



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:41 PM GMT

PDB ID : 1LFD  
Title : CRYSTAL STRUCTURE OF THE ACTIVE RAS PROTEIN COMPLEXED  
WITH THE RAS-INTERACTING DOMAIN OF RALGDS  
Authors : Huang, L.; Hofer, F.; Martin, G.S.; Kim, S.-H.  
Deposited on : 1998-04-29  
Resolution : 2.10 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](mailto: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.

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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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

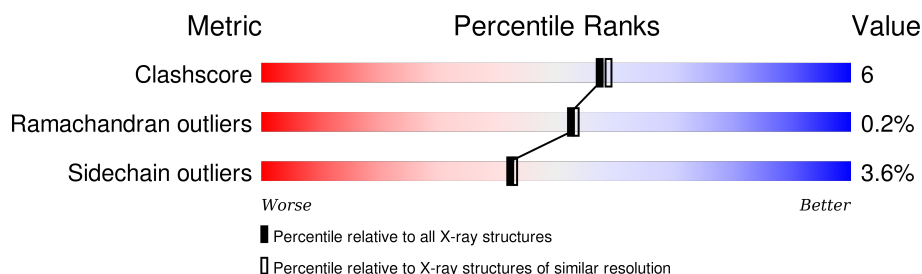
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)

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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	87	
1	C	87	
2	B	167	
2	D	167	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4414 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 RALGDS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	87	Total	C	N	O	S	0	0	0
			697	436	117	139	5			
1	C	87	Total	C	N	O	S	0	0	0
			697	436	117	139	5			

- Molecule 2 is a protein called RAS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	167	Total	C	N	O	S	0	0	0
			1331	831	231	262	7			
2	D	167	Total	C	N	O	S	0	0	0
			1331	831	231	262	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	231	LYS	GLU	MUTATION	UNP P01112
D	231	LYS	GLU	MUTATION	UNP P01112

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total 32	C 10	N 6	O 13	P 3	0	0
4	D	1	Total 32	C 10	N 6	O 13	P 3	0	0

- Molecule 5 is water.

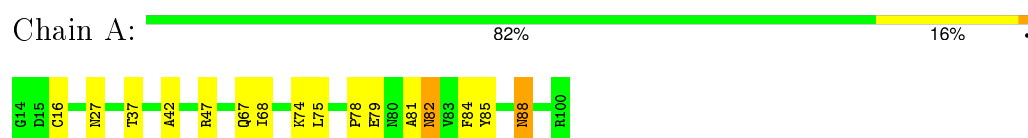
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	38	Total O 38 38	0	0
5	B	103	Total O 103 103	0	0
5	C	54	Total O 54 54	0	0
5	D	97	Total O 97 97	0	0

### 3 Residue-property plots [i](#)

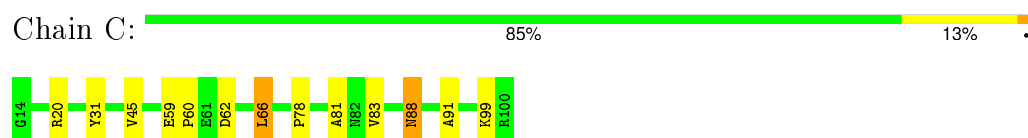
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

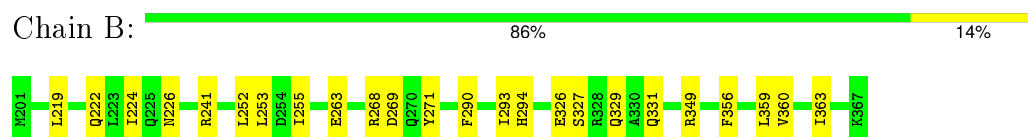
- Molecule 1: RALGDS



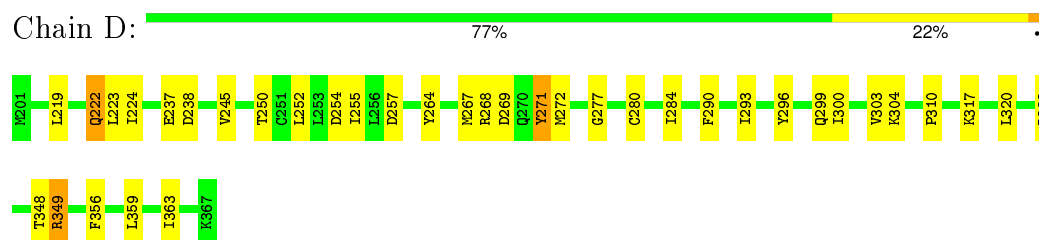
- Molecule 1: RALGDS



- Molecule 2: RAS



- Molecule 2: RAS



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.65Å 78.26Å 87.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.10	Depositor
% Data completeness (in resolution range)	98.0 (6.00-2.10)	Depositor
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
Refinement program	X-PLOR 3.85	Depositor
R, $R_{free}$	0.206 , 0.282	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4414	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

## 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: GNP, MG

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	0.30	0/707	0.54	0/953
1	C	0.30	0/707	0.57	0/953
2	B	0.29	0/1350	0.53	0/1820
2	D	0.29	0/1350	0.55	0/1820
All	All	0.29	0/4114	0.55	0/5546

There are no bond length outliers.

There are no bond angle outliers.

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	697	0	687	10	0
1	C	697	0	687	7	0
2	B	1331	0	1313	14	0
2	D	1331	0	1313	18	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
4	B	32	0	13	0	0
4	D	32	0	13	0	0
5	A	38	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	103	0	0	1	0
5	C	54	0	0	0	0
5	D	97	0	0	0	0
All	All	4414	0	4026	49	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 (49) 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)
2:D:284:ILE:HD13	2:D:323:ARG:HG2	1.62	0.80
2:D:317:LYS:HB3	2:D:320:LEU:HD23	1.74	0.68
2:D:222:GLN:NE2	2:D:349:ARG:HG2	2.13	0.64
2:D:245:VAL:HG22	2:D:250:THR:HG22	1.81	0.62
1:C:78:PRO:HB2	1:C:81:ALA:HB2	1.83	0.61
2:B:326:GLU:H	2:B:329:GLN:HE21	1.50	0.60
1:A:16:CYS:HB3	1:A:37:THR:HA	1.84	0.59
1:A:82:ASN:HD21	1:A:84:PHE:HB3	1.69	0.58
1:A:82:ASN:HD22	1:A:85:TYR:H	1.51	0.57
2:D:290:PHE:O	2:D:293:ILE:HG22	2.06	0.56
2:D:224:ILE:HD11	2:D:255:ILE:HD12	1.88	0.56
2:B:224:ILE:HD11	2:B:255:ILE:HD12	1.89	0.55
1:A:88:ASN:C	1:A:88:ASN:HD22	2.10	0.54
2:D:277:GLY:HA2	2:D:310:PRO:HB2	1.91	0.52
2:B:222:GLN:O	2:B:226:ASN:HA	2.10	0.52
1:C:45:VAL:HG21	1:C:83:VAL:HG21	1.90	0.52
2:B:349:ARG:HG2	2:B:349:ARG:O	2.11	0.51
1:C:88:ASN:HD22	1:C:91:ALA:H	1.57	0.51
2:D:238:ASP:HB2	2:D:257:ASP:HB3	1.92	0.51
2:D:280:CYS:HB3	2:D:293:ILE:HD11	1.93	0.50
2:D:272:MET:O	2:D:304:LYS:HE2	2.11	0.50
2:D:299:GLN:O	2:D:303:VAL:HG23	2.11	0.49
1:A:78:PRO:HB2	1:A:81:ALA:HB2	1.95	0.49
1:C:20:ARG:HG3	1:C:31:TYR:CD1	2.48	0.47
2:D:326:GLU:HB3	2:D:329:GLN:HG3	1.97	0.47
1:A:42:ALA:HB3	1:A:79:GLU:O	2.15	0.47
2:D:268:ARG:HA	2:D:271:TYR:CE2	2.50	0.46
2:B:326:GLU:HG2	2:B:329:GLN:HE21	1.80	0.45
1:C:62:ASP:O	1:C:99:LYS:HG3	2.16	0.44
2:B:268:ARG:HA	2:B:271:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:241:ARG:HA	2:B:253:LEU:O	2.17	0.44
2:B:326:GLU:HG3	2:B:329:GLN:H	1.83	0.44
2:B:327:SER:O	2:B:331:GLN:HG3	2.18	0.44
2:B:263:GLU:HG2	5:B:1401:HOH:O	2.16	0.43
2:D:223:LEU:HD22	2:D:356:PHE:CG	2.54	0.43
1:A:47:ARG:O	1:A:47:ARG:HD3	2.18	0.43
2:D:296:TYR:O	2:D:300:ILE:HG13	2.18	0.43
2:D:359:LEU:O	2:D:363:ILE:HG13	2.18	0.43
1:C:66:LEU:HD22	1:C:66:LEU:N	2.33	0.43
1:A:68:ILE:HG12	1:A:74:LYS:NZ	2.33	0.43
2:B:290:PHE:O	2:B:293:ILE:HG22	2.20	0.42
2:D:264:TYR:HB3	2:D:267:MET:SD	2.59	0.42
2:D:348:THR:O	2:D:349:ARG:HB2	2.19	0.42
1:A:67:GLN:HB3	1:A:75:LEU:HB3	2.03	0.41
2:B:356:PHE:O	2:B:360:VAL:HG23	2.19	0.41
2:B:359:LEU:O	2:B:363:ILE:HG13	2.21	0.41
2:B:326:GLU:HG2	2:B:329:GLN:NE2	2.36	0.41
1:C:59:GLU:HA	1:C:60:PRO:HD3	1.92	0.40
1:A:82:ASN:ND2	1:A:84:PHE:HB3	2.36	0.40

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	85/87 (98%)	80 (94%)	5 (6%)	0	100	100
1	C	85/87 (98%)	85 (100%)	0	0	100	100
2	B	165/167 (99%)	160 (97%)	5 (3%)	0	100	100
2	D	165/167 (99%)	160 (97%)	4 (2%)	1 (1%)	30	24
All	All	500/508 (98%)	485 (97%)	14 (3%)	1 (0%)	52	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	237	GLU

### 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	79/79 (100%)	76 (96%)	3 (4%)	40	40
1	C	79/79 (100%)	77 (98%)	2 (2%)	55	59
2	B	145/145 (100%)	141 (97%)	4 (3%)	51	55
2	D	145/145 (100%)	138 (95%)	7 (5%)	31	29
All	All	448/448 (100%)	432 (96%)	16 (4%)	42	43

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	82	ASN
1	A	88	ASN
2	B	219	LEU
2	B	252	LEU
2	B	269	ASP
2	B	294	HIS
1	C	66	LEU
1	C	88	ASN
2	D	219	LEU
2	D	222	GLN
2	D	252	LEU
2	D	254	ASP
2	D	269	ASP
2	D	271	TYR
2	D	349	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	82	ASN
1	A	88	ASN
2	B	299	GLN
2	B	329	GLN
2	B	350	GLN
1	C	29	ASN
1	C	73	HIS
1	C	88	ASN
2	D	222	GLN
2	D	225	GLN
2	D	261	GLN
2	D	329	GLN
2	D	365	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z  > 2$
4	GNP	B	1380	3	28,34,34	2.98	7 (25%)	33,54,54	2.49	7 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GNP	D	2380	3	28,34,34	2.86	6 (21%)	33,54,54	2.47	8 (24%)

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
4	GNP	B	1380	3	-	0/12/38/38	0/3/3/3
4	GNP	D	2380	3	-	0/12/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	2380	GNP	PG-O3G	-4.36	1.44	1.56
4	B	1380	GNP	PG-O3G	-4.32	1.44	1.56
4	B	1380	GNP	PG-O2G	-3.95	1.45	1.56
4	B	1380	GNP	C8-N7	-3.63	1.27	1.34
4	D	2380	GNP	PG-O2G	-3.58	1.46	1.56
4	D	2380	GNP	C8-N7	-3.52	1.27	1.34
4	B	1380	GNP	PB-O2B	-3.04	1.48	1.56
4	D	2380	GNP	PB-O2B	-2.96	1.48	1.56
4	B	1380	GNP	PB-O1B	2.40	1.48	1.46
4	D	2380	GNP	C6-N1	4.03	1.40	1.33
4	B	1380	GNP	C6-N1	4.11	1.40	1.33
4	D	2380	GNP	PG-O1G	11.89	1.59	1.46
4	B	1380	GNP	PG-O1G	12.44	1.60	1.46

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1380	GNP	C5-C6-N1	-8.97	111.32	123.59
4	D	2380	GNP	C5-C6-N1	-8.89	111.43	123.59
4	D	2380	GNP	O3G-PG-O1G	-4.94	100.36	113.49
4	B	1380	GNP	O3G-PG-O1G	-4.55	101.41	113.49
4	D	2380	GNP	N3-C2-N1	-3.40	122.27	127.44
4	B	1380	GNP	N3-C2-N1	-3.28	122.45	127.44
4	B	1380	GNP	PA-O3A-PB	-2.36	124.74	132.67
4	B	1380	GNP	O2G-PG-O1G	-2.25	107.52	113.49
4	D	2380	GNP	PA-O3A-PB	-2.24	125.16	132.67
4	D	2380	GNP	O3A-PA-O5'	-2.20	97.10	102.94
4	D	2380	GNP	O2G-PG-O1G	-2.11	107.89	113.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1380	GNP	O1B-PB-N3B	2.15	115.19	111.90
4	D	2380	GNP	O2B-PB-O1B	2.18	114.54	110.00
4	D	2380	GNP	C6-N1-C2	6.78	125.35	115.94
4	B	1380	GNP	C6-N1-C2	7.03	125.70	115.94

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](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.