



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:39 PM UTC

PDB ID : 3ASK / pdb_00003ask
Title : Structure of UHRF1 in complex with histone tail
Authors : Arita, K.; Sugita, K.; Unoki, M.; Hamamoto, R.; Sekiyama, N.; Tochio, H.;
Ariyoshi, M.; Shirakawa, M.
Deposited on : 2010-12-16
Resolution : 2.90 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

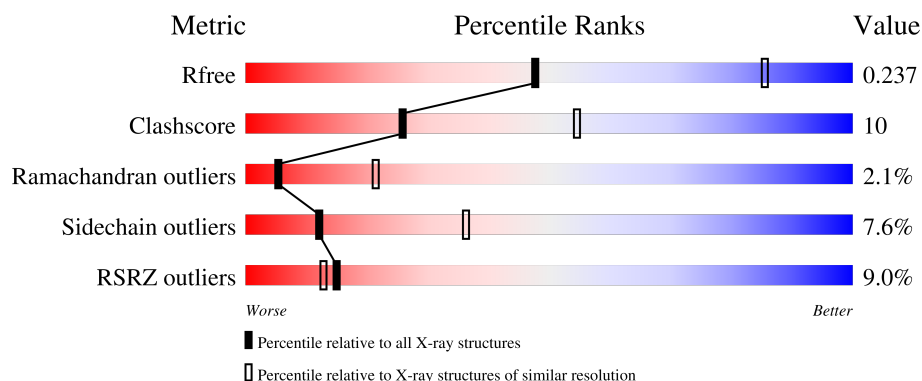
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	226	
1	B	226	
1	C	226	
1	D	226	
2	P	13	

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Mol	Chain	Length	Quality of chain
2	Q	13	 46% 23% 31%
2	R	13	 8% 92%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6141 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	213	Total	C	N	O	S	55	0	0
			1731	1070	313	329	19			
1	B	211	Total	C	N	O	S	88	0	0
			1714	1060	308	327	19			
1	C	154	Total	C	N	O	S	30	0	0
			1265	787	229	242	7			
1	D	152	Total	C	N	O	S	42	0	0
			1250	778	226	240	6			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	133	SER	-	expression tag	UNP Q96T88
A	?	-	SER	deletion	UNP Q96T88
A	?	-	ARG	deletion	UNP Q96T88
A	?	-	ASP	deletion	UNP Q96T88
A	?	-	GLU	deletion	UNP Q96T88
A	?	-	PRO	deletion	UNP Q96T88
A	?	-	CYS	deletion	UNP Q96T88
A	?	-	SER	deletion	UNP Q96T88
A	?	-	SER	deletion	UNP Q96T88
A	?	-	THR	deletion	UNP Q96T88
B	133	SER	-	expression tag	UNP Q96T88
B	?	-	SER	deletion	UNP Q96T88
B	?	-	ARG	deletion	UNP Q96T88
B	?	-	ASP	deletion	UNP Q96T88
B	?	-	GLU	deletion	UNP Q96T88
B	?	-	PRO	deletion	UNP Q96T88
B	?	-	CYS	deletion	UNP Q96T88
B	?	-	SER	deletion	UNP Q96T88
B	?	-	SER	deletion	UNP Q96T88
B	?	-	THR	deletion	UNP Q96T88
C	133	SER	-	expression tag	UNP Q96T88

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	SER	deletion	UNP Q96T88
C	?	-	ARG	deletion	UNP Q96T88
C	?	-	ASP	deletion	UNP Q96T88
C	?	-	GLU	deletion	UNP Q96T88
C	?	-	PRO	deletion	UNP Q96T88
C	?	-	CYS	deletion	UNP Q96T88
C	?	-	SER	deletion	UNP Q96T88
C	?	-	SER	deletion	UNP Q96T88
C	?	-	THR	deletion	UNP Q96T88
D	133	SER	-	expression tag	UNP Q96T88
D	?	-	SER	deletion	UNP Q96T88
D	?	-	ARG	deletion	UNP Q96T88
D	?	-	ASP	deletion	UNP Q96T88
D	?	-	GLU	deletion	UNP Q96T88
D	?	-	PRO	deletion	UNP Q96T88
D	?	-	CYS	deletion	UNP Q96T88
D	?	-	SER	deletion	UNP Q96T88
D	?	-	SER	deletion	UNP Q96T88
D	?	-	THR	deletion	UNP Q96T88

- Molecule 2 is a protein called Histone H3.3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	10	Total	C	N	O	6	0	0
			82	49	19	14			
2	Q	9	Total	C	N	O	4	0	0
			76	46	18	12			
2	R	1	Total	C	N	O	4	0	0
			12	9	2	1			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Zn	0	0
			3	3		
3	B	3	Total	Zn	0	0
			3	3		
3	C	1	Total	Zn	0	0
			1	1		

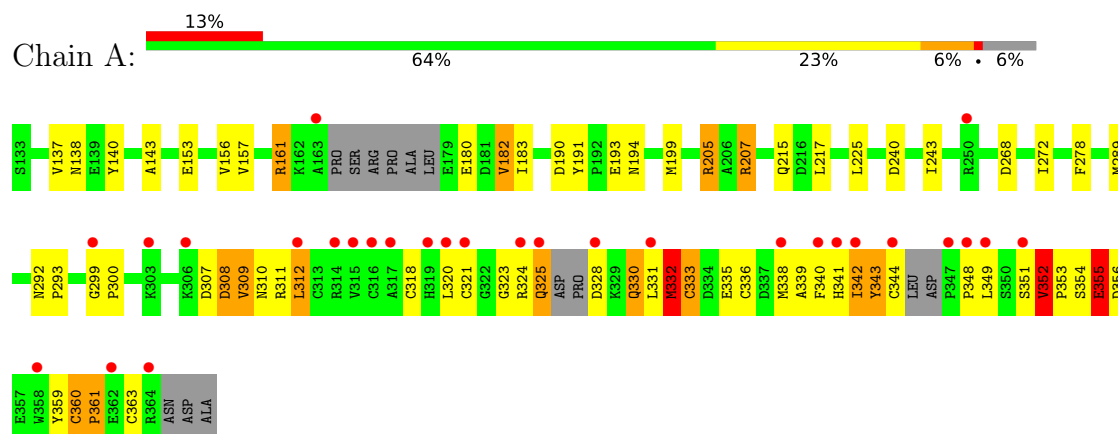
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total 3	O 3	0	0
4	C	1	Total 1	O 1	0	0

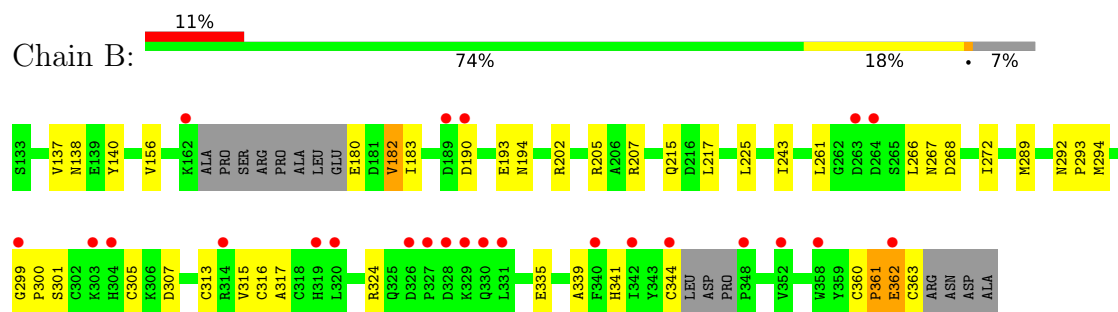
3 Residue-property plots [i](#)

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.

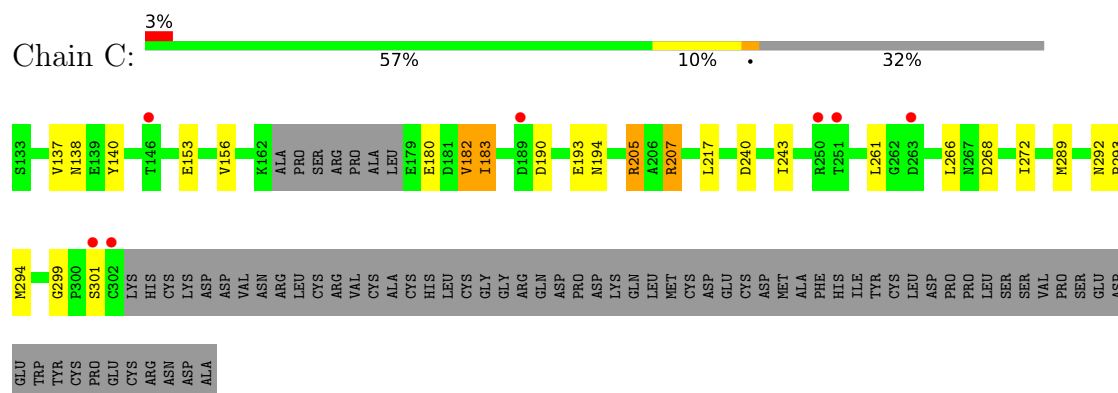
- Molecule 1: E3 ubiquitin-protein ligase UHRF1



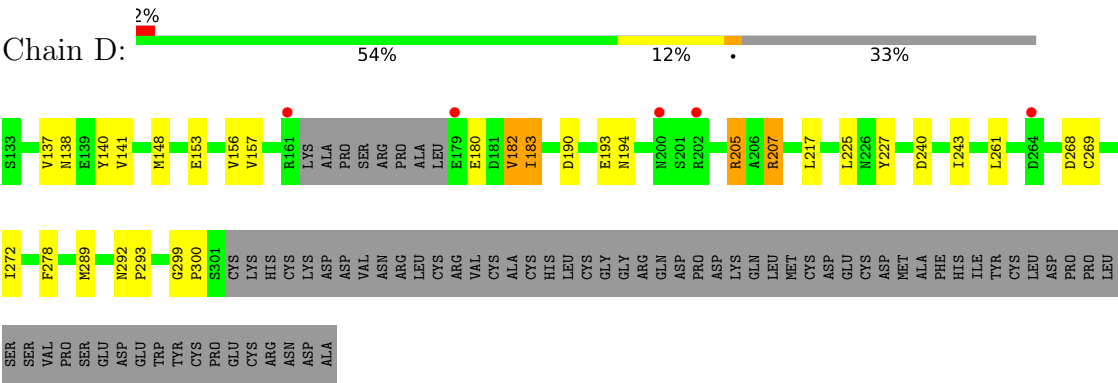
- Molecule 1: E3 ubiquitin-protein ligase UHRF1



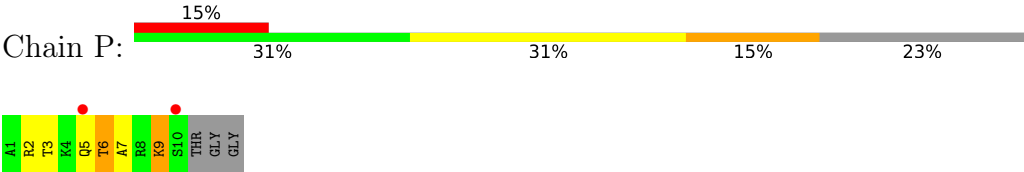
- Molecule 1: E3 ubiquitin-protein ligase UHRF1



● Molecule 1: E3 ubiquitin-protein ligase UHRF1



● Molecule 2: Histone H3.3



● Molecule 2: Histone H3.3



● Molecule 2: Histone H3.3



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	145.18Å 145.18Å 125.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.27 – 2.90 40.27 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.6 (40.27-2.90) 97.6 (40.27-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.243 , 0.286 0.242 , 0.237	Depositor DCC
R_{free} test set	2965 reflections (9.83%)	wwPDB-VP
Wilson B-factor (Å ²)	70.5	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6141	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% 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: ZN, M3L

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	0.60	0/1765	1.06	12/2380 (0.5%)
1	B	0.54	0/1749	0.91	4/2361 (0.2%)
1	C	0.52	0/1289	0.89	4/1741 (0.2%)
1	D	0.49	0/1274	0.89	3/1722 (0.2%)
2	P	0.52	0/69	0.96	0/90
2	Q	0.41	0/63	0.86	0/82
All	All	0.54	0/6209	0.95	23/8376 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	202	ARG	NE-CZ-NH2	11.58	129.62	119.20
1	A	360	CYS	CA-C-N	11.29	133.95	119.84
1	A	360	CYS	C-N-CA	11.29	133.95	119.84
1	B	202	ARG	NE-CZ-NH1	-11.14	110.36	121.50
1	A	161	ARG	NE-CZ-NH2	10.73	128.85	119.20
1	A	161	ARG	NE-CZ-NH1	-10.10	111.40	121.50
1	A	207	ARG	NE-CZ-NH2	-9.78	110.39	119.20
1	B	202	ARG	CD-NE-CZ	8.77	136.67	124.40
1	A	207	ARG	CD-NE-CZ	8.20	135.88	124.40
1	A	161	ARG	CD-NE-CZ	7.88	135.44	124.40
1	C	299	GLY	CA-C-N	-7.79	112.81	120.52
1	C	299	GLY	C-N-CA	-7.79	112.81	120.52
1	A	207	ARG	NE-CZ-NH1	7.23	128.73	121.50
1	D	299	GLY	CA-C-N	6.95	128.53	119.84
1	D	299	GLY	C-N-CA	6.95	128.53	119.84
1	A	359	TYR	N-CA-C	6.52	119.09	109.24
1	C	301	SER	N-CA-C	-5.87	106.16	113.55
1	A	352	VAL	CA-C-N	5.46	126.67	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	352	VAL	C-N-CA	5.46	126.67	119.84
1	A	332	MET	N-CA-C	5.36	118.18	108.69
1	D	207	ARG	NE-CZ-NH1	-5.15	116.35	121.50
1	B	267	ASN	N-CA-C	5.14	117.68	110.23
1	C	207	ARG	NE-CZ-NH1	-5.09	116.41	121.50

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	1731	0	1645	55	1
1	B	1714	0	1626	24	0
1	C	1265	0	1224	13	0
1	D	1250	0	1206	16	0
2	P	82	0	96	6	0
2	Q	76	0	91	0	0
2	R	12	0	20	2	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	1	0	0	0	0
4	A	3	0	0	0	0
4	C	1	0	0	0	0
All	All	6141	0	5908	109	1

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 10.

All (109) 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:336:CYS:HB3	1:A:363:CYS:SG	2.10	0.90
1:B:360:CYS:HB2	1:B:361:PRO:HD2	1.62	0.81
1:A:333:CYS:HB2	1:A:360:CYS:SG	2.28	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:PHE:CE2	1:A:360:CYS:SG	2.82	0.73
1:B:317:ALA:O	1:B:339:ALA:HB3	1.89	0.72
1:C:193:GLU:HG2	1:C:194:ASN:HD22	1.55	0.71
1:D:193:GLU:HG2	1:D:194:ASN:HD22	1.57	0.70
1:B:193:GLU:HG2	1:B:194:ASN:HD22	1.57	0.70
1:B:140:TYR:HB2	1:B:289:MET:HG2	1.75	0.67
1:A:193:GLU:HG2	1:A:194:ASN:HD22	1.59	0.67
1:C:140:TYR:HB2	1:C:289:MET:HG2	1.76	0.67
1:A:193:GLU:HG2	1:A:194:ASN:ND2	2.11	0.66
1:B:305:CYS:HB3	1:B:313:CYS:SG	2.34	0.66
1:D:193:GLU:HG2	1:D:194:ASN:ND2	2.10	0.66
1:A:341:HIS:CD2	1:A:343:TYR:CE1	2.84	0.66
1:C:193:GLU:HG2	1:C:194:ASN:ND2	2.12	0.65
1:A:321:CYS:SG	1:A:341:HIS:HE1	2.19	0.65
1:D:140:TYR:HB2	1:D:289:MET:HG2	1.77	0.65
1:A:328:ASP:N	1:A:330:GLN:HE22	1.95	0.64
1:B:193:GLU:HG2	1:B:194:ASN:ND2	2.12	0.64
1:B:207:ARG:HG2	1:B:289:MET:HE3	1.80	0.63
1:A:335:GLU:HB3	1:A:363:CYS:HB3	1.79	0.63
1:A:137:VAL:O	1:A:138:ASN:HB2	1.99	0.63
1:A:191:TYR:CE1	2:P:9:M3L:HE2	2.33	0.62
1:A:341:HIS:CD2	1:A:343:TYR:OH	2.52	0.62
1:A:341:HIS:CG	1:A:343:TYR:CE1	2.87	0.62
1:A:341:HIS:CD2	1:A:343:TYR:CZ	2.88	0.62
1:B:299:GLY:HA2	1:B:300:PRO:O	2.00	0.61
1:D:182:VAL:HG23	1:D:183:ILE:N	2.16	0.61
1:D:137:VAL:O	1:D:138:ASN:HB2	2.00	0.61
1:A:308:ASP:C	1:A:310:ASN:H	2.09	0.60
1:B:301:SER:HB2	1:B:307:ASP:OD2	2.00	0.60
1:A:182:VAL:HG23	1:A:183:ILE:N	2.16	0.60
1:A:140:TYR:HB2	1:A:289:MET:HG2	1.82	0.60
1:A:330:GLN:O	2:P:3:THR:HA	2.02	0.60
1:B:137:VAL:O	1:B:138:ASN:HB2	2.01	0.60
1:B:313:CYS:O	1:B:317:ALA:HB3	2.01	0.60
1:B:335:GLU:HB3	1:B:363:CYS:SG	2.42	0.59
1:A:330:GLN:HE21	1:A:330:GLN:N	2.00	0.59
1:B:360:CYS:HB2	1:B:361:PRO:CD	2.31	0.59
1:A:309:VAL:HG22	1:A:309:VAL:O	2.04	0.57
1:C:207:ARG:HG2	1:C:289:MET:HE3	1.85	0.57
1:A:336:CYS:SG	1:A:338:MET:HB2	2.45	0.57
1:A:323:GLY:C	1:A:325:GLN:H	2.12	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:VAL:HA	1:D:156:VAL:HG12	1.87	0.56
1:A:330:GLN:C	1:A:331:LEU:HD12	2.30	0.56
1:A:343:TYR:HD1	1:A:344:CYS:H	1.54	0.56
1:B:182:VAL:HG23	1:B:183:ILE:N	2.19	0.56
1:C:137:VAL:HA	1:C:156:VAL:HG12	1.88	0.56
1:B:313:CYS:HB3	1:B:316:CYS:SG	2.46	0.54
1:C:182:VAL:HG23	1:C:183:ILE:N	2.21	0.54
1:A:318:CYS:HB2	1:A:341:HIS:CE1	2.42	0.54
1:D:207:ARG:HG2	1:D:289:MET:HE3	1.90	0.54
1:A:307:ASP:O	1:A:309:VAL:N	2.40	0.53
1:B:361:PRO:O	1:B:362:GLU:C	2.52	0.53
1:A:339:ALA:C	1:A:340:PHE:CD2	2.87	0.53
1:C:137:VAL:O	1:C:138:ASN:HB2	2.08	0.52
1:B:341:HIS:HB2	1:B:344:CYS:SG	2.49	0.52
1:A:137:VAL:HA	1:A:156:VAL:HG12	1.92	0.51
1:A:193:GLU:HG3	2:P:5:GLN:OE1	2.10	0.51
1:A:335:GLU:HB3	1:A:363:CYS:CB	2.41	0.51
1:B:217:LEU:HD22	1:B:243:ILE:CD1	2.41	0.50
1:A:340:PHE:CZ	1:A:360:CYS:SG	3.05	0.49
1:D:205:ARG:NH2	1:D:240:ASP:OD2	2.46	0.48
2:P:7:ALA:C	2:P:9:M3L:N	2.69	0.48
1:A:323:GLY:HA3	1:A:325:GLN:HG2	1.95	0.48
1:B:289:MET:SD	1:B:293:PRO:HG3	2.54	0.48
1:A:205:ARG:NH2	1:A:240:ASP:OD2	2.47	0.47
1:A:308:ASP:O	1:A:310:ASN:N	2.46	0.47
1:C:217:LEU:HD22	1:C:243:ILE:CD1	2.45	0.47
1:C:205:ARG:NH2	1:C:240:ASP:OD2	2.48	0.47
1:B:137:VAL:HA	1:B:156:VAL:HG12	1.96	0.46
1:D:292:ASN:HA	1:D:293:PRO:HD3	1.77	0.46
1:D:148:MET:SD	2:R:9:M3L:HB2	2.56	0.46
1:B:294:MET:HE3	1:D:278:PHE:HZ	1.81	0.46
2:P:6:THR:HG22	2:P:7:ALA:N	2.31	0.46
1:A:217:LEU:HD22	1:A:243:ILE:CD1	2.47	0.45
1:D:217:LEU:HD22	1:D:243:ILE:CD1	2.47	0.45
1:A:323:GLY:C	1:A:325:GLN:N	2.75	0.45
1:C:261:LEU:HD11	1:C:266:LEU:HD22	1.99	0.45
1:A:340:PHE:CZ	1:A:360:CYS:HB2	2.53	0.44
1:A:205:ARG:HH22	1:A:240:ASP:CG	2.26	0.44
1:A:318:CYS:CB	1:A:341:HIS:CE1	3.00	0.44
1:D:205:ARG:HH22	1:D:240:ASP:CG	2.26	0.44
1:A:312:LEU:C	1:A:312:LEU:HD23	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:TYR:O	1:A:344:CYS:C	2.61	0.43
1:C:205:ARG:HH22	1:C:240:ASP:CG	2.26	0.43
1:A:354:SER:O	1:A:355:GLU:C	2.61	0.43
1:A:339:ALA:O	1:A:340:PHE:CD2	2.72	0.43
1:A:308:ASP:CG	1:A:311:ARG:HB2	2.44	0.43
1:A:325:GLN:NE2	1:A:343:TYR:OH	2.52	0.43
1:A:341:HIS:HD2	1:A:343:TYR:OH	1.98	0.43
1:B:292:ASN:HA	1:B:293:PRO:HD3	1.75	0.43
1:A:292:ASN:HA	1:A:293:PRO:HD3	1.76	0.42
1:A:332:MET:HE3	1:A:332:MET:HB3	1.62	0.42
1:B:261:LEU:HD11	1:B:266:LEU:HD22	2.02	0.42
2:R:9:M3L:HD3	2:R:9:M3L:HM22	1.89	0.42
1:A:299:GLY:HA3	1:A:300:PRO:HA	1.67	0.42
1:D:227:TYR:CE2	1:D:269:CYS:SG	3.05	0.42
1:C:292:ASN:HA	1:C:293:PRO:HD3	1.76	0.41
2:P:7:ALA:O	2:P:9:M3L:N	2.53	0.41
1:A:215:GLN:HB3	1:B:215:GLN:O	2.20	0.41
1:A:278:PHE:HZ	1:C:294:MET:HE3	1.85	0.41
1:A:143:ALA:HB1	1:A:199:MET:HE2	2.02	0.41
1:A:336:CYS:CB	1:A:363:CYS:SG	2.88	0.41
1:D:141:VAL:HB	1:D:205:ARG:O	2.21	0.41
1:D:227:TYR:HE2	1:D:269:CYS:HG	1.57	0.41
1:A:352:VAL:O	1:A:352:VAL:HG23	2.22	0.40
1:A:318:CYS:SG	1:A:320:LEU:HB2	2.62	0.40

All (1) 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:194:ASN:OD1	1:A:356:ASP:OD1[8_556]	2.17	0.03

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	205/226 (91%)	171 (83%)	23 (11%)	11 (5%)	1	5
1	B	205/226 (91%)	182 (89%)	20 (10%)	3 (2%)	8	28
1	C	150/226 (66%)	142 (95%)	8 (5%)	0	100	100
1	D	148/226 (66%)	139 (94%)	8 (5%)	1 (1%)	18	48
2	P	7/13 (54%)	6 (86%)	1 (14%)	0	100	100
2	Q	7/13 (54%)	6 (86%)	1 (14%)	0	100	100
All	All	722/930 (78%)	646 (90%)	61 (8%)	15 (2%)	5	21

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	308	ASP
1	A	355	GLU
1	A	309	VAL
1	A	342	ILE
1	A	361	PRO
1	A	343	TYR
1	A	351	SER
1	B	362	GLU
1	A	352	VAL
1	B	324	ARG
1	B	361	PRO
1	A	333	CYS
1	A	348	PRO
1	A	353	PRO
1	D	300	PRO

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	A	192/203 (95%)	172 (90%)	20 (10%)	7	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	191/203 (94%)	183 (96%)	8 (4%)	26	60
1	C	138/203 (68%)	130 (94%)	8 (6%)	18	49
1	D	136/203 (67%)	125 (92%)	11 (8%)	11	33
2	P	7/8 (88%)	5 (71%)	2 (29%)	0	1
2	Q	6/8 (75%)	4 (67%)	2 (33%)	0	0
All	All	670/828 (81%)	619 (92%)	51 (8%)	12	36

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	GLU
1	A	157	VAL
1	A	161	ARG
1	A	180	GLU
1	A	182	VAL
1	A	190	ASP
1	A	205	ARG
1	A	207	ARG
1	A	225	LEU
1	A	268	ASP
1	A	272	ILE
1	A	312	LEU
1	A	324	ARG
1	A	325	GLN
1	A	330	GLN
1	A	332	MET
1	A	342	ILE
1	A	349	LEU
1	A	355	GLU
1	A	361	PRO
1	B	180	GLU
1	B	182	VAL
1	B	190	ASP
1	B	205	ARG
1	B	225	LEU
1	B	268	ASP
1	B	272	ILE
1	B	315	VAL
1	C	153	GLU
1	C	180	GLU

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Mol	Chain	Res	Type
1	C	182	VAL
1	C	183	ILE
1	C	190	ASP
1	C	205	ARG
1	C	268	ASP
1	C	272	ILE
1	D	153	GLU
1	D	157	VAL
1	D	180	GLU
1	D	182	VAL
1	D	183	ILE
1	D	190	ASP
1	D	205	ARG
1	D	225	LEU
1	D	261	LEU
1	D	268	ASP
1	D	272	ILE
2	P	2	ARG
2	P	6	THR
2	Q	3	THR
2	Q	4	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	ASN
1	A	194	ASN
1	A	231	ASN
1	A	325	GLN
1	A	330	GLN
1	A	341	HIS
1	B	147	ASN
1	B	194	ASN
1	B	231	ASN
1	B	267	ASN
1	C	147	ASN
1	C	194	ASN
1	C	200	ASN
1	C	231	ASN
1	C	267	ASN
1	D	147	ASN
1	D	194	ASN

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Mol	Chain	Res	Type
1	D	231	ASN
1	D	267	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	M3L	P	9	2	10,11,12	1.17	0	9,14,16	1.99	3 (33%)
2	M3L	R	9	-	10,11,12	0.93	0	9,14,16	1.97	2 (22%)
2	M3L	Q	9	2	10,11,12	1.01	0	9,14,16	2.28	3 (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	M3L	P	9	2	-	2/9/10/12	-
2	M3L	R	9	-	-	0/9/10/12	-
2	M3L	Q	9	2	-	0/9/10/12	-

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	9	M3L	CM3-NZ-CE	4.83	129.09	109.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	9	M3L	CM3-NZ-CE	4.11	126.23	109.91
2	P	9	M3L	CM3-NZ-CE	3.87	125.31	109.91
2	R	9	M3L	CM3-NZ-CM2	-2.93	101.29	108.98
2	Q	9	M3L	CM3-NZ-CM2	-2.88	101.40	108.98
2	P	9	M3L	CM3-NZ-CM2	-2.47	102.48	108.98
2	Q	9	M3L	CM2-NZ-CE	-2.12	101.47	109.91
2	P	9	M3L	CD-CG-CB	2.08	121.44	113.62

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	P	9	M3L	O-C-CA-CB
2	P	9	M3L	CD-CE-NZ-CM3

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	9	M3L	3	0
2	R	9	M3L	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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	213/226 (94%)	0.70	29 (13%) 7 5	32, 63, 111, 126	14 (6%)
1	B	211/226 (93%)	0.72	24 (11%) 10 8	31, 63, 132, 151	22 (10%)
1	C	154/226 (68%)	0.43	7 (4%) 38 30	35, 62, 94, 133	7 (4%)
1	D	152/226 (67%)	0.49	5 (3%) 49 40	33, 64, 93, 131	10 (6%)
2	P	9/13 (69%)	1.03	2 (22%) 2 2	52, 71, 90, 105	2 (22%)
2	Q	8/13 (61%)	0.80	0 100 100	55, 92, 113, 119	1 (12%)
2	R	0/13	-	-	-	-
All	All	747/943 (79%)	0.61	67 (8%) 15 13	31, 64, 119, 151	56 (7%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	348	PRO	5.4
1	A	315	VAL	4.8
1	A	351	SER	4.6
1	B	328	ASP	4.4
1	A	342	ILE	4.1
1	A	314	ARG	4.1
1	A	299	GLY	4.0
1	C	302	CYS	3.9
1	D	202	ARG	3.8
1	A	358	TRP	3.7
1	D	179	GLU	3.7
1	B	327	PRO	3.6
1	B	326	ASP	3.6
1	C	263	ASP	3.5
1	A	331	LEU	3.4
1	B	263	ASP	3.4
1	A	348	PRO	3.3
1	D	200	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	264	ASP	3.2
1	A	347	PRO	3.2
1	B	342	ILE	3.1
1	B	330	GLN	3.1
1	B	162	LYS	3.0
1	D	161	ARG	3.0
1	B	329	LYS	2.8
1	C	301	SER	2.8
1	A	324	ARG	2.7
1	A	340	PHE	2.7
1	A	317	ALA	2.7
1	B	358	TRP	2.7
1	D	264	ASP	2.7
1	B	331	LEU	2.7
1	A	341	HIS	2.6
1	A	303	LYS	2.6
2	P	10	SER	2.6
1	B	304	HIS	2.5
1	A	328	ASP	2.5
1	A	338	MET	2.5
1	B	319	HIS	2.5
1	B	340	PHE	2.5
1	C	251	THR	2.5
1	A	320	LEU	2.5
1	A	344	CYS	2.4
1	B	352	VAL	2.4
1	A	312	LEU	2.4
1	C	189	ASP	2.3
1	A	364	ARG	2.3
1	A	250	ARG	2.3
1	B	344	CYS	2.3
1	B	362	GLU	2.2
1	B	189	ASP	2.2
1	B	190	ASP	2.2
1	B	314	ARG	2.2
1	A	319	HIS	2.2
1	A	316	CYS	2.2
1	A	349	LEU	2.2
2	P	5	GLN	2.1
1	B	320	LEU	2.1
1	A	163	ALA	2.1
1	B	303	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	325	GLN	2.1
1	C	250	ARG	2.1
1	A	321	CYS	2.1
1	C	146	THR	2.1
1	A	306	LYS	2.0
1	B	299	GLY	2.0
1	A	362	GLU	2.0

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. 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(Å ²)	Q<0.9
2	M3L	R	9	12/13	0.87	0.28	57,98,102,116	4
2	M3L	Q	9	12/13	0.89	0.16	48,72,85,88	0
2	M3L	P	9	12/13	0.95	0.08	37,51,102,109	0

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(Å ²)	Q<0.9
3	ZN	C	501	1/1	0.93	0.11	106,106,106,106	0
3	ZN	A	502	1/1	0.96	0.06	122,122,122,122	0
3	ZN	A	503	1/1	0.96	0.07	109,109,109,109	0
3	ZN	A	501	1/1	0.96	0.06	130,130,130,130	0
3	ZN	B	506	1/1	0.97	0.08	150,150,150,150	0
3	ZN	B	504	1/1	0.98	0.06	125,125,125,125	0
3	ZN	B	505	1/1	0.98	0.08	135,135,135,135	0

6.5 Other polymers [i](#)

There are no such residues in this entry.