



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:04 AM UTC

PDB ID : 3BFT / pdb\_00003bft  
Title : Structure of the ligand-binding core of GluR2 in complex with the agonist (S)-TDPA at 2.25 Å resolution  
Authors : Beich-Frandsen, M.; Mirza, O.; Vestergaard, B.; Gajhede, M.; Kastrop, J.S.  
Deposited on : 2007-11-23  
Resolution : 2.27 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

<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>.

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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

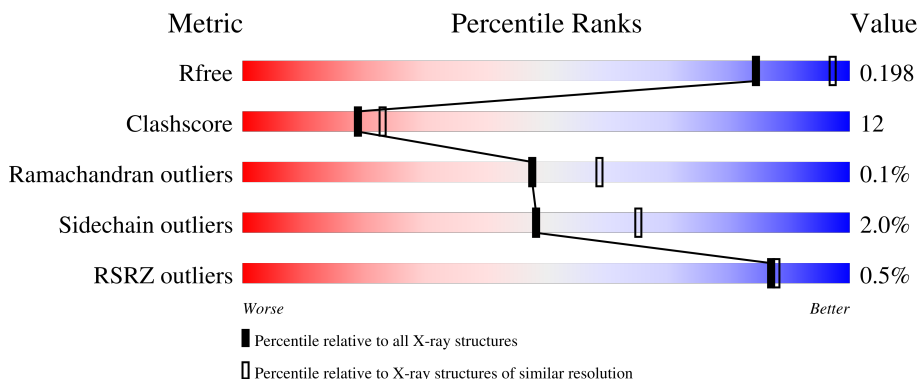
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.27 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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  $\geq 3$ , 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	 73% 24% ..
1	B	263	 71% 27% ..
1	C	263	 79% 18% ..

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6695 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.

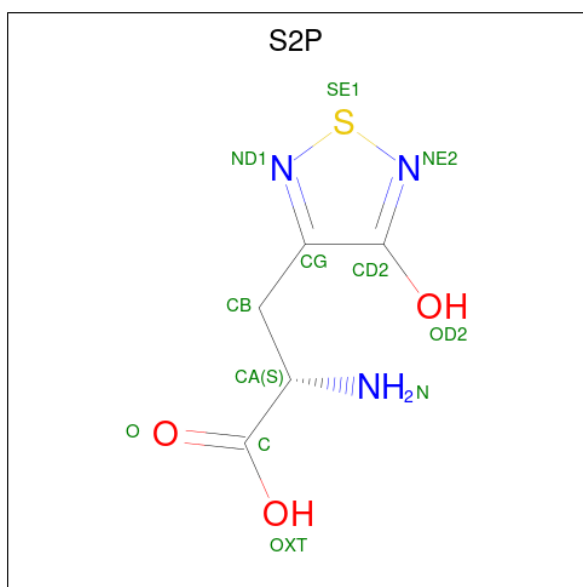
- Molecule 1 is a protein called Glutamate receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	259	2026	1290	338	384	14	0	0	0
1	B	259	2026	1290	338	384	14	0	0	0
1	C	259	2026	1290	338	384	14	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP P19491
A	2	ALA	-	expression tag	UNP P19491
B	1	GLY	-	expression tag	UNP P19491
B	2	ALA	-	expression tag	UNP P19491
C	1	GLY	-	expression tag	UNP P19491
C	2	ALA	-	expression tag	UNP P19491
A	118	GLY	-	linker	UNP P19491
A	119	THR	-	linker	UNP P19491
B	118	GLY	-	linker	UNP P19491
B	119	THR	-	linker	UNP P19491
C	118	GLY	-	linker	UNP P19491
C	119	THR	-	linker	UNP P19491

- Molecule 2 is (2S)-2-amino-3-(4-hydroxy-1,2,5-thiadiazol-3-yl)propanoic acid (CCD ID: S2P) (formula: C<sub>5</sub>H<sub>7</sub>N<sub>3</sub>O<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	12	5	3	3	1	0	0
2	B	1	12	5	3	3	1	0	0
2	C	1	12	5	3	3	1	0	0

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
3	A	2	2	2	0	0
3	B	3	3	3	0	0

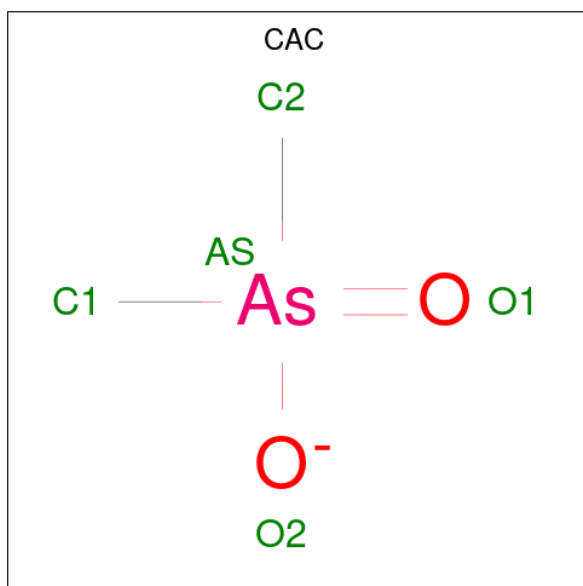
- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

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

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0

- Molecule 6 is CACODYLATE ION (CCD ID: CAC) (formula:  $C_2H_6AsO_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total As C O 5 1 2 2	0	0

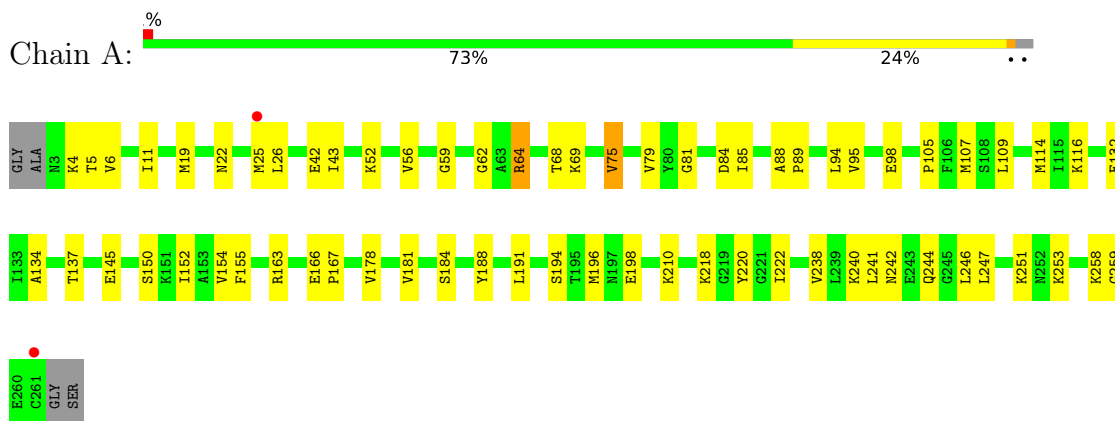
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	154	Total O 154 154	0	0
7	B	222	Total O 222 222	0	0
7	C	192	Total O 192 192	0	0

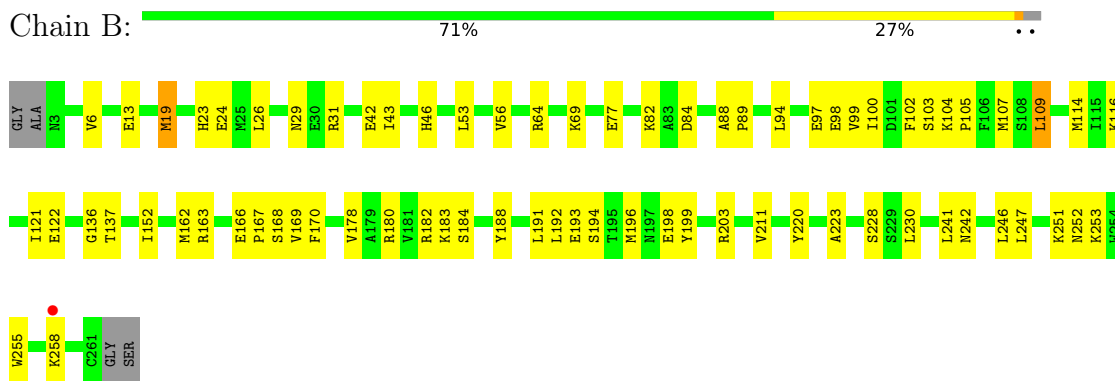
### 3 Residue-property plots

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.

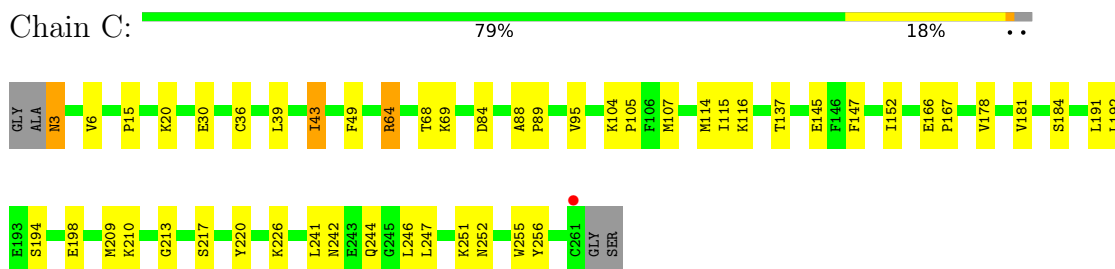
- Molecule 1: Glutamate receptor 2



- Molecule 1: Glutamate receptor 2



- Molecule 1: Glutamate receptor 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	113.82Å 163.33Å 47.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.27 20.00 – 2.27	Depositor EDS
% Data completeness (in resolution range)	98.8 (20.00-2.27) 98.7 (20.00-2.27)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.03 (at 2.26Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.200 , 0.250 0.196 , 0.198	Depositor DCC
$R_{free}$ test set	2079 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.0	Xtrriage
Anisotropy	0.226	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 39.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6695	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.90% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: NA, CL, CAC, S2P, ZN

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	0.38	0/2062	0.83	2/2773 (0.1%)
1	B	0.40	0/2062	0.86	5/2773 (0.2%)
1	C	0.40	0/2062	0.85	3/2773 (0.1%)
All	All	0.39	0/6186	0.84	10/8319 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	116	LYS	N-CA-C	-5.99	100.24	109.76
1	B	182	ARG	N-CA-C	5.70	117.49	111.28
1	B	169	VAL	N-CA-C	-5.67	107.21	112.83
1	B	116	LYS	N-CA-C	-5.52	101.47	110.20
1	A	210	LYS	N-CA-C	-5.45	100.30	108.96
1	B	104	LYS	N-CA-C	-5.32	102.44	109.84
1	C	104	LYS	N-CA-C	-5.29	102.78	110.08
1	B	103	SER	N-CA-C	-5.22	103.80	110.53
1	A	116	LYS	N-CA-C	-5.12	102.10	110.20
1	C	256	TYR	N-CA-C	5.02	119.40	113.12

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	2026	0	2056	47	0
1	B	2026	0	2056	60	0
1	C	2026	0	2056	37	0
2	A	12	0	5	1	0
2	B	12	0	5	1	0
2	C	12	0	5	0	0
3	A	2	0	0	0	0
3	B	3	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
6	B	5	0	0	1	0
7	A	154	0	0	11	0
7	B	222	0	0	13	0
7	C	192	0	0	13	0
All	All	6695	0	6183	145	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 12.

All (145) 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:B:136:GLY:HA2	1:B:162:MET:HE3	1.43	0.97
1:B:137:THR:HG23	1:B:162:MET:HE1	1.51	0.93
1:A:62:GLY:HA3	1:A:75:VAL:HG12	1.54	0.89
1:A:22:ASN:HB2	1:A:25:MET:HE2	1.54	0.88
1:C:107:MET:HE2	1:C:251:LYS:HD2	1.58	0.84
1:A:241:LEU:HD22	1:A:246:LEU:HD22	1.64	0.79
1:B:241:LEU:HD22	1:B:246:LEU:HD22	1.65	0.79
1:A:62:GLY:HA3	1:A:75:VAL:CG1	2.13	0.78
1:B:253:LYS:O	1:B:258:LYS:HD3	1.84	0.78
1:B:137:THR:HG22	1:B:191:LEU:HB2	1.66	0.78
1:A:114:MET:HE1	1:A:178:VAL:HA	1.69	0.75
1:A:42:GLU:HG3	1:A:246:LEU:HD21	1.70	0.73
1:C:252:ASN:HB3	7:C:284:HOH:O	1.87	0.73
1:C:114:MET:HE1	1:C:178:VAL:HA	1.71	0.73
1:B:184:SER:HB2	7:B:1617:HOH:O	1.86	0.73
1:B:194:SER:O	1:B:198:GLU:HG3	1.88	0.72
1:C:210:LYS:NZ	7:C:435:HOH:O	2.23	0.71
1:A:253:LYS:O	1:A:258:LYS:HD3	1.91	0.70
1:B:64:ARG:CZ	7:B:1661:HOH:O	2.40	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:MET:HE2	1:B:251:LYS:HD3	1.75	0.69
1:C:115:ILE:HD11	1:C:209:MET:HE2	1.75	0.69
1:B:19:MET:HE1	1:B:31:ARG:HB3	1.76	0.68
1:B:42:GLU:HG3	1:B:246:LEU:HD21	1.77	0.67
1:C:3:ASN:HA	7:C:404:HOH:O	1.95	0.66
1:A:194:SER:O	1:A:198:GLU:HG3	1.96	0.66
1:A:184:SER:HB2	7:A:1150:HOH:O	1.96	0.65
1:A:81:GLY:HA2	7:A:1147:HOH:O	1.97	0.64
1:A:88:ALA:HB1	1:A:89:PRO:HD2	1.79	0.63
1:C:184:SER:HB2	7:C:272:HOH:O	1.99	0.63
1:B:122:GLU:HA	1:B:211:VAL:CG2	2.29	0.63
1:A:137:THR:HG22	1:A:191:LEU:HB2	1.82	0.62
1:B:122:GLU:HA	1:B:211:VAL:HG21	1.81	0.62
1:A:134:ALA:O	1:A:188:TYR:HA	2.00	0.61
1:C:114:MET:HE2	1:C:181:VAL:HG21	1.83	0.61
1:A:105:PRO:HA	1:A:220:TYR:O	2.00	0.61
1:A:75:VAL:O	1:A:79:VAL:HG13	2.02	0.60
1:C:88:ALA:HB1	1:C:89:PRO:HD2	1.83	0.60
1:A:94:LEU:O	1:A:98:GLU:HG3	2.02	0.60
1:B:29:ASN:HD21	1:B:53:LEU:H	1.48	0.60
1:B:183:LYS:HB3	7:B:1474:HOH:O	2.02	0.59
1:A:64:ARG:HG3	1:A:64:ARG:HH11	1.68	0.58
1:A:43:ILE:HD13	1:A:222:ILE:HD13	1.87	0.57
1:C:137:THR:HG22	1:C:191:LEU:HB2	1.85	0.57
1:C:226:LYS:HE2	7:C:269:HOH:O	2.03	0.56
1:B:162:MET:HE2	1:B:170:PHE:CE1	2.40	0.56
1:A:154:VAL:HG23	7:A:1051:HOH:O	2.05	0.56
1:A:240:LYS:O	1:A:244:GLN:HG3	2.05	0.56
1:B:29:ASN:ND2	1:B:53:LEU:H	2.04	0.55
1:C:194:SER:O	1:C:198:GLU:HG3	2.07	0.55
1:C:145:GLU:HG2	7:C:376:HOH:O	2.07	0.55
1:A:196:MET:HE2	2:A:1009:S2P:SE1	2.57	0.55
1:A:145:GLU:HG2	7:A:1089:HOH:O	2.07	0.54
1:B:168:SER:HB2	6:B:1439:CAC:O1	2.08	0.54
1:A:163:ARG:HD2	7:A:1146:HOH:O	2.08	0.54
1:C:147:PHE:CZ	1:C:191:LEU:HD13	2.44	0.53
1:C:241:LEU:HD22	1:C:246:LEU:HD22	1.90	0.53
1:A:19:MET:HE3	7:A:1161:HOH:O	2.09	0.53
1:B:99:VAL:HG23	1:B:100:ILE:HG23	1.92	0.52
1:A:152:ILE:HD12	1:A:152:ILE:N	2.24	0.52
1:B:24:GLU:CD	1:B:24:GLU:H	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:ASN:HB3	7:B:1502:HOH:O	2.10	0.52
1:C:107:MET:HE3	7:C:266:HOH:O	2.10	0.52
1:B:180:ARG:HD3	7:B:1646:HOH:O	2.10	0.51
1:B:241:LEU:HD22	1:B:246:LEU:CD2	2.38	0.51
1:B:105:PRO:HA	1:B:220:TYR:O	2.11	0.51
1:B:152:ILE:HD12	1:B:152:ILE:N	2.25	0.51
1:C:152:ILE:H	1:C:152:ILE:HD12	1.76	0.51
1:C:192:LEU:H	1:C:192:LEU:HD23	1.76	0.51
1:B:88:ALA:HB1	1:B:89:PRO:HD2	1.93	0.50
1:B:23:HIS:CD2	1:B:26:LEU:HD12	2.47	0.50
1:C:68:THR:O	1:C:69:LYS:HB2	2.11	0.50
1:C:213:GLY:C	7:C:435:HOH:O	2.55	0.50
1:A:4:LYS:HG2	1:A:5:THR:N	2.26	0.49
1:C:39:LEU:HG	1:C:43:ILE:CD1	2.41	0.49
1:B:180:ARG:HD2	1:B:188:TYR:CD2	2.48	0.49
1:B:242:ASN:HB2	1:B:247:LEU:HD12	1.94	0.49
1:B:97:GLU:HG2	1:B:102:PHE:HD2	1.78	0.49
1:A:218:LYS:HB3	7:C:443:HOH:O	2.12	0.49
2:B:1440:S2P:HBA	7:B:1649:HOH:O	2.13	0.49
1:B:136:GLY:CA	1:B:162:MET:HE3	2.29	0.48
1:A:242:ASN:HB2	1:A:247:LEU:HD12	1.95	0.48
1:A:152:ILE:HD12	1:A:152:ILE:H	1.77	0.48
1:B:69:LYS:CD	7:B:1661:HOH:O	2.62	0.48
1:A:52:LYS:HE3	7:A:1105:HOH:O	2.13	0.48
1:C:242:ASN:HB2	1:C:247:LEU:HD12	1.96	0.47
1:B:69:LYS:HE3	7:B:1661:HOH:O	2.13	0.47
1:B:180:ARG:HD2	1:B:188:TYR:CG	2.49	0.47
1:B:82:LYS:HG3	7:B:1638:HOH:O	2.13	0.47
1:C:152:ILE:HD12	1:C:152:ILE:N	2.29	0.47
1:A:43:ILE:HD11	1:A:222:ILE:HD11	1.97	0.47
1:B:64:ARG:HD3	1:B:69:LYS:HA	1.97	0.47
1:C:251:LYS:O	1:C:255:TRP:HB2	2.15	0.47
1:B:109:LEU:HD13	1:B:193:GLU:HB3	1.97	0.46
1:A:68:THR:O	1:A:69:LYS:HB2	2.14	0.46
1:C:6:VAL:HG13	1:C:84:ASP:HB2	1.97	0.46
1:B:13:GLU:HG2	1:B:196:MET:HE3	1.97	0.46
1:B:168:SER:HA	7:B:1476:HOH:O	2.16	0.45
1:B:64:ARG:HG3	1:B:64:ARG:NH1	2.31	0.45
1:B:114:MET:HE1	1:B:178:VAL:HG22	1.98	0.45
1:A:43:ILE:CD1	1:A:222:ILE:CD1	2.95	0.45
1:A:218:LYS:HD3	7:C:443:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:MET:HE2	1:B:170:PHE:HE1	1.81	0.45
1:B:64:ARG:HG3	1:B:64:ARG:HH11	1.81	0.45
1:C:166:GLU:HA	1:C:167:PRO:C	2.41	0.45
1:C:192:LEU:HD23	1:C:192:LEU:N	2.31	0.45
1:B:199:TYR:O	1:B:203:ARG:HG2	2.17	0.44
1:B:94:LEU:O	1:B:98:GLU:HG3	2.17	0.44
1:B:163:ARG:NH1	7:B:1662:HOH:O	2.41	0.44
1:A:95:VAL:HG22	7:A:1026:HOH:O	2.18	0.44
1:B:69:LYS:HD2	7:B:1661:HOH:O	2.16	0.44
1:B:121:ILE:O	1:B:211:VAL:HG21	2.17	0.44
1:C:105:PRO:HA	1:C:220:TYR:O	2.17	0.44
1:A:107:MET:HE2	1:A:251:LYS:HD3	1.98	0.44
1:C:95:VAL:HG22	7:C:296:HOH:O	2.17	0.43
1:A:150:SER:OG	1:A:155:PHE:HB2	2.18	0.43
1:B:19:MET:CE	1:B:31:ARG:HB3	2.46	0.43
1:B:42:GLU:O	1:B:46:HIS:HD2	2.01	0.43
1:C:244:GLN:NE2	7:C:437:HOH:O	2.34	0.42
1:A:56:VAL:HG13	1:A:59:GLY:HA2	2.01	0.42
1:B:56:VAL:O	1:B:56:VAL:HG13	2.20	0.42
1:A:6:VAL:HG13	1:A:84:ASP:HB2	2.01	0.42
1:A:85:ILE:HD11	1:A:222:ILE:CG2	2.49	0.42
1:B:100:ILE:HD12	1:B:223:ALA:HB1	2.01	0.42
1:A:11:ILE:HG13	7:A:1060:HOH:O	2.19	0.42
1:A:222:ILE:HD12	1:A:238:VAL:CG2	2.50	0.42
1:C:217:SER:CB	7:C:438:HOH:O	2.67	0.42
1:C:39:LEU:HG	1:C:43:ILE:HD12	2.01	0.41
1:C:49:PHE:CD1	1:C:49:PHE:C	2.99	0.41
1:A:166:GLU:HA	1:A:167:PRO:C	2.44	0.41
1:A:64:ARG:NH1	7:A:1083:HOH:O	2.46	0.41
1:B:166:GLU:HA	1:B:167:PRO:C	2.46	0.41
1:B:180:ARG:HD2	1:B:188:TYR:CD1	2.56	0.41
1:B:64:ARG:NH2	7:B:1661:HOH:O	2.51	0.41
1:B:251:LYS:HD2	1:B:255:TRP:CE3	2.56	0.41
1:A:163:ARG:NH1	7:A:1146:HOH:O	2.49	0.41
1:C:114:MET:CE	1:C:178:VAL:HA	2.46	0.41
1:B:137:THR:HG23	1:B:162:MET:CE	2.37	0.40
1:C:36:CYS:SG	1:C:88:ALA:HA	2.61	0.40
1:C:64:ARG:HH11	1:C:64:ARG:HG3	1.86	0.40
1:B:192:LEU:HD23	1:B:192:LEU:H	1.87	0.40
1:B:228:SER:C	1:B:230:LEU:H	2.29	0.40
1:A:43:ILE:HD11	1:A:222:ILE:CD1	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:MET:HE2	1:A:181:VAL:HG21	2.02	0.40
1:B:6:VAL:HG13	1:B:84:ASP:HB2	2.03	0.40
1:C:20:LYS:HD2	1:C:30:GLU:O	2.21	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	257/263 (98%)	248 (96%)	8 (3%)	1 (0%)	30	36
1	B	257/263 (98%)	252 (98%)	5 (2%)	0	100	100
1	C	257/263 (98%)	249 (97%)	8 (3%)	0	100	100
All	All	771/789 (98%)	749 (97%)	21 (3%)	1 (0%)	48	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	259	GLY

### 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	218/219 (100%)	213 (98%)	5 (2%)	44	60

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	218/219 (100%)	214 (98%)	4 (2%)	51	68
1	C	218/219 (100%)	214 (98%)	4 (2%)	51	68
All	All	654/657 (100%)	641 (98%)	13 (2%)	48	65

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LEU
1	A	64	ARG
1	A	75	VAL
1	A	109	LEU
1	A	132	GLU
1	B	19	MET
1	B	43	ILE
1	B	77	GLU
1	B	109	LEU
1	C	3	ASN
1	C	15	PRO
1	C	43	ILE
1	C	64	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	22	ASN
1	B	29	ASN
1	B	214	ASN
1	B	232	ASN
1	B	242	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 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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
2	S2P	B	1440	-	10,12,12	4.01	4 (40%)	8,16,16	2.74	3 (37%)
6	CAC	B	1439	3	2,4,4	2.23	1 (50%)	4,6,6	0.86	0
2	S2P	A	1009	-	10,12,12	4.36	4 (40%)	8,16,16	3.01	4 (50%)
2	S2P	C	264	-	10,12,12	4.50	3 (30%)	8,16,16	3.27	4 (50%)

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
2	S2P	B	1440	-	-	1/8/8/8	0/1/1/1
2	S2P	A	1009	-	-	3/8/8/8	0/1/1/1
2	S2P	C	264	-	-	1/8/8/8	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	264	S2P	SE1-NE2	10.26	1.79	1.62
2	A	1009	S2P	SE1-ND1	9.27	1.77	1.62
2	A	1009	S2P	SE1-NE2	8.73	1.76	1.62
2	B	1440	S2P	SE1-ND1	8.69	1.76	1.62
2	C	264	S2P	SE1-ND1	8.32	1.75	1.62
2	B	1440	S2P	SE1-NE2	8.31	1.75	1.62
2	C	264	S2P	CD2-NE2	4.14	1.38	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1009	S2P	CD2-NE2	3.60	1.37	1.30
2	A	1009	S2P	OD2-CD2	-2.79	1.25	1.33
6	B	1439	CAC	AS-C1	2.76	1.96	1.90
2	B	1440	S2P	OD2-CD2	-2.50	1.26	1.33
2	B	1440	S2P	CD2-NE2	2.00	1.34	1.30

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	264	S2P	NE2-SE1-ND1	-6.83	92.92	100.70
2	A	1009	S2P	NE2-SE1-ND1	-6.65	93.12	100.70
2	B	1440	S2P	NE2-SE1-ND1	-6.38	93.43	100.70
2	C	264	S2P	OD2-CD2-NE2	3.86	130.55	122.92
2	C	264	S2P	OD2-CD2-CG	-3.51	115.28	124.33
2	A	1009	S2P	OD2-CD2-CG	-3.31	115.78	124.33
2	A	1009	S2P	OD2-CD2-NE2	2.99	128.85	122.92
2	C	264	S2P	OXT-C-O	2.73	130.27	124.08
2	B	1440	S2P	OD2-CD2-CG	-2.59	117.66	124.33
2	A	1009	S2P	OXT-C-O	2.29	129.28	124.08
2	B	1440	S2P	OD2-CD2-NE2	2.29	127.45	122.92

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1009	S2P	CA-CB-CG-ND1
2	A	1009	S2P	O-C-CA-N
2	B	1440	S2P	O-C-CA-N
2	C	264	S2P	O-C-CA-N
2	A	1009	S2P	O-C-CA-CB

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1440	S2P	1	0
6	B	1439	CAC	1	0
2	A	1009	S2P	1	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	259/263 (98%)	-0.14	2 (0%) 82 84	18, 31, 52, 59	0
1	B	259/263 (98%)	-0.39	1 (0%) 88 89	15, 26, 44, 58	0
1	C	259/263 (98%)	-0.29	1 (0%) 88 89	16, 28, 44, 57	0
All	All	777/789 (98%)	-0.27	4 (0%) 87 88	15, 28, 48, 59	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	261	CYS	2.4
1	A	25	MET	2.3
1	A	261	CYS	2.2
1	B	258	LYS	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, 95<sup>th</sup> 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
5	CL	A	1008	1/1	0.60	0.24	84,84,84,84	0
2	S2P	A	1009	12/12	0.89	0.11	25,34,46,50	0
2	S2P	C	264	12/12	0.90	0.09	24,34,43,49	0
4	NA	B	1006	1/1	0.91	0.18	44,44,44,44	0
4	NA	A	1007	1/1	0.92	0.18	50,50,50,50	0
3	ZN	A	1003	1/1	0.93	0.08	40,40,40,40	0
2	S2P	B	1440	12/12	0.94	0.08	19,26,34,39	0
3	ZN	B	1004	1/1	0.97	0.03	31,31,31,31	0
3	ZN	B	1001	1/1	0.97	0.04	32,32,32,32	0
6	CAC	B	1439	5/5	0.98	0.10	46,47,49,50	0
3	ZN	A	1002	1/1	0.99	0.03	61,61,61,61	0
3	ZN	B	1005	1/1	0.99	0.02	25,25,25,25	0

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

There are no such residues in this entry.