width: 8%;"></div> <div style="width: 4%;"></div> <div style="width: 0%;"></div> </div> <div>88% 8% .</div>
1	D	212	<div> <div style="width: 88%;"></div> <div style="width: 9%;"></div> <div style="width: 3%;"></div> <div style="width: 0%;"></div> </div> <div>88% 9% ..</div>
2	E	12	<div> <div style="width: 42%;"></div> <div style="width: 50%;"></div> <div style="width: 8%;"></div> <div style="width: 0%;"></div> </div> <div>42% 50% 8%</div>
2	G	12	<div> <div style="width: 42%;"></div> <div style="width: 58%;"></div> <div style="width: 0%;"></div> <div style="width: 0%;"></div> </div> <div>42% 58%</div>

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Mol	Chain	Length	Quality of chain
2	I	12	
2	K	12	
3	F	12	
3	H	12	
3	J	12	
3	L	12	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9090 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 E3 ubiquitin-protein ligase UHRF1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	204	Total	C	N	O	S	0	1	0
			1609	1007	299	299	4			
1	B	210	Total	C	N	O	S	0	1	0
			1646	1028	307	307	4			
1	C	204	Total	C	N	O	S	0	1	0
			1597	1000	294	299	4			
1	D	209	Total	C	N	O	S	0	0	0
			1638	1025	306	303	4			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	413	MET	-	INITIATING METHIONINE	UNP Q96T88
A	618	ALA	-	EXPRESSION TAG	UNP Q96T88
A	619	HIS	-	EXPRESSION TAG	UNP Q96T88
A	620	HIS	-	EXPRESSION TAG	UNP Q96T88
A	621	HIS	-	EXPRESSION TAG	UNP Q96T88
A	622	HIS	-	EXPRESSION TAG	UNP Q96T88
A	623	HIS	-	EXPRESSION TAG	UNP Q96T88
A	624	HIS	-	EXPRESSION TAG	UNP Q96T88
B	413	MET	-	INITIATING METHIONINE	UNP Q96T88
B	618	ALA	-	EXPRESSION TAG	UNP Q96T88
B	619	HIS	-	EXPRESSION TAG	UNP Q96T88
B	620	HIS	-	EXPRESSION TAG	UNP Q96T88
B	621	HIS	-	EXPRESSION TAG	UNP Q96T88
B	622	HIS	-	EXPRESSION TAG	UNP Q96T88
B	623	HIS	-	EXPRESSION TAG	UNP Q96T88
B	624	HIS	-	EXPRESSION TAG	UNP Q96T88
C	413	MET	-	INITIATING METHIONINE	UNP Q96T88
C	618	ALA	-	EXPRESSION TAG	UNP Q96T88
C	619	HIS	-	EXPRESSION TAG	UNP Q96T88
C	620	HIS	-	EXPRESSION TAG	UNP Q96T88
C	621	HIS	-	EXPRESSION TAG	UNP Q96T88

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Chain	Residue	Modelled	Actual	Comment	Reference
C	622	HIS	-	EXPRESSION TAG	UNP Q96T88
C	623	HIS	-	EXPRESSION TAG	UNP Q96T88
C	624	HIS	-	EXPRESSION TAG	UNP Q96T88
D	413	MET	-	INITIATING METHIONINE	UNP Q96T88
D	618	ALA	-	EXPRESSION TAG	UNP Q96T88
D	619	HIS	-	EXPRESSION TAG	UNP Q96T88
D	620	HIS	-	EXPRESSION TAG	UNP Q96T88
D	621	HIS	-	EXPRESSION TAG	UNP Q96T88
D	622	HIS	-	EXPRESSION TAG	UNP Q96T88
D	623	HIS	-	EXPRESSION TAG	UNP Q96T88
D	624	HIS	-	EXPRESSION TAG	UNP Q96T88

- Molecule 2 is a DNA chain called 5'-D(*DGP*DGP*DGP*DCP*DCP*(5CM)P*DGP*DCP*DAP*DGP*DGP*DG)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	12	Total	C	N	O	P	0	0	0
			249	117	52	69	11			
2	G	12	Total	C	N	O	P	0	0	0
			249	117	52	69	11			
2	I	12	Total	C	N	O	P	0	0	0
			249	117	52	69	11			
2	K	12	Total	C	N	O	P	0	0	0
			249	117	52	69	11			

- Molecule 3 is a DNA chain called 5'-D(*DCP*DCP*DCP*DTP*DGP*DCP*DGP*DGP*DP*DCP*DCP*DC)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	12	Total	C	N	O	P	0	0	0
			238	113	43	71	11			
3	H	12	Total	C	N	O	P	0	0	0
			238	113	43	71	11			
3	J	12	Total	C	N	O	P	0	0	0
			238	113	43	71	11			
3	L	12	Total	C	N	O	P	0	0	0
			238	113	43	71	11			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	126	Total 127	O 127	0	1
4	B	104	Total 110	O 110	0	6
4	C	90	Total 90	O 90	0	0
4	D	109	Total 110	O 110	0	1
4	E	36	Total 37	O 37	0	1
4	F	19	Total 19	O 19	0	0
4	G	29	Total 29	O 29	0	0
4	H	20	Total 20	O 20	0	0
4	I	38	Total 40	O 40	0	2
4	J	19	Total 19	O 19	0	0
4	K	25	Total 26	O 26	0	1
4	L	25	Total 25	O 25	0	0

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 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: E3 ubiquitin-protein ligase UHRF1

Chain A: 




- Molecule 1: E3 ubiquitin-protein ligase UHRF1

Chain B: 




- Molecule 1: E3 ubiquitin-protein ligase UHRF1

Chain C: 



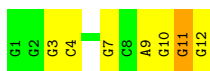
- Molecule 1: E3 ubiquitin-protein ligase UHRF1

Chain D: 



- Molecule 2: 5'-D(*DGP*DGP*DGP*DCP*DCP*(5CM)P*DGP*DCP*DAP*DGP*DGP*DG)-3'

Chain E: 



- Molecule 2: 5'-D(*DGP*DGP*DGP*DCP*DCP*(5CM)P*DGP*DCP*DAP*DGP*DGP*DG)-3'

Chain G:  42% 58%



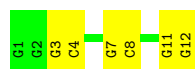
- Molecule 2: 5'-D(*DGP*DGP*DGP*DCP*DCP*(5CM)P*DGP*DCP*DAP*DGP*DGP*DG)-3'

Chain I:  25% 58% 17%



- Molecule 2: 5'-D(*DGP*DGP*DGP*DCP*DCP*(5CM)P*DGP*DCP*DAP*DGP*DGP*DG)-3'

Chain K:  50% 50%



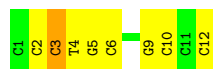
- Molecule 3: 5'-D(*DCP*DCP*DCP*DTP*DGP*DCP*DGP*DGP*DGP*DCP*DCP*DC)-3',

Chain F:  75% 25%



- Molecule 3: 5'-D(*DCP*DCP*DCP*DTP*DGP*DCP*DGP*DGP*DGP*DCP*DCP*DC)-3',

Chain H:  33% 58% 8%



- Molecule 3: 5'-D(*DCP*DCP*DCP*DTP*DGP*DCP*DGP*DGP*DGP*DCP*DCP*DC)-3',

Chain J:  33% 50% 17%



- Molecule 3: 5'-D(*DCP*DCP*DCP*DTP*DGP*DCP*DGP*DGP*DGP*DCP*DCP*DC)-3',

Chain L:  42% 33% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.20Å 111.36Å 97.40Å 90.00° 90.53° 90.00°	Depositor
Resolution (Å)	38.63 – 2.20 38.10 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (38.63-2.20) 96.7 (38.10-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.189 , 0.231 0.191 , 0.231	Depositor DCC
R_{free} test set	4085 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.389 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9090	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.02 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 9.1282e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: 5CM

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	0.62	1/1651 (0.1%)	0.66	1/2233 (0.0%)
1	B	0.59	0/1691	0.64	0/2290
1	C	0.62	0/1639	0.68	1/2220 (0.0%)
1	D	0.60	0/1685	0.66	0/2284
2	E	1.15	0/257	1.88	11/394 (2.8%)
2	G	1.14	0/257	2.08	13/394 (3.3%)
2	I	1.28	0/257	2.19	16/394 (4.1%)
2	K	1.13	0/257	1.83	9/394 (2.3%)
3	F	0.99	0/265	1.87	4/406 (1.0%)
3	H	1.01	0/265	1.88	10/406 (2.5%)
3	J	1.07	0/265	1.98	10/406 (2.5%)
3	L	1.12	1/265 (0.4%)	1.82	9/406 (2.2%)
All	All	0.76	2/8754 (0.0%)	1.14	84/12227 (0.7%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	619	HIS	C-O	5.81	1.34	1.23
3	L	8	DG	C3'-O3'	-5.01	1.37	1.44

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	4	DC	O4'-C1'-N1	-12.57	99.20	108.00
2	E	7	DG	O4'-C1'-N9	-10.33	100.77	108.00
2	K	11	DG	O4'-C1'-N9	9.99	114.99	108.00
2	G	7	DG	O4'-C1'-N9	-9.86	101.10	108.00
2	G	8	DC	OP1-P-O3'	-9.64	83.98	105.20
2	I	11	DG	O4'-C1'-N9	9.62	114.73	108.00
3	F	3	DC	O4'-C1'-N1	-9.23	101.54	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	3	DC	O4'-C1'-N1	-9.20	101.56	108.00
2	I	8	DC	OP1-P-O3'	-9.15	85.07	105.20
3	L	9	DG	O4'-C1'-N9	-8.81	101.83	108.00
3	F	4	DT	O4'-C1'-N1	-8.64	101.95	108.00
3	H	12	DC	O4'-C1'-N1	8.59	114.01	108.00
3	J	3	DC	O4'-C1'-N1	-8.57	102.00	108.00
2	G	8	DC	O3'-P-O5'	-8.24	88.35	104.00
2	G	9	DA	OP1-P-OP2	8.18	131.88	119.60
2	I	8	DC	O3'-P-O5'	-8.11	88.59	104.00
3	H	4	DT	O4'-C1'-N1	-8.10	102.33	108.00
2	G	11	DG	O4'-C1'-N9	7.92	113.55	108.00
3	J	4	DT	O4'-C1'-N1	-7.92	102.46	108.00
2	G	8	DC	OP2-P-O3'	-7.90	87.82	105.20
2	I	7	DG	O4'-C1'-N9	-7.89	102.48	108.00
3	F	9	DG	O4'-C1'-N9	-7.65	102.64	108.00
2	I	5	DC	O4'-C1'-N1	7.65	113.35	108.00
2	I	12	DG	O4'-C4'-C3'	-7.52	101.49	106.00
3	L	12	DC	O4'-C1'-N1	7.51	113.26	108.00
2	E	4	DC	O4'-C1'-N1	-7.51	102.74	108.00
3	L	4	DT	O4'-C1'-N1	-7.49	102.76	108.00
2	I	8	DC	OP2-P-O3'	-7.46	88.78	105.20
2	I	3	DG	O4'-C1'-N9	-7.33	102.87	108.00
2	I	9	DA	OP1-P-OP2	7.27	130.51	119.60
2	K	7	DG	O4'-C1'-N9	-7.27	102.91	108.00
2	G	3	DG	O4'-C1'-N9	-7.12	103.02	108.00
2	K	4	DC	O4'-C1'-N1	-7.09	103.03	108.00
3	H	9	DG	O4'-C1'-N9	-6.78	103.25	108.00
3	J	7	DG	C8-N9-C4	-6.69	103.72	106.40
3	J	5	DG	O4'-C1'-N9	6.49	112.54	108.00
2	E	11	DG	O4'-C1'-N9	6.33	112.43	108.00
2	K	3	DG	O4'-C1'-N9	-6.28	103.60	108.00
3	J	7	DG	O4'-C1'-N9	6.26	112.39	108.00
3	L	8	DG	P-O5'-C5'	-6.25	110.89	120.90
2	E	10	DG	O4'-C1'-N9	-6.21	103.65	108.00
3	J	7	DG	N9-C4-C5	6.19	107.88	105.40
2	E	9	DA	O4'-C1'-N9	6.15	112.31	108.00
2	K	12	DG	O4'-C4'-C3'	-6.07	102.07	104.50
3	J	2	DC	N1-C2-O2	6.01	122.51	118.90
3	H	10	DC	P-O3'-C3'	5.92	126.81	119.70
2	I	10	DG	O4'-C1'-N9	-5.91	103.86	108.00
2	G	4	DC	P-O3'-C3'	5.84	126.71	119.70
3	J	12	DC	O4'-C1'-N1	5.83	112.08	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	8	DC	O4'-C1'-N1	-5.83	103.92	108.00
2	I	4	DC	P-O3'-C3'	5.78	126.63	119.70
2	G	8	DC	P-O3'-C3'	5.74	126.59	119.70
3	L	7	DG	C8-N9-C4	-5.74	104.10	106.40
2	G	10	DG	O4'-C1'-N9	-5.73	103.99	108.00
3	J	4	DT	N3-C2-O2	-5.69	118.88	122.30
2	I	4	DC	O4'-C1'-N1	-5.66	104.04	108.00
2	E	7	DG	N1-C6-O6	5.65	123.29	119.90
3	H	5	DG	C5-C6-N1	5.64	114.32	111.50
2	K	11	DG	C2-N3-C4	5.54	114.67	111.90
2	E	11	DG	C5-C6-O6	-5.53	125.28	128.60
1	C	433	ARG	NE-CZ-NH2	-5.52	117.54	120.30
2	E	4	DC	P-O3'-C3'	5.38	126.16	119.70
2	I	8	DC	P-O3'-C3'	5.33	126.10	119.70
3	H	3	DC	P-O3'-C3'	5.32	126.08	119.70
3	H	9	DG	P-O3'-C3'	5.30	126.06	119.70
3	H	6	DC	O4'-C1'-N1	-5.29	104.30	108.00
2	I	9	DA	N1-C6-N6	5.29	121.77	118.60
2	G	11	DG	P-O3'-C3'	5.26	126.01	119.70
3	H	5	DG	C5-C6-O6	-5.24	125.45	128.60
2	E	12	DG	O4'-C4'-C3'	-5.23	102.41	104.50
2	G	9	DA	O4'-C1'-N9	5.22	111.66	108.00
3	L	5	DG	C3'-C2'-C1'	-5.21	96.24	102.50
2	I	11	DG	P-O3'-C3'	5.17	125.90	119.70
2	E	3	DG	O4'-C1'-N9	-5.17	104.38	108.00
3	F	9	DG	C6-N1-C2	-5.16	122.01	125.10
1	A	433	ARG	NE-CZ-NH2	-5.16	117.72	120.30
2	K	11	DG	P-O3'-C3'	5.14	125.87	119.70
2	E	11	DG	O4'-C1'-C2'	-5.13	101.80	105.90
3	L	9	DG	P-O3'-C3'	5.12	125.85	119.70
3	J	8	DG	P-O5'-C5'	-5.07	112.78	120.90
3	L	3	DC	N1-C1'-C2'	5.05	122.20	112.60
2	I	8	DC	O4'-C1'-N1	-5.04	104.47	108.00
3	L	8	DG	C5'-C4'-C3'	-5.03	105.05	114.10
2	K	11	DG	C5-C6-N1	5.00	114.00	111.50

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	1609	0	1536	12	0
1	B	1646	0	1534	9	0
1	C	1597	0	1507	8	0
1	D	1638	0	1527	10	0
2	E	249	0	136	1	0
2	G	249	0	136	0	0
2	I	249	0	136	2	0
2	K	249	0	136	0	0
3	F	238	0	135	0	0
3	H	238	0	135	1	0
3	J	238	0	135	4	0
3	L	238	0	135	2	0
4	A	127	0	0	0	0
4	B	110	0	0	0	0
4	C	90	0	0	0	0
4	D	110	0	0	1	0
4	E	37	0	0	1	0
4	F	19	0	0	0	0
4	G	29	0	0	0	0
4	H	20	0	0	0	0
4	I	40	0	0	2	0
4	J	19	0	0	0	0
4	K	26	0	0	1	0
4	L	25	0	0	0	0
All	All	9090	0	7188	48	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 3.

All (48) 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)
3:L:7:DG:H2''	3:L:8:DG:H5''	1.31	1.11
3:J:7:DG:H2''	3:J:8:DG:H5''	1.52	0.92
1:A:416:ASN:N	1:A:416:ASN:HD22	1.73	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:589:PRO:HB3	1:D:598:ILE:HD11	1.71	0.72
1:B:429:MET:O	1:B:606:GLN:HG2	1.93	0.69
3:J:7:DG:C2'	3:J:8:DG:H5''	2.22	0.69
1:A:416:ASN:ND2	1:A:416:ASN:N	2.39	0.66
3:J:7:DG:H2''	3:J:8:DG:C5'	2.24	0.66
2:I:12:DG:N7	4:I:374:HOH:O	2.29	0.65
3:L:8:DG:H2''	3:L:9:DG:C8	2.34	0.62
1:A:589:PRO:HB3	1:A:598:ILE:HD11	1.84	0.60
1:A:537:ARG:HD2	1:A:556:ASP:OD2	2.04	0.57
1:B:589:PRO:HB3	1:B:598:ILE:HD11	1.88	0.55
1:D:537:ARG:HD2	1:D:556:ASP:OD2	2.09	0.53
1:A:460:LEU:HD12	1:A:460:LEU:C	2.30	0.52
1:C:460:LEU:HD12	1:C:460:LEU:C	2.31	0.51
1:C:448:GLY:O	1:C:460:LEU:HA	2.12	0.50
1:B:434:VAL:HG23	1:B:445:HIS:CE1	2.48	0.48
1:A:598:ILE:HG23	1:A:603:LEU:HB2	1.96	0.48
1:A:615:LEU:C	1:A:617:ASN:H	2.17	0.47
1:C:444:PRO:CD	1:C:450:HIS:HB2	2.44	0.47
1:D:460:LEU:HD12	1:D:460:LEU:C	2.35	0.47
1:D:555:TYR:CZ	1:D:557:GLY:HA2	2.50	0.47
3:J:8:DG:H2''	3:J:9:DG:C8	2.50	0.47
3:H:2:DC:H1'	3:H:3:DC:H5'	1.97	0.47
1:D:598:ILE:HG23	1:D:603:LEU:HB2	1.98	0.46
1:B:622:HIS:C	1:B:624:HIS:H	2.20	0.46
1:D:619:HIS:O	1:D:622:HIS:HB2	2.16	0.46
1:A:418:TYR:CZ	1:A:531:LYS:HD3	2.52	0.45
1:B:460:LEU:HD12	1:B:460:LEU:C	2.37	0.45
1:C:589:PRO:HB3	1:C:598:ILE:HD11	1.98	0.45
1:A:546:LYS:HD2	1:C:488:GLY:HA2	1.98	0.44
1:D:482:GLY:HA2	4:K:14:HOH:O	2.17	0.44
1:D:620:HIS:C	1:D:622:HIS:H	2.22	0.43
1:A:448:GLY:O	1:A:460:LEU:HA	2.20	0.42
1:C:429:MET:HA	1:C:553:ASN:O	2.19	0.42
1:B:588:GLY:HA2	1:B:589:PRO:HD3	1.91	0.42
1:C:598:ILE:HG23	1:C:603:LEU:HB2	2.02	0.42
1:B:537:ARG:HD2	1:B:556:ASP:OD2	2.20	0.42
1:B:448:GLY:O	1:B:460:LEU:HA	2.20	0.42
1:C:418:TYR:CZ	1:C:531:LYS:HD3	2.55	0.41
2:E:11:DG:N3	4:E:195:HOH:O	2.37	0.41
1:D:564:TYR:HA	1:D:577:ARG:O	2.21	0.41
2:I:9:DA:N3	4:I:57:HOH:O	2.37	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:SER:O	1:A:506[A]:ARG:NH1	2.42	0.41
1:D:504:THR:HG22	4:D:689:HOH:O	2.21	0.41
1:A:568:LYS:HB3	1:A:568:LYS:HE3	1.81	0.40
1:B:589:PRO:HD2	1:B:590:TRP:CE3	2.57	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	203/212 (96%)	195 (96%)	8 (4%)	0	100	100
1	B	209/212 (99%)	201 (96%)	7 (3%)	1 (0%)	32	34
1	C	203/212 (96%)	193 (95%)	10 (5%)	0	100	100
1	D	207/212 (98%)	197 (95%)	9 (4%)	1 (0%)	32	34
All	All	822/848 (97%)	786 (96%)	34 (4%)	2 (0%)	51	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	621	HIS
1	B	623	HIS

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	162/172 (94%)	158 (98%)	4 (2%)	53	65
1	B	162/172 (94%)	157 (97%)	5 (3%)	45	57
1	C	159/172 (92%)	153 (96%)	6 (4%)	38	47
1	D	162/172 (94%)	153 (94%)	9 (6%)	25	29
All	All	645/688 (94%)	621 (96%)	24 (4%)	40	49

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	416	ASN
1	A	516	ILE
1	A	560	LYS
1	A	567	GLU
1	B	470	VAL
1	B	474	ASN
1	B	516	ILE
1	B	606	GLN
1	B	620	HIS
1	C	438[A]	GLU
1	C	438[B]	GLU
1	C	470	VAL
1	C	474	ASN
1	C	516	ILE
1	C	560	LYS
1	D	470	VAL
1	D	474	ASN
1	D	504	THR
1	D	520	GLU
1	D	560	LYS
1	D	596	ASP
1	D	615	LEU
1	D	619	HIS
1	D	620	HIS

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

Mol	Chain	Res	Type
1	A	445	HIS
1	A	544	ASN
1	B	445	HIS

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Mol	Chain	Res	Type
1	D	416	ASN
1	D	445	HIS
1	D	617	ASN
1	D	620	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	5CM	E	6	2	14,21,22	1.83	1 (7%)	18,30,33	1.42	4 (22%)
2	5CM	G	6	2	14,21,22	1.91	2 (14%)	18,30,33	1.67	4 (22%)
2	5CM	I	6	2	14,21,22	1.79	1 (7%)	18,30,33	1.55	4 (22%)
2	5CM	K	6	2	14,21,22	1.65	1 (7%)	18,30,33	1.46	4 (22%)

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	5CM	E	6	2	-	0/3/21/22	0/2/2/2
2	5CM	G	6	2	-	0/3/21/22	0/2/2/2
2	5CM	I	6	2	-	0/3/21/22	0/2/2/2
2	5CM	K	6	2	-	0/3/21/22	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	6	5CM	O5'-C5'	-2.24	1.41	1.44
2	K	6	5CM	C5-C4	5.41	1.49	1.41
2	I	6	5CM	C5-C4	5.64	1.49	1.41
2	E	6	5CM	C5-C4	5.98	1.50	1.41
2	G	6	5CM	C5-C4	6.16	1.50	1.41

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	6	5CM	O4'-C1'-N1	-3.12	102.53	107.78
2	G	6	5CM	O3'-C3'-C4'	-3.08	97.96	110.13
2	E	6	5CM	O4'-C1'-N1	-2.98	102.75	107.78
2	K	6	5CM	O3'-C3'-C4'	-2.97	98.41	110.13
2	I	6	5CM	O3'-C3'-C4'	-2.91	98.62	110.13
2	E	6	5CM	O3'-C3'-C4'	-2.88	98.77	110.13
2	G	6	5CM	O4'-C1'-N1	-2.65	103.31	107.78
2	K	6	5CM	O4'-C1'-N1	-2.16	104.14	107.78
2	I	6	5CM	C5A-C5-C4	2.18	123.89	121.65
2	E	6	5CM	C2'-C3'-C4'	2.21	107.44	102.73
2	K	6	5CM	C5A-C5-C4	2.32	124.04	121.65
2	G	6	5CM	C2'-C3'-C4'	2.46	107.98	102.73
2	E	6	5CM	C2'-C1'-N1	2.55	120.26	114.23
2	K	6	5CM	C2'-C1'-N1	3.07	121.49	114.23
2	I	6	5CM	C2'-C1'-N1	3.34	122.11	114.23
2	G	6	5CM	C2'-C1'-N1	3.98	123.64	114.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	204/212 (96%)	-0.63	0 100 100	28, 41, 70, 163	0
1	B	210/212 (99%)	-0.62	0 100 100	28, 43, 71, 117	0
1	C	204/212 (96%)	-0.64	0 100 100	29, 41, 73, 144	0
1	D	209/212 (98%)	-0.60	0 100 100	28, 43, 72, 143	0
2	E	11/12 (91%)	-1.13	0 100 100	35, 45, 52, 53	0
2	G	11/12 (91%)	-1.14	0 100 100	34, 46, 51, 53	0
2	I	11/12 (91%)	-1.10	0 100 100	35, 44, 50, 54	0
2	K	11/12 (91%)	-1.12	0 100 100	35, 45, 51, 54	0
3	F	12/12 (100%)	-1.10	0 100 100	35, 44, 51, 53	0
3	H	12/12 (100%)	-1.06	0 100 100	37, 46, 52, 52	0
3	J	12/12 (100%)	-1.14	0 100 100	36, 46, 51, 52	0
3	L	12/12 (100%)	-1.05	0 100 100	34, 47, 53, 54	0
All	All	919/944 (97%)	-0.67	0 100 100	28, 43, 72, 163	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	5CM	G	6	20/21	0.99	0.09	-	27,30,33,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	5CM	I	6	20/21	0.98	0.10	-	25,28,35,38	0
2	5CM	K	6	20/21	0.99	0.11	-	29,34,35,36	0
2	5CM	E	6	20/21	0.98	0.09	-	25,30,34,36	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.