



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 10:05 AM EDT

PDB ID : 6KUY / pdb_00006kuy
Title : Crystal structure of the alpha2A adrenergic receptor in complex with a partial agonist
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Deposited on : 2019-09-02
Resolution : 3.20 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

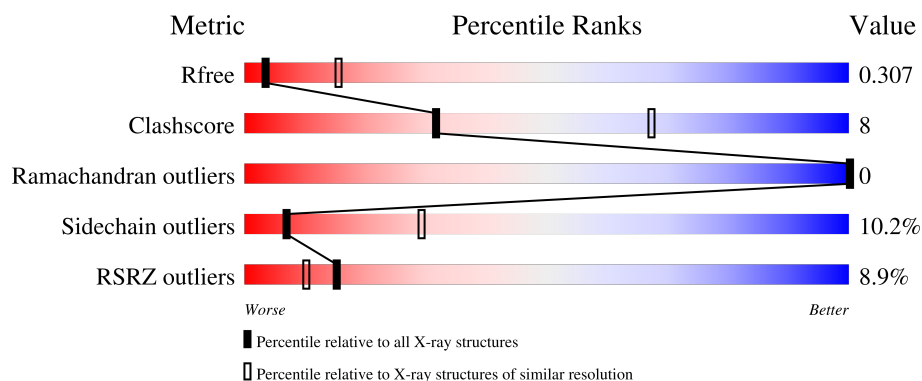
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	395	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2786 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 Alpha2A adrenergic receptor.

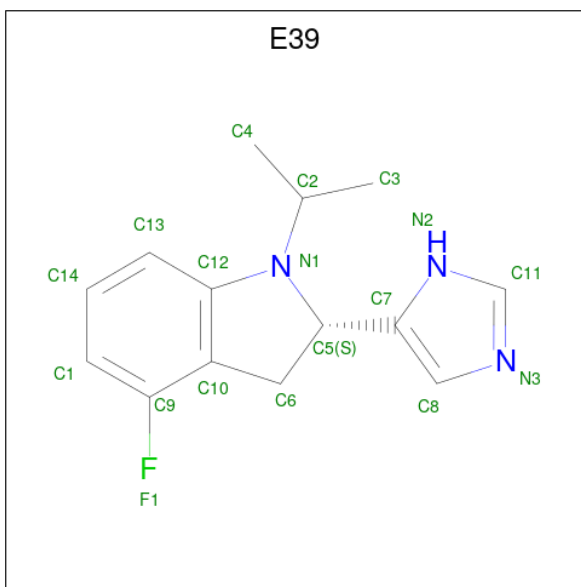
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	370	2761	1796	452	498	15	0	0	0

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	7	4	3	0	0

- Molecule 3 is (2 {S})-4-fluoranyl-2-(1 {H}-imidazol-5-yl)-1-propan-2-yl-2,3-dihydroindole (CCD ID: E39) (formula: C₁₄H₁₆FN₃).

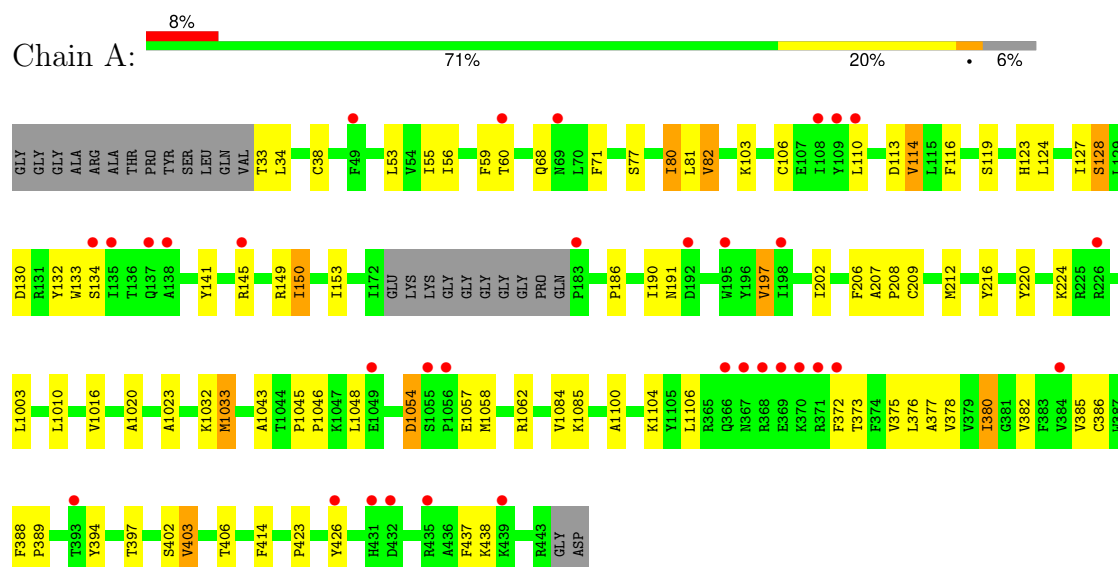


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	N		
3	A	1	18	14	1	3	0	0

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: Alpha2A adrenergic receptor



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	71.21Å 71.62Å 286.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.62 – 3.20 47.62 – 3.20	Depositor EDS
% Data completeness (in resolution range)	89.7 (47.62-3.20) 89.3 (47.62-3.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 3.12Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.267 , 0.298 0.272 , 0.307	Depositor DCC
R_{free} test set	601 reflections (3.25%)	wwPDB-VP
Wilson B-factor (Å ²)	100.6	Xtriage
Anisotropy	0.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 121.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.108 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2786	wwPDB-VP
Average B, all atoms (Å ²)	160.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.90% 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: E39, PEG

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	0.17	0/2822	0.40	0/3857

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	2761	0	2651	44	0
2	A	7	0	10	0	0
3	A	18	0	0	0	0
All	All	2786	0	2661	44	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 8.

All (44) 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:127:ILE:HD12	1:A:376:LEU:HD13	1.59	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:PHE:HD2	1:A:127:ILE:HG22	1.51	0.76
1:A:1023:ALA:HB2	1:A:1084:VAL:HG22	1.76	0.68
1:A:202:ILE:HA	1:A:206:PHE:HB2	1.78	0.63
1:A:1054:ASP:OD1	1:A:1054:ASP:N	2.31	0.62
1:A:103:LYS:HG3	1:A:186:PRO:HG3	1.82	0.61
1:A:132:TYR:HD1	1:A:133:TRP:HD1	1.50	0.60
1:A:77:SER:HA	1:A:80:ILE:HG22	1.84	0.59
1:A:1100:ALA:O	1:A:1104:LYS:NZ	2.38	0.56
1:A:1085:LYS:HD2	1:A:1085:LYS:H	1.73	0.52
1:A:134:SER:HA	1:A:141:TYR:HD2	1.74	0.52
1:A:1046:PRO:HD2	1:A:1106:LEU:HA	1.92	0.52
1:A:1048:LEU:HD21	1:A:1057:GLU:HB3	1.92	0.52
1:A:1048:LEU:HD13	1:A:1058:MET:HA	1.91	0.51
1:A:423:PRO:HA	1:A:426:TYR:HB3	1.93	0.51
1:A:1045:PRO:HG2	1:A:1048:LEU:HB2	1.93	0.51
1:A:56:ILE:HA	1:A:59:PHE:HB2	1.92	0.50
1:A:191:ASN:OD1	1:A:191:ASN:N	2.44	0.50
1:A:82:VAL:HG21	1:A:116:PHE:CG	2.47	0.50
1:A:1010:LEU:HD23	1:A:1033:MET:HB3	1.94	0.49
1:A:1003:LEU:HD13	1:A:1043:ALA:HB3	1.96	0.47
1:A:82:VAL:HG21	1:A:116:PHE:CD1	2.51	0.46
1:A:212:MET:HE3	1:A:380:ILE:HG13	1.97	0.45
1:A:53:LEU:HD21	1:A:437:PHE:HA	1.99	0.45
1:A:33:THR:OG1	1:A:34:LEU:N	2.42	0.44
1:A:124:LEU:O	1:A:128:SER:HB2	2.18	0.44
1:A:197:VAL:HG13	1:A:394:TYR:CE2	2.54	0.43
1:A:150:ILE:H	1:A:150:ILE:HG12	1.48	0.43
1:A:77:SER:O	1:A:81:LEU:HG	2.18	0.43
1:A:388:PHE:N	1:A:389:PRO:HD2	2.34	0.42
1:A:119:SER:O	1:A:123:HIS:ND1	2.53	0.42
1:A:1048:LEU:HD23	1:A:1048:LEU:HA	1.72	0.42
1:A:207:ALA:HB3	1:A:208:PRO:HD3	2.02	0.42
1:A:1016:VAL:O	1:A:1020:ALA:HB2	2.20	0.42
1:A:103:LYS:O	1:A:106:CYS:N	2.53	0.41
1:A:71:PHE:CD2	1:A:127:ILE:HG22	2.42	0.41
1:A:141:TYR:O	1:A:145:ARG:N	2.54	0.41
1:A:402:SER:OG	1:A:403:VAL:N	2.52	0.41
1:A:216:TYR:CE2	1:A:377:ALA:HB2	2.56	0.41
1:A:110:LEU:O	1:A:114:VAL:HG12	2.21	0.41
1:A:149:ARG:O	1:A:153:ILE:HG13	2.20	0.41
1:A:220:TYR:CZ	1:A:224:LYS:HE3	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ILE:O	1:A:59:PHE:HB2	2.21	0.40
1:A:386:CYS:HA	1:A:414:PHE:HD2	1.86	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	366/395 (93%)	347 (95%)	19 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	274/335 (82%)	246 (90%)	28 (10%)	7	29

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	CYS
1	A	60	THR
1	A	68	GLN
1	A	80	ILE

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Mol	Chain	Res	Type
1	A	82	VAL
1	A	113	ASP
1	A	114	VAL
1	A	128	SER
1	A	130	ASP
1	A	150	ILE
1	A	190	ILE
1	A	197	VAL
1	A	209	CYS
1	A	1032	LYS
1	A	1033	MET
1	A	1054	ASP
1	A	1062	ARG
1	A	372	PHE
1	A	373	THR
1	A	375	VAL
1	A	378	VAL
1	A	380	ILE
1	A	382	VAL
1	A	385	VAL
1	A	397	THR
1	A	403	VAL
1	A	406	THR
1	A	438	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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

2 ligands are modelled in this entry.

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
3	E39	A	1202	-	18,20,20	2.02	4 (22%)	19,29,29	1.55	3 (15%)
2	PEG	A	1201	-	6,6,6	0.50	0	5,5,5	0.36	0

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
3	E39	A	1202	-	-	4/8/20/20	0/3/3/3
2	PEG	A	1201	-	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1202	E39	C7-N2	-5.28	1.32	1.37
3	A	1202	E39	C12-N1	3.59	1.49	1.41
3	A	1202	E39	C8-N3	-2.91	1.32	1.37
3	A	1202	E39	C6-C5	-2.46	1.49	1.54

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1202	E39	C6-C5-C7	-4.28	108.47	114.22
3	A	1202	E39	C13-C12-C10	-2.12	119.12	121.73
3	A	1202	E39	C1-C9-C10	-2.02	120.33	123.99

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1202	E39	C6-C5-C7-N2
3	A	1202	E39	N1-C5-C7-N2
2	A	1201	PEG	C1-C2-O2-C3
3	A	1202	E39	C6-C5-C7-C8
2	A	1201	PEG	O2-C3-C4-O4
3	A	1202	E39	N1-C5-C7-C8

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 ⓘ

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	370/395 (93%)	0.53	33 (8%) 15 10	56, 162, 254, 326	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	371	ARG	12.6
1	A	432	ASP	11.8
1	A	369	GLU	11.4
1	A	109	TYR	7.9
1	A	431	HIS	7.2
1	A	135	ILE	6.7
1	A	134	SER	6.6
1	A	366	GLN	5.7
1	A	370	LYS	5.3
1	A	435	ARG	4.5
1	A	60	THR	4.4
1	A	1056	PRO	4.1
1	A	393	THR	3.9
1	A	198	ILE	3.9
1	A	1049	GLU	3.8
1	A	108	ILE	3.6
1	A	367	ASN	3.5
1	A	1055	SER	3.4
1	A	368	ARG	3.3
1	A	195	TRP	3.2
1	A	439	LYS	3.2
1	A	110	LEU	3.0
1	A	69	ASN	2.7
1	A	137	GLN	2.7
1	A	183	PRO	2.6
1	A	138	ALA	2.6
1	A	426	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	226	ARG	2.2
1	A	49	PHE	2.2
1	A	145	ARG	2.1
1	A	192	ASP	2.1
1	A	372	PHE	2.1
1	A	384	VAL	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](#)

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	PEG	A	1201	7/7	0.84	0.13	119,127,134,134	0
3	E39	A	1202	18/18	0.95	0.14	150,163,169,169	0

6.5 Other polymers [i](#)

There are no such residues in this entry.