



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:20 AM UTC

PDB ID : 3B6T / pdb_00003b6t
Title : Crystal Structure of the GLUR2 Ligand Binding Core (S1S2J) T686A Mutant
in Complex with Quisqualate at 2.1 Resolution
Authors : Cho, Y.; Lolis, E.; Howe, J.R.
Deposited on : 2007-10-29
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

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

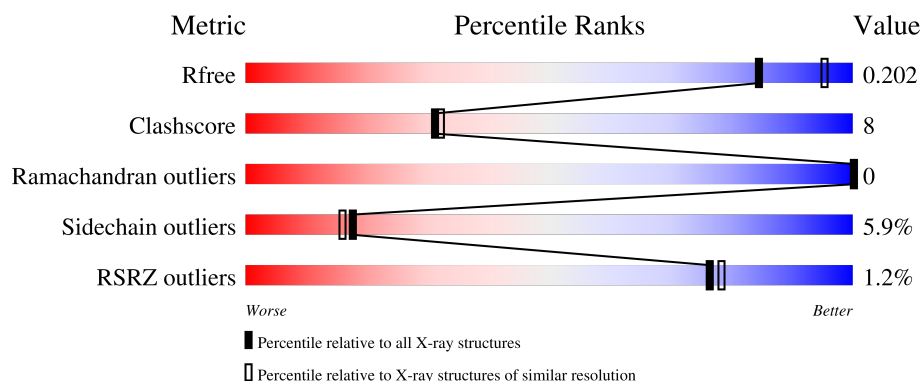
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
R_{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (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 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	263	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2136 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.

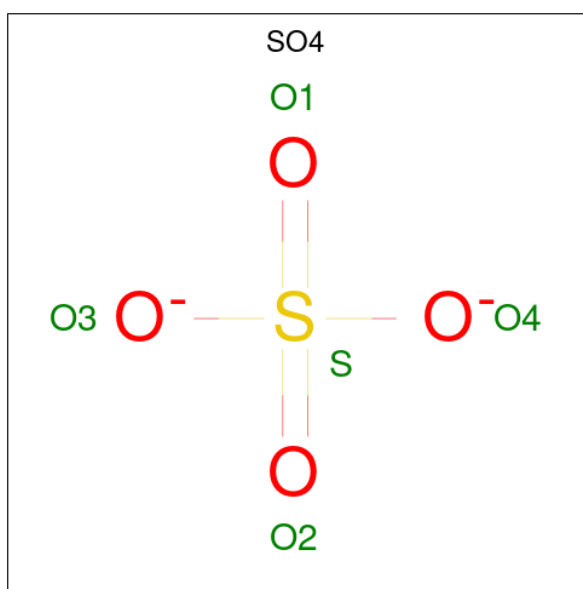
- Molecule 1 is a protein called Glutamate receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	258	1961	1247	322	378	14	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

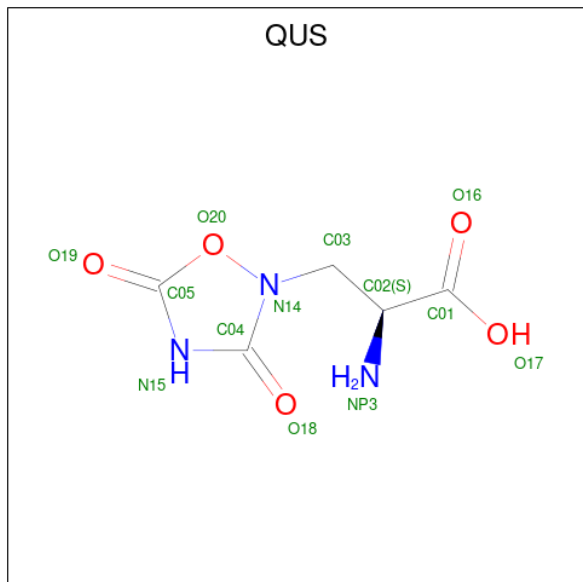
Chain	Residue	Modelled	Actual	Comment	Reference
A	390	GLY	-	expression tag	UNP P19491
A	391	ALA	-	expression tag	UNP P19491
A	630	GLY	-	linker	UNP P19491
A	631	THR	-	linker	UNP P19491
A	686	ALA	THR	engineered mutation	UNP P19491

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

- Molecule 3 is (S)-2-AMINO-3-(3,5-DIOXO-[1,2,4]OXADIAZOLIDIN-2-YL)-PROPIONIC ACID (CCD ID: QUS) (formula: C₅H₇N₃O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	5	3	5		

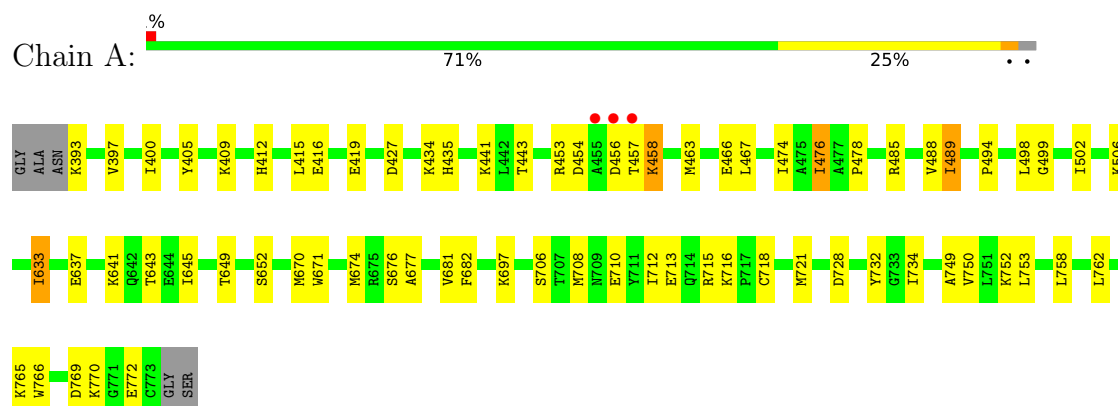
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	157	Total	O	0	0
			157	157		

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: Glutamate receptor 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	123.00Å 47.21Å 49.94Å 90.00° 110.83° 90.00°	Depositor
Resolution (Å)	57.45 – 2.10 57.45 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.0 (57.45-2.10) 98.0 (57.45-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.61 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.3.0038	Depositor
R, R_{free}	0.206 , 0.274 0.209 , 0.202	Depositor DCC
R_{free} test set	770 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	17.1	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2136	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.36% 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: SO4, QUS

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	1.40	8/1996 (0.4%)	1.38	10/2697 (0.4%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	697	LYS	CA-CB	8.46	1.66	1.53
1	A	489	ILE	CA-CB	6.58	1.64	1.54
1	A	645	ILE	CA-CB	6.36	1.61	1.54
1	A	633	ILE	CA-CB	5.44	1.61	1.54
1	A	677	ALA	N-CA	5.28	1.52	1.46
1	A	750	VAL	CA-CB	5.24	1.60	1.54
1	A	502	ILE	CA-CB	5.10	1.60	1.54
1	A	434	LYS	CA-CB	5.06	1.61	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	427	ASP	CA-CB-CG	8.21	120.81	112.60
1	A	706	SER	CA-C-N	6.41	128.77	120.44
1	A	706	SER	C-N-CA	6.41	128.77	120.44
1	A	671	TRP	N-CA-C	5.67	117.91	111.11
1	A	710	GLU	N-CA-C	5.64	117.42	111.28
1	A	681	VAL	CA-C-O	5.31	124.30	119.20
1	A	466	GLU	N-CA-C	5.22	116.97	111.28
1	A	485	ARG	N-CA-C	5.13	116.56	111.07
1	A	670	MET	N-CA-C	5.10	116.53	111.07
1	A	674	MET	N-CA-C	5.06	116.80	111.28

There are no chirality outliers.

There are no planarity outliers.

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	1961	0	1927	32	0
2	A	5	0	0	0	0
3	A	13	0	6	0	0
4	A	157	0	0	4	0
All	All	2136	0	1933	32	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 (32) 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:454:ASP:HB3	1:A:457:THR:HG22	1.72	0.70
1:A:637:GLU:O	1:A:641:LYS:HG3	1.97	0.65
1:A:716:LYS:HE3	1:A:772:GLU:CD	2.22	0.64
1:A:409:LYS:HE3	1:A:419:GLU:O	2.00	0.61
1:A:506:LYS:HE2	1:A:718:CYS:O	2.02	0.58
1:A:454:ASP:CB	1:A:457:THR:HG22	2.36	0.55
1:A:474:ILE:HD11	1:A:734:ILE:HG21	1.88	0.54
1:A:474:ILE:HD11	1:A:734:ILE:CG2	2.38	0.54
1:A:715:ARG:NH1	1:A:770:LYS:O	2.38	0.52
1:A:494:PRO:HA	1:A:732:TYR:O	2.10	0.50
1:A:765:LYS:HE2	1:A:766:TRP:CZ2	2.46	0.50
1:A:393:LYS:N	4:A:87:HOH:O	2.45	0.49
1:A:435:HIS:CE1	1:A:752:LYS:HD2	2.48	0.49
1:A:749:ALA:O	1:A:753:LEU:HG	2.14	0.48
1:A:716:LYS:H	1:A:772:GLU:HG3	1.78	0.47
1:A:758:LEU:O	1:A:758:LEU:HD22	2.15	0.47
1:A:405:TYR:CD2	1:A:478:PRO:HG3	2.49	0.46
1:A:441:LYS:HD2	4:A:102:HOH:O	2.15	0.46
1:A:721:MET:HE3	1:A:721:MET:HB2	1.65	0.46
1:A:488:VAL:HG23	1:A:489:ILE:HG23	1.98	0.45
1:A:409:LYS:O	1:A:412:HIS:HD2	2.00	0.43
1:A:409:LYS:O	1:A:412:HIS:CD2	2.70	0.43
1:A:405:TYR:CE2	1:A:708:MET:HE3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:397:VAL:HG13	1:A:476:ILE:HG13	2.00	0.42
1:A:400:ILE:HB	1:A:463:MET:HE2	2.02	0.42
1:A:499:GLY:HA3	1:A:728:ASP:O	2.20	0.42
1:A:649:THR:O	1:A:682:PHE:HA	2.21	0.41
1:A:676:SER:HB2	4:A:155:HOH:O	2.21	0.41
1:A:412:HIS:HA	1:A:415:LEU:HD12	2.04	0.40
1:A:453:ARG:CZ	1:A:458:LYS:HD2	2.50	0.40
1:A:443:THR:CG2	4:A:71:HOH:O	2.69	0.40
1:A:713:GLU:HG3	1:A:721:MET:HA	2.03	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	256/263 (97%)	247 (96%)	9 (4%)	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	204/218 (94%)	192 (94%)	12 (6%)	18	16

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	416	GLU
1	A	456	ASP
1	A	458	LYS
1	A	467	LEU
1	A	476	ILE
1	A	498	LEU
1	A	633	ILE
1	A	643	THR
1	A	652	SER
1	A	712	ILE
1	A	762	LEU
1	A	769	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	412	HIS
1	A	435	HIS
1	A	709	ASN
1	A	714	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
2	SO4	A	901	-	4,4,4	0.56	0	6,6,6	0.75	0
3	QUS	A	801	-	9,13,13	4.95	5 (55%)	10,18,18	5.13	7 (70%)

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	QUS	A	801	-	-	2/6/8/8	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	QUS	O19-C05	13.33	1.46	1.21
3	A	801	QUS	C04-N14	-4.53	1.31	1.39
3	A	801	QUS	O17-C01	-2.75	1.21	1.30
3	A	801	QUS	C03-C02	2.58	1.56	1.53
3	A	801	QUS	O18-C04	2.46	1.27	1.23

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	QUS	N15-C04-N14	9.75	115.44	106.10
3	A	801	QUS	O20-C05-N15	8.25	113.57	106.14
3	A	801	QUS	C04-N15-C05	-6.95	100.85	112.62
3	A	801	QUS	O18-C04-N14	-4.61	120.33	126.37
3	A	801	QUS	C03-N14-C04	3.34	137.14	129.17
3	A	801	QUS	O19-C05-N15	-2.87	124.99	129.68
3	A	801	QUS	O20-C05-O19	-2.17	121.07	122.95

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	801	QUS	C01-C02-C03-N14
3	A	801	QUS	NP3-C02-C03-N14

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	258/263 (98%)	-0.15	3 (1%) 76 78	5, 14, 26, 43	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	457	THR	2.9
1	A	456	ASP	2.1
1	A	455	ALA	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(Å ²)	Q<0.9
2	SO4	A	901	5/5	0.95	0.08	26,27,29,29	0
3	QUS	A	801	13/13	0.95	0.06	5,10,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.