



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:22 PM GMT

PDB ID : 4LCJ
Title : CtBP2 in complex with substrate MTOB
Authors : Hilbert, B.J.; Schiffer, C.A.; Royer Jr., W.E.
Deposited on : 2013-06-21
Resolution : 2.86 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

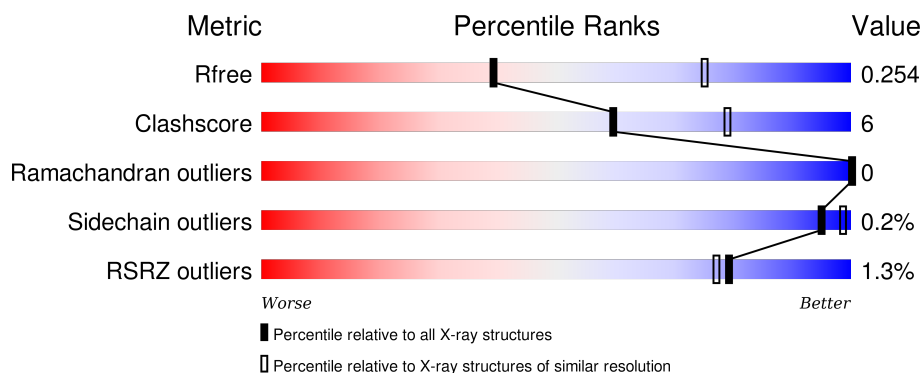
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.86 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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	349	<div> <div>83%</div> <div>11%</div> <div>5%</div> </div>
1	B	349	<div> <div>83%</div> <div>11%</div> <div>5%</div> </div>
1	C	349	<div> <div>85%</div> <div>9%</div> <div>5%</div> </div>
1	D	349	<div> <div>81%</div> <div>13%</div> <div>6%</div> </div>
1	E	349	<div> <div>83%</div> <div>11%</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	349	 % 83% 11% 6%
1	G	349	 3% 82% 11% 7%
1	H	349	 % 84% 11% 5%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	KMT	A	402	-	-	-	X
3	KMT	B	402	-	-	-	X
3	KMT	C	402	-	-	-	X
3	KMT	D	402	-	-	-	X
3	KMT	E	402	-	-	-	X
3	KMT	F	402	-	-	-	X
3	KMT	G	402	-	-	-	X
3	KMT	H	402	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 20058 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 C-terminal-binding protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	1	0
			2493	1568	442	471	12			
1	B	330	Total	C	N	O	S	0	1	0
			2463	1552	433	465	13			
1	C	330	Total	C	N	O	S	0	1	0
			2502	1574	445	470	13			
1	D	329	Total	C	N	O	S	0	1	0
			2403	1509	431	452	11			
1	E	327	Total	C	N	O	S	0	0	0
			2371	1483	425	450	13			
1	F	329	Total	C	N	O	S	0	0	0
			2443	1540	430	460	13			
1	G	325	Total	C	N	O	S	0	0	0
			2291	1424	409	445	13			
1	H	330	Total	C	N	O	S	0	1	0
			2463	1554	433	463	13			

There are 136 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	GLY	-	EXPRESSION TAG	UNP P56545
A	15	SER	-	EXPRESSION TAG	UNP P56545
A	16	HIS	-	EXPRESSION TAG	UNP P56545
A	17	MET	-	EXPRESSION TAG	UNP P56545
A	18	ALA	-	EXPRESSION TAG	UNP P56545
A	19	SER	-	EXPRESSION TAG	UNP P56545
A	20	MET	-	EXPRESSION TAG	UNP P56545
A	21	THR	-	EXPRESSION TAG	UNP P56545
A	22	GLY	-	EXPRESSION TAG	UNP P56545
A	23	GLY	-	EXPRESSION TAG	UNP P56545
A	24	GLN	-	EXPRESSION TAG	UNP P56545
A	25	GLN	-	EXPRESSION TAG	UNP P56545
A	26	MET	-	EXPRESSION TAG	UNP P56545

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Chain	Residue	Modelled	Actual	Comment	Reference
A	27	GLY	-	EXPRESSION TAG	UNP P56545
A	28	ARG	-	EXPRESSION TAG	UNP P56545
A	29	GLY	-	EXPRESSION TAG	UNP P56545
A	30	SER	-	EXPRESSION TAG	UNP P56545
B	14	GLY	-	EXPRESSION TAG	UNP P56545
B	15	SER	-	EXPRESSION TAG	UNP P56545
B	16	HIS	-	EXPRESSION TAG	UNP P56545
B	17	MET	-	EXPRESSION TAG	UNP P56545
B	18	ALA	-	EXPRESSION TAG	UNP P56545
B	19	SER	-	EXPRESSION TAG	UNP P56545
B	20	MET	-	EXPRESSION TAG	UNP P56545
B	21	THR	-	EXPRESSION TAG	UNP P56545
B	22	GLY	-	EXPRESSION TAG	UNP P56545
B	23	GLY	-	EXPRESSION TAG	UNP P56545
B	24	GLN	-	EXPRESSION TAG	UNP P56545
B	25	GLN	-	EXPRESSION TAG	UNP P56545
B	26	MET	-	EXPRESSION TAG	UNP P56545
B	27	GLY	-	EXPRESSION TAG	UNP P56545
B	28	ARG	-	EXPRESSION TAG	UNP P56545
B	29	GLY	-	EXPRESSION TAG	UNP P56545
B	30	SER	-	EXPRESSION TAG	UNP P56545
C	14	GLY	-	EXPRESSION TAG	UNP P56545
C	15	SER	-	EXPRESSION TAG	UNP P56545
C	16	HIS	-	EXPRESSION TAG	UNP P56545
C	17	MET	-	EXPRESSION TAG	UNP P56545
C	18	ALA	-	EXPRESSION TAG	UNP P56545
C	19	SER	-	EXPRESSION TAG	UNP P56545
C	20	MET	-	EXPRESSION TAG	UNP P56545
C	21	THR	-	EXPRESSION TAG	UNP P56545
C	22	GLY	-	EXPRESSION TAG	UNP P56545
C	23	GLY	-	EXPRESSION TAG	UNP P56545
C	24	GLN	-	EXPRESSION TAG	UNP P56545
C	25	GLN	-	EXPRESSION TAG	UNP P56545
C	26	MET	-	EXPRESSION TAG	UNP P56545
C	27	GLY	-	EXPRESSION TAG	UNP P56545
C	28	ARG	-	EXPRESSION TAG	UNP P56545
C	29	GLY	-	EXPRESSION TAG	UNP P56545
C	30	SER	-	EXPRESSION TAG	UNP P56545
D	14	GLY	-	EXPRESSION TAG	UNP P56545
D	15	SER	-	EXPRESSION TAG	UNP P56545
D	16	HIS	-	EXPRESSION TAG	UNP P56545
D	17	MET	-	EXPRESSION TAG	UNP P56545

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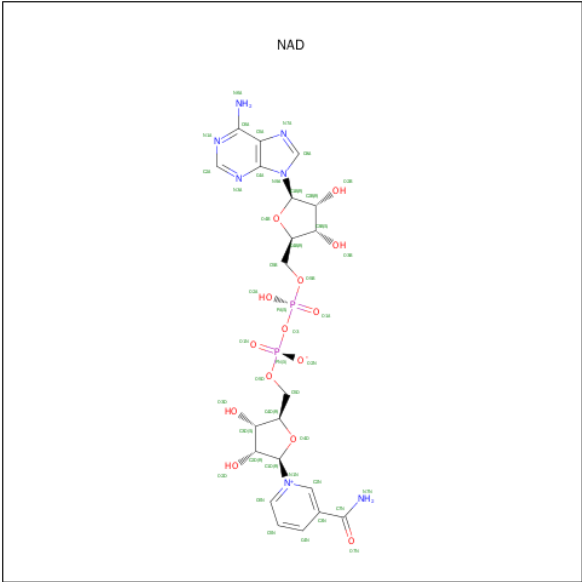
Chain	Residue	Modelled	Actual	Comment	Reference
D	18	ALA	-	EXPRESSION TAG	UNP P56545
D	19	SER	-	EXPRESSION TAG	UNP P56545
D	20	MET	-	EXPRESSION TAG	UNP P56545
D	21	THR	-	EXPRESSION TAG	UNP P56545
D	22	GLY	-	EXPRESSION TAG	UNP P56545
D	23	GLY	-	EXPRESSION TAG	UNP P56545
D	24	GLN	-	EXPRESSION TAG	UNP P56545
D	25	GLN	-	EXPRESSION TAG	UNP P56545
D	26	MET	-	EXPRESSION TAG	UNP P56545
D	27	GLY	-	EXPRESSION TAG	UNP P56545
D	28	ARG	-	EXPRESSION TAG	UNP P56545
D	29	GLY	-	EXPRESSION TAG	UNP P56545
D	30	SER	-	EXPRESSION TAG	UNP P56545
E	14	GLY	-	EXPRESSION TAG	UNP P56545
E	15	SER	-	EXPRESSION TAG	UNP P56545
E	16	HIS	-	EXPRESSION TAG	UNP P56545
E	17	MET	-	EXPRESSION TAG	UNP P56545
E	18	ALA	-	EXPRESSION TAG	UNP P56545
E	19	SER	-	EXPRESSION TAG	UNP P56545
E	20	MET	-	EXPRESSION TAG	UNP P56545
E	21	THR	-	EXPRESSION TAG	UNP P56545
E	22	GLY	-	EXPRESSION TAG	UNP P56545
E	23	GLY	-	EXPRESSION TAG	UNP P56545
E	24	GLN	-	EXPRESSION TAG	UNP P56545
E	25	GLN	-	EXPRESSION TAG	UNP P56545
E	26	MET	-	EXPRESSION TAG	UNP P56545
E	27	GLY	-	EXPRESSION TAG	UNP P56545
E	28	ARG	-	EXPRESSION TAG	UNP P56545
E	29	GLY	-	EXPRESSION TAG	UNP P56545
E	30	SER	-	EXPRESSION TAG	UNP P56545
F	14	GLY	-	EXPRESSION TAG	UNP P56545
F	15	SER	-	EXPRESSION TAG	UNP P56545
F	16	HIS	-	EXPRESSION TAG	UNP P56545
F	17	MET	-	EXPRESSION TAG	UNP P56545
F	18	ALA	-	EXPRESSION TAG	UNP P56545
F	19	SER	-	EXPRESSION TAG	UNP P56545
F	20	MET	-	EXPRESSION TAG	UNP P56545
F	21	THR	-	EXPRESSION TAG	UNP P56545
F	22	GLY	-	EXPRESSION TAG	UNP P56545
F	23	GLY	-	EXPRESSION TAG	UNP P56545
F	24	GLN	-	EXPRESSION TAG	UNP P56545
F	25	GLN	-	EXPRESSION TAG	UNP P56545

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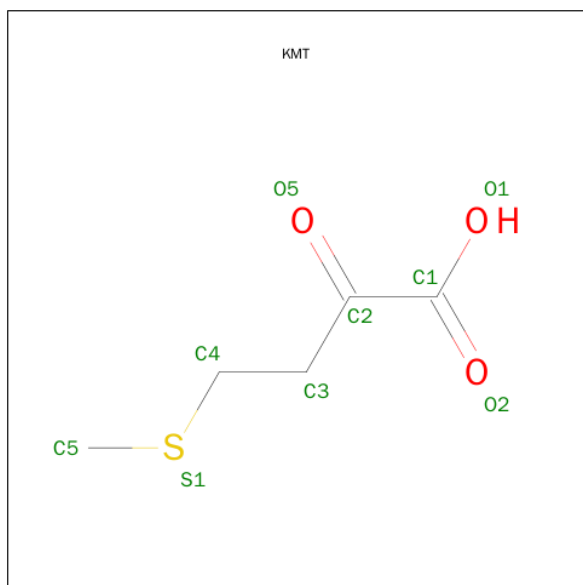
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F	26	MET	-	EXPRESSION TAG	UNP P56545
F	27	GLY	-	EXPRESSION TAG	UNP P56545
F	28	ARG	-	EXPRESSION TAG	UNP P56545
F	29	GLY	-	EXPRESSION TAG	UNP P56545
F	30	SER	-	EXPRESSION TAG	UNP P56545
G	14	GLY	-	EXPRESSION TAG	UNP P56545
G	15	SER	-	EXPRESSION TAG	UNP P56545
G	16	HIS	-	EXPRESSION TAG	UNP P56545
G	17	MET	-	EXPRESSION TAG	UNP P56545
G	18	ALA	-	EXPRESSION TAG	UNP P56545
G	19	SER	-	EXPRESSION TAG	UNP P56545
G	20	MET	-	EXPRESSION TAG	UNP P56545
G	21	THR	-	EXPRESSION TAG	UNP P56545
G	22	GLY	-	EXPRESSION TAG	UNP P56545
G	23	GLY	-	EXPRESSION TAG	UNP P56545
G	24	GLN	-	EXPRESSION TAG	UNP P56545
G	25	GLN	-	EXPRESSION TAG	UNP P56545
G	26	MET	-	EXPRESSION TAG	UNP P56545
G	27	GLY	-	EXPRESSION TAG	UNP P56545
G	28	ARG	-	EXPRESSION TAG	UNP P56545
G	29	GLY	-	EXPRESSION TAG	UNP P56545
G	30	SER	-	EXPRESSION TAG	UNP P56545
H	14	GLY	-	EXPRESSION TAG	UNP P56545
H	15	SER	-	EXPRESSION TAG	UNP P56545
H	16	HIS	-	EXPRESSION TAG	UNP P56545
H	17	MET	-	EXPRESSION TAG	UNP P56545
H	18	ALA	-	EXPRESSION TAG	UNP P56545
H	19	SER	-	EXPRESSION TAG	UNP P56545
H	20	MET	-	EXPRESSION TAG	UNP P56545
H	21	THR	-	EXPRESSION TAG	UNP P56545
H	22	GLY	-	EXPRESSION TAG	UNP P56545
H	23	GLY	-	EXPRESSION TAG	UNP P56545
H	24	GLN	-	EXPRESSION TAG	UNP P56545
H	25	GLN	-	EXPRESSION TAG	UNP P56545
H	26	MET	-	EXPRESSION TAG	UNP P56545
H	27	GLY	-	EXPRESSION TAG	UNP P56545
H	28	ARG	-	EXPRESSION TAG	UNP P56545
H	29	GLY	-	EXPRESSION TAG	UNP P56545
H	30	SER	-	EXPRESSION TAG	UNP P56545

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	G	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	H	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is 4-(METHYLSULFANYL)-2-OXOBUTANOIC ACID (three-letter code: KMT) (formula: C₅H₈O₃S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			9	5	3	1		
3	B	1	Total	C	O	S	0	0
			9	5	3	1		
3	C	1	Total	C	O	S	0	0
			9	5	3	1		
3	D	1	Total	C	O	S	0	0
			9	5	3	1		
3	E	1	Total	C	O	S	0	0
			9	5	3	1		
3	F	1	Total	C	O	S	0	0
			9	5	3	1		
3	G	1	Total	C	O	S	0	0
			9	5	3	1		
3	H	1	Total	C	O	S	0	0
			9	5	3	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	33	Total	O	0	0
			33	33		
4	B	29	Total	O	0	0
			29	29		
4	C	30	Total	O	0	0
			30	30		
4	D	18	Total	O	0	0
			18	18		

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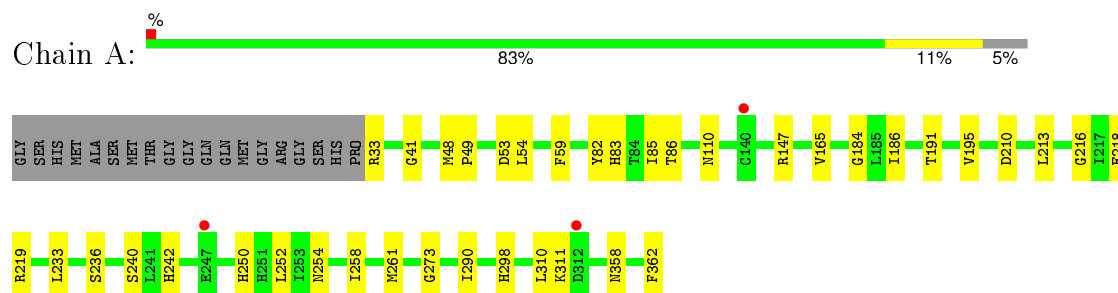
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	22	Total 22	O 22	0	0
4	F	32	Total 32	O 32	0	0
4	G	18	Total 18	O 18	0	0
4	H	23	Total 23	O 23	0	0

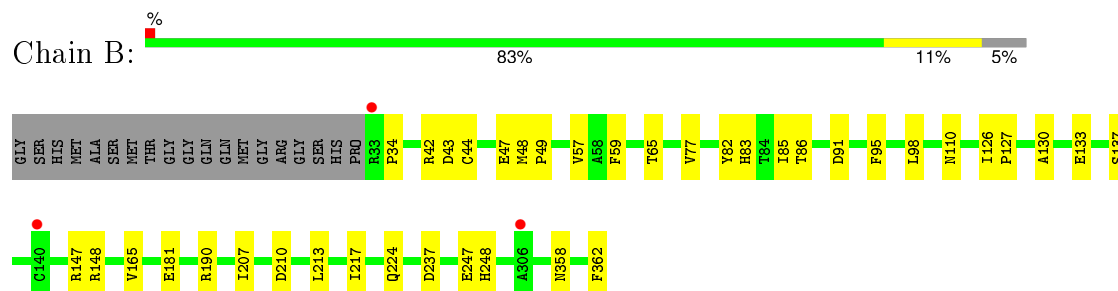
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

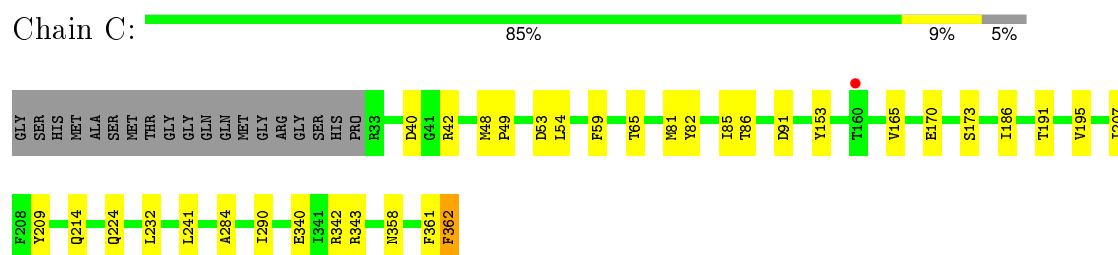
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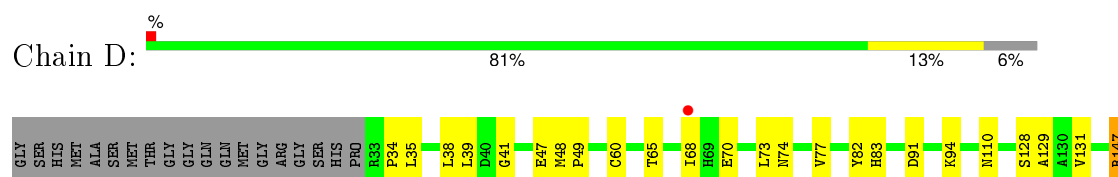
• Molecule 1: C-terminal-binding protein 2



• Molecule 1: C-terminal-binding protein 2

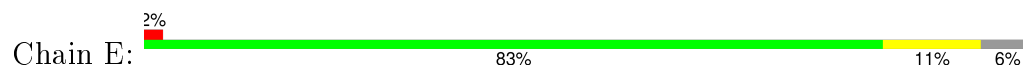


• Molecule 1: C-terminal-binding protein 2

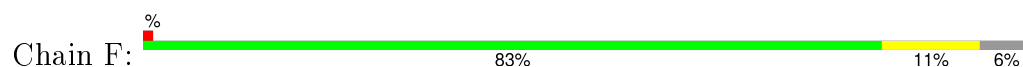




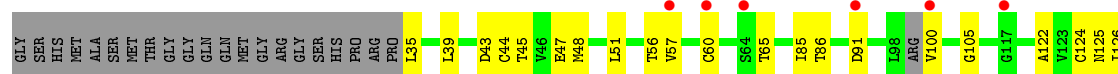
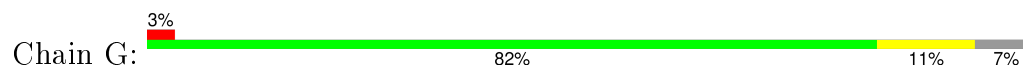
● Molecule 1: C-terminal-binding protein 2



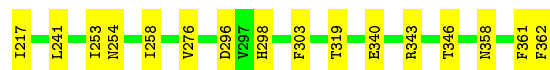
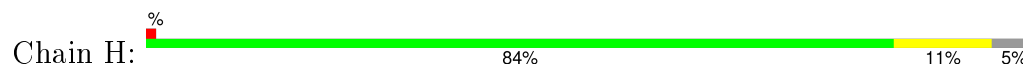
● Molecule 1: C-terminal-binding protein 2



● Molecule 1: C-terminal-binding protein 2



● Molecule 1: C-terminal-binding protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.15Å 140.60Å 135.13Å 90.00° 97.87° 90.00°	Depositor
Resolution (Å)	32.30 – 2.86 32.30 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.2 (32.30-2.86) 99.2 (32.30-2.86)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 2.85Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.214 , 0.251 0.224 , 0.254	Depositor DCC
R_{free} test set	3689 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	54.1	Xtriage
Anisotropy	0.476	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 27.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	1 of 73327 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20058	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: KMT, NAD

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	0.45	0/2535	0.62	0/3447
1	B	0.44	0/2508	0.62	1/3415 (0.0%)
1	C	0.48	0/2547	0.62	0/3460
1	D	0.44	0/2445	0.59	1/3333 (0.0%)
1	E	0.44	0/2409	0.58	0/3284
1	F	0.43	0/2484	0.59	0/3380
1	G	0.43	0/2325	0.60	0/3172
1	H	0.45	0/2509	0.63	2/3413 (0.1%)
All	All	0.45	0/19762	0.61	4/26904 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	147	ARG	NE-CZ-NH1	5.90	123.25	120.30
1	B	147	ARG	NE-CZ-NH1	5.71	123.15	120.30
1	H	147	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	H	147	ARG	NE-CZ-NH1	5.09	122.84	120.30

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	2493	0	2402	25	0
1	B	2463	0	2360	27	0
1	C	2502	0	2431	28	0
1	D	2403	0	2266	37	0
1	E	2371	0	2222	32	0
1	F	2443	0	2342	29	0
1	G	2291	0	2050	26	0
1	H	2463	0	2367	27	0
2	A	44	0	26	0	0
2	B	44	0	26	1	0
2	C	44	0	26	0	0
2	D	44	0	26	0	0
2	E	44	0	26	1	0
2	F	44	0	26	0	0
2	G	44	0	26	0	0
2	H	44	0	26	0	0
3	A	9	0	7	0	0
3	B	9	0	7	1	0
3	C	9	0	7	1	0
3	D	9	0	7	1	0
3	E	9	0	7	0	0
3	F	9	0	7	0	0
3	G	9	0	7	0	0
3	H	9	0	7	1	0
4	A	33	0	0	1	0
4	B	29	0	0	1	0
4	C	30	0	0	1	0
4	D	18	0	0	0	0
4	E	22	0	0	1	0
4	F	32	0	0	0	0
4	G	18	0	0	0	0
4	H	23	0	0	0	0
All	All	20058	0	18704	218	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 6.

All (218) 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:D:73:LEU:HD21	1:D:94:LYS:O	1.86	0.74
1:D:283:GLN:NE2	1:D:287:GLU:OE2	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:402:KMT:H41	3:C:402:KMT:O1	1.90	0.71
1:D:65:THR:HG21	1:D:91:ASP:OD1	1.89	0.71
1:D:65:THR:HA	1:D:68:ILE:HG13	1.73	0.70
1:B:48:MET:HB3	1:B:49:PRO:CD	2.24	0.67
1:C:165:VAL:HG21	1:D:82:TYR:CG	2.30	0.66
1:E:47:GLU:N	1:E:47:GLU:OE2	2.29	0.66
1:H:89:ARG:NH2	1:H:118:GLU:OE1	2.28	0.65
1:F:216:GLY:HA2	1:G:128:SER:HB2	1.78	0.65
1:E:48:MET:SD	1:E:57:VAL:CG2	2.85	0.65
1:E:48:MET:SD	1:E:57:VAL:HG21	2.36	0.64
1:A:53:ASP:OD2	1:A:54:LEU:CD1	2.46	0.64
1:F:298:HIS:CD2	1:F:303:PHE:CD2	2.86	0.64
1:G:47:GLU:OE1	1:G:47:GLU:N	2.30	0.64
1:G:132:GLU:OE1	1:H:179:ARG:NH1	2.30	0.63
3:B:402:KMT:H41	3:B:402:KMT:O2	1.99	0.63
1:E:48:MET:CE	1:E:57:VAL:CG2	2.76	0.63
1:E:210:ASP:OD2	2:E:401:NAD:H1B	2.00	0.62
1:B:48:MET:HB3	1:B:49:PRO:HD3	1.83	0.61
1:C:65:THR:HA	1:C:85:ILE:HD11	1.81	0.61
1:A:53:ASP:OD2	1:A:54:LEU:HD13	2.01	0.60
1:E:35:LEU:O	1:E:77:VAL:HG12	2.02	0.60
1:D:38:LEU:HD21	1:D:41:GLY:HA3	1.84	0.60
1:G:165:VAL:HG21	1:H:82:TYR:CG	2.36	0.60
1:F:81:MET:HE3	1:F:85:ILE:HG23	1.85	0.58
1:D:38:LEU:HD21	1:D:41:GLY:CA	2.33	0.58
1:F:48:MET:HG3	1:F:57:VAL:HG21	1.85	0.57
1:E:48:MET:CE	1:E:57:VAL:HG23	2.34	0.56
1:C:284:ALA:HB1	1:C:290:ILE:HG12	1.87	0.56
1:G:44:CYS:SG	1:G:57:VAL:CG1	2.94	0.56
1:D:47:GLU:N	1:D:47:GLU:OE2	2.38	0.56
1:E:82:TYR:CG	1:F:165:VAL:HG21	2.41	0.56
1:H:44:CYS:SG	1:H:57:VAL:HG12	2.44	0.56
1:E:65:THR:HG21	1:E:91:ASP:OD1	2.05	0.56
1:A:165:VAL:HG21	1:B:82:TYR:CG	2.40	0.56
1:F:188:PHE:HB2	1:F:208:PHE:CD2	2.41	0.56
1:D:65:THR:OG1	1:D:68:ILE:HD12	2.05	0.55
1:E:147:ARG:NH2	1:F:325:TYR:O	2.30	0.55
1:E:346:THR:HG22	1:E:346:THR:O	2.07	0.55
1:A:186:ILE:HG21	1:A:252:LEU:HD21	1.87	0.55
1:H:38:LEU:HD21	1:H:41:GLY:HA3	1.88	0.55
1:C:53:ASP:OD2	1:C:54:LEU:CD1	2.55	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:126:ILE:HD11	1:G:337:ALA:HA	1.88	0.55
1:G:35:LEU:HD12	1:G:56:THR:O	2.08	0.54
1:G:44:CYS:SG	1:G:57:VAL:HG11	2.49	0.53
1:C:82:TYR:CG	1:D:165:VAL:HG21	2.43	0.53
1:B:247:GLU:HG3	1:B:248:HIS:CE1	2.43	0.53
1:G:207:ILE:HA	1:G:224:GLN:O	2.08	0.53
1:G:295:LEU:O	1:G:318:CYS:HA	2.08	0.53
1:A:33:ARG:NH1	1:A:53:ASP:O	2.42	0.53
1:B:210:ASP:OD2	2:B:401:NAD:H1B	2.09	0.53
1:D:284:ALA:HB1	1:D:290:ILE:HG12	1.91	0.52
1:B:48:MET:SD	1:B:57:VAL:HG21	2.50	0.52
1:A:83:HIS:HA	1:A:110:ASN:ND2	2.25	0.52
1:F:48:MET:N	1:F:49:PRO:HD2	2.24	0.52
1:E:48:MET:HB3	1:E:49:PRO:CD	2.40	0.52
1:E:48:MET:HE2	1:E:57:VAL:CG2	2.40	0.52
1:B:34:PRO:HB3	1:B:77:VAL:HG11	1.90	0.52
1:D:48:MET:N	1:D:49:PRO:HD2	2.25	0.52
1:E:282:ALA:HB2	1:E:310:LEU:HD23	1.91	0.51
1:D:65:THR:HG21	1:D:91:ASP:CG	2.31	0.51
1:C:53:ASP:OD2	1:C:54:LEU:HD13	2.10	0.51
1:A:82:TYR:CG	1:B:165:VAL:HG21	2.45	0.51
1:B:213:LEU:HD22	1:B:217:ILE:HD12	1.92	0.50
1:D:266:PHE:CZ	1:D:291:ARG:HD2	2.47	0.50
1:F:346:THR:HG22	1:F:346:THR:O	2.10	0.50
1:B:190:ARG:HH22	1:C:214:GLN:CD	2.15	0.50
1:C:340:GLU:OE1	1:C:343:ARG:NH1	2.44	0.50
1:B:95:PHE:HB3	1:B:98:LEU:HB2	1.93	0.50
1:B:65:THR:HG21	1:B:91:ASP:OD1	2.11	0.49
1:C:186:ILE:HB	1:C:241:LEU:HD23	1.93	0.49
1:H:44:CYS:SG	1:H:57:VAL:CG1	3.01	0.49
1:C:186:ILE:HD12	1:C:241:LEU:CD2	2.43	0.49
1:G:65:THR:CB	1:G:91:ASP:OD1	2.60	0.49
1:C:361:PHE:O	1:C:362:PHE:HB2	2.12	0.49
1:F:207:ILE:HA	1:F:224:GLN:O	2.12	0.49
1:A:261:MET:CE	1:A:290:ILE:HD11	2.42	0.49
1:F:83:HIS:HA	1:F:110:ASN:ND2	2.27	0.49
1:F:233:LEU:HD22	1:F:261:MET:HG2	1.93	0.49
1:E:48:MET:HE2	1:E:57:VAL:HG23	1.94	0.48
1:A:147:ARG:HD3	1:B:137:SER:OG	2.14	0.48
1:H:254:ASN:O	1:H:258:ILE:HG13	2.13	0.48
1:F:117:GLY:O	1:F:361:PHE:O	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:38:LEU:HD21	1:F:41:GLY:HA3	1.95	0.48
1:H:33:ARG:O	1:H:34:PRO:C	2.52	0.48
1:H:59:PHE:C	1:H:59:PHE:CD1	2.87	0.48
1:E:302:PRO:HB3	1:F:162:VAL:O	2.13	0.48
1:D:73:LEU:CD2	1:D:94:LYS:O	2.57	0.48
1:E:354:ARG:NH2	1:H:215:ASP:OD1	2.47	0.48
1:E:128:SER:OG	1:H:216:GLY:HA2	2.14	0.48
1:C:342:ARG:NH2	4:C:526:HOH:O	2.47	0.48
1:H:358:ASN:CG	1:H:358:ASN:O	2.52	0.48
1:E:43:ASP:OD1	1:E:45:THR:HG23	2.14	0.48
1:F:287:GLU:OE1	1:F:289:ARG:NE	2.47	0.48
1:G:43:ASP:OD1	1:G:45:THR:HG23	2.14	0.47
1:A:358:ASN:CG	1:A:358:ASN:O	2.53	0.47
1:C:53:ASP:CG	1:C:54:LEU:HD13	2.35	0.47
1:D:147:ARG:O	1:D:148:ARG:HB2	2.14	0.47
1:D:34:PRO:HB2	1:D:77:VAL:HG11	1.96	0.47
1:D:83:HIS:HA	1:D:110:ASN:ND2	2.30	0.47
1:D:246:ASN:O	1:D:250[B]:HIS:CE1	2.68	0.47
1:F:124:CYS:SG	1:F:353:LEU:HD13	2.54	0.47
1:A:218:GLU:HG2	1:A:219:ARG:N	2.29	0.47
1:D:255:ASP:O	1:D:259:LYS:HG3	2.15	0.47
1:G:144:ASN:ND2	1:G:149:ASN:HB2	2.30	0.47
1:F:209:TYR:HB3	1:F:232:LEU:HD13	1.97	0.47
1:H:54:LEU:HD21	1:H:346:THR:CG2	2.44	0.46
1:F:140:CYS:O	1:F:144:ASN:OD1	2.33	0.46
1:G:163:GLN:N	1:G:167:GLN:OE1	2.41	0.46
1:F:38:LEU:CD2	1:F:41:GLY:HA3	2.45	0.46
1:A:240:SER:OG	1:A:242:HIS:CE1	2.69	0.46
1:D:73:LEU:HD23	1:D:73:LEU:C	2.36	0.46
1:H:33:ARG:HB2	1:H:56:THR:OG1	2.15	0.46
1:D:230:GLN:HG2	1:D:234:TYR:CE2	2.51	0.46
1:D:233:LEU:HD22	1:D:261:MET:HG2	1.97	0.46
1:C:207:ILE:HG22	1:C:224:GLN:HB3	1.97	0.46
1:H:213:LEU:HD13	1:H:217:ILE:HD12	1.98	0.46
1:B:181:GLU:OE1	1:B:237:ASP:HB3	2.16	0.45
1:E:257:THR:O	1:E:260:GLN:HB2	2.15	0.45
3:D:402:KMT:O2	3:D:402:KMT:H41	2.16	0.45
1:F:81:MET:CE	1:F:85:ILE:HG23	2.46	0.45
1:F:344:ALA:HB1	1:F:349:ILE:HD11	1.98	0.45
1:F:48:MET:HB3	1:F:49:PRO:CD	2.47	0.45
1:H:38:LEU:CD2	1:H:41:GLY:HA3	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:83:HIS:HA	1:E:110:ASN:ND2	2.32	0.45
1:E:267:LEU:O	1:E:293:ALA:HA	2.17	0.45
1:F:216:GLY:HA2	1:G:128:SER:CB	2.46	0.45
1:H:38:LEU:CD2	1:H:41:GLY:CA	2.95	0.45
1:B:207:ILE:HG22	1:B:224:GLN:HB3	1.99	0.45
1:A:273:GLY:HA3	1:A:298:HIS:C	2.37	0.45
1:C:48:MET:HB3	1:C:49:PRO:CD	2.46	0.45
1:C:209:TYR:HB3	1:C:232:LEU:HD13	2.00	0.44
1:E:346:THR:O	1:E:346:THR:CG2	2.65	0.44
1:H:253:ILE:HB	1:H:276:VAL:HG22	2.00	0.44
1:G:105:GLY:O	1:G:125:ASN:HB2	2.18	0.44
1:G:48:MET:O	1:G:51:LEU:N	2.45	0.44
1:A:191:THR:O	1:A:195:VAL:HG23	2.18	0.44
1:G:100:VAL:HA	1:G:122:ALA:O	2.18	0.44
1:C:186:ILE:HD12	1:C:241:LEU:HD21	2.00	0.44
1:F:358:ASN:O	1:F:358:ASN:CG	2.56	0.44
1:H:340:GLU:OE1	1:H:343:ARG:NH1	2.50	0.43
1:D:73:LEU:HD23	1:D:73:LEU:O	2.17	0.43
1:H:42:ARG:O	1:H:59:PHE:CD2	2.71	0.43
1:A:261:MET:HE1	1:A:290:ILE:HD11	2.00	0.43
1:D:129:ALA:CB	1:D:333:MET:HB3	2.48	0.43
1:G:43:ASP:OD1	1:G:44:CYS:N	2.52	0.43
1:E:209:TYR:HB3	1:E:232:LEU:HD13	2.01	0.43
1:C:170:GLU:O	1:C:173:SER:HB2	2.19	0.43
1:H:296:ASP:OD1	1:H:319:THR:OG1	2.31	0.43
1:E:38:LEU:HD23	1:E:44:CYS:SG	2.59	0.43
1:C:81:MET:CE	1:C:85:ILE:HG23	2.49	0.43
1:G:328:GLN:OE1	1:G:328:GLN:N	2.42	0.43
1:A:85:ILE:HG12	1:A:86:THR:N	2.33	0.43
1:A:254:ASN:O	1:A:258:ILE:HG13	2.18	0.43
1:D:70:GLU:O	1:D:74:ASN:HB2	2.18	0.43
1:G:85:ILE:HG12	1:G:86:THR:N	2.33	0.43
1:B:85:ILE:HG12	1:B:86:THR:N	2.33	0.43
1:A:250:HIS:HB2	4:A:513:HOH:O	2.19	0.43
1:C:358:ASN:O	1:C:358:ASN:CG	2.57	0.43
1:G:338:ALA:O	1:G:341:ILE:HB	2.18	0.43
1:C:81:MET:HE3	1:C:85:ILE:HG23	2.00	0.43
1:A:233:LEU:HD22	1:A:261:MET:HG2	2.00	0.43
1:D:65:THR:HA	1:D:68:ILE:CG1	2.45	0.42
1:H:298:HIS:CD2	1:H:303:PHE:CD1	3.07	0.42
1:B:57:VAL:HG23	1:B:57:VAL:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:ARG:NH1	4:B:517:HOH:O	2.51	0.42
1:C:85:ILE:HG12	1:C:86:THR:N	2.34	0.42
1:C:65:THR:HG21	1:C:91:ASP:CG	2.39	0.42
1:B:207:ILE:HA	1:B:224:GLN:O	2.20	0.42
1:H:48:MET:N	1:H:49:PRO:HD2	2.34	0.42
1:D:213:LEU:HD22	1:D:217:ILE:HD12	2.01	0.42
1:F:33:ARG:CZ	1:F:53:ASP:O	2.67	0.42
1:G:358:ASN:O	1:G:358:ASN:CG	2.58	0.42
1:G:184:GLY:HA3	1:G:236:SER:OG	2.20	0.42
1:B:126:ILE:HA	1:B:127:PRO:HD3	1.89	0.42
1:B:48:MET:SD	1:B:57:VAL:CG2	3.08	0.42
1:H:53:ASP:C	1:H:54:LEU:HD12	2.40	0.42
1:H:186:ILE:HB	1:H:241:LEU:HD23	2.02	0.42
1:G:124:CYS:SG	1:G:353:LEU:HD13	2.60	0.42
1:B:130:ALA:HA	1:B:133:GLU:HB2	2.01	0.42
1:B:47:GLU:N	1:B:47:GLU:OE2	2.52	0.42
1:A:210:ASP:CG	1:A:213:LEU:HG	2.40	0.42
1:E:48:MET:HB3	1:E:49:PRO:HD3	2.01	0.41
1:C:40:ASP:N	1:C:40:ASP:OD1	2.53	0.41
1:E:358:ASN:CG	1:E:358:ASN:O	2.57	0.41
1:A:41:GLY:O	1:A:59:PHE:HB2	2.20	0.41
1:B:42:ARG:HA	1:B:59:PHE:CD1	2.55	0.41
1:C:191:THR:O	1:C:195:VAL:HG23	2.21	0.41
1:F:40:ASP:OD2	1:F:82:TYR:HB2	2.20	0.41
1:D:128:SER:HA	1:D:131:VAL:HG23	2.02	0.41
1:D:38:LEU:O	1:D:38:LEU:HD23	2.20	0.41
1:A:48:MET:N	1:A:49:PRO:HD2	2.35	0.41
1:D:358:ASN:O	1:D:358:ASN:CG	2.59	0.41
1:B:358:ASN:O	1:B:358:ASN:CG	2.59	0.41
1:H:48:MET:HB3	1:H:49:PRO:CD	2.51	0.41
1:G:39:LEU:O	1:G:60:CYS:HB2	2.20	0.41
1:D:161:ARG:NH2	1:D:163:GLN:HG3	2.36	0.41
1:D:39:LEU:O	1:D:60:CYS:HB2	2.21	0.41
1:F:207:ILE:HG22	1:F:224:GLN:HB2	2.01	0.41
1:E:38:LEU:CD2	1:E:44:CYS:SG	3.09	0.41
1:C:42:ARG:O	1:C:59:PHE:CD2	2.74	0.41
1:D:35:LEU:O	1:D:77:VAL:HG12	2.21	0.41
1:C:340:GLU:HA	1:C:340:GLU:OE1	2.21	0.40
1:E:207:ILE:HA	1:E:224:GLN:O	2.20	0.40
3:H:402:KMT:O2	3:H:402:KMT:H41	2.20	0.40
1:C:153:TYR:CD2	1:D:150:THR:CG2	3.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:266:PHE:CZ	1:D:291:ARG:CD	3.04	0.40
1:F:170:GLU:CG	1:F:171:VAL:N	2.84	0.40
1:B:43:ASP:OD1	1:B:44:CYS:N	2.53	0.40
1:E:138:THR:HG21	1:E:242:HIS:CE1	2.55	0.40
1:B:83:HIS:HA	1:B:110:ASN:ND2	2.36	0.40
1:A:216:GLY:HA2	1:D:128:SER:CB	2.51	0.40
1:E:146:TYR:CE1	1:E:266:PHE:HD2	2.40	0.40
1:H:361:PHE:O	1:H:362:PHE:O	2.39	0.40
1:A:310:LEU:O	1:A:311:LYS:C	2.60	0.40
1:E:291:ARG:NH2	4:E:503:HOH:O	2.54	0.40
1:A:184:GLY:HA3	1:A:236:SER:OG	2.21	0.40

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	329/349 (94%)	318 (97%)	11 (3%)	0	100	100
1	B	329/349 (94%)	317 (96%)	12 (4%)	0	100	100
1	C	329/349 (94%)	317 (96%)	12 (4%)	0	100	100
1	D	328/349 (94%)	316 (96%)	12 (4%)	0	100	100
1	E	325/349 (93%)	313 (96%)	12 (4%)	0	100	100
1	F	327/349 (94%)	314 (96%)	13 (4%)	0	100	100
1	G	321/349 (92%)	307 (96%)	14 (4%)	0	100	100
1	H	329/349 (94%)	317 (96%)	12 (4%)	0	100	100
All	All	2617/2792 (94%)	2519 (96%)	98 (4%)	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	247/283 (87%)	246 (100%)	1 (0%)	93	98
1	B	243/283 (86%)	242 (100%)	1 (0%)	93	98
1	C	251/283 (89%)	250 (100%)	1 (0%)	93	98
1	D	226/283 (80%)	226 (100%)	0	100	100
1	E	224/283 (79%)	224 (100%)	0	100	100
1	F	238/283 (84%)	237 (100%)	1 (0%)	93	98
1	G	202/283 (71%)	202 (100%)	0	100	100
1	H	240/283 (85%)	240 (100%)	0	100	100
All	All	1871/2264 (83%)	1867 (100%)	4 (0%)	95	98

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	362	PHE
1	B	362	PHE
1	C	362	PHE
1	F	361	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	242	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

16 ligands 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	NAD	A	401	-	38,48,48	0.80	1 (2%)	47,73,73	1.77	6 (12%)
3	KMT	A	402	-	5,8,8	0.69	0	4,9,9	2.06	2 (50%)
2	NAD	B	401	-	38,48,48	0.87	2 (5%)	47,73,73	1.58	5 (10%)
3	KMT	B	402	-	5,8,8	0.97	0	4,9,9	2.37	2 (50%)
2	NAD	C	401	-	38,48,48	0.84	1 (2%)	47,73,73	1.70	5 (10%)
3	KMT	C	402	-	5,8,8	0.99	0	4,9,9	1.64	1 (25%)
2	NAD	D	401	-	38,48,48	0.79	1 (2%)	47,73,73	1.65	5 (10%)
3	KMT	D	402	-	5,8,8	1.18	0	4,9,9	1.84	1 (25%)
2	NAD	E	401	-	38,48,48	0.82	1 (2%)	47,73,73	1.64	6 (12%)
3	KMT	E	402	-	5,8,8	0.67	0	4,9,9	1.67	1 (25%)
2	NAD	F	401	-	38,48,48	0.94	2 (5%)	47,73,73	1.96	9 (19%)
3	KMT	F	402	-	5,8,8	0.62	0	4,9,9	1.67	1 (25%)
2	NAD	G	401	-	38,48,48	0.82	1 (2%)	47,73,73	1.85	7 (14%)
3	KMT	G	402	-	5,8,8	0.70	0	4,9,9	1.69	2 (50%)
2	NAD	H	401	-	38,48,48	0.73	1 (2%)	47,73,73	1.99	9 (19%)
3	KMT	H	402	-	5,8,8	0.93	0	4,9,9	1.50	1 (25%)

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	NAD	A	401	-	-	0/22/62/62	0/5/5/5
3	KMT	A	402	-	-	1/4/8/8	0/0/0/0
2	NAD	B	401	-	-	0/22/62/62	0/5/5/5
3	KMT	B	402	-	-	0/4/8/8	0/0/0/0
2	NAD	C	401	-	-	0/22/62/62	0/5/5/5
3	KMT	C	402	-	-	1/4/8/8	0/0/0/0
2	NAD	D	401	-	-	0/22/62/62	0/5/5/5
3	KMT	D	402	-	-	1/4/8/8	0/0/0/0
2	NAD	E	401	-	-	0/22/62/62	0/5/5/5
3	KMT	E	402	-	-	1/4/8/8	0/0/0/0
2	NAD	F	401	-	-	0/22/62/62	0/5/5/5
3	KMT	F	402	-	-	1/4/8/8	0/0/0/0
2	NAD	G	401	-	-	0/22/62/62	0/5/5/5
3	KMT	G	402	-	-	0/4/8/8	0/0/0/0
2	NAD	H	401	-	-	0/22/62/62	0/5/5/5
3	KMT	H	402	-	-	1/4/8/8	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAD	O4B-C1B	2.18	1.44	1.41
2	F	401	NAD	O4D-C1D	2.64	1.44	1.41
2	D	401	NAD	C5A-C4A	2.68	1.46	1.40
2	A	401	NAD	C5A-C4A	2.71	1.46	1.40
2	H	401	NAD	C5A-C4A	2.74	1.46	1.40
2	B	401	NAD	C5A-C4A	2.74	1.46	1.40
2	E	401	NAD	C5A-C4A	2.78	1.46	1.40
2	G	401	NAD	C5A-C4A	2.84	1.46	1.40
2	C	401	NAD	C5A-C4A	3.13	1.47	1.40
2	F	401	NAD	C5A-C4A	3.58	1.48	1.40

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	401	NAD	N3A-C2A-N1A	-9.08	121.94	128.89
2	D	401	NAD	N3A-C2A-N1A	-8.56	122.34	128.89
2	A	401	NAD	N3A-C2A-N1A	-8.51	122.38	128.89
2	F	401	NAD	N3A-C2A-N1A	-8.25	122.58	128.89
2	G	401	NAD	N3A-C2A-N1A	-7.99	122.77	128.89
2	C	401	NAD	N3A-C2A-N1A	-7.66	123.03	128.89
2	E	401	NAD	N3A-C2A-N1A	-7.64	123.05	128.89
2	B	401	NAD	N3A-C2A-N1A	-7.05	123.50	128.89
2	H	401	NAD	PN-O3-PA	-5.16	118.22	132.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	NAD	C4B-O4B-C1B	-4.90	104.34	109.72
2	C	401	NAD	PN-O3-PA	-4.38	120.43	132.73
2	H	401	NAD	O4D-C1D-N1N	-4.04	103.70	108.13
2	F	401	NAD	C4B-O4B-C1B	-3.80	105.54	109.72
2	A	401	NAD	C4A-C5A-N7A	-3.70	106.07	109.48
2	H	401	NAD	C4B-O4B-C1B	-3.58	105.78	109.72
2	C	401	NAD	C4A-C5A-N7A	-3.51	106.25	109.48
2	G	401	NAD	PN-O3-PA	-3.22	123.69	132.73
2	A	401	NAD	C1B-N9A-C4A	-3.18	122.14	126.94
2	F	401	NAD	PN-O3-PA	-3.16	123.86	132.73
2	G	401	NAD	C1B-N9A-C4A	-2.98	122.45	126.94
2	A	401	NAD	PN-O3-PA	-2.97	124.40	132.73
2	G	401	NAD	C4A-C5A-N7A	-2.81	106.90	109.48
2	E	401	NAD	C4A-C5A-N7A	-2.69	107.00	109.48
2	A	401	NAD	C4B-O4B-C1B	-2.66	106.80	109.72
2	H	401	NAD	C4A-C5A-N7A	-2.58	107.11	109.48
2	H	401	NAD	C1B-N9A-C4A	-2.57	123.07	126.94
2	D	401	NAD	C4A-C5A-N7A	-2.56	107.12	109.48
2	B	401	NAD	O3-PN-O5D	-2.48	96.35	102.94
2	D	401	NAD	C4B-O4B-C1B	-2.46	107.02	109.72
2	E	401	NAD	C1B-N9A-C4A	-2.44	123.27	126.94
2	E	401	NAD	O3-PN-O5D	-2.40	96.58	102.94
2	D	401	NAD	O3-PN-O5D	-2.20	97.09	102.94
2	F	401	NAD	C4A-C5A-N7A	-2.13	107.52	109.48
2	B	401	NAD	O7N-C7N-C3N	-2.12	117.28	119.59
2	F	401	NAD	C1B-N9A-C4A	-2.06	123.83	126.94
2	A	401	NAD	C2A-N1A-C6A	2.06	122.44	118.77
2	B	401	NAD	N6A-C6A-N1A	2.10	123.72	119.20
3	G	402	KMT	C3-C4-S1	2.12	118.08	112.88
2	H	401	NAD	C3N-C7N-N7N	2.12	120.14	117.82
2	F	401	NAD	C2A-N1A-C6A	2.13	122.58	118.77
2	H	401	NAD	C2A-N1A-C6A	2.14	122.58	118.77
2	E	401	NAD	O4D-C1D-N1N	2.16	110.50	108.13
2	G	401	NAD	C2A-N1A-C6A	2.18	122.66	118.77
2	B	401	NAD	C3N-C7N-N7N	2.20	120.23	117.82
2	F	401	NAD	O4B-C1B-N9A	2.22	112.75	108.10
2	H	401	NAD	O4B-C1B-N9A	2.24	112.78	108.10
3	B	402	KMT	C3-C4-S1	2.24	118.37	112.88
2	D	401	NAD	C2A-N1A-C6A	2.26	122.81	118.77
2	C	401	NAD	O4B-C1B-N9A	2.29	112.89	108.10
3	G	402	KMT	C5-S1-C4	2.32	108.30	100.37
3	H	402	KMT	C3-C4-S1	2.39	118.75	112.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	NAD	C3N-C7N-N7N	2.46	120.50	117.82
2	E	401	NAD	O2A-PA-O3	2.46	116.27	105.09
3	F	402	KMT	C5-S1-C4	2.66	109.44	100.37
3	D	402	KMT	C5-S1-C4	2.67	109.47	100.37
3	E	402	KMT	C5-S1-C4	2.67	109.47	100.37
3	A	402	KMT	C5-S1-C4	2.68	109.52	100.37
2	C	401	NAD	C3N-C7N-N7N	2.81	120.90	117.82
3	C	402	KMT	C5-S1-C4	2.83	110.02	100.37
3	A	402	KMT	C3-C4-S1	2.83	119.83	112.88
2	G	401	NAD	C3N-C7N-N7N	3.25	121.37	117.82
3	B	402	KMT	C5-S1-C4	3.70	112.98	100.37
2	F	401	NAD	O4D-C1D-N1N	6.05	114.78	108.13

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	KMT	O5-C2-C3-C4
3	E	402	KMT	O5-C2-C3-C4
3	C	402	KMT	O5-C2-C3-C4
3	H	402	KMT	O5-C2-C3-C4
3	D	402	KMT	O5-C2-C3-C4
3	F	402	KMT	O5-C2-C3-C4

There are no ring outliers.

6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	NAD	1	0
3	B	402	KMT	1	0
3	C	402	KMT	1	0
3	D	402	KMT	1	0
2	E	401	NAD	1	0
3	H	402	KMT	1	0

5.7 Other polymers ⓘ

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 ⓘ

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	330/349 (94%)	-0.14	3 (0%)	85 84	41, 49, 59, 65	1 (0%)
1	B	330/349 (94%)	-0.10	3 (0%)	85 84	39, 51, 69, 73	0
1	C	330/349 (94%)	-0.21	1 (0%)	94 93	40, 48, 55, 62	1 (0%)
1	D	329/349 (94%)	-0.08	3 (0%)	85 84	41, 53, 73, 77	0
1	E	327/349 (93%)	0.03	7 (2%)	67 62	41, 56, 76, 82	1 (0%)
1	F	329/349 (94%)	-0.13	4 (1%)	81 78	42, 52, 60, 66	1 (0%)
1	G	325/349 (93%)	0.04	12 (3%)	45 38	40, 55, 74, 77	1 (0%)
1	H	330/349 (94%)	-0.13	2 (0%)	90 89	42, 50, 58, 62	0
All	All	2630/2792 (94%)	-0.09	35 (1%)	79 77	39, 51, 71, 82	5 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	354	ARG	3.6
1	G	60	CYS	3.3
1	B	140	CYS	3.1
1	G	91	ASP	2.8
1	B	306	ALA	2.8
1	G	306	ALA	2.6
1	G	117	GLY	2.6
1	E	58	ALA	2.5
1	A	140	CYS	2.5
1	G	159	GLY	2.5
1	E	346	THR	2.4
1	C	160	THR	2.4
1	B	33	ARG	2.4
1	A	247	GLU	2.3
1	G	100	VAL	2.3
1	H	137	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	G	350	PRO	2.3
1	G	140	CYS	2.2
1	E	352	SER	2.2
1	D	306	ALA	2.2
1	G	137	SER	2.1
1	E	63	GLN	2.1
1	E	353	LEU	2.1
1	D	294	ALA	2.1
1	F	234	TYR	2.1
1	E	74	ASN	2.1
1	F	227	TYR	2.1
1	G	64	SER	2.1
1	H	140	CYS	2.1
1	G	323	ALA	2.1
1	G	57	VAL	2.1
1	A	312	ASP	2.0
1	F	158	GLU	2.0
1	D	68	ILE	2.0
1	F	140	CYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	KMT	G	402	9/9	0.76	0.47	10.17	56,58,59,62	9
3	KMT	B	402	9/9	0.81	0.40	9.76	50,51,53,54	9

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	KMT	F	402	9/9	0.87	0.34	6.56	49,49,50,50	9
3	KMT	D	402	9/9	0.87	0.35	4.80	53,54,56,58	9
3	KMT	H	402	9/9	0.83	0.42	4.18	46,47,49,49	9
3	KMT	A	402	9/9	0.88	0.31	3.71	45,46,48,48	9
3	KMT	E	402	9/9	0.88	0.35	3.19	57,58,60,61	9
3	KMT	C	402	9/9	0.87	0.25	2.42	44,44,45,45	9
2	NAD	G	401	44/44	0.94	0.20	0.36	47,51,55,55	0
2	NAD	F	401	44/44	0.95	0.19	0.17	45,48,49,49	0
2	NAD	B	401	44/44	0.96	0.17	-0.09	42,43,48,48	0
2	NAD	A	401	44/44	0.95	0.18	-0.10	40,44,44,44	0
2	NAD	E	401	44/44	0.93	0.19	-0.25	46,48,55,56	0
2	NAD	H	401	44/44	0.94	0.18	-0.35	42,45,46,46	0
2	NAD	D	401	44/44	0.96	0.16	-0.55	44,45,51,51	0
2	NAD	C	401	44/44	0.97	0.15	-0.63	40,42,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.