



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 23, 2021 – 02:55 PM EDT

PDB ID : 1DKG
Title : CRYSTAL STRUCTURE OF THE NUCLEOTIDE EXCHANGE FACTOR GRPE BOUND TO THE ATPASE DOMAIN OF THE MOLECULAR CHAPERONE DNAK
Authors : Harrison, C.J.; Kuriyan, J.
Deposited on : 1997-02-13
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

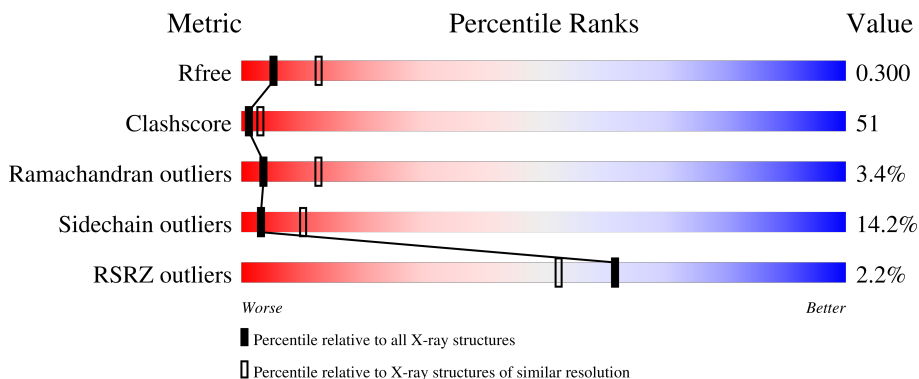
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	197	 3% 34% 38% 9% 20%
1	B	197	 2% 34% 37% 6% 23%
2	D	383	 3% 37% 50% 10% ..

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5183 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 NUCLEOTIDE EXCHANGE FACTOR GRPE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	158	1203	750	211	235	7	0	0	0
1	B	151	1144	716	204	217	7	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	122	ASP	GLY	engineered mutation	UNP P09372
B	122	ASP	GLY	engineered mutation	UNP P09372

- Molecule 2 is a protein called MOLECULAR CHAPERONE DNAK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	376	2809	1763	483	554	9	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	319	LEU	PRO	engineered mutation	UNP P04475

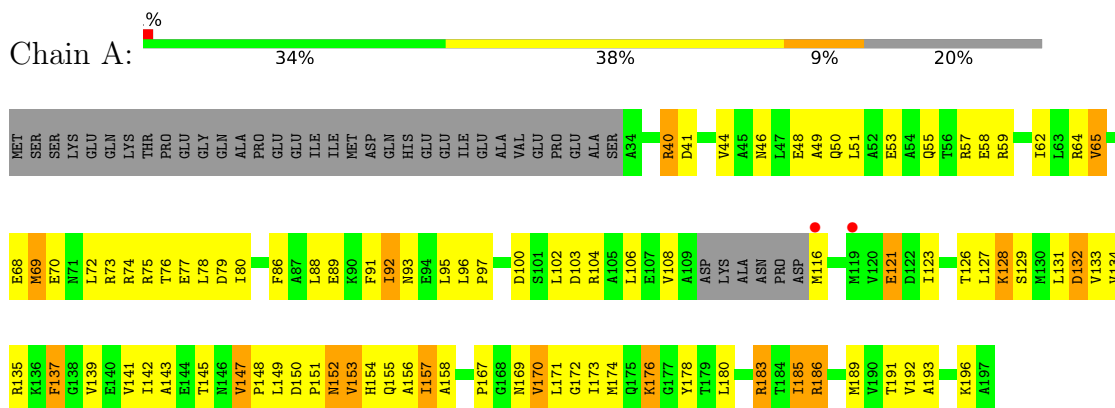
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	7	Total O 7 7	0	0
3	B	4	Total O 4 4	0	0
3	D	16	Total O 16 16	0	0

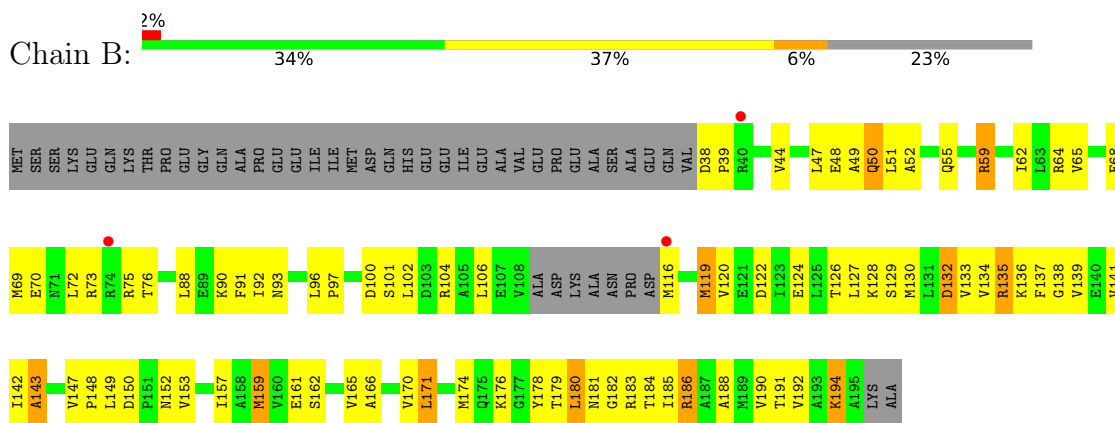
3 Residue-property plots [i](#)

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

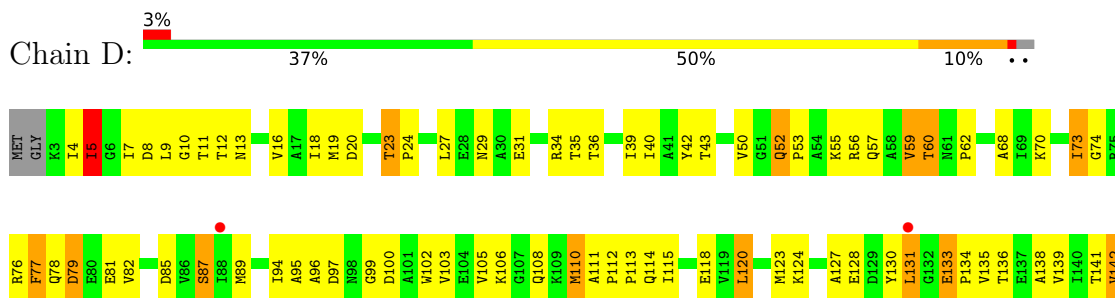
- Molecule 1: NUCLEOTIDE EXCHANGE FACTOR GRPE

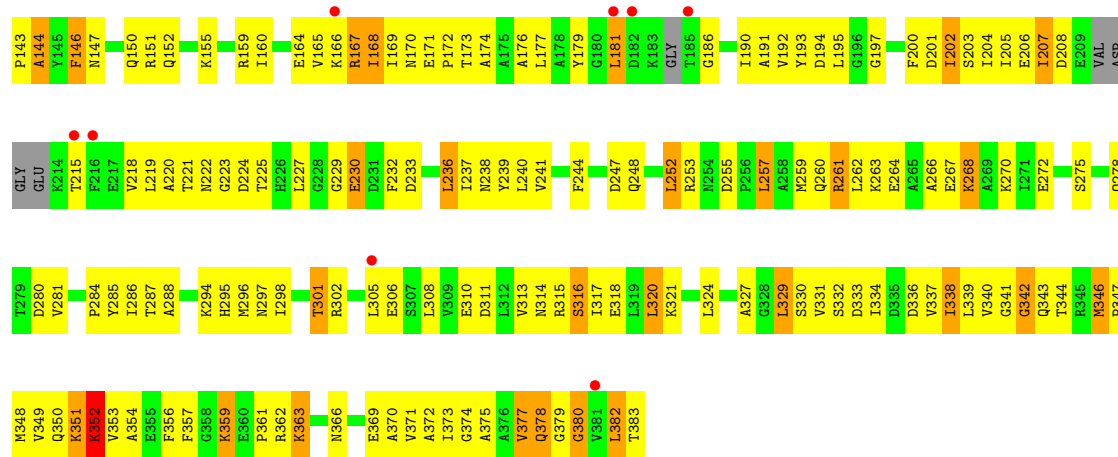


- Molecule 1: NUCLEOTIDE EXCHANGE FACTOR GRPE



- Molecule 2: MOLECULAR CHAPERONE DNAK





4 Data and refinement statistics i

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	149.43Å 149.43Å 49.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 29.89 – 2.80	Depositor EDS
% Data completeness (in resolution range)	87.9 (30.00-2.80) 93.9 (29.89-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.50 (at 2.80Å)	Xtrriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.223 , 0.317 0.246 , 0.300	Depositor DCC
R_{free} test set	1968 reflections (7.70%)	wwPDB-VP
Wilson B-factor (Å ²)	70.3	Xtrriage
Anisotropy	0.039	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.036 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5183	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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](#)

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	0.62	0/1212	0.81	0/1639
1	B	0.82	2/1153 (0.2%)	0.84	3/1557 (0.2%)
2	D	0.55	0/2843	0.75	0/3853
All	All	0.63	2/5208 (0.0%)	0.78	3/7049 (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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	132	ASP	CG-OD2	13.64	1.56	1.25
1	B	132	ASP	CG-OD1	12.77	1.54	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	ASP	CB-CG-OD2	-9.92	109.37	118.30
1	B	132	ASP	CB-CG-OD1	-8.79	110.39	118.30
1	B	132	ASP	OD1-CG-OD2	-5.72	112.44	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	132	ASP	Sidechain

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	1203	0	1213	166	0
1	B	1144	0	1158	120	0
2	D	2809	0	2807	287	0
3	A	7	0	0	0	0
3	B	4	0	0	0	0
3	D	16	0	0	2	0
All	All	5183	0	5178	525	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 51.

The worst 5 of 525 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ARG:HH11	1:A:40:ARG:HG3	1.03	1.12
1:A:142:ILE:HG13	1:A:153:VAL:CG2	1.77	1.12
2:D:202:ILE:HD12	2:D:203:SER:N	1.66	1.11
2:D:18:ILE:HD13	2:D:131:LEU:HD21	1.23	1.10
2:D:202:ILE:HD12	2:D:203:SER:H	1.12	1.10

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	154/197 (78%)	134 (87%)	17 (11%)	3 (2%)	8 26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	147/197 (75%)	127 (86%)	18 (12%)	2 (1%)	11	34
2	D	370/383 (97%)	307 (83%)	45 (12%)	18 (5%)	2	7
All	All	671/777 (86%)	568 (85%)	80 (12%)	23 (3%)	3	13

5 of 23 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	LYS
2	D	167	ARG
2	D	342	GLY
2	D	352	LYS
1	B	143	ALA

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	126/166 (76%)	109 (86%)	17 (14%)	4	11
1	B	118/166 (71%)	105 (89%)	13 (11%)	6	19
2	D	292/312 (94%)	246 (84%)	46 (16%)	2	8
All	All	536/644 (83%)	460 (86%)	76 (14%)	3	10

5 of 76 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	252	LEU
2	D	359	LYS
2	D	261	ARG
2	D	316	SER
2	D	378	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	98	ASN
2	D	150	GLN
2	D	378	GLN
2	D	297	ASN
2	D	343	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	158/197 (80%)	-0.13	2 (1%) 77 72	27, 44, 64, 71	0
1	B	151/197 (76%)	-0.32	3 (1%) 65 56	30, 51, 72, 87	0
2	D	376/383 (98%)	-0.10	10 (2%) 54 44	27, 52, 79, 98	0
All	All	685/777 (88%)	-0.15	15 (2%) 62 52	27, 51, 73, 98	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	185	THR	4.6
1	B	116	MET	3.4
2	D	381	VAL	3.1
2	D	88	ILE	2.9
2	D	216	PHE	2.6

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 monosaccharides 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.