



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:31 PM UTC

PDB ID : 3U5N / pdb_00003u5n
Title : Crystal structure of the complex of TRIM33 PHD-Bromo and H3(1-20)K9me3K14ac histone peptide
Authors : Wang, Z.; Patel, D.J.
Deposited on : 2011-10-11
Resolution : 1.95 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

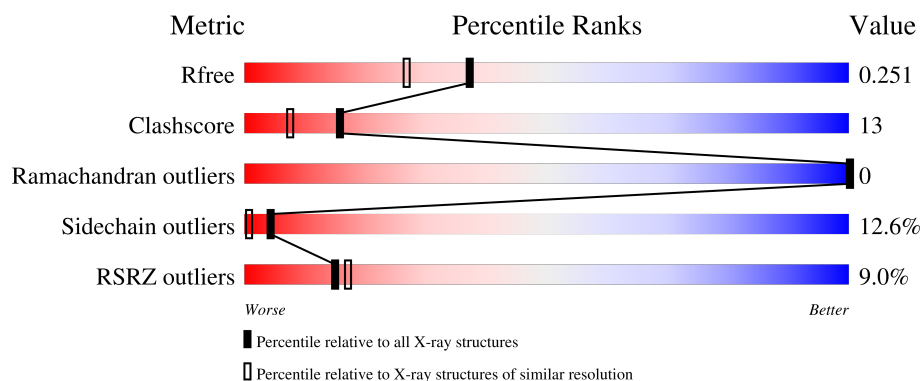
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	207	4% 67% 17% 5% 11%
1	B	207	10% 62% 21% 6% 10%
2	C	20	10% 40% 5% 5% 50%
2	D	20	20% 55% 35% 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3443 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 TRIM33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	0	0
			1489	950	246	278	15			
1	B	186	Total	C	N	O	S	0	0	0
			1498	957	245	281	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	881	SER	-	expression tag	UNP Q9UPN9
B	881	SER	-	expression tag	UNP Q9UPN9

- Molecule 2 is a protein called Histone H3.1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	0	0	0
			82	49	19	14			
2	D	20	Total	C	N	O	0	0	0
			158	96	35	27			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	2	Total	Zn	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	121	Total	O	0	0
			121	121		

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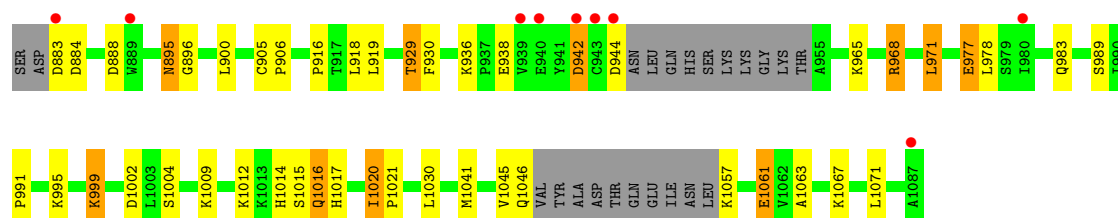
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	86	Total 86	O 86	0	0
4	D	5	Total 5	O 5	0	0

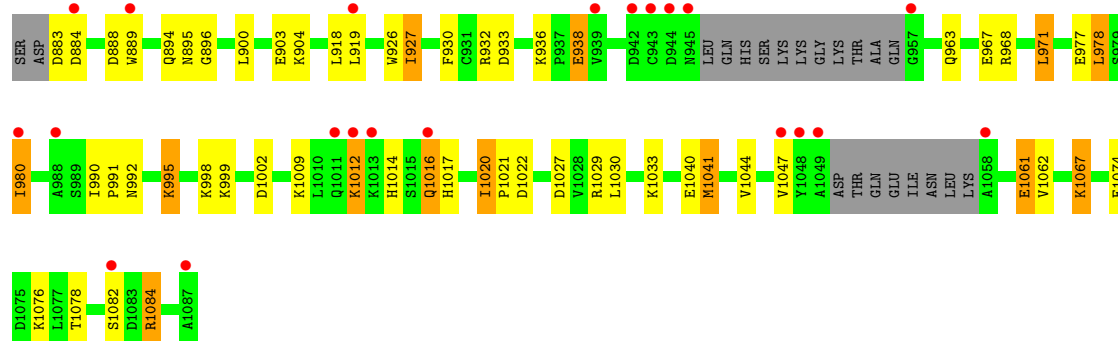
3 Residue-property plots [i](#)

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.

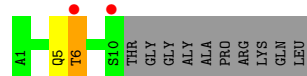
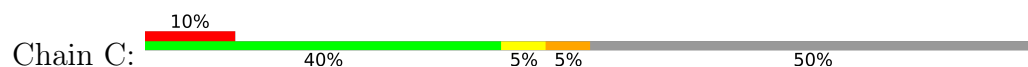
• Molecule 1: E3 ubiquitin-protein ligase TRIM33



• Molecule 1: E3 ubiquitin-protein ligase TRIM33

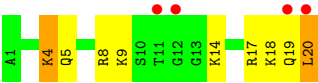


• Molecule 2: Histone H3.1



• Molecule 2: Histone H3.1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.59Å 113.74Å 39.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.79 – 1.95 29.79 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.4 (29.79-1.95) 92.3 (29.79-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.59 (at 1.95Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.202 , 0.256 0.198 , 0.251	Depositor DCC
R_{free} test set	1652 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.910	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3443	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ALY, 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	0.49	0/1523	0.77	2/2057 (0.1%)
1	B	0.41	0/1533	0.79	2/2073 (0.1%)
2	C	0.43	0/69	0.69	0/90
2	D	0.36	0/133	0.75	0/174
All	All	0.45	0/3258	0.78	4/4394 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	884	ASP	CA-C-N	-8.04	113.53	121.65
1	B	884	ASP	C-N-CA	-8.04	113.53	121.65
1	A	936	LYS	CA-C-N	5.08	125.32	119.83
1	A	936	LYS	C-N-CA	5.08	125.32	119.83

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1489	0	1450	38	0
1	B	1498	0	1451	45	0
2	C	82	0	97	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	158	0	181	11	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	121	0	0	5	0
4	B	86	0	0	1	0
4	D	5	0	0	0	0
All	All	3443	0	3179	85	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 13.

All (85) 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:B:895:ASN:HD22	1:B:896:GLY:H	1.13	0.96
1:A:977:GLU:CD	1:A:977:GLU:H	1.77	0.91
1:B:1002:ASP:OD1	4:B:174:HOH:O	1.96	0.84
1:B:995:LYS:O	1:B:998:LYS:NZ	2.10	0.83
1:B:904:LYS:HB2	1:B:927:ILE:HD11	1.61	0.82
1:B:895:ASN:ND2	1:B:896:GLY:H	1.81	0.79
1:A:1016:GLN:OE1	2:D:20:LEU:N	2.14	0.79
1:A:1061:GLU:OE1	1:A:1061:GLU:N	2.15	0.78
1:B:895:ASN:HD21	2:D:9:M3L:H	1.30	0.77
1:A:1061:GLU:H	1:A:1061:GLU:CD	1.90	0.77
1:A:1020:ILE:HG13	1:A:1021:PRO:HD2	1.66	0.76
1:A:883:ASP:CG	1:A:884:ASP:H	1.96	0.72
1:B:895:ASN:HD22	1:B:896:GLY:N	1.86	0.70
1:B:1014:HIS:HD2	1:B:1016:GLN:H	1.40	0.68
1:A:905:CYS:HB2	1:A:906:PRO:HD2	1.76	0.68
1:B:977:GLU:HG3	1:B:978:LEU:HD13	1.76	0.68
1:B:1040:GLU:HG2	1:B:1041:MET:HE3	1.76	0.67
1:A:1041:MET:SD	4:A:83:HOH:O	2.51	0.67
1:A:977:GLU:CD	1:A:977:GLU:N	2.53	0.66
1:A:999:LYS:HD2	1:A:1030:LEU:HD21	1.78	0.66
1:B:1020:ILE:HG23	1:B:1021:PRO:HD2	1.76	0.66
1:B:999:LYS:HG3	1:B:1030:LEU:HD21	1.76	0.65
1:A:965:LYS:NZ	4:A:80:HOH:O	2.32	0.62
1:B:932:ARG:O	1:B:968:ARG:NH2	2.34	0.61
1:A:916:PRO:HG3	1:A:944:ASP:HA	1.82	0.61
1:B:1061:GLU:OE1	1:B:1061:GLU:N	2.25	0.60
1:A:883:ASP:CG	1:A:884:ASP:N	2.58	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:980:ILE:HD13	1:B:980:ILE:H	1.68	0.59
1:B:980:ILE:HD13	1:B:980:ILE:N	2.17	0.59
1:A:930:PHE:HA	1:A:968:ARG:HD2	1.85	0.58
1:A:895:ASN:HD22	1:A:896:GLY:H	1.51	0.58
1:A:1015:SER:HB3	2:D:17:ARG:NH2	2.18	0.58
1:B:980:ILE:H	1:B:980:ILE:CD1	2.16	0.58
1:B:1029:ARG:O	1:B:1033:LYS:HG3	2.05	0.57
1:B:977:GLU:O	1:B:980:ILE:HD11	2.05	0.57
1:B:895:ASN:ND2	2:D:9:M3L:H	2.03	0.55
1:A:1067:LYS:NZ	4:A:149:HOH:O	2.40	0.55
1:A:1067:LYS:O	1:A:1071:LEU:HD23	2.07	0.55
1:B:1061:GLU:H	1:B:1061:GLU:CD	2.03	0.55
1:A:888:ASP:O	2:C:6:THR:HG21	2.07	0.54
1:B:1012:LYS:HA	1:B:1017:HIS:CD2	2.42	0.54
1:B:904:LYS:HB2	1:B:927:ILE:CD1	2.33	0.54
1:B:894:GLN:O	2:D:9:M3L:HM22	2.08	0.54
1:A:1020:ILE:HD12	4:A:98:HOH:O	2.06	0.54
1:A:1014:HIS:CD2	1:A:1016:GLN:H	2.27	0.53
1:A:1020:ILE:HD11	4:A:198:HOH:O	2.08	0.52
1:A:895:ASN:ND2	1:A:896:GLY:H	2.07	0.52
1:A:942:ASP:OD2	1:A:942:ASP:N	2.43	0.52
1:A:929:THR:HB	1:A:971:LEU:HD12	1.93	0.51
1:B:903:GLU:HG3	1:B:926:TRP:O	2.12	0.50
1:A:1014:HIS:HD2	1:A:1016:GLN:H	1.60	0.49
1:B:991:PRO:O	1:B:992:ASN:HB2	2.12	0.49
1:B:1067:LYS:HE3	1:B:1067:LYS:HB2	1.71	0.49
1:B:1020:ILE:HG22	1:B:1022:ASP:H	1.77	0.48
1:B:1044:VAL:O	1:B:1047:VAL:HG12	2.12	0.48
1:A:1016:GLN:O	1:A:1017:HIS:C	2.57	0.47
1:B:1009:LYS:HE2	1:B:1027:ASP:OD2	2.14	0.47
1:B:938:GLU:H	1:B:938:GLU:HG2	1.43	0.47
1:B:889:TRP:CE3	2:D:9:M3L:HM23	2.50	0.46
1:B:888:ASP:OD1	2:D:4:LYS:NZ	2.48	0.46
1:B:933:ASP:HB3	1:B:936:LYS:O	2.15	0.46
2:D:14:ALY:HE2	2:D:14:ALY:HH31	1.66	0.46
1:A:989:SER:O	1:A:991:PRO:HD3	2.16	0.46
1:B:1014:HIS:CD2	1:B:1016:GLN:H	2.27	0.46
2:D:9:M3L:HM22	2:D:9:M3L:HD3	1.46	0.46
1:B:904:LYS:HB2	1:B:927:ILE:CG1	2.47	0.45
1:A:895:ASN:HD22	1:A:896:GLY:N	2.15	0.45
1:A:905:CYS:CB	1:A:906:PRO:HD2	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1063:ALA:O	1:A:1067:LYS:HG3	2.17	0.45
1:B:930:PHE:CE2	1:B:971:LEU:HD13	2.52	0.45
1:A:1009:LYS:HE3	1:A:1016:GLN:HB3	1.99	0.45
1:A:919:LEU:N	1:A:919:LEU:HD12	2.33	0.44
1:B:889:TRP:CD2	2:D:9:M3L:HM23	2.52	0.44
1:A:1002:ASP:OD2	1:A:1004:SER:OG	2.35	0.43
1:B:1076:LYS:HD3	1:B:1076:LYS:HA	1.88	0.43
1:A:896:GLY:HA3	2:C:6:THR:HG22	2.00	0.42
1:A:938:GLU:OE1	1:A:938:GLU:N	2.48	0.42
1:A:983:GLN:O	1:A:1004:SER:HB3	2.19	0.42
1:B:895:ASN:ND2	1:B:896:GLY:N	2.56	0.42
1:A:916:PRO:CG	1:A:944:ASP:HA	2.47	0.42
1:B:963:GLN:O	1:B:967:GLU:HG3	2.20	0.42
1:B:1022:ASP:OD1	1:B:1084:ARG:NH2	2.50	0.42
1:B:896:GLY:O	2:D:8:ARG:NH1	2.53	0.42
1:B:1020:ILE:HG23	1:B:1021:PRO:CD	2.47	0.41
1:B:1074:GLU:O	1:B:1078:THR:HG23	2.21	0.41

There are no symmetry-related clashes.

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	179/207 (86%)	177 (99%)	2 (1%)	0	100	100
1	B	180/207 (87%)	175 (97%)	5 (3%)	0	100	100
2	C	7/20 (35%)	7 (100%)	0	0	100	100
2	D	16/20 (80%)	16 (100%)	0	0	100	100
All	All	382/454 (84%)	375 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	168/189 (89%)	150 (89%)	18 (11%)	6	1
1	B	169/189 (89%)	149 (88%)	20 (12%)	5	1
2	C	7/13 (54%)	5 (71%)	2 (29%)	0	0
2	D	13/13 (100%)	8 (62%)	5 (38%)	0	0
All	All	357/404 (88%)	312 (87%)	45 (13%)	4	1

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	895	ASN
1	A	900	LEU
1	A	918	LEU
1	A	929	THR
1	A	942	ASP
1	A	968	ARG
1	A	971	LEU
1	A	977	GLU
1	A	978	LEU
1	A	995	LYS
1	A	999	LYS
1	A	1012	LYS
1	A	1016	GLN
1	A	1020	ILE
1	A	1045	VAL
1	A	1046	GLN
1	A	1057	LYS
1	A	1061	GLU
2	C	5	GLN
2	C	6	THR
1	B	883	ASP
1	B	900	LEU
1	B	918	LEU
1	B	919	LEU

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Mol	Chain	Res	Type
1	B	927	ILE
1	B	938	GLU
1	B	971	LEU
1	B	978	LEU
1	B	980	ILE
1	B	990	ILE
1	B	995	LYS
1	B	1012	LYS
1	B	1016	GLN
1	B	1020	ILE
1	B	1041	MET
1	B	1061	GLU
1	B	1062	VAL
1	B	1067	LYS
1	B	1082	SER
1	B	1084	ARG
2	D	4	LYS
2	D	5	GLN
2	D	18	LYS
2	D	19	GLN
2	D	20	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	894	GLN
1	A	895	ASN
1	A	983	GLN
1	A	992	ASN
1	A	1014	HIS
1	A	1017	HIS
1	A	1046	GLN
1	B	895	ASN
1	B	983	GLN
1	B	1014	HIS
1	B	1016	GLN
1	B	1064	GLN
2	D	5	GLN

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	ALY	D	14	2	10,11,12	0.40	0	7,12,14	0.54	0
2	M3L	C	9	2	10,11,12	0.83	0	9,14,16	0.64	0
2	M3L	D	9	2	10,11,12	0.83	0	9,14,16	0.57	0

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	ALY	D	14	2	-	3/9/10/12	-
2	M3L	C	9	2	-	7/9/10/12	-
2	M3L	D	9	2	-	5/9/10/12	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	9	M3L	N-CA-CB-CG
2	C	9	M3L	C-CA-CB-CG
2	D	14	ALY	OH-CH-NZ-CE
2	D	14	ALY	CH3-CH-NZ-CE

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Mol	Chain	Res	Type	Atoms
2	D	9	M3L	CG-CD-CE-NZ
2	C	9	M3L	CG-CD-CE-NZ
2	D	9	M3L	CE-CD-CG-CB
2	D	14	ALY	CE-CD-CG-CB
2	D	9	M3L	CD-CE-NZ-CM2
2	C	9	M3L	CD-CE-NZ-CM2
2	D	9	M3L	CD-CE-NZ-CM3
2	C	9	M3L	CD-CE-NZ-CM1
2	D	9	M3L	CD-CE-NZ-CM1
2	C	9	M3L	CD-CE-NZ-CM3
2	C	9	M3L	CA-CB-CG-CD

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	14	ALY	1	0
2	D	9	M3L	6	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 ⓘ

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	185/207 (89%)	0.21	9 (4%) 35 41	17, 29, 51, 78	0
1	B	186/207 (89%)	0.67	21 (11%) 10 11	21, 38, 67, 81	0
2	C	9/20 (45%)	1.36	2 (22%) 2 2	36, 49, 69, 85	0
2	D	18/20 (90%)	1.18	4 (22%) 2 2	31, 45, 62, 63	0
All	All	398/454 (87%)	0.49	36 (9%) 15 17	17, 35, 62, 85	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1087	ALA	5.3
1	B	945	ASN	4.0
1	A	942	ASP	4.0
1	B	943	CYS	3.9
1	A	943	CYS	3.8
2	C	10	SER	3.7
2	D	20	LEU	3.6
1	B	980	ILE	3.2
1	B	1087	ALA	3.0
1	B	1048	TYR	3.0
2	C	6	THR	2.9
1	A	980	ILE	2.8
1	B	889	TRP	2.8
1	B	1012	LYS	2.6
1	B	944	ASP	2.6
1	B	1049	ALA	2.6
1	B	1016	GLN	2.6
1	B	1047	VAL	2.6
1	B	919	LEU	2.5
1	B	1082	SER	2.5
1	A	883	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	939	VAL	2.4
2	D	19	GLN	2.3
1	B	1013	LYS	2.3
1	B	1011	GLN	2.3
1	B	988	ALA	2.2
1	B	942	ASP	2.2
1	B	884	ASP	2.2
1	B	957	GLY	2.2
1	A	889	TRP	2.2
1	A	944	ASP	2.2
1	B	1058	ALA	2.2
1	A	939	VAL	2.1
2	D	11	THR	2.1
2	D	12	GLY	2.1
1	A	940	GLU	2.1

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	C	9	12/13	0.69	0.20	68,76,84,87	0
2	M3L	D	9	12/13	0.91	0.11	27,38,69,69	0
2	ALY	D	14	12/13	0.94	0.09	25,35,41,44	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($\text{\AA}^2$)	Q<0.9
3	ZN	B	2	1/1	0.95	0.06	40,40,40,40	0
3	ZN	A	2	1/1	0.99	0.03	36,36,36,36	0
3	ZN	B	1	1/1	0.99	0.04	37,37,37,37	0
3	ZN	A	1	1/1	0.99	0.06	35,35,35,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.