



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 05:24 PM UTC

PDB ID : 2WYR / pdb_00002wyr
Title : 3-D structure of PhTET1-12s, dodecamer in the asymmetric unit
Authors : Vellieux, F.M.D.; Dura, M.A.; Franzetti, B.
Deposited on : 2009-11-20
Resolution : 2.25 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

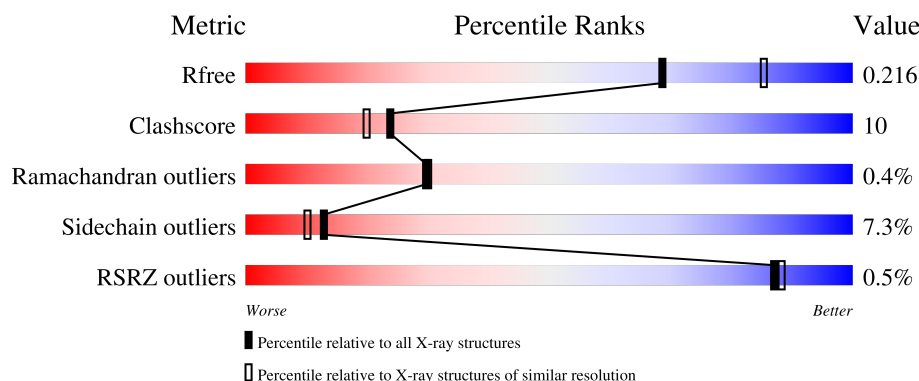
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	332	% 79% 17% ..
1	B	332	 78% 17% ..
1	C	332	% 77% 19% ..
1	D	332	 71% 21% 5% .
1	E	332	 77% 18% ..

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Mol	Chain	Length	Quality of chain
1	F	332	% 74%21% ..
1	G	332	79%18% ..
1	H	332	72%21%5% .
1	I	332	2% 67%25%5% ..
1	J	332	% 79%16% ..
1	K	332	76%19% ..
1	L	332	% 75%19% ...

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 32303 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 COBALT-ACTIVATED PEPTIDASE TET1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	5	0
			2582	1665	426	485	6			
1	B	325	Total	C	N	O	S	0	5	0
			2579	1666	426	480	7			
1	C	326	Total	C	N	O	S	0	2	0
			2568	1656	428	478	6			
1	D	325	Total	C	N	O	S	0	2	0
			2560	1653	423	478	6			
1	E	326	Total	C	N	O	S	0	3	0
			2574	1662	425	481	6			
1	F	324	Total	C	N	O	S	0	2	0
			2554	1648	424	477	5			
1	G	325	Total	C	N	O	S	0	1	0
			2554	1648	422	478	6			
1	H	326	Total	C	N	O	S	0	4	0
			2579	1664	430	479	6			
1	I	325	Total	C	N	O	S	0	0	0
			2545	1640	422	477	6			
1	J	326	Total	C	N	O	S	0	1	0
			2560	1651	425	478	6			
1	K	325	Total	C	N	O	S	0	3	0
			2565	1656	423	480	6			
1	L	325	Total	C	N	O	S	0	1	0
			2553	1645	425	477	6			

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Co	0	0
			2	2		
2	B	2	Total	Co	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	2	Total 2	Co 2	0	0
2	D	2	Total 2	Co 2	0	0
2	E	2	Total 2	Co 2	0	0
2	F	2	Total 2	Co 2	0	0
2	G	2	Total 2	Co 2	0	0
2	H	2	Total 2	Co 2	0	0
2	I	2	Total 2	Co 2	0	0
2	J	2	Total 2	Co 2	0	0
2	K	2	Total 2	Co 2	0	0
2	L	2	Total 2	Co 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	140	Total 140	O 140	0	0
3	B	122	Total 122	O 122	0	0
3	C	146	Total 146	O 146	0	0
3	D	125	Total 125	O 125	0	0
3	E	143	Total 143	O 143	0	0
3	F	110	Total 110	O 110	0	0
3	G	142	Total 142	O 142	0	0
3	H	110	Total 110	O 110	0	0
3	I	84	Total 84	O 84	0	0

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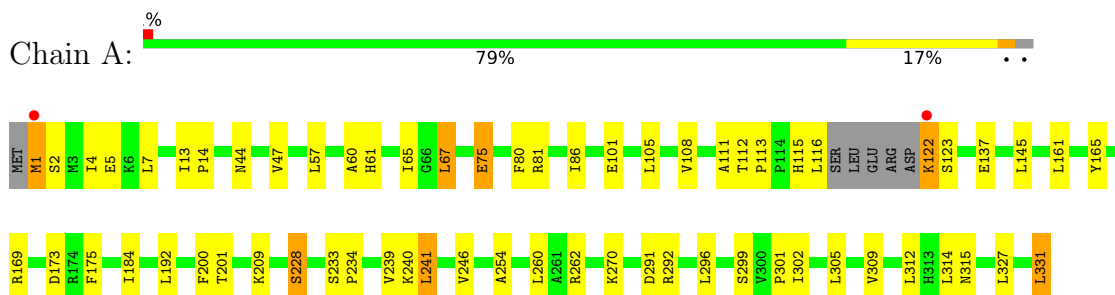
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	J	135	Total 135	O 135	0	0
3	K	122	Total 122	O 122	0	0
3	L	127	Total 127	O 127	0	0

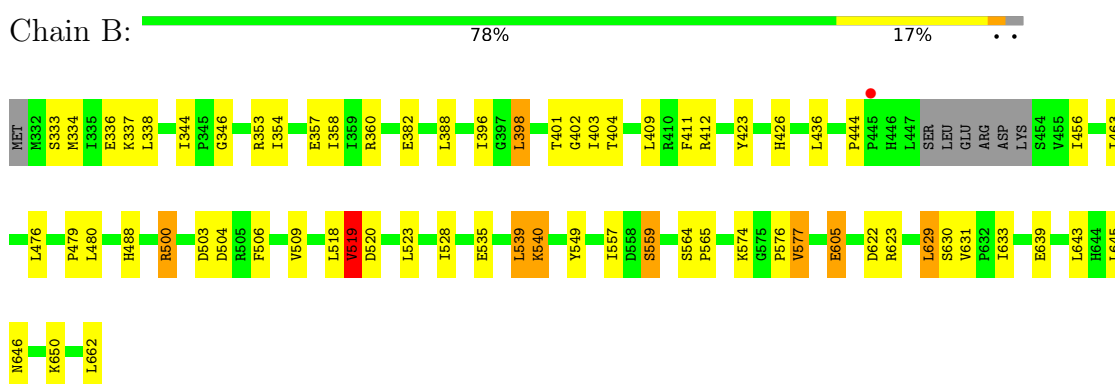
3 Residue-property plots

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.

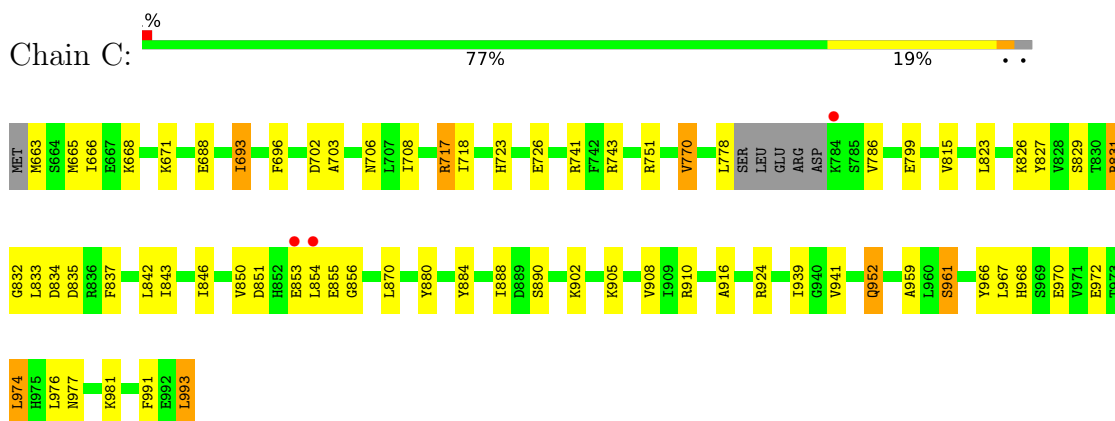
• Molecule 1: COBAL-T-ACTIVATED PEPTIDASE TET1



• Molecule 1: COBAL-T-ACTIVATED PEPTIDASE TET1

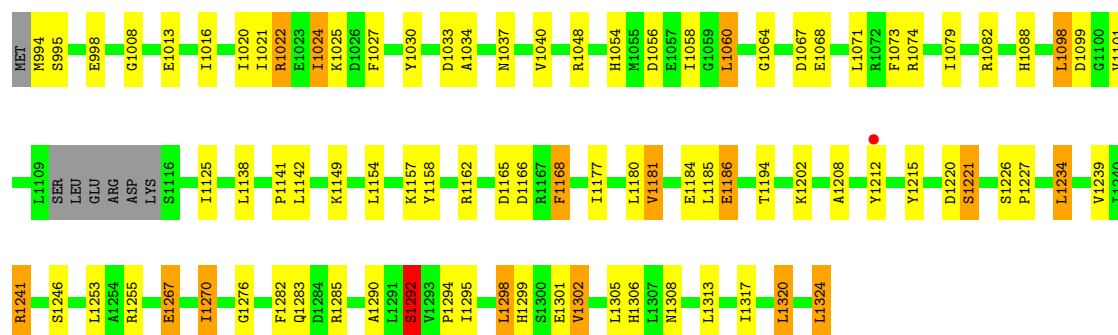


• Molecule 1: COBAL-T-ACTIVATED PEPTIDASE TET1



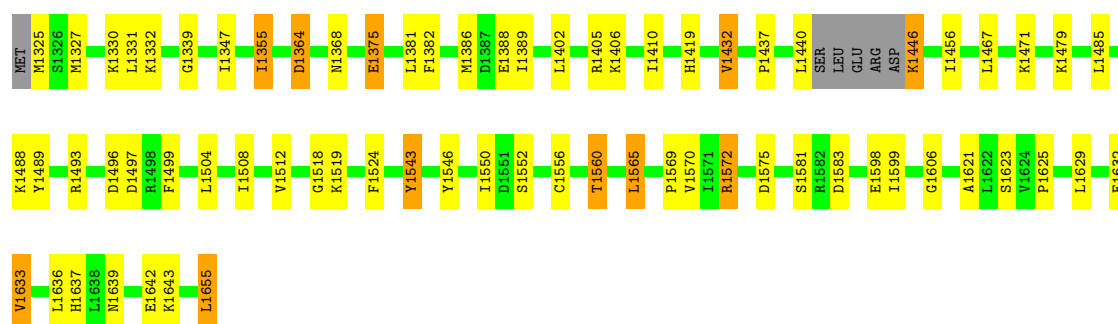
• Molecule 1: COBAL-T-ACTIVATED PEPTIDASE TET1

Chain D:  71% 21% 5% .



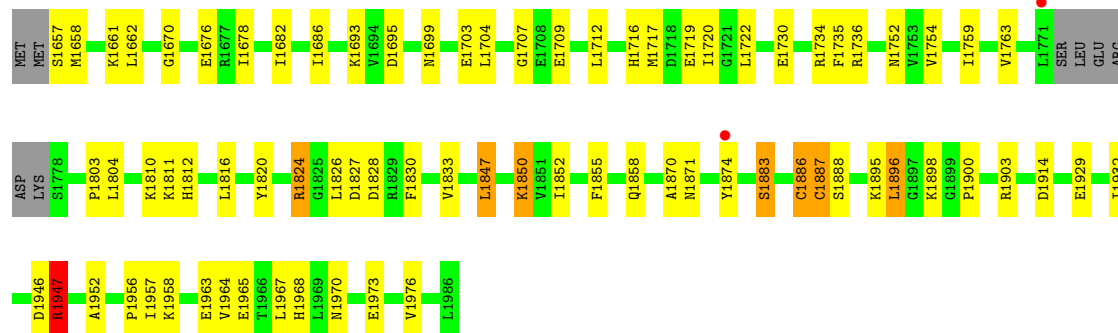
• Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1

Chain E:  77% 18% . .



• Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1

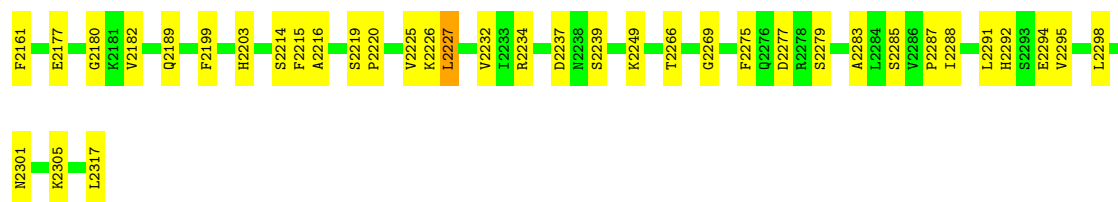
Chain F:  74% 21% . .



• Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1

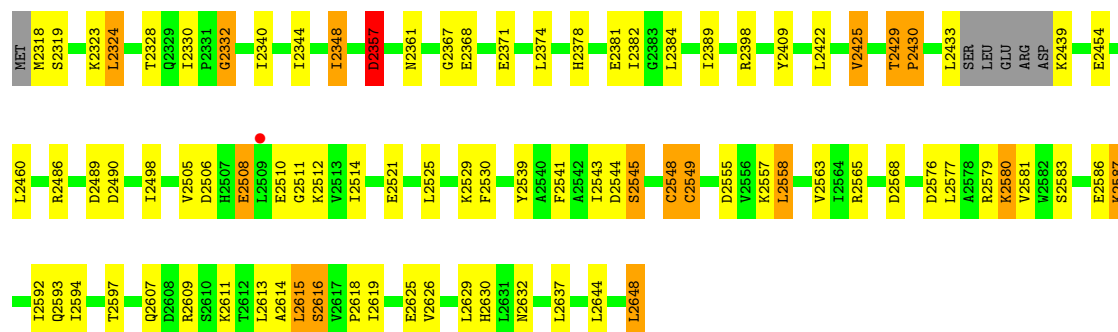
Chain G:  79% 18% . .





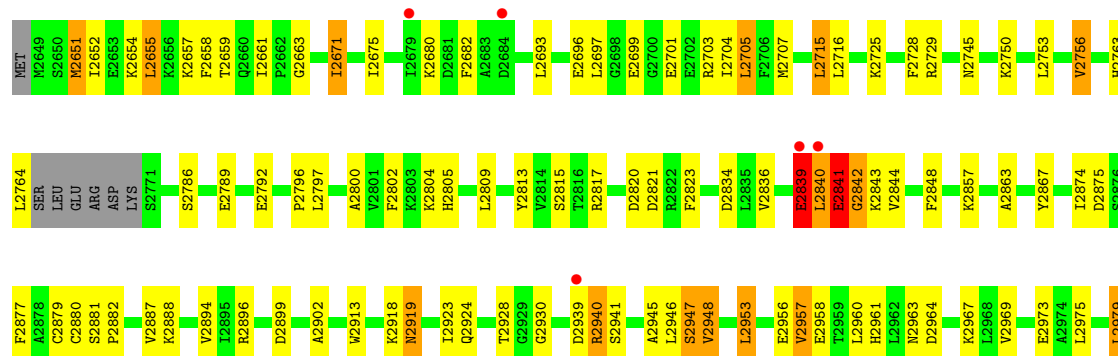
• Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1

Chain H: 72% 21% 5% .



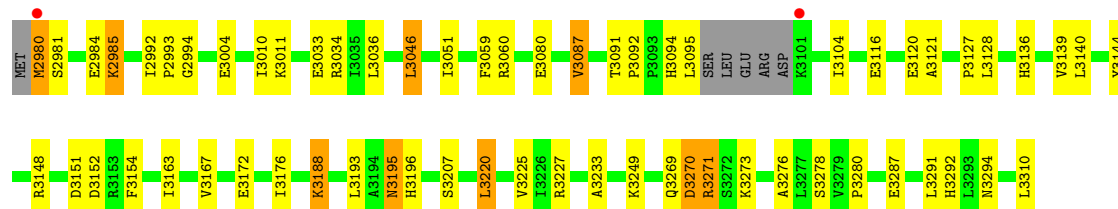
• Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1

Chain I: 67% 25% 5% . .



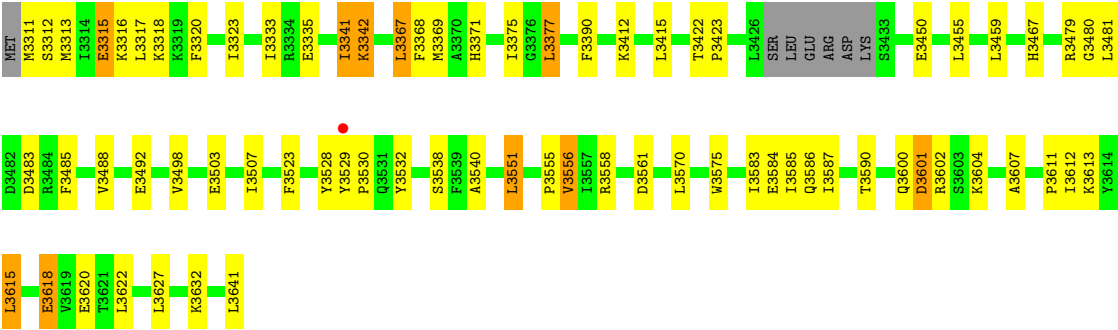
• Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1

Chain J: 79% 16% . .

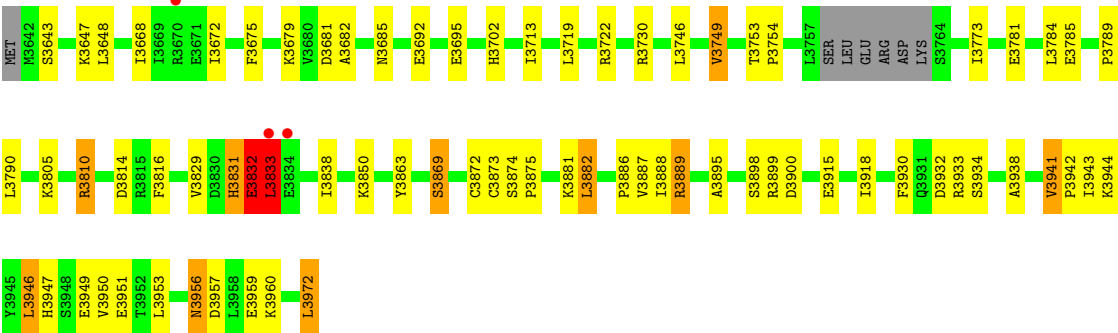


• Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1

Chain K: 76% 19% . .



● Molecule 1: COBALT-ACTIVATED PEPTIDASE TET1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	113.26Å 205.22Å 113.49Å 90.00° 100.78° 90.00°	Depositor
Resolution (Å)	49.68 – 2.25 49.68 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.68-2.25) 99.8 (49.68-2.25)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 2.25Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.170 , 0.220 0.167 , 0.216	Depositor DCC
R_{free} test set	12098 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.199	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.038 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	32303	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.48	0/2641	0.80	1/3559 (0.0%)
1	B	0.50	0/2639	0.84	0/3558
1	C	0.52	0/2618	0.84	3/3527 (0.1%)
1	D	0.50	0/2611	0.86	1/3521 (0.0%)
1	E	0.51	0/2628	0.84	1/3543 (0.0%)
1	F	0.46	0/2605	0.84	1/3514 (0.0%)
1	G	0.53	0/2602	0.88	0/3510
1	H	0.49	0/2635	0.86	3/3549 (0.1%)
1	I	0.46	0/2589	0.84	3/3491 (0.1%)
1	J	0.52	0/2607	0.88	5/3513 (0.1%)
1	K	0.47	0/2619	0.85	2/3532 (0.1%)
1	L	0.52	0/2600	0.84	1/3505 (0.0%)
All	All	0.50	0/31394	0.85	21/42322 (0.0%)

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
1	C	0	1
1	F	0	1
1	H	0	1
1	J	0	1
1	L	0	1
All	All	0	6

There are no bond length outliers.

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	2580	LYS	N-CA-C	-11.09	95.90	110.53
1	F	1887	CYS	N-CA-C	9.53	124.25	110.24
1	H	2549	CYS	N-CA-C	9.18	123.78	110.10
1	J	3195	ASN	N-CA-C	-7.87	99.93	110.55
1	K	3422	THR	CA-C-N	6.54	124.37	119.66

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	299	SER	Peptide
1	C	961	SER	Peptide
1	F	1886	CYS	Peptide
1	H	2548	CYS	Peptide
1	J	3278	SER	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	2582	0	2642	40	0
1	B	2579	0	2642	60	0
1	C	2568	0	2631	52	0
1	D	2560	0	2614	82	0
1	E	2574	0	2631	54	0
1	F	2554	0	2605	58	0
1	G	2554	0	2601	42	0
1	H	2579	0	2650	73	0
1	I	2545	0	2592	77	0
1	J	2560	0	2618	42	0
1	K	2565	0	2618	49	0
1	L	2553	0	2605	60	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0
2	K	2	0	0	0	0
2	L	2	0	0	0	0
3	A	140	0	0	4	0
3	B	122	0	0	3	0
3	C	146	0	0	5	0
3	D	125	0	0	8	0
3	E	143	0	0	2	0
3	F	110	0	0	5	0
3	G	142	0	0	4	0
3	H	110	0	0	6	0
3	I	84	0	0	4	0
3	J	135	0	0	1	0
3	K	122	0	0	2	0
3	L	127	0	0	5	0
All	All	32303	0	31449	629	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 10.

The worst 5 of 629 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:1220:ASP:O	1:D:1292:SER:HB3	1.43	1.18
1:I:2841:GLU:N	1:I:2842:GLY:HA3	1.56	1.10
1:B:559:SER:HB3	1:B:633:ILE:HD12	1.40	1.04
1:I:2875:ASP:O	1:I:2947:SER:HB3	1.57	1.03
1:L:3810:ARG:HH12	1:L:3951:GLU:CD	1.69	1.01

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	326/332 (98%)	316 (97%)	10 (3%)	0	100	100
1	B	326/332 (98%)	308 (94%)	17 (5%)	1 (0%)	36	38
1	C	324/332 (98%)	310 (96%)	14 (4%)	0	100	100
1	D	323/332 (97%)	307 (95%)	15 (5%)	1 (0%)	36	38
1	E	325/332 (98%)	314 (97%)	11 (3%)	0	100	100
1	F	322/332 (97%)	311 (97%)	10 (3%)	1 (0%)	36	38
1	G	322/332 (97%)	313 (97%)	8 (2%)	1 (0%)	36	38
1	H	326/332 (98%)	309 (95%)	14 (4%)	3 (1%)	14	11
1	I	321/332 (97%)	297 (92%)	18 (6%)	6 (2%)	6	2
1	J	323/332 (97%)	312 (97%)	10 (3%)	1 (0%)	36	38
1	K	324/332 (98%)	308 (95%)	15 (5%)	1 (0%)	36	38
1	L	322/332 (97%)	305 (95%)	16 (5%)	1 (0%)	36	38
All	All	3884/3984 (98%)	3710 (96%)	158 (4%)	16 (0%)	30	30

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	519	VAL
1	I	2841	GLU
1	K	3342	LYS
1	F	1947	ARG
1	I	2842	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	278/279 (100%)	260 (94%)	18 (6%)	15	13
1	B	277/279 (99%)	261 (94%)	16 (6%)	18	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	275/279 (99%)	259 (94%)	16 (6%)	18	17
1	D	274/279 (98%)	251 (92%)	23 (8%)	10	7
1	E	276/279 (99%)	254 (92%)	22 (8%)	11	8
1	F	273/279 (98%)	259 (95%)	14 (5%)	21	21
1	G	273/279 (98%)	255 (93%)	18 (7%)	15	13
1	H	277/279 (99%)	254 (92%)	23 (8%)	10	7
1	I	272/279 (98%)	244 (90%)	28 (10%)	7	4
1	J	274/279 (98%)	254 (93%)	20 (7%)	13	10
1	K	275/279 (99%)	253 (92%)	22 (8%)	11	8
1	L	273/279 (98%)	250 (92%)	23 (8%)	10	7
All	All	3297/3348 (98%)	3054 (93%)	243 (7%)	13	10

5 of 243 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	2086	THR
1	L	3679	LYS
1	H	2586	GLU
1	K	3641	LEU
1	L	3900	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 53 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	2538	GLN
1	I	2869	GLN
1	L	3653	GLN
1	H	2607	GLN
1	I	2805	HIS

5.3.3 RNA ⓘ

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 [i](#)

Of 24 ligands modelled in this entry, 24 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	326/332 (98%)	-0.48	2 (0%) 85 87	17, 29, 57, 83	5 (1%)
1	B	325/332 (97%)	-0.37	1 (0%) 90 91	16, 31, 60, 85	5 (1%)
1	C	326/332 (98%)	-0.53	3 (0%) 81 82	15, 28, 53, 82	2 (0%)
1	D	325/332 (97%)	-0.26	1 (0%) 90 91	20, 35, 60, 83	2 (0%)
1	E	326/332 (98%)	-0.49	0 100 100	19, 29, 51, 73	3 (0%)
1	F	324/332 (97%)	-0.29	2 (0%) 85 87	20, 35, 62, 81	2 (0%)
1	G	325/332 (97%)	-0.56	0 100 100	17, 27, 48, 70	1 (0%)
1	H	326/332 (98%)	-0.21	1 (0%) 90 91	22, 36, 62, 85	4 (1%)
1	I	325/332 (97%)	-0.09	5 (1%) 72 73	25, 41, 70, 94	0
1	J	326/332 (98%)	-0.49	2 (0%) 85 87	19, 28, 53, 74	1 (0%)
1	K	325/332 (97%)	-0.36	1 (0%) 90 91	20, 32, 57, 81	3 (0%)
1	L	325/332 (97%)	-0.48	3 (0%) 81 82	16, 29, 53, 78	2 (0%)
All	All	3904/3984 (97%)	-0.38	21 (0%) 87 88	15, 32, 59, 94	30 (0%)

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	2509	LEU	3.9
1	D	1212[A]	TYR	3.7
1	L	3833	LEU	3.4
1	A	122[A]	LYS	3.3
1	I	2840	LEU	3.1

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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CO	B	1663	1/1	0.98	0.04	30,30,30,30	0
2	CO	E	2656	1/1	0.98	0.03	30,30,30,30	0
2	CO	B	1664	1/1	0.99	0.03	25,25,25,25	0
2	CO	D	2325	1/1	0.99	0.02	35,35,35,35	0
2	CO	D	2326	1/1	0.99	0.02	28,28,28,28	0
2	CO	A	1333	1/1	0.99	0.02	24,24,24,24	0
2	CO	F	2987	1/1	0.99	0.02	35,35,35,35	0
2	CO	G	3318	1/1	0.99	0.02	27,27,27,27	0
2	CO	H	3649	1/1	0.99	0.04	39,39,39,39	0
2	CO	H	3650	1/1	0.99	0.02	29,29,29,29	0
2	CO	I	3980	1/1	0.99	0.03	38,38,38,38	0
2	CO	I	3981	1/1	0.99	0.03	33,33,33,33	0
2	CO	J	4311	1/1	0.99	0.02	28,28,28,28	0
2	CO	K	4642	1/1	0.99	0.03	35,35,35,35	0
2	CO	C	1995	1/1	1.00	0.03	22,22,22,22	0
2	CO	E	2657	1/1	1.00	0.03	27,27,27,27	0
2	CO	A	1332	1/1	1.00	0.01	28,28,28,28	0
2	CO	F	2988	1/1	1.00	0.03	29,29,29,29	0
2	CO	C	1994	1/1	1.00	0.01	30,30,30,30	0
2	CO	J	4312	1/1	1.00	0.03	23,23,23,23	0
2	CO	G	3319	1/1	1.00	0.03	24,24,24,24	0
2	CO	K	4643	1/1	1.00	0.03	26,26,26,26	0
2	CO	L	4973	1/1	1.00	0.01	29,29,29,29	0
2	CO	L	4974	1/1	1.00	0.02	23,23,23,23	0

6.5 Other polymers ⓘ

There are no such residues in this entry.