



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:09 PM UTC

PDB ID : 6Q0V / pdb_00006q0v
Title : Structure of DDB1-DDA1-DCAF15 complex bound to tasisulam and RBM39
Authors : Faust, T.; Yoon, H.; Nowak, R.P.; Donovan, K.A.; Li, Z.; Cai, Q.; Eleuteri, N.A.; Zhang, T.; Gray, N.S.; Fischer, E.S.
Deposited on : 2019-08-02
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

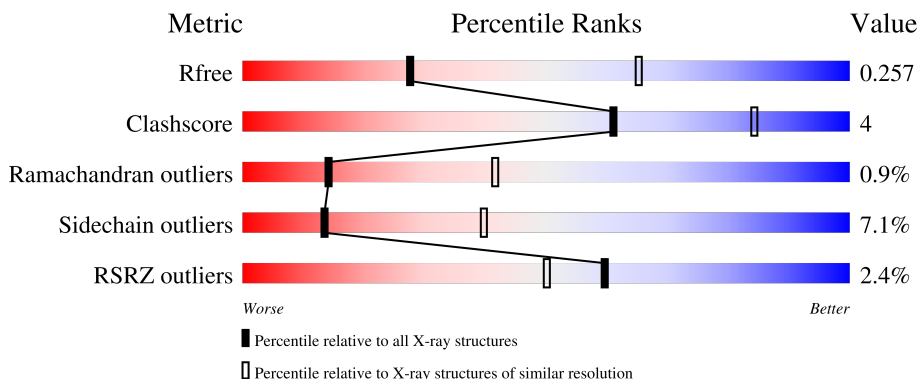
1 Overall quality at a glance

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	864	2% 75% 12% 11%
2	B	276	3% 61% 12% 25%
3	C	263	2% 57% 14% 27%
4	D	107	% 56% 16% 26%
5	E	126	2% 34% 11% 51%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10494 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 DNA damage-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	772	Total	C	N	O	S	0	0	0
			6082	3862	1023	1163	34			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-27	MET	-	initiating methionine	UNP Q16531
A	-26	GLY	-	expression tag	UNP Q16531
A	-25	SER	-	expression tag	UNP Q16531
A	-24	SER	-	expression tag	UNP Q16531
A	-23	HIS	-	expression tag	UNP Q16531
A	-22	HIS	-	expression tag	UNP Q16531
A	-21	HIS	-	expression tag	UNP Q16531
A	-20	HIS	-	expression tag	UNP Q16531
A	-19	HIS	-	expression tag	UNP Q16531
A	-18	HIS	-	expression tag	UNP Q16531
A	-17	SER	-	expression tag	UNP Q16531
A	-16	ALA	-	expression tag	UNP Q16531
A	-15	ALA	-	expression tag	UNP Q16531
A	-14	HIS	-	expression tag	UNP Q16531
A	-13	ILE	-	expression tag	UNP Q16531
A	-12	VAL	-	expression tag	UNP Q16531
A	-11	MET	-	expression tag	UNP Q16531
A	-10	VAL	-	expression tag	UNP Q16531
A	-9	ASP	-	expression tag	UNP Q16531
A	-8	ALA	-	expression tag	UNP Q16531
A	-7	TYR	-	expression tag	UNP Q16531
A	-6	LYS	-	expression tag	UNP Q16531
A	-5	PRO	-	expression tag	UNP Q16531
A	-4	THR	-	expression tag	UNP Q16531
A	-3	LYS	-	expression tag	UNP Q16531
A	-2	GLY	-	expression tag	UNP Q16531
A	-1	GLY	-	expression tag	UNP Q16531

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ARG	-	expression tag	UNP Q16531
A	700	GLY	-	linker	UNP Q16531
A	701	ASN	-	linker	UNP Q16531
A	702	GLY	-	linker	UNP Q16531
A	703	ASN	-	linker	UNP Q16531
A	704	SER	-	linker	UNP Q16531
A	705	GLY	-	linker	UNP Q16531

- Molecule 2 is a protein called DDB1- and CUL4-associated factor 15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	208	Total 1690	C 1091	N 291	O 298	S 10	0	0	0

There are 49 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	MET	-	initiating methionine	UNP Q66K64
B	-14	ASP	-	expression tag	UNP Q66K64
B	-13	TRP	-	expression tag	UNP Q66K64
B	-12	SER	-	expression tag	UNP Q66K64
B	-11	HIS	-	expression tag	UNP Q66K64
B	-10	PRO	-	expression tag	UNP Q66K64
B	-9	GLN	-	expression tag	UNP Q66K64
B	-8	PHE	-	expression tag	UNP Q66K64
B	-7	GLU	-	expression tag	UNP Q66K64
B	-6	LYS	-	expression tag	UNP Q66K64
B	-5	SER	-	expression tag	UNP Q66K64
B	-4	ALA	-	expression tag	UNP Q66K64
B	-3	VAL	-	expression tag	UNP Q66K64
B	-2	GLY	-	expression tag	UNP Q66K64
B	-1	LEU	-	expression tag	UNP Q66K64
B	0	ASN	-	expression tag	UNP Q66K64
B	1	ASP	-	expression tag	UNP Q66K64
B	2	ILE	-	expression tag	UNP Q66K64
B	3	PHE	-	expression tag	UNP Q66K64
B	4	GLU	-	expression tag	UNP Q66K64
B	5	ALA	-	expression tag	UNP Q66K64
B	6	GLN	-	expression tag	UNP Q66K64
B	7	LYS	-	expression tag	UNP Q66K64
B	8	ILE	-	expression tag	UNP Q66K64
B	9	GLU	-	expression tag	UNP Q66K64

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Chain	Residue	Modelled	Actual	Comment	Reference
B	10	TRP	-	expression tag	UNP Q66K64
B	11	HIS	-	expression tag	UNP Q66K64
B	12	GLU	-	expression tag	UNP Q66K64
B	13	GLY	-	expression tag	UNP Q66K64
B	14	GLY	-	expression tag	UNP Q66K64
B	15	GLY	-	expression tag	UNP Q66K64
B	16	GLY	-	expression tag	UNP Q66K64
B	17	SER	-	expression tag	UNP Q66K64
B	18	GLY	-	expression tag	UNP Q66K64
B	19	GLU	-	expression tag	UNP Q66K64
B	20	ASN	-	expression tag	UNP Q66K64
B	21	LEU	-	expression tag	UNP Q66K64
B	22	TYR	-	expression tag	UNP Q66K64
B	23	PHE	-	expression tag	UNP Q66K64
B	24	GLN	-	expression tag	UNP Q66K64
B	25	GLY	-	expression tag	UNP Q66K64
B	26	GLY	-	expression tag	UNP Q66K64
B	27	GLY	-	expression tag	UNP Q66K64
B	28	ARG	-	expression tag	UNP Q66K64
B	29	MET	-	expression tag	UNP Q66K64
B	30	GLY	-	expression tag	UNP Q66K64
B	31	ARG	-	expression tag	UNP Q66K64
B	32	ARG	-	expression tag	UNP Q66K64
B	33	ARG	-	expression tag	UNP Q66K64

- Molecule 3 is a protein called DDB1- and CUL4-associated factor 15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	192	Total	C	N	O	S	0	0	0
			1545	992	256	290	7			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	338	MET	-	initiating methionine	UNP Q66K64
C	339	ASP	-	expression tag	UNP Q66K64
C	340	TRP	-	expression tag	UNP Q66K64
C	341	SER	-	expression tag	UNP Q66K64
C	342	HIS	-	expression tag	UNP Q66K64
C	343	PRO	-	expression tag	UNP Q66K64
C	344	GLN	-	expression tag	UNP Q66K64
C	345	PHE	-	expression tag	UNP Q66K64

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Chain	Residue	Modelled	Actual	Comment	Reference
C	346	GLU	-	expression tag	UNP Q66K64
C	347	LYS	-	expression tag	UNP Q66K64
C	348	SER	-	expression tag	UNP Q66K64
C	349	ALA	-	expression tag	UNP Q66K64
C	350	VAL	-	expression tag	UNP Q66K64
C	351	GLY	-	expression tag	UNP Q66K64
C	352	LEU	-	expression tag	UNP Q66K64
C	353	ASN	-	expression tag	UNP Q66K64
C	354	ASP	-	expression tag	UNP Q66K64
C	355	ILE	-	expression tag	UNP Q66K64
C	356	PHE	-	expression tag	UNP Q66K64
C	357	GLU	-	expression tag	UNP Q66K64
C	358	ALA	-	expression tag	UNP Q66K64
C	359	GLN	-	expression tag	UNP Q66K64
C	360	LYS	-	expression tag	UNP Q66K64
C	361	ILE	-	expression tag	UNP Q66K64
C	362	GLU	-	expression tag	UNP Q66K64
C	363	TRP	-	expression tag	UNP Q66K64
C	364	HIS	-	expression tag	UNP Q66K64
C	365	GLU	-	expression tag	UNP Q66K64
C	366	GLY	-	expression tag	UNP Q66K64
C	367	GLY	-	expression tag	UNP Q66K64
C	368	GLY	-	expression tag	UNP Q66K64
C	369	GLY	-	expression tag	UNP Q66K64
C	370	SER	-	expression tag	UNP Q66K64
C	371	GLY	-	expression tag	UNP Q66K64
C	372	GLU	-	expression tag	UNP Q66K64
C	373	ASN	-	expression tag	UNP Q66K64
C	374	LEU	-	expression tag	UNP Q66K64
C	375	TYR	-	expression tag	UNP Q66K64
C	376	PHE	-	expression tag	UNP Q66K64
C	377	GLN	-	expression tag	UNP Q66K64
C	378	GLY	-	expression tag	UNP Q66K64
C	379	GLY	-	expression tag	UNP Q66K64
C	380	GLY	-	expression tag	UNP Q66K64
C	381	ARG	-	expression tag	UNP Q66K64
C	382	MET	-	expression tag	UNP Q66K64

- Molecule 4 is a protein called RNA-binding protein 39.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	79	Total	C	N	O	S	0	0	0
			619	392	106	115	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	226	MET	-	initiating methionine	UNP Q14498
D	227	GLY	-	expression tag	UNP Q14498
D	228	SER	-	expression tag	UNP Q14498
D	229	SER	-	expression tag	UNP Q14498
D	230	HIS	-	expression tag	UNP Q14498
D	231	HIS	-	expression tag	UNP Q14498
D	232	HIS	-	expression tag	UNP Q14498
D	233	HIS	-	expression tag	UNP Q14498
D	234	HIS	-	expression tag	UNP Q14498
D	235	HIS	-	expression tag	UNP Q14498
D	236	SER	-	expression tag	UNP Q14498
D	237	ALA	-	expression tag	UNP Q14498
D	238	VAL	-	expression tag	UNP Q14498
D	239	ASP	-	expression tag	UNP Q14498
D	240	GLU	-	expression tag	UNP Q14498
D	241	ASN	-	expression tag	UNP Q14498
D	242	LEU	-	expression tag	UNP Q14498
D	243	TYR	-	expression tag	UNP Q14498
D	244	PHE	-	expression tag	UNP Q14498
D	245	GLN	-	expression tag	UNP Q14498
D	246	GLY	-	expression tag	UNP Q14498
D	247	GLY	-	expression tag	UNP Q14498
D	248	GLY	-	expression tag	UNP Q14498
D	249	ARG	-	expression tag	UNP Q14498

- Molecule 5 is a protein called DET1- and DDB1-associated protein 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	E	62	Total	C	N	O	0	0	0
			520	336	91	93			

There are 24 discrepancies between the modelled and reference sequences:

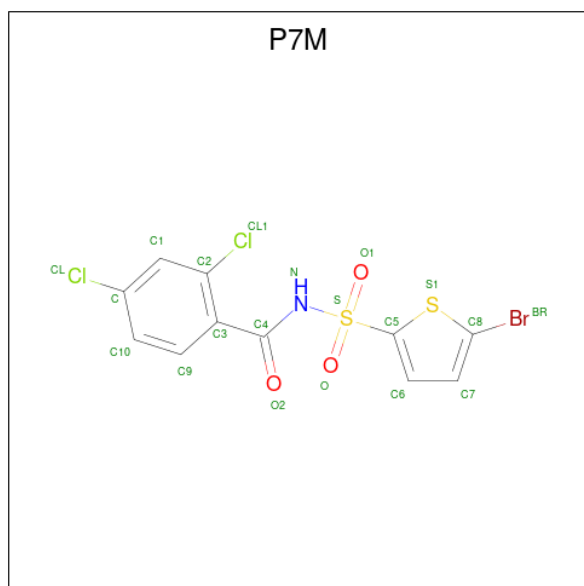
Chain	Residue	Modelled	Actual	Comment	Reference
E	-23	MET	-	initiating methionine	UNP Q9BW61
E	-22	GLY	-	expression tag	UNP Q9BW61
E	-21	SER	-	expression tag	UNP Q9BW61
E	-20	SER	-	expression tag	UNP Q9BW61
E	-19	HIS	-	expression tag	UNP Q9BW61
E	-18	HIS	-	expression tag	UNP Q9BW61
E	-17	HIS	-	expression tag	UNP Q9BW61

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-16	HIS	-	expression tag	UNP Q9BW61
E	-15	HIS	-	expression tag	UNP Q9BW61
E	-14	HIS	-	expression tag	UNP Q9BW61
E	-13	SER	-	expression tag	UNP Q9BW61
E	-12	ALA	-	expression tag	UNP Q9BW61
E	-11	VAL	-	expression tag	UNP Q9BW61
E	-10	ASP	-	expression tag	UNP Q9BW61
E	-9	GLU	-	expression tag	UNP Q9BW61
E	-8	ASN	-	expression tag	UNP Q9BW61
E	-7	LEU	-	expression tag	UNP Q9BW61
E	-6	TYR	-	expression tag	UNP Q9BW61
E	-5	PHE	-	expression tag	UNP Q9BW61
E	-4	GLN	-	expression tag	UNP Q9BW61
E	-3	GLY	-	expression tag	UNP Q9BW61
E	-2	GLY	-	expression tag	UNP Q9BW61
E	-1	GLY	-	expression tag	UNP Q9BW61
E	0	ARG	-	expression tag	UNP Q9BW61

- Molecule 6 is N-[(5-bromothiophen-2-yl)sulfonyl]-2,4-dichlorobenzamide (CCD ID: P7M) (formula: $C_{11}H_6BrCl_2NO_3S_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	Br	C	Cl	N	O	S		
6	B	1	20	1	11	2	1	3	2	0	0

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total 1	Zn 1	0	0

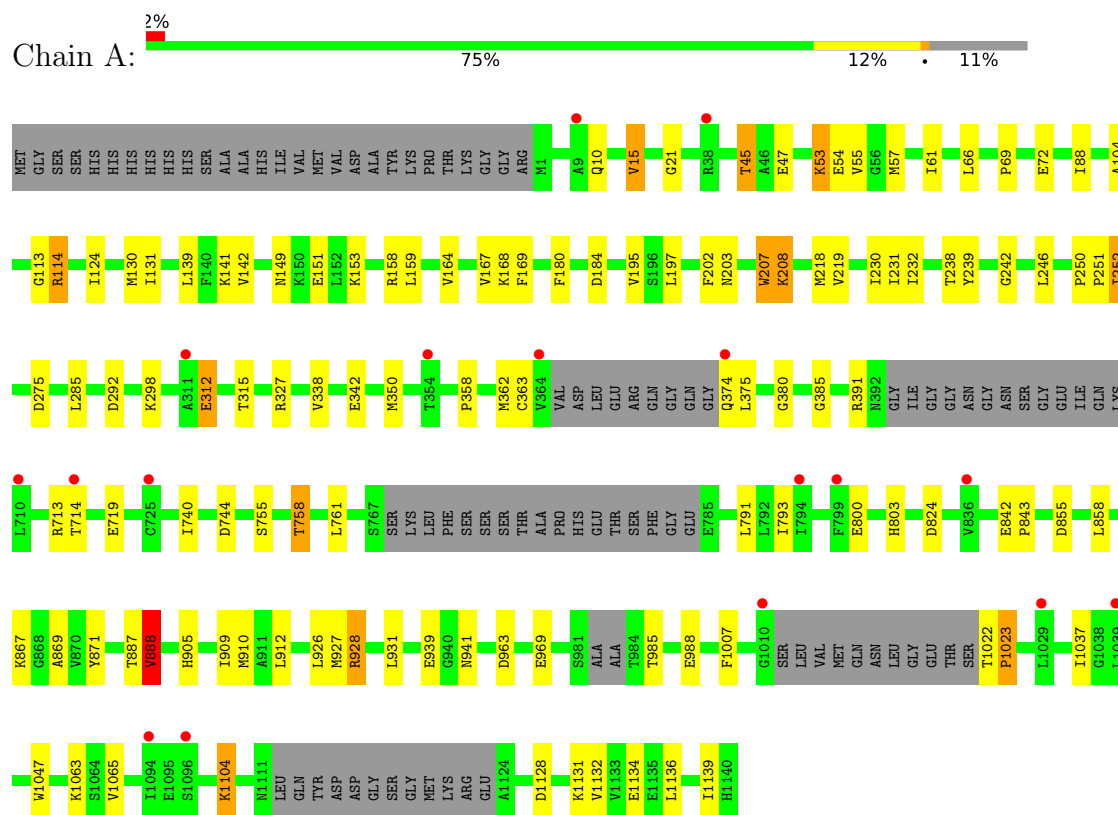
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	7	Total 7	O 7	0	0
8	B	6	Total 6	O 6	0	0
8	C	2	Total 2	O 2	0	0
8	D	1	Total 1	O 1	0	0
8	E	1	Total 1	O 1	0	0

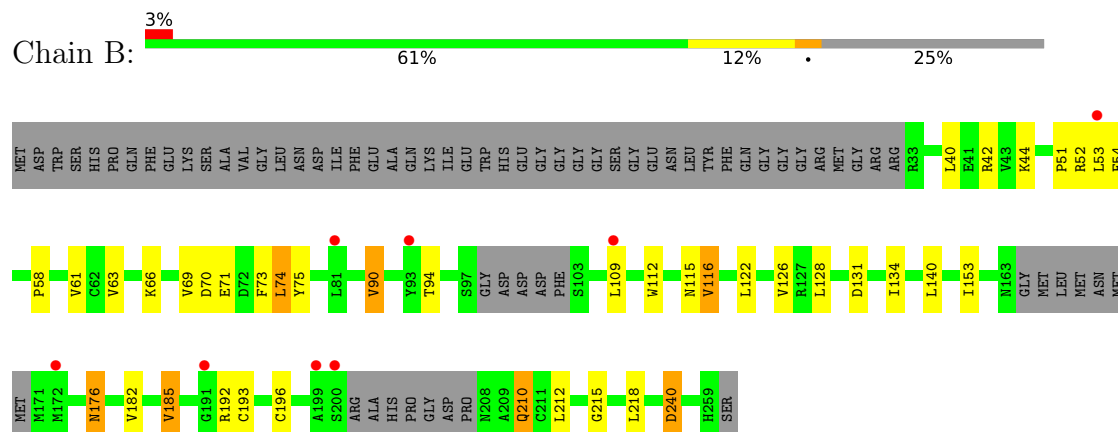
3 Residue-property plots [i](#)

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.

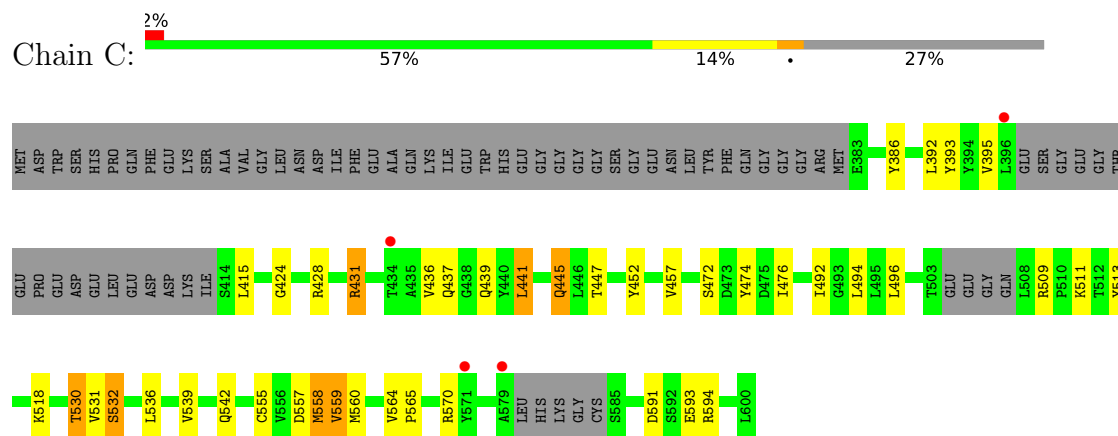
• Molecule 1: DNA damage-binding protein 1



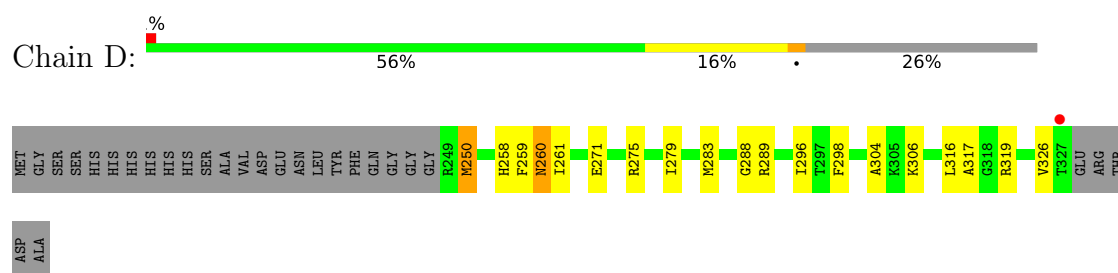
• Molecule 2: DDB1- and CUL4-associated factor 15



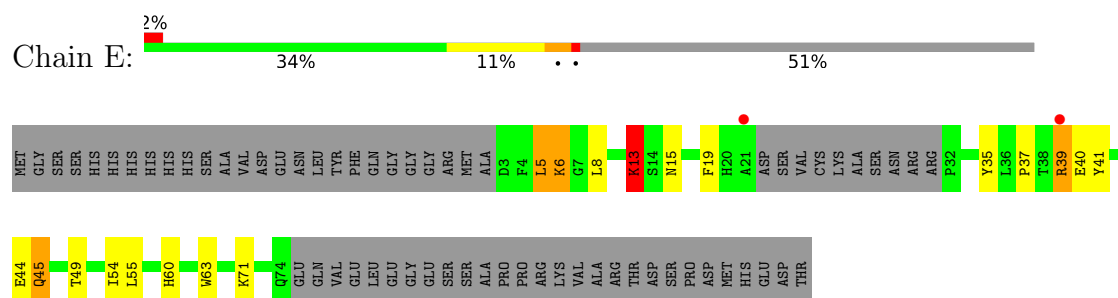
- Molecule 3: DDB1- and CUL4-associated factor 15



- Molecule 4: RNA-binding protein 39



- Molecule 5: DET1- and DDB1-associated protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.77Å 94.61Å 260.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.55 – 2.90 46.55 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.55-2.90) 99.7 (46.55-2.90)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.91Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.206 , 0.242 0.218 , 0.257	Depositor DCC
R_{free} test set	2249 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	85.7	Xtriage
Anisotropy	0.597	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 84.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10494	wwPDB-VP
Average B, all atoms (Å ²)	120.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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: ZN, P7M

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	0.87	1/6189 (0.0%)	1.24	23/8369 (0.3%)
2	B	0.91	0/1734	1.27	7/2349 (0.3%)
3	C	0.89	0/1579	1.27	8/2150 (0.4%)
4	D	0.90	0/630	1.36	5/841 (0.6%)
5	E	0.90	0/533	1.52	6/718 (0.8%)
All	All	0.88	1/10665 (0.0%)	1.27	49/14427 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	15	VAL	CA-CB	5.49	1.60	1.54

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	TRP	CA-C-N	8.54	134.03	121.31
1	A	207	TRP	C-N-CA	8.54	134.03	121.31
1	A	888	VAL	N-CA-CB	7.67	120.20	110.99
5	E	15	ASN	CA-C-N	-7.39	110.77	122.29
5	E	15	ASN	C-N-CA	-7.39	110.77	122.29
1	A	941	ASN	CA-CB-CG	6.63	119.23	112.60
1	A	15	VAL	N-CA-CB	6.60	118.93	111.21
5	E	13	LYS	CA-C-N	6.38	128.84	120.28
5	E	13	LYS	C-N-CA	6.38	128.84	120.28
1	A	164	VAL	N-CA-CB	6.28	118.52	110.99
3	C	591	ASP	CA-C-N	6.27	131.16	121.19
3	C	591	ASP	C-N-CA	6.27	131.16	121.19
2	B	54	PHE	CA-C-N	6.08	128.75	120.54
2	B	54	PHE	C-N-CA	6.08	128.75	120.54
1	A	184	ASP	CA-CB-CG	6.03	118.63	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	312	GLU	N-CA-C	-6.02	104.09	112.45
4	D	260	ASN	CA-CB-CG	5.94	118.54	112.60
1	A	203	ASN	CA-CB-CG	5.90	118.50	112.60
1	A	362	MET	CA-C-N	5.88	131.80	122.62
1	A	362	MET	C-N-CA	5.88	131.80	122.62
2	B	176	ASN	CA-CB-CG	5.72	118.32	112.60
1	A	197	LEU	CA-C-N	5.63	127.76	120.44
1	A	197	LEU	C-N-CA	5.63	127.76	120.44
3	C	570	ARG	CA-C-N	5.61	128.83	120.87
3	C	570	ARG	C-N-CA	5.61	128.83	120.87
2	B	240	ASP	N-CA-C	5.58	118.89	111.75
1	A	824	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	292	ASP	CA-CB-CG	5.51	118.11	112.60
2	B	58	PRO	N-CA-C	5.50	117.41	110.70
1	A	871	TYR	N-CA-C	5.49	117.35	111.36
1	A	275	ASP	CA-CB-CG	5.46	118.06	112.60
5	E	5	LEU	CA-C-N	5.45	128.58	120.90
5	E	5	LEU	C-N-CA	5.45	128.58	120.90
1	A	909	ILE	N-CA-C	-5.27	106.53	111.48
3	C	555	CYS	N-CA-C	-5.23	105.27	110.97
3	C	564	VAL	CB-CA-C	5.19	115.16	110.13
1	A	149	ASN	CA-C-N	5.18	127.48	120.38
1	A	149	ASN	C-N-CA	5.18	127.48	120.38
3	C	557	ASP	CA-C-N	5.15	127.14	120.44
3	C	557	ASP	C-N-CA	5.15	127.14	120.44
4	D	316	LEU	CA-C-N	5.07	131.23	121.54
4	D	316	LEU	C-N-CA	5.07	131.23	121.54
1	A	1007	PHE	CA-CB-CG	5.07	118.87	113.80
1	A	744	ASP	CA-CB-CG	5.03	117.63	112.60
2	B	42	ARG	CA-C-N	5.03	126.89	120.56
2	B	42	ARG	C-N-CA	5.03	126.89	120.56
1	A	202	PHE	CA-CB-CG	5.02	118.82	113.80
4	D	326	VAL	CA-C-N	5.01	130.72	121.70
4	D	326	VAL	C-N-CA	5.01	130.72	121.70

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	6082	0	6066	52	0
2	B	1690	0	1665	13	0
3	C	1545	0	1520	17	0
4	D	619	0	609	7	0
5	E	520	0	517	12	0
6	B	20	0	0	2	0
7	B	1	0	0	0	0
8	A	7	0	0	0	0
8	B	6	0	0	0	0
8	C	2	0	0	1	0
8	D	1	0	0	0	0
8	E	1	0	0	0	0
All	All	10494	0	10377	88	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 4.

All (88) 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:207:TRP:HB3	1:A:242:GLY:HA2	1.71	0.72
1:A:55:VAL:HG21	1:A:1065:VAL:HG21	1.74	0.70
1:A:53:LYS:HG2	5:E:35:TYR:HB2	1.80	0.62
2:B:70:ASP:HB3	2:B:73:PHE:HD2	1.64	0.62
1:A:130:MET:HE1	1:A:195:VAL:HG11	1.83	0.60
3:C:518:LYS:HB2	3:C:532:SER:HB3	1.83	0.60
1:A:130:MET:HE2	1:A:169:PHE:HZ	1.66	0.60
1:A:57:MET:HG3	1:A:61:ILE:HD11	1.84	0.59
1:A:219:VAL:HG12	1:A:231:ILE:HG22	1.86	0.58
3:C:560:MET:HE1	4:D:260:ASN:HD22	1.69	0.58
2:B:126:VAL:HG21	2:B:185:VAL:HG11	1.87	0.57
1:A:843:PRO:HG2	1:A:869:ALA:HB2	1.88	0.56
1:A:755:SER:H	1:A:758:THR:HG22	1.70	0.56
1:A:910:MET:HE2	1:A:912:LEU:HD21	1.87	0.55
1:A:113:GLY:HA2	8:C:701:HOH:O	2.07	0.55
6:B:301:P7M:CL1	3:C:559:VAL:HG11	2.44	0.55
2:B:193:CYS:HB3	2:B:196:CYS:HB2	1.89	0.55
1:A:374:GLN:HG2	1:A:391:ARG:HB3	1.88	0.54
1:A:793:ILE:HD13	1:A:803:HIS:HB3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:301:P7M:CL1	6:B:301:P7M:O2	2.64	0.53
5:E:35:TYR:CE2	5:E:37:PRO:HG3	2.44	0.53
4:D:250:MET:HE2	4:D:304:ALA:HB3	1.91	0.52
1:A:141:LYS:HZ1	5:E:49:THR:HG21	1.76	0.51
1:A:231:ILE:HG13	1:A:238:THR:OG1	2.10	0.51
1:A:374:GLN:HE21	1:A:391:ARG:HD3	1.77	0.51
1:A:312:GLU:HB2	1:A:327:ARG:HB2	1.93	0.50
1:A:358:PRO:HD2	1:A:380:GLY:HA2	1.94	0.50
1:A:928:ARG:HH22	2:B:115:ASN:HA	1.77	0.50
2:B:61:VAL:HG11	2:B:116:VAL:HG13	1.94	0.49
1:A:363:CYS:O	1:A:375:LEU:HD23	2.13	0.49
2:B:70:ASP:HB3	2:B:73:PHE:CD2	2.45	0.48
3:C:494:LEU:HD23	3:C:496:LEU:HD23	1.96	0.48
1:A:910:MET:HE1	2:B:40:LEU:HD11	1.95	0.48
3:C:560:MET:HE1	4:D:260:ASN:CB	2.44	0.48
1:A:985:THR:HB	1:A:988:GLU:HG2	1.96	0.47
1:A:888:VAL:HG22	1:A:905:HIS:HB3	1.95	0.47
1:A:385:GLY:HA3	1:A:719:GLU:O	2.14	0.47
1:A:793:ILE:HD11	1:A:858:LEU:HD22	1.96	0.47
1:A:1047:TRP:HZ3	1:A:1132:VAL:HG13	1.80	0.47
1:A:1136:LEU:O	1:A:1139:ILE:HG12	2.15	0.46
2:B:71:GLU:HA	2:B:74:LEU:HD13	1.97	0.46
1:A:114:ARG:HG2	3:C:565:PRO:HB2	1.97	0.46
1:A:791:LEU:HD23	1:A:858:LEU:HD21	1.97	0.46
1:A:230:ILE:HD11	1:A:285:LEU:HD21	1.98	0.46
1:A:1128:ASP:O	1:A:1131:LYS:HG3	2.15	0.46
1:A:69:PRO:HD2	1:A:72:GLU:HG3	1.97	0.45
3:C:539:VAL:HG13	3:C:542:GLN:HG3	1.98	0.45
4:D:298:PHE:HB2	4:D:304:ALA:HB2	1.97	0.45
5:E:6:LYS:HD3	5:E:6:LYS:H	1.82	0.45
1:A:239:TYR:HB3	1:A:246:LEU:HB2	1.98	0.44
1:A:1104:LYS:HE2	5:E:39:ARG:H	1.82	0.44
2:B:51:PRO:C	2:B:53:LEU:H	2.25	0.44
3:C:393:TYR:HD2	3:C:431:ARG:HA	1.82	0.44
1:A:167:VAL:HG23	1:A:180:PHE:HB3	1.99	0.44
1:A:88:ILE:HG12	1:A:104:ALA:HB3	1.99	0.44
1:A:21:GLY:HA3	1:A:66:LEU:HD13	1.99	0.44
1:A:45:THR:HG23	1:A:47:GLU:H	1.82	0.44
1:A:124:ILE:HG12	1:A:131:ILE:HG12	2.00	0.43
3:C:558:MET:HE2	5:E:55:LEU:HD11	2.00	0.43
5:E:44:GLU:O	5:E:45:GLN:HB2	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:196:CYS:SG	2:B:210:GLN:HG3	2.59	0.43
3:C:474:TYR:HB3	3:C:494:LEU:HD12	2.00	0.43
3:C:476:ILE:HD12	3:C:492:ILE:HG12	2.00	0.43
1:A:250:PRO:HA	1:A:251:PRO:HD3	1.94	0.43
2:B:212:LEU:HD21	3:C:441:LEU:HB2	2.01	0.42
3:C:511:LYS:HB2	3:C:513:TYR:CE1	2.54	0.42
4:D:258:HIS:HB3	4:D:261:ILE:HG12	2.00	0.42
1:A:926:LEU:HD23	1:A:927:MET:HE3	2.02	0.42
1:A:130:MET:HE2	1:A:169:PHE:CZ	2.51	0.42
1:A:208:LYS:HD2	1:A:208:LYS:H	1.84	0.42
1:A:151:GLU:HG3	1:A:153:LYS:HE2	2.01	0.42
1:A:350:MET:HE3	5:E:19:PHE:CZ	2.55	0.41
1:A:1063:LYS:HB2	5:E:37:PRO:HB2	2.02	0.41
1:A:252:ILE:H	1:A:252:ILE:HG13	1.55	0.41
1:A:218:MET:HG2	1:A:232:ILE:HB	2.03	0.41
1:A:10:GLN:HB3	1:A:1037:ILE:HB	2.02	0.41
3:C:415:LEU:H	3:C:445:GLN:HE22	1.68	0.41
4:D:279:ILE:HG12	4:D:296:ILE:HG12	2.02	0.41
2:B:90:VAL:HG13	2:B:112:TRP:HB2	2.03	0.41
3:C:386:TYR:HB3	3:C:452:TYR:CD1	2.56	0.41
1:A:740:ILE:N	1:A:740:ILE:HD12	2.36	0.41
5:E:13:LYS:HE2	5:E:13:LYS:H	1.86	0.40
1:A:1022:THR:HA	1:A:1023:PRO:HD3	1.94	0.40
3:C:457:VAL:HG12	3:C:530:THR:CG2	2.52	0.40
2:B:109:LEU:HB2	2:B:128:LEU:HD11	2.03	0.40
3:C:531:VAL:O	5:E:63:TRP:HB2	2.22	0.40
1:A:158:ARG:HB2	5:E:54:ILE:HD12	2.03	0.40
4:D:259:PHE:HB3	4:D:289:ARG:HD3	2.03	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	758/864 (88%)	726 (96%)	31 (4%)	1 (0%)	48	77
2	B	200/276 (72%)	186 (93%)	11 (6%)	3 (2%)	8	28
3	C	184/263 (70%)	162 (88%)	19 (10%)	3 (2%)	7	27
4	D	77/107 (72%)	73 (95%)	2 (3%)	2 (3%)	4	17
5	E	58/126 (46%)	50 (86%)	6 (10%)	2 (3%)	3	12
All	All	1277/1636 (78%)	1197 (94%)	69 (5%)	11 (1%)	14	41

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	436	VAL
4	D	288	GLY
4	D	317	ALA
5	E	39	ARG
5	E	45	GLN
2	B	116	VAL
3	C	424	GLY
3	C	439	GLN
1	A	1023	PRO
2	B	176	ASN
2	B	215	GLY

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	677/749 (90%)	645 (95%)	32 (5%)	23	56
2	B	189/242 (78%)	169 (89%)	20 (11%)	6	22
3	C	174/232 (75%)	157 (90%)	17 (10%)	7	25
4	D	66/89 (74%)	60 (91%)	6 (9%)	9	28
5	E	57/112 (51%)	49 (86%)	8 (14%)	3	11
All	All	1163/1424 (82%)	1080 (93%)	83 (7%)	13	40

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	VAL
1	A	45	THR
1	A	53	LYS
1	A	54	GLU
1	A	114	ARG
1	A	139	LEU
1	A	142	VAL
1	A	159	LEU
1	A	168	LYS
1	A	208	LYS
1	A	252	ILE
1	A	298	LYS
1	A	315	THR
1	A	338	VAL
1	A	342	GLU
1	A	713	ARG
1	A	714	THR
1	A	758	THR
1	A	761	LEU
1	A	800	GLU
1	A	842	GLU
1	A	855	ASP
1	A	867	LYS
1	A	887	THR
1	A	888	VAL
1	A	928	ARG
1	A	931	LEU
1	A	939	GLU
1	A	963	ASP
1	A	969	GLU
1	A	1104	LYS
1	A	1134	GLU
2	B	44	LYS
2	B	52	ARG
2	B	63	VAL
2	B	66	LYS
2	B	69	VAL
2	B	74	LEU
2	B	75	TYR
2	B	90	VAL
2	B	94	THR
2	B	122	LEU

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Mol	Chain	Res	Type
2	B	131	ASP
2	B	134	ILE
2	B	140	LEU
2	B	153	ILE
2	B	182	VAL
2	B	185	VAL
2	B	192	ARG
2	B	210	GLN
2	B	218	LEU
2	B	240	ASP
3	C	392	LEU
3	C	395	VAL
3	C	428	ARG
3	C	431	ARG
3	C	437	GLN
3	C	441	LEU
3	C	445	GLN
3	C	447	THR
3	C	472	SER
3	C	509	ARG
3	C	530	THR
3	C	532	SER
3	C	536	LEU
3	C	558	MET
3	C	559	VAL
3	C	593	GLU
3	C	594	ARG
4	D	250	MET
4	D	271	GLU
4	D	275	ARG
4	D	283	MET
4	D	306	LYS
4	D	319	ARG
5	E	5	LEU
5	E	6	LYS
5	E	8	LEU
5	E	13	LYS
5	E	40	GLU
5	E	41	TYR
5	E	60	HIS
5	E	71	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	149	ASN
1	A	156	ASN
1	A	203	ASN
1	A	209	GLN
1	A	319	ASN
1	A	337	ASN
1	A	374	GLN
1	A	759	GLN
1	A	797	HIS
1	A	852	GLN
1	A	859	GLN
1	A	1005	ASN
2	B	39	GLN
2	B	125	GLN
2	B	158	ASN
2	B	176	ASN
2	B	197	GLN
3	C	388	ASN
3	C	455	ASN
3	C	467	GLN
3	C	514	HIS
4	D	260	ASN
4	D	312	ASN
4	D	325	HIS
5	E	12	ASN
5	E	61	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
6	P7M	B	301	-	21,21,21	3.12	1 (4%)	29,31,31	1.11	4 (13%)

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
6	P7M	B	301	-	-	1/15/15/15	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	301	P7M	C5-S	-14.02	1.53	1.74

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	301	P7M	C2-C3-C4	-3.67	116.62	122.56
6	B	301	P7M	C4-N-S	-2.32	120.33	123.38
6	B	301	P7M	C9-C3-C4	2.28	124.61	118.53
6	B	301	P7M	O1-S-N	2.09	111.76	106.66

There are no chirality outliers.

All (1) torsion outliers are listed below:

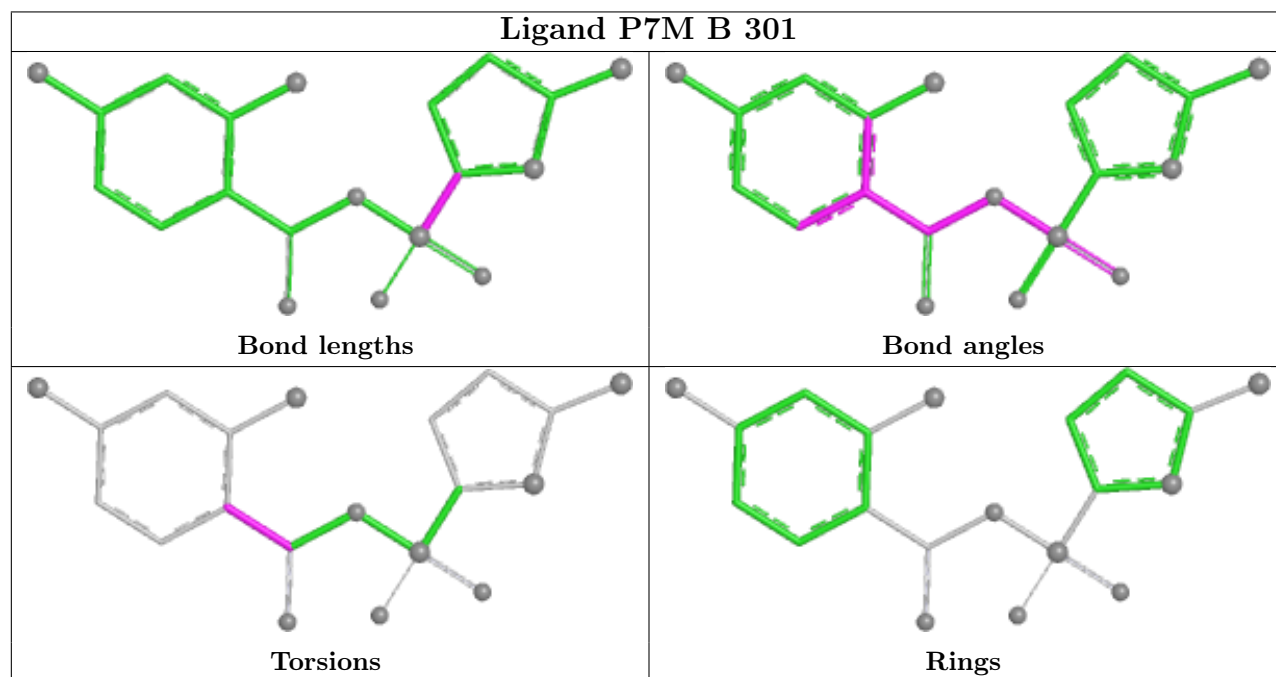
Mol	Chain	Res	Type	Atoms
6	B	301	P7M	C2-C3-C4-N

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	301	P7M	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	772/864 (89%)	0.07	17 (2%) 62 53	69, 120, 174, 190	0
2	B	208/276 (75%)	0.08	8 (3%) 44 36	67, 100, 156, 201	0
3	C	192/263 (73%)	0.05	4 (2%) 63 54	69, 102, 183, 208	0
4	D	79/107 (73%)	-0.02	1 (1%) 75 67	89, 114, 145, 155	0
5	E	62/126 (49%)	0.34	2 (3%) 50 41	88, 152, 193, 223	0
All	All	1313/1636 (80%)	0.08	32 (2%) 59 50	67, 114, 174, 223	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	E	21	ALA	4.3
1	A	9	ALA	4.1
2	B	200	SER	3.4
1	A	1029	LEU	3.3
2	B	109	LEU	3.2
1	A	1010	GLY	3.0
4	D	327	THR	3.0
1	A	799	PHE	2.9
3	C	571	TYR	2.9
1	A	714	THR	2.9
1	A	1039	LEU	2.7
2	B	172	MET	2.7
1	A	710	LEU	2.7
1	A	1096	SER	2.6
3	C	579	ALA	2.6
1	A	38	ARG	2.3
1	A	725	CYS	2.3
1	A	374	GLN	2.3
3	C	396	LEU	2.3
1	A	794	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	53	LEU	2.2
2	B	199	ALA	2.2
1	A	1094	ILE	2.2
5	E	39	ARG	2.2
1	A	354	THR	2.2
2	B	191	GLY	2.1
2	B	81	LEU	2.1
3	C	434	THR	2.1
1	A	311	ALA	2.0
2	B	93	TYR	2.0
1	A	836	VAL	2.0
1	A	364	VAL	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](#)

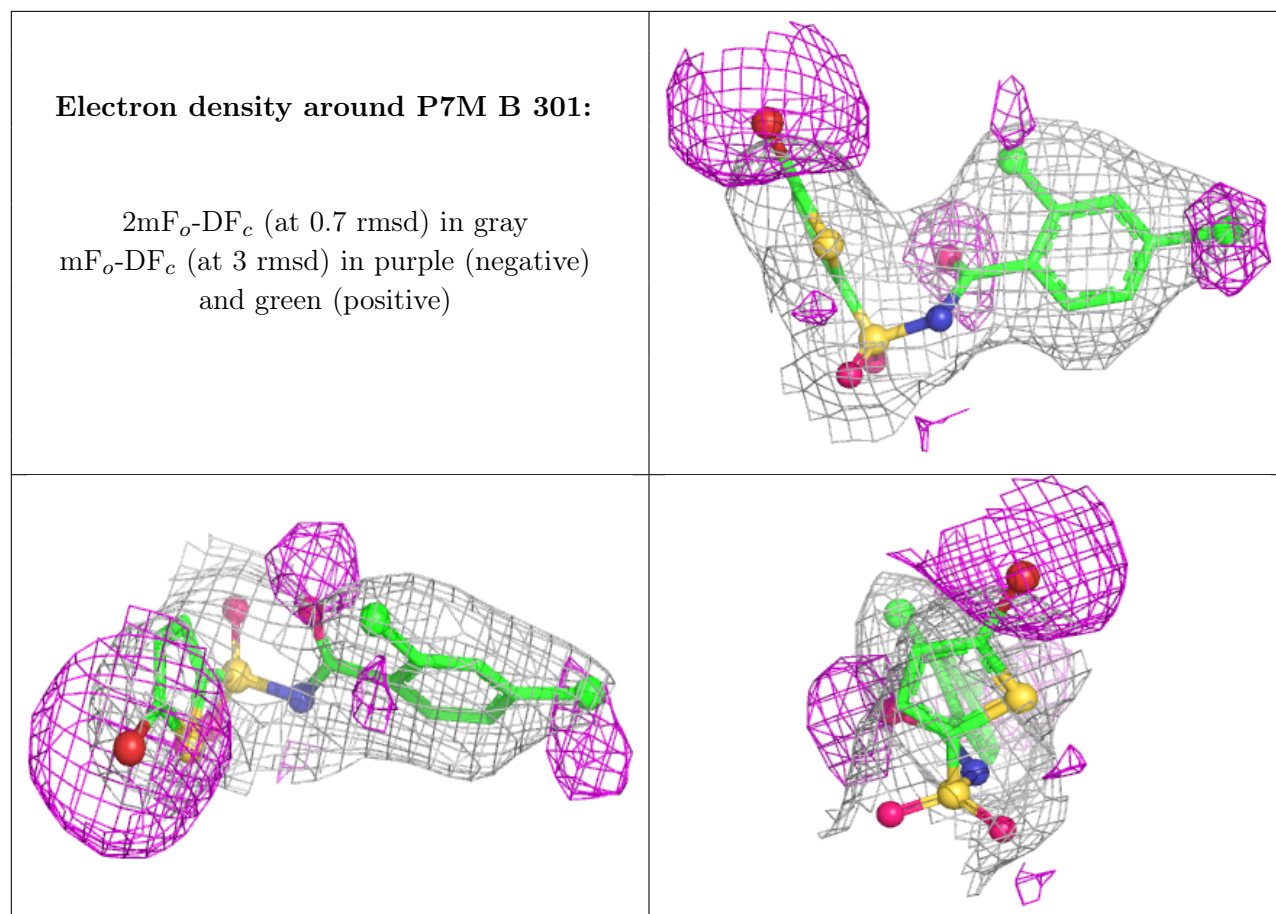
There are no oligosaccharides in this entry.

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
6	P7M	B	301	20/20	0.84	0.14	85,102,155,184	0
7	ZN	B	302	1/1	0.99	0.07	140,140,140,140	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.