



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 2XRO / pdb_00002xro
Title : Crystal structure of TtgV in complex with its DNA operator
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Deposited on : 2010-09-17
Resolution : 3.40 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

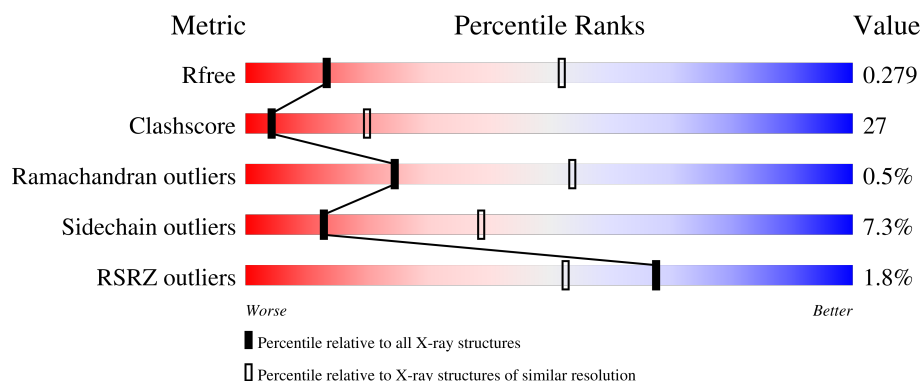
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	241	3% 53% 41% 5%
1	B	241	% 58% 36% 5%
1	E	241	3% 51% 43% 5%
1	F	241	% 57% 37% 5%
2	X	42	26% 60% 14%

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Mol	Chain	Length	Quality of chain
3	Y	43	16% 79% 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8840 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 HTH-TYPE TRANSCRIPTIONAL REGULATOR TTGV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1768	1123	309	334	2			
1	B	240	Total	C	N	O	S	0	0	0
			1780	1130	313	335	2			
1	E	239	Total	C	N	O	S	0	0	0
			1774	1126	312	334	2			
1	F	240	Total	C	N	O	S	0	0	0
			1780	1130	313	335	2			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	SER	-	expression tag	UNP Q93PU6
A	109	SER	CYS	engineered mutation	UNP Q93PU6
A	205	SER	CYS	engineered mutation	UNP Q93PU6
B	13	SER	-	expression tag	UNP Q93PU6
B	109	SER	CYS	engineered mutation	UNP Q93PU6
B	205	SER	CYS	engineered mutation	UNP Q93PU6
E	13	SER	-	expression tag	UNP Q93PU6
E	109	SER	CYS	engineered mutation	UNP Q93PU6
E	205	SER	CYS	engineered mutation	UNP Q93PU6
F	13	SER	-	expression tag	UNP Q93PU6
F	109	SER	CYS	engineered mutation	UNP Q93PU6
F	205	SER	CYS	engineered mutation	UNP Q93PU6

- Molecule 2 is a DNA chain called TTGV OPERATOR DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	X	42	Total	C	N	O	P	0	0	0
			850	408	153	248	41			

- Molecule 3 is a DNA chain called TTGV OPERATOR DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Y	43	Total	C	N	O	P	0	0	0
			886	424	161	259	42			

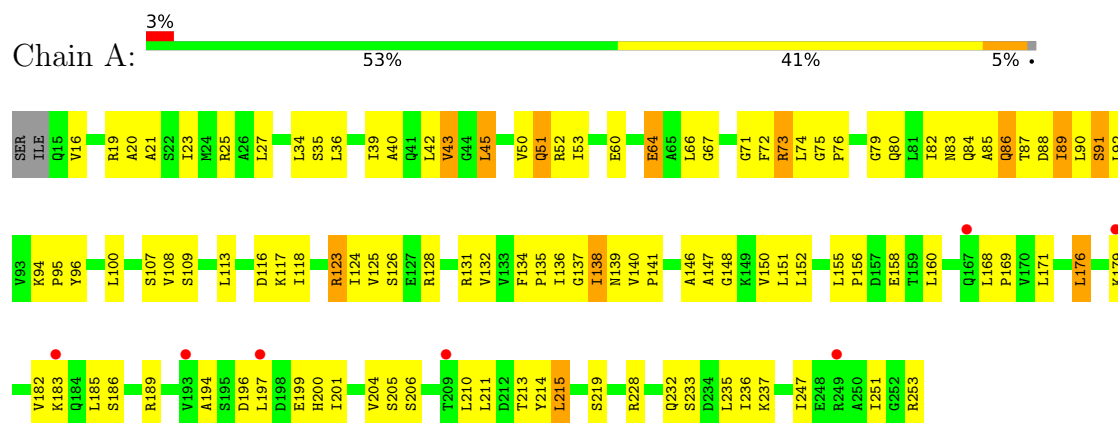
- Molecule 4 is OSMIUM ION (CCD ID: OS) (formula: Os).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Os	0	0
			1	1		
4	B	1	Total	Os	0	0
			1	1		

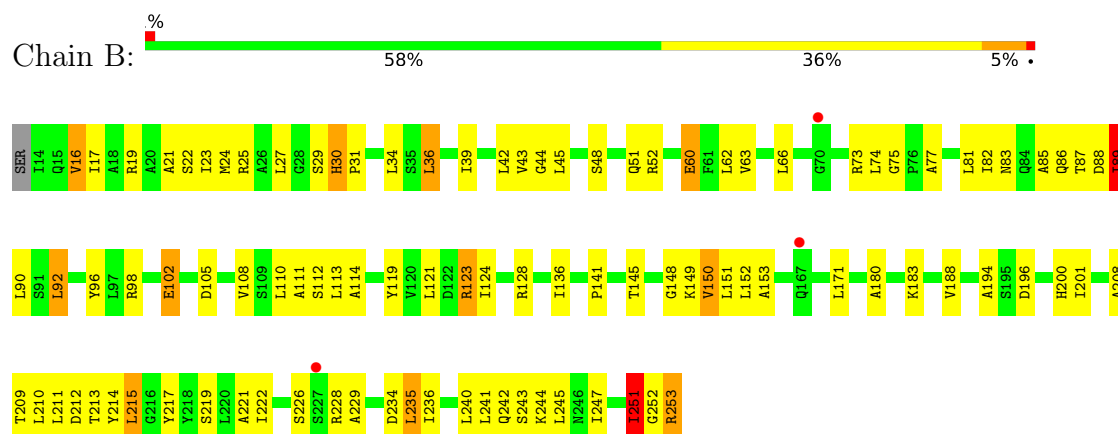
3 Residue-property plots [i](#)

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.

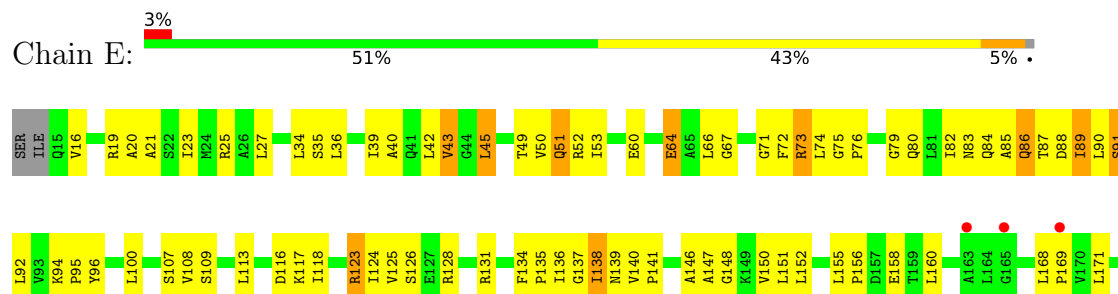
• Molecule 1: HTH-TYPE TRANSCRIPTIONAL REGULATOR TTGV

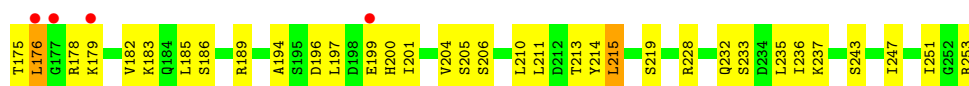


• Molecule 1: HTH-TYPE TRANSCRIPTIONAL REGULATOR TTGV



• Molecule 1: HTH-TYPE TRANSCRIPTIONAL REGULATOR TTGV





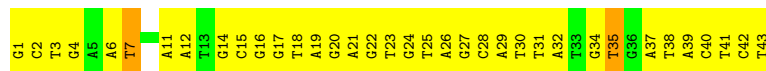
• Molecule 1: HTH-TYPE TRANSCRIPTIONAL REGULATOR TTGV



• Molecule 2: TTGV OPERATOR DNA



• Molecule 3: TTGV OPERATOR DNA



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	89.58Å 89.58Å 416.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	62.23 – 3.40 62.23 – 3.40	Depositor EDS
% Data completeness (in resolution range)	92.6 (62.23-3.40) 96.5 (62.23-3.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.71 (at 3.41Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.214 , 0.282 0.209 , 0.279	Depositor DCC
R_{free} test set	1203 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	116.6	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 345.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.447 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8840	wwPDB-VP
Average B, all atoms (Å ²)	152.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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: OS

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	0.34	0/1789	0.79	6/2429 (0.2%)
1	B	0.35	0/1801	0.87	6/2444 (0.2%)
1	E	0.34	0/1795	0.80	6/2436 (0.2%)
1	F	0.35	0/1801	0.88	6/2444 (0.2%)
2	X	0.60	0/952	1.28	8/1465 (0.5%)
3	Y	0.59	0/994	1.29	2/1535 (0.1%)
All	All	0.41	0/9132	0.96	34/12753 (0.3%)

There are no bond length outliers.

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	83	ASN	N-CA-C	7.91	119.70	108.00
1	F	16	VAL	N-CA-C	-7.63	103.84	113.22
1	B	16	VAL	N-CA-C	-7.56	103.93	113.22
2	X	12	DA	C3'-C2'-C1'	-6.86	91.31	101.60
1	F	75	GLY	CA-C-N	6.33	126.25	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1768	0	1852	108	0
1	B	1780	0	1867	84	0
1	E	1774	0	1863	110	0
1	F	1780	0	1867	85	0
2	X	850	0	475	64	0
3	Y	886	0	489	65	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	8840	0	8413	464	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 27.

The worst 5 of 464 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:F:89:ILE:HD13	1:F:89:ILE:H	1.28	0.97
1:B:89:ILE:HD13	1:B:89:ILE:H	1.31	0.94
2:X:1:DG:H5''	2:X:1:DG:H8	1.32	0.92
1:A:135:PRO:HD2	1:A:138:ILE:HD12	1.56	0.87
1:E:135:PRO:HD2	1:E:138:ILE:HD12	1.56	0.87

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	237/241 (98%)	196 (83%)	40 (17%)	1 (0%)	30 59
1	B	238/241 (99%)	206 (87%)	30 (13%)	2 (1%)	16 44
1	E	237/241 (98%)	197 (83%)	39 (16%)	1 (0%)	30 59
1	F	238/241 (99%)	207 (87%)	30 (13%)	1 (0%)	30 59

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	950/964 (98%)	806 (85%)	139 (15%)	5 (0%)	24 54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	89	ILE
1	F	89	ILE
1	B	251	ILE
1	A	138	ILE
1	E	138	ILE

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	187/193 (97%)	176 (94%)	11 (6%)	18 44
1	B	188/193 (97%)	171 (91%)	17 (9%)	9 30
1	E	188/193 (97%)	177 (94%)	11 (6%)	18 44
1	F	188/193 (97%)	172 (92%)	16 (8%)	10 33
All	All	751/772 (97%)	696 (93%)	55 (7%)	13 39

5 of 55 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	43	VAL
1	E	108	VAL
1	F	253	ARG
1	F	150	VAL
1	E	51	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	55	ASN
1	F	232	GLN
1	B	232	GLN
1	E	80	GLN
1	E	83	ASN

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 [i](#)

Of 2 ligands modelled in this entry, 2 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	239/241 (99%)	0.01	7 (2%) 53 39	63, 154, 257, 392	0
1	B	240/241 (99%)	-0.20	3 (1%) 75 61	77, 138, 233, 292	0
1	E	239/241 (99%)	0.05	7 (2%) 53 39	68, 151, 245, 403	0
1	F	240/241 (99%)	-0.23	2 (0%) 82 71	79, 136, 228, 303	0
2	X	42/42 (100%)	0.12	0 100 100	102, 162, 183, 196	0
3	Y	43/43 (100%)	0.16	0 100 100	108, 153, 190, 227	0
All	All	1043/1049 (99%)	-0.08	19 (1%) 67 53	63, 146, 241, 403	0

The worst 5 of 19 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	179	LYS	4.3
1	A	179	LYS	4.1
1	F	227	SER	3.5
1	E	177	GLY	3.5
1	B	227	SER	3.2

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($\text{\AA}^2$)	Q<0.9
4	OS	A	1001	1/1	0.99	0.10	370,370,370,370	0
4	OS	B	1001	1/1	0.99	0.11	373,373,373,373	0

6.5 Other polymers [i](#)

There are no such residues in this entry.