



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 01:31 PM UTC

PDB ID : 1WN0 / pdb_00001wn0
Title : Crystal Structure of Histidine-containing Phosphotransfer Protein, ZmHP2, from maize
Authors : Sugawara, H.; Kawano, Y.; Hatakeyama, T.; Yamaya, T.; Kamiya, N.; Sakakibara, H.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2004-07-24
Resolution : 2.20 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

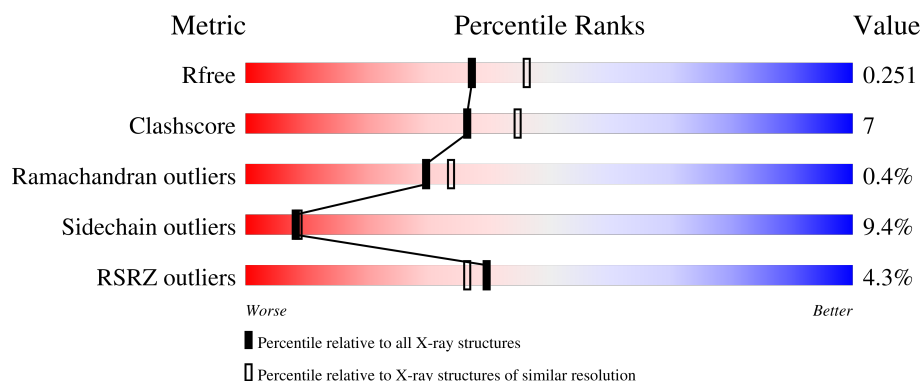
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	145	
1	B	145	
1	C	145	
1	D	145	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4278 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 histidine-containing phosphotransfer protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	0	0
			1027	646	173	199	9			
1	B	124	Total	C	N	O	S	0	0	0
			976	617	164	186	9			
1	C	138	Total	C	N	O	S	0	0	0
			1086	682	185	210	9			
1	D	138	Total	C	N	O	S	0	0	0
			1086	682	185	210	9			

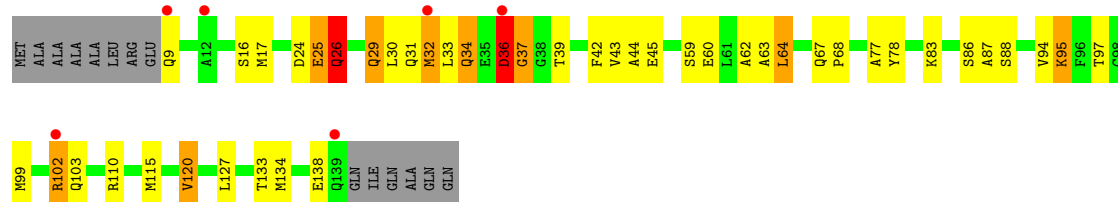
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	27	Total	O	0	0
			27	27		
2	B	12	Total	O	0	0
			12	12		
2	C	37	Total	O	0	0
			37	37		
2	D	27	Total	O	0	0
			27	27		

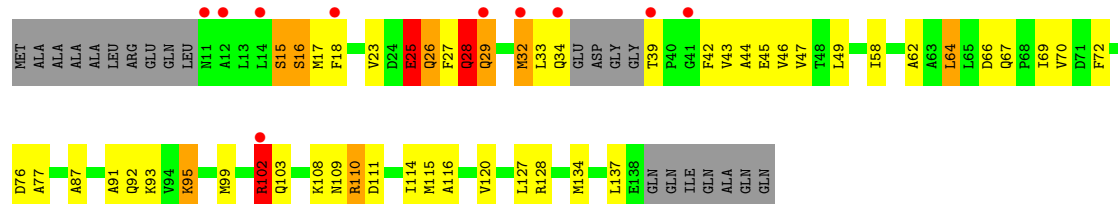
3 Residue-property plots [i](#)

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.

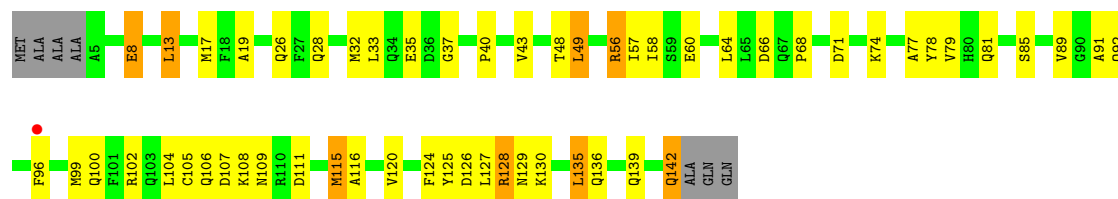
- Molecule 1: histidine-containing phosphotransfer protein



- Molecule 1: histidine-containing phosphotransfer protein

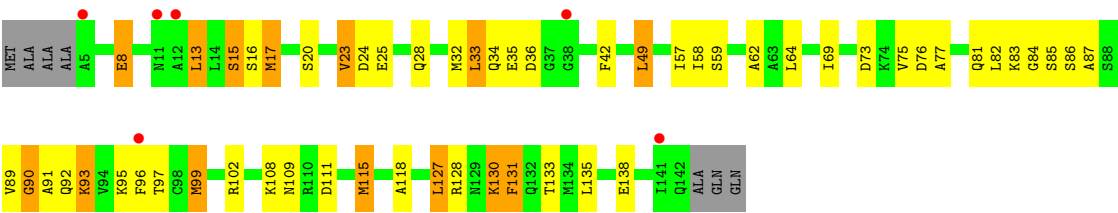


- Molecule 1: histidine-containing phosphotransfer protein



- Molecule 1: histidine-containing phosphotransfer protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	148.80Å 81.41Å 89.50Å 90.00° 123.42° 90.00°	Depositor
Resolution (Å)	19.84 – 2.20 19.84 – 2.20	Depositor EDS
% Data completeness (in resolution range)	94.2 (19.84-2.20) 94.1 (19.84-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.25 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.209 , 0.248 0.212 , 0.251	Depositor DCC
R_{free} test set	2141 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	39.7	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 23.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.010 for $1/2^*h-3/2^*k,-1/2^*h-1/2^*k,-1/2^*h+1/2^*k-l$ 0.008 for $1/2^*h+3/2^*k,1/2^*h-1/2^*k,-1/2^*h-1/2^*k-l$	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4278	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	2.11	28/1040 (2.7%)	1.65	16/1400 (1.1%)
1	B	2.13	31/988 (3.1%)	1.60	10/1329 (0.8%)
1	C	2.37	44/1099 (4.0%)	1.70	18/1479 (1.2%)
1	D	2.39	52/1099 (4.7%)	1.59	9/1479 (0.6%)
All	All	2.26	155/4226 (3.7%)	1.63	53/5687 (0.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (155) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	32	MET	SD-CE	16.84	2.21	1.79
1	A	110	ARG	CD-NE	-11.91	1.29	1.46
1	C	115	MET	SD-CE	-11.87	1.49	1.79
1	D	99	MET	SD-CE	11.17	2.07	1.79
1	C	102	ARG	NE-CZ	9.46	1.43	1.33
1	A	110	ARG	NE-CZ	-9.20	1.23	1.33
1	D	111	ASP	C-O	-8.77	1.13	1.24
1	B	77	ALA	CA-C	8.70	1.64	1.52
1	D	32	MET	SD-CE	8.59	2.01	1.79
1	D	58	ILE	N-CA	-8.51	1.36	1.46
1	D	131	PHE	C-O	-8.38	1.14	1.24
1	B	77	ALA	CA-CB	8.31	1.66	1.53
1	D	35	GLU	C-O	-8.25	1.14	1.24
1	A	95	LYS	CE-NZ	8.24	1.74	1.49
1	B	108	LYS	C-O	-8.24	1.13	1.23
1	A	99	MET	SD-CE	8.12	1.99	1.79

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	32	MET	SD-CE	8.06	1.99	1.79
1	B	110	ARG	CD-NE	-8.04	1.35	1.46
1	C	99	MET	SD-CE	8.00	1.99	1.79
1	B	69	ILE	C-O	-7.96	1.15	1.24
1	A	44	ALA	CA-CB	7.95	1.65	1.53
1	D	102	ARG	NE-CZ	7.95	1.41	1.33
1	B	95	LYS	CE-NZ	7.61	1.72	1.49
1	C	129	ASN	C-O	-7.46	1.15	1.24
1	B	43	VAL	CA-CB	-7.42	1.46	1.54
1	D	16	SER	CA-C	7.42	1.62	1.52
1	B	134	MET	SD-CE	-7.39	1.61	1.79
1	B	91	ALA	CA-CB	-7.34	1.43	1.52
1	C	19	ALA	N-CA	7.34	1.55	1.46
1	D	97	THR	CA-C	7.29	1.62	1.52
1	D	115	MET	SD-CE	-7.27	1.61	1.79
1	C	32	MET	SD-CE	7.26	1.97	1.79
1	D	82	LEU	N-CA	7.24	1.55	1.46
1	C	74	LYS	C-O	-7.20	1.15	1.24
1	C	106	GLN	C-O	-7.16	1.15	1.24
1	B	47	VAL	CA-CB	-7.14	1.44	1.54
1	D	23	VAL	C-O	7.12	1.31	1.24
1	C	60	GLU	CA-CB	7.02	1.64	1.53
1	C	13	LEU	CA-C	6.99	1.61	1.52
1	C	28	GLN	C-O	-6.86	1.16	1.24
1	C	85	SER	CA-C	6.72	1.61	1.52
1	A	68	PRO	C-O	-6.72	1.15	1.24
1	D	17	MET	SD-CE	6.70	1.96	1.79
1	A	138	GLU	CA-C	6.69	1.61	1.52
1	A	115	MET	SD-CE	-6.69	1.62	1.79
1	B	114	ILE	CA-C	-6.59	1.44	1.52
1	B	62	ALA	C-O	-6.57	1.16	1.24
1	D	76	ASP	CA-C	6.53	1.61	1.52
1	B	114	ILE	CA-CB	6.49	1.62	1.54
1	C	17	MET	C-O	-6.42	1.16	1.24
1	A	103	GLN	C-O	-6.41	1.16	1.24
1	D	130	LYS	C-O	6.39	1.31	1.24
1	B	110	ARG	NE-CZ	-6.37	1.26	1.33
1	C	128	ARG	N-CA	-6.36	1.38	1.46
1	D	8	GLU	CA-C	6.36	1.61	1.52
1	C	35	GLU	C-O	-6.35	1.15	1.24
1	D	108	LYS	CA-C	-6.30	1.45	1.53
1	A	110	ARG	CA-CB	-6.25	1.43	1.53

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	120	VAL	C-O	6.25	1.31	1.24
1	B	49	LEU	C-O	-6.25	1.16	1.24
1	C	107	ASP	N-CA	-6.24	1.37	1.46
1	C	108	LYS	C-O	-6.22	1.15	1.23
1	D	15	SER	CA-CB	6.21	1.63	1.53
1	C	77	ALA	CA-CB	6.18	1.63	1.53
1	C	126	ASP	N-CA	6.15	1.53	1.46
1	D	77	ALA	C-O	-6.14	1.16	1.24
1	D	128	ARG	C-O	-6.12	1.17	1.24
1	B	44	ALA	CA-CB	6.12	1.63	1.53
1	B	109	ASN	CA-C	6.12	1.60	1.52
1	A	67	GLN	CA-C	-6.09	1.44	1.52
1	D	84	GLY	C-O	6.08	1.31	1.23
1	D	69	ILE	CB-CG2	6.06	1.72	1.52
1	C	124	PHE	N-CA	-5.99	1.38	1.46
1	A	59	SER	CA-C	-5.96	1.45	1.52
1	D	17	MET	C-O	-5.95	1.17	1.24
1	C	74	LYS	CA-CB	-5.89	1.44	1.53
1	D	57	ILE	C-O	-5.88	1.17	1.24
1	C	13	LEU	CG-CD2	-5.86	1.33	1.52
1	A	24	ASP	CA-C	5.83	1.61	1.53
1	A	26	GLN	CD-OE1	5.81	1.34	1.23
1	A	62	ALA	C-O	-5.78	1.17	1.24
1	C	92	GLN	N-CA	-5.78	1.39	1.46
1	D	83	LYS	CD-CE	5.75	1.69	1.52
1	A	83	LYS	C-O	5.75	1.30	1.24
1	D	75	VAL	CA-CB	-5.75	1.47	1.54
1	C	79	VAL	CA-CB	-5.71	1.47	1.54
1	C	136	GLN	C-O	-5.71	1.17	1.24
1	A	32	MET	CG-SD	5.71	1.95	1.80
1	A	77	ALA	CA-CB	5.71	1.62	1.53
1	C	99	MET	CG-SD	5.66	1.95	1.80
1	D	87	ALA	CA-CB	5.65	1.62	1.53
1	B	87	ALA	N-CA	5.65	1.53	1.46
1	B	17	MET	SD-CE	5.63	1.93	1.79
1	D	62	ALA	CA-CB	5.62	1.62	1.53
1	A	99	MET	CB-CG	5.57	1.69	1.52
1	C	111	ASP	CA-CB	5.57	1.62	1.53
1	A	64	LEU	N-CA	5.56	1.53	1.46
1	D	85	SER	CA-C	5.54	1.60	1.52
1	D	28	GLN	C-O	-5.54	1.17	1.24
1	D	81	GLN	CA-C	5.54	1.59	1.52

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	115	MET	CG-SD	5.53	1.94	1.80
1	C	92	GLN	CA-C	-5.53	1.45	1.52
1	B	58	ILE	CG1-CD1	5.51	1.73	1.51
1	C	26	GLN	CD-NE2	5.51	1.44	1.33
1	C	57	ILE	CA-C	5.51	1.59	1.52
1	D	36	ASP	CA-C	5.49	1.60	1.52
1	C	43	VAL	C-O	-5.49	1.18	1.24
1	C	68	PRO	C-O	5.48	1.31	1.24
1	C	58	ILE	C-O	5.44	1.30	1.24
1	C	57	ILE	CA-CB	-5.42	1.47	1.54
1	C	78	TYR	CG-CD2	-5.39	1.28	1.39
1	D	62	ALA	C-O	-5.38	1.17	1.24
1	D	90	GLY	C-O	5.38	1.30	1.24
1	B	102	ARG	CB-CG	5.37	1.68	1.52
1	D	42	PHE	C-O	-5.37	1.17	1.24
1	D	20	SER	CA-C	5.36	1.61	1.52
1	B	28	GLN	CA-CB	5.36	1.61	1.53
1	C	66	ASP	C-O	-5.36	1.16	1.24
1	D	130	LYS	CB-CG	5.33	1.68	1.52
1	D	128	ARG	CZ-NH1	5.33	1.40	1.32
1	C	104	LEU	CA-C	5.32	1.59	1.52
1	C	58	ILE	CA-CB	-5.32	1.48	1.54
1	B	93	LYS	CD-CE	5.31	1.68	1.52
1	B	16	SER	CA-C	5.29	1.59	1.52
1	D	91	ALA	CA-CB	-5.28	1.45	1.52
1	D	102	ARG	CB-CG	5.27	1.68	1.52
1	A	87	ALA	N-CA	5.26	1.52	1.46
1	C	81	GLN	CA-CB	5.25	1.61	1.53
1	D	81	GLN	CD-OE1	5.25	1.33	1.23
1	C	40	PRO	CA-CB	5.24	1.60	1.53
1	B	103	GLN	CA-C	5.23	1.59	1.52
1	B	66	ASP	N-CA	-5.23	1.39	1.46
1	A	94	VAL	N-CA	5.21	1.52	1.46
1	A	17	MET	SD-CE	5.18	1.92	1.79
1	A	110	ARG	CZ-NH1	-5.16	1.25	1.32
1	A	77	ALA	CA-C	5.16	1.59	1.52
1	B	110	ARG	CZ-NH1	-5.16	1.25	1.32
1	C	116	ALA	C-O	-5.16	1.18	1.24
1	C	48	THR	CA-CB	5.14	1.61	1.53
1	D	24	ASP	CG-OD1	-5.14	1.15	1.25
1	D	135	LEU	CA-C	5.14	1.59	1.52
1	B	58	ILE	CA-C	5.13	1.59	1.52

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	34	GLN	CA-C	5.12	1.59	1.52
1	A	63	ALA	CA-CB	5.12	1.61	1.53
1	D	86	SER	C-O	5.12	1.30	1.24
1	D	73	ASP	CA-CB	5.11	1.61	1.53
1	D	83	LYS	C-O	5.10	1.30	1.24
1	D	93	LYS	C-O	-5.09	1.18	1.24
1	B	99	MET	CB-CG	5.08	1.67	1.52
1	B	116	ALA	N-CA	5.06	1.52	1.46
1	D	82	LEU	CA-C	5.06	1.59	1.52
1	D	59	SER	C-O	-5.03	1.18	1.24
1	D	69	ILE	CA-C	5.01	1.59	1.52
1	C	37	GLY	C-O	-5.00	1.20	1.24
1	C	125	TYR	C-O	-5.00	1.18	1.24

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	60	GLU	CB-CG-CD	9.15	128.15	112.60
1	C	135	LEU	N-CA-C	-8.94	100.70	111.69
1	B	25	GLU	N-CA-C	7.71	119.76	111.36
1	C	89	VAL	CB-CA-C	-7.37	102.06	111.20
1	D	49	LEU	CA-CB-CG	-7.31	90.73	116.30
1	C	49	LEU	CA-CB-CG	-7.16	91.24	116.30
1	C	49	LEU	N-CA-C	6.36	118.21	111.28
1	A	88	SER	N-CA-C	6.21	118.85	111.71
1	A	110	ARG	NE-CZ-NH2	-6.14	113.67	119.20
1	A	110	ARG	CG-CD-NE	-6.10	98.58	112.00
1	B	72	PHE	N-CA-C	6.09	119.97	112.54
1	A	110	ARG	CB-CG-CD	-5.96	97.58	111.30
1	D	89	VAL	CB-CA-C	-5.94	104.31	111.85
1	A	67	GLN	N-CA-C	-5.93	102.42	110.36
1	C	136	GLN	CB-CA-C	-5.88	101.03	110.79
1	B	16	SER	N-CA-C	5.85	117.33	111.07
1	D	99	MET	CG-SD-CE	5.83	113.73	100.90
1	C	71	ASP	N-CA-CB	-5.76	101.96	110.49
1	C	56	ARG	N-CA-C	-5.75	105.15	111.82
1	A	97	THR	N-CA-C	5.67	117.92	111.11
1	D	25	GLU	CB-CG-CD	5.64	122.20	112.60
1	B	42	PHE	N-CA-C	-5.64	104.77	111.03
1	B	70	VAL	CB-CA-C	-5.59	103.97	110.91
1	A	59	SER	CA-CB-OG	-5.56	99.98	111.10
1	A	36	ASP	CB-CA-C	5.52	118.85	109.75

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	15	SER	CA-C-N	5.51	127.61	120.44
1	B	15	SER	C-N-CA	5.51	127.61	120.44
1	A	34	GLN	CA-CB-CG	5.46	125.02	114.10
1	C	99	MET	CG-SD-CE	5.46	112.90	100.90
1	A	25	GLU	N-CA-C	5.43	118.71	111.75
1	D	92	GLN	N-CA-C	5.42	117.27	111.36
1	A	63	ALA	CA-C-N	-5.38	113.07	120.28
1	A	63	ALA	C-N-CA	-5.38	113.07	120.28
1	B	26	GLN	N-CA-CB	5.32	119.49	110.49
1	D	25	GLU	CA-CB-CG	5.25	124.60	114.10
1	A	86	SER	N-CA-C	-5.22	105.76	111.82
1	C	91	ALA	CA-C-N	5.21	127.22	120.44
1	C	91	ALA	C-N-CA	5.21	127.22	120.44
1	C	105	CYS	CA-C-N	5.20	127.25	120.28
1	C	105	CYS	C-N-CA	5.20	127.25	120.28
1	C	89	VAL	N-CA-C	5.20	117.89	112.90
1	A	26	GLN	N-CA-CB	5.19	117.53	110.01
1	C	136	GLN	N-CA-CB	5.18	117.73	110.12
1	A	25	GLU	CA-C-N	-5.16	113.73	120.44
1	A	25	GLU	C-N-CA	-5.16	113.73	120.44
1	B	46	VAL	CB-CA-C	-5.16	105.17	112.14
1	D	138	GLU	CA-C-N	5.12	127.40	120.38
1	D	138	GLU	C-N-CA	5.12	127.40	120.38
1	D	133	THR	N-CA-C	-5.09	105.81	111.36
1	C	129	ASN	CB-CA-C	-5.08	102.35	110.79
1	C	8	GLU	CB-CA-C	5.06	120.70	110.38
1	B	64	LEU	CB-CG-CD1	5.06	125.87	110.70
1	C	126	ASP	N-CA-C	-5.05	105.69	111.14

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	37	GLY	Peptide

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	1027	0	1013	22	0
1	B	976	0	969	23	0
1	C	1086	0	1075	20	0
1	D	1086	0	1075	20	0
2	A	27	0	0	0	0
2	B	12	0	0	0	0
2	C	37	0	0	2	0
2	D	27	0	0	0	0
All	All	4278	0	4132	60	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 7.

All (60) 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:95:LYS:CE	1:B:95:LYS:NZ	1.72	1.51
1:A:95:LYS:CE	1:A:95:LYS:NZ	1.74	1.49
1:D:99:MET:CE	1:D:99:MET:SD	2.07	1.43
1:B:32:MET:CE	1:B:32:MET:SD	2.21	1.28
1:A:26:GLN:HE21	1:D:99:MET:CE	1.71	1.01
1:A:26:GLN:NE2	1:D:99:MET:HE2	1.80	0.96
1:A:26:GLN:HE21	1:D:99:MET:HE2	1.31	0.96
1:A:45:GLU:OE1	1:D:109:ASN:ND2	2.08	0.86
1:A:26:GLN:NE2	1:D:99:MET:CE	2.41	0.78
1:D:33:LEU:C	1:D:33:LEU:HD23	2.16	0.69
1:B:26:GLN:HG2	1:C:96:PHE:CD1	2.30	0.67
1:C:142:GLN:HE21	1:C:142:GLN:HA	1.61	0.65
1:B:67:GLN:O	1:B:110:ARG:NH2	2.26	0.64
1:B:26:GLN:OE1	1:C:100:GLN:NE2	2.30	0.64
1:B:45:GLU:OE1	1:C:109:ASN:ND2	2.34	0.60
1:B:76:ASP:OD1	1:B:102:ARG:HD2	2.01	0.60
1:A:34:GLN:HB2	1:A:43:VAL:HG21	1.84	0.59
1:B:26:GLN:NE2	1:C:96:PHE:HB3	2.17	0.58
1:B:29:GLN:NE2	1:C:96:PHE:HD1	2.02	0.57
1:B:26:GLN:HG2	1:C:96:PHE:CG	2.40	0.56
1:A:102:ARG:N	1:A:102:ARG:HD3	2.19	0.56
1:B:25:GLU:O	1:B:26:GLN:C	2.49	0.56
1:A:29:GLN:HE21	1:D:96:PHE:HD1	1.53	0.55
1:B:111:ASP:O	1:B:115:MET:HG3	2.07	0.55
1:B:29:GLN:NE2	1:C:96:PHE:CD1	2.75	0.55
1:C:142:GLN:HE21	1:C:142:GLN:CA	2.19	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:LEU:C	1:C:33:LEU:HD23	2.33	0.54
1:A:29:GLN:NE2	1:D:96:PHE:HD1	2.06	0.52
1:C:128:ARG:NH2	2:C:173:HOH:O	2.35	0.51
1:A:29:GLN:NE2	1:D:93:LYS:HA	2.25	0.51
1:A:26:GLN:NE2	1:D:96:PHE:HB3	2.25	0.51
1:C:128:ARG:NE	2:C:173:HOH:O	2.31	0.49
1:A:36:ASP:O	1:A:37:GLY:C	2.54	0.48
1:B:26:GLN:CD	1:C:100:GLN:HE22	2.21	0.48
1:D:23:VAL:HA	1:D:90:GLY:O	2.14	0.47
1:A:39:THR:HG22	1:A:42:PHE:HB2	1.97	0.46
1:C:142:GLN:HA	1:C:142:GLN:NE2	2.29	0.46
1:A:26:GLN:CD	1:D:96:PHE:HB3	2.41	0.46
1:A:60:GLU:OE1	1:A:78:TYR:CE1	2.69	0.46
1:A:39:THR:HG23	1:D:115:MET:HE2	1.99	0.45
1:D:115:MET:HE3	1:D:118:ALA:HB3	1.98	0.45
1:A:39:THR:HG23	1:D:115:MET:SD	2.56	0.45
1:B:18:PHE:CE2	1:B:28:GLN:HG2	2.52	0.44
1:B:39:THR:HG22	1:C:115:MET:HE2	1.99	0.44
1:B:26:GLN:HE21	1:C:96:PHE:HB3	1.83	0.44
1:C:135:LEU:O	1:C:139:GLN:HG3	2.17	0.44
1:B:110:ARG:O	1:B:111:ASP:C	2.60	0.43
1:B:76:ASP:OD1	1:B:102:ARG:CD	2.66	0.43
1:A:39:THR:HG22	1:A:42:PHE:CB	2.49	0.43
1:A:30:LEU:HD22	1:A:43:VAL:HG22	2.01	0.42
1:D:13:LEU:O	1:D:17:MET:HG3	2.19	0.42
1:A:29:GLN:NE2	1:D:96:PHE:CD1	2.86	0.42
1:B:26:GLN:NE2	1:C:100:GLN:HE22	2.18	0.42
1:B:23:VAL:HG21	1:B:27:PHE:CD1	2.54	0.41
1:D:127:LEU:HD13	1:D:131:PHE:CE2	2.55	0.41
1:B:39:THR:HG22	1:C:115:MET:CE	2.50	0.41
1:C:115:MET:HE3	1:C:115:MET:HA	2.03	0.41
1:D:33:LEU:HD23	1:D:33:LEU:O	2.20	0.41
1:A:133:THR:O	1:A:134:MET:C	2.60	0.41
1:B:102:ARG:HA	1:B:102:ARG:HD3	1.84	0.41

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	129/145 (89%)	126 (98%)	3 (2%)	0	100	100
1	B	120/145 (83%)	116 (97%)	2 (2%)	2 (2%)	7	5
1	C	136/145 (94%)	135 (99%)	1 (1%)	0	100	100
1	D	136/145 (94%)	135 (99%)	1 (1%)	0	100	100
All	All	521/580 (90%)	512 (98%)	7 (1%)	2 (0%)	30	34

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	137	LEU
1	B	15	SER

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	113/122 (93%)	100 (88%)	13 (12%)	5	5
1	B	108/122 (88%)	96 (89%)	12 (11%)	6	6
1	C	119/122 (98%)	110 (92%)	9 (8%)	12	14
1	D	119/122 (98%)	110 (92%)	9 (8%)	12	14
All	All	459/488 (94%)	416 (91%)	43 (9%)	8	9

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	16	SER
1	A	25	GLU
1	A	26	GLN
1	A	29	GLN
1	A	31	GLN
1	A	32	MET
1	A	33	LEU
1	A	36	ASP
1	A	64	LEU
1	A	102	ARG
1	A	120	VAL
1	A	127	LEU
1	B	16	SER
1	B	25	GLU
1	B	28	GLN
1	B	29	GLN
1	B	33	LEU
1	B	34	GLN
1	B	64	LEU
1	B	92	GLN
1	B	102	ARG
1	B	120	VAL
1	B	127	LEU
1	B	128	ARG
1	C	8	GLU
1	C	13	LEU
1	C	49	LEU
1	C	56	ARG
1	C	64	LEU
1	C	120	VAL
1	C	127	LEU
1	C	130	LYS
1	C	142	GLN
1	D	8	GLU
1	D	13	LEU
1	D	15	SER
1	D	33	LEU
1	D	49	LEU
1	D	64	LEU
1	D	95	LYS
1	D	127	LEU
1	D	130	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	26	GLN
1	A	28	GLN
1	A	29	GLN
1	A	92	GLN
1	A	100	GLN
1	A	132	GLN
1	B	26	GLN
1	B	29	GLN
1	B	92	GLN
1	C	11	ASN
1	C	100	GLN
1	C	106	GLN
1	C	122	ASN
1	C	132	GLN
1	C	142	GLN
1	D	11	ASN
1	D	28	GLN
1	D	81	GLN
1	D	100	GLN
1	D	106	GLN
1	D	122	ASN
1	D	132	GLN
1	D	140	GLN
1	D	142	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	131/145 (90%)	0.15	6 (4%) 37 34	30, 41, 70, 80	0
1	B	124/145 (85%)	0.28	10 (8%) 18 15	32, 42, 73, 78	0
1	C	138/145 (95%)	0.01	1 (0%) 84 82	29, 37, 56, 77	0
1	D	138/145 (95%)	0.19	6 (4%) 40 36	31, 40, 70, 83	0
All	All	531/580 (91%)	0.15	23 (4%) 40 36	29, 40, 71, 83	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	9	GLN	5.5
1	B	32	MET	4.3
1	B	41	GLY	3.0
1	D	96	PHE	2.9
1	D	38	GLY	2.8
1	D	5	ALA	2.8
1	D	141	ILE	2.6
1	B	12	ALA	2.6
1	A	36	ASP	2.5
1	B	102	ARG	2.4
1	D	11	ASN	2.3
1	B	34	GLN	2.3
1	B	11	ASN	2.3
1	A	102	ARG	2.3
1	B	39	THR	2.2
1	B	18	PHE	2.2
1	A	139	GLN	2.2
1	D	12	ALA	2.2
1	B	14	LEU	2.2
1	A	32	MET	2.2
1	B	29	GLN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	12	ALA	2.1
1	C	96	PHE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.