



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:43 PM UTC

PDB ID : 2X11 / pdb_00002x11
Title : Crystal structure of the complete EphA2 ectodomain in complex with ephrin A5 receptor binding domain
Authors : Seiradake, E.; Harlos, K.; Sutton, G.; Aricescu, A.R.; Jones, E.Y.
Deposited on : 2009-12-21
Resolution : 4.83 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
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

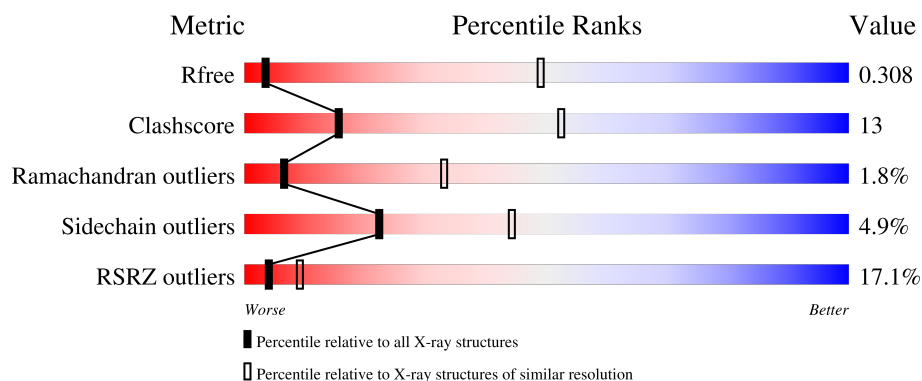
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.83 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	1024 (5.74-3.94)
Clashscore	190562	1007 (5.70-3.98)
Ramachandran outliers	187476	1110 (5.78-3.90)
Sidechain outliers	187428	1091 (5.78-3.90)
RSRZ outliers	180081	1019 (5.74-3.94)

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	545	14% 64% 23% • 10%
2	B	177	16% 53% 22% • • 21%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4889 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 EPHRIN TYPE-A RECEPTOR 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	2	0
			3738	2358	628	724	28			

- Molecule 2 is a protein called EPHRIN-A5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	139	Total	C	N	O	S	0	0	1
			1151	736	199	208	8			

4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	173.63Å 59.63Å 112.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.48 – 4.83 40.48 – 4.83	Depositor EDS
% Data completeness (in resolution range)	93.7 (40.48-4.83) 99.6 (40.48-4.83)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 4.86Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.312 , 0.314 0.304 , 0.308	Depositor DCC
R_{free} test set	275 reflections (4.57%)	wwPDB-VP
Wilson B-factor (Å ²)	180.0	Xtriage
Anisotropy	0.319	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 257.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.78	EDS
Total number of atoms	4889	wwPDB-VP
Average B, all atoms (Å ²)	247.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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	0.26	0/3834	0.75	7/5226 (0.1%)
2	B	0.32	1/1190 (0.1%)	0.78	0/1612
All	All	0.28	1/5024 (0.0%)	0.76	7/6838 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	170	MET	C-N	-7.01	1.23	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	237	PRO	N-CA-C	6.59	118.74	110.70
1	A	236	VAL	CA-C-N	6.02	126.58	120.38
1	A	236	VAL	C-N-CA	6.02	126.58	120.38
1	A	201	CYS	CA-C-N	5.30	125.21	119.85
1	A	201	CYS	C-N-CA	5.30	125.21	119.85
1	A	403	GLU	CA-C-N	5.15	125.09	119.78
1	A	403	GLU	C-N-CA	5.15	125.09	119.78

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	3738	0	3490	94	0
2	B	1151	0	1068	49	0
All	All	4889	0	4558	127	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 13.

All (127) 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:B:126:THR:HG21	2:B:131:GLY:HA3	1.24	1.11
1:A:159:ARG:HA	2:B:130:LEU:CD1	1.89	1.02
1:A:159:ARG:HA	2:B:130:LEU:HD13	1.45	0.98
1:A:92:GLU:HB3	1:A:218:SER:HB2	1.51	0.93
1:A:156:PHE:O	2:B:130:LEU:HD21	1.70	0.91
2:B:154:ARG:HH11	2:B:154:ARG:HB2	1.37	0.90
2:B:87:ASN:OD1	2:B:99:LYS:HB3	1.80	0.81
1:A:293:CYS:HB3	1:A:297:THR:HG21	1.63	0.78
1:A:237:PRO:HD2	1:A:238:PRO:HD3	1.67	0.76
1:A:68:SER:HB2	2:B:127:PRO:HB2	1.69	0.75
2:B:126:THR:HG22	2:B:128:PHE:H	1.53	0.74
1:A:190:ALA:HB3	2:B:128:PHE:HD2	1.53	0.73
2:B:35:ARG:HG2	2:B:61:TYR:HB2	1.69	0.73
1:A:156:PHE:HA	2:B:130:LEU:HG	1.71	0.72
1:A:295:GLU:HB2	1:A:324:PRO:HB3	1.71	0.72
1:A:370:TRP:HB3	1:A:373:SER:HB3	1.71	0.71
2:B:170:MET:H	2:B:170:MET:HE2	1.59	0.67
1:A:103:ARG:HB3	1:A:188:CYS:HB3	1.76	0.65
2:B:141:PHE:CE2	2:B:160:LYS:HG3	2.32	0.65
2:B:126:THR:HG21	2:B:131:GLY:CA	2.16	0.64
1:A:215:ILE:HD13	1:A:215:ILE:N	2.11	0.64
2:B:88:PHE:O	2:B:91:TYR:N	2.32	0.62
1:A:237:PRO:CD	1:A:238:PRO:HD3	2.29	0.61
1:A:269:VAL:HB	1:A:274:GLN:HG3	1.82	0.61
1:A:335:LEU:HD22	1:A:412:VAL:HG12	1.84	0.59
2:B:80:ARG:HB3	2:B:147:ILE:HD12	1.83	0.59
1:A:153:SER:O	1:A:156:PHE:HB2	2.02	0.59
1:A:246:HIS:ND1	1:A:255:VAL:HG23	2.18	0.59
2:B:151:ASN:HB3	2:B:154:ARG:HH22	1.68	0.58
1:A:411:THR:HG22	1:A:429:THR:HG22	1.86	0.57
1:A:460:PRO:N	1:A:461:PRO:HD2	2.19	0.57
1:A:58:ILE:HG21	2:B:102:LYS:NZ	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ARG:HA	2:B:130:LEU:HD11	1.83	0.56
1:A:99:LYS:HB2	1:A:193:SER:HB3	1.85	0.56
1:A:380:GLU:O	1:A:383:VAL:HG22	2.05	0.56
2:B:170:MET:HE2	2:B:170:MET:N	2.21	0.56
2:B:154:ARG:HB2	2:B:154:ARG:NH1	2.15	0.56
1:A:223:LEU:HD21	1:A:254:LEU:HD12	1.87	0.56
1:A:448:SER:O	1:A:449:THR:HB	2.05	0.56
1:A:482:TYR:CD2	1:A:482:TYR:N	2.74	0.55
2:B:58:ILE:O	2:B:59:ASN:HB2	2.06	0.55
1:A:473:TYR:CE2	1:A:483:ASN:HB2	2.41	0.55
2:B:102:LYS:NZ	2:B:102:LYS:HB3	2.21	0.55
1:A:508:GLN:HA	1:A:518:GLY:HA2	1.88	0.55
1:A:344:VAL:CG2	1:A:399:VAL:HB	2.37	0.54
1:A:159:ARG:CA	2:B:130:LEU:CD1	2.76	0.54
1:A:102:VAL:HG21	1:A:147:PRO:HG3	1.90	0.53
1:A:245:MET:HB3	1:A:253:TRP:CE3	2.44	0.53
1:A:293:CYS:HB3	1:A:297:THR:CG2	2.38	0.52
2:B:97:THR:O	2:B:99:LYS:HD3	2.08	0.52
1:A:112:ALA:C	1:A:114:SER:H	2.17	0.52
2:B:122:PHE:CE1	2:B:161:VAL:HG11	2.45	0.51
1:A:152:VAL:O	1:A:155:ASP:HB2	2.10	0.51
1:A:70:CYS:HB2	2:B:127:PRO:HG3	1.93	0.51
1:A:474:ARG:HG2	1:A:475:LYS:N	2.27	0.50
2:B:101:PHE:CD1	2:B:101:PHE:N	2.79	0.50
1:A:108:PHE:HZ	1:A:188:CYS:SG	2.34	0.50
2:B:126:THR:CG2	2:B:131:GLY:HA3	2.17	0.50
1:A:85:TRP:CZ2	1:A:177:GLY:HA3	2.47	0.49
1:A:123:TYR:HB3	1:A:142:ILE:HD11	1.94	0.49
1:A:444:LEU:HD11	1:A:524:PHE:HB3	1.95	0.49
1:A:440:PRO:HA	1:A:458:ILE:HG23	1.94	0.49
1:A:449:THR:OG1	1:A:499:PRO:HA	2.13	0.49
1:A:471:VAL:O	1:A:484:VAL:HA	2.13	0.48
1:A:344:VAL:HG22	1:A:402:LEU:HD11	1.96	0.48
1:A:328:PRO:HG2	1:A:422:VAL:HG11	1.96	0.48
1:A:190:ALA:HB3	2:B:128:PHE:CD2	2.42	0.48
1:A:46:HIS:HA	1:A:47:PRO:C	2.39	0.47
1:A:281:PHE:C	1:A:281:PHE:CD1	2.93	0.47
1:A:236:VAL:HA	1:A:237:PRO:HD3	1.58	0.47
1:A:237:PRO:HG2	1:A:261:LEU:HD12	1.97	0.47
2:B:83:LEU:HD22	2:B:144:SER:HB3	1.97	0.46
1:A:44:LEU:C	1:A:44:LEU:HD23	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:PHE:HB3	1:A:325:CYS:HB3	1.96	0.46
1:A:327:ARG:O	1:A:355:GLY:HA3	2.16	0.46
1:A:392:LEU:HD21	1:A:397:VAL:HG11	1.97	0.46
1:A:468:LYS:O	1:A:510:LEU:HB2	2.16	0.46
2:B:88:PHE:O	2:B:89:ASP:C	2.59	0.46
2:B:103:ARG:HG3	2:B:104:TRP:CE2	2.51	0.46
1:A:70:CYS:HB2	2:B:127:PRO:CA	2.46	0.45
1:A:456:TRP:HZ3	1:A:492:VAL:HG22	1.80	0.45
2:B:73:VAL:HA	2:B:74:PRO:HD3	1.82	0.45
1:A:220:ALA:N	1:A:221:PRO:HD2	2.31	0.45
2:B:128:PHE:CD1	2:B:129:SER:O	2.70	0.45
1:A:447:ARG:HG3	1:A:527:LEU:CD1	2.47	0.45
1:A:294:PRO:HG3	1:A:321:ALA:O	2.17	0.45
1:A:188:CYS:SG	2:B:127:PRO:HA	2.57	0.45
1:A:385:TYR:C	1:A:387:GLU:H	2.25	0.44
1:A:369:CYS:O	1:A:371:PRO:HD3	2.18	0.44
1:A:387:GLU:O	1:A:388:PRO:C	2.61	0.44
1:A:440:PRO:HA	1:A:458:ILE:CG2	2.47	0.44
1:A:264:ALA:HB1	1:A:305:THR:HG22	1.99	0.43
1:A:348:TRP:CZ3	1:A:397:VAL:HG22	2.53	0.43
1:A:344:VAL:HG23	1:A:399:VAL:HB	1.99	0.43
1:A:74:SER:HB2	1:A:77:GLN:NE2	2.33	0.43
1:A:96:ILE:O	1:A:167:GLU:HA	2.18	0.43
1:A:348:TRP:CH2	1:A:397:VAL:HG22	2.53	0.43
1:A:220:ALA:C	1:A:222:SER:H	2.27	0.43
2:B:108:ARG:HB2	2:B:116:LEU:HD12	2.00	0.42
1:A:337:ALA:HB1	1:A:432:VAL:CG2	2.49	0.42
1:A:215:ILE:N	1:A:215:ILE:CD1	2.74	0.42
1:A:262:CYS:O	1:A:282:LYS:NZ	2.52	0.42
2:B:73:VAL:C	2:B:75:GLU:H	2.28	0.42
1:A:189:VAL:HG12	1:A:190:ALA:N	2.33	0.42
1:A:472:THR:HG23	1:A:482:TYR:HD1	1.85	0.42
2:B:68:HIS:HE1	2:B:111:SER:O	2.03	0.42
2:B:116:LEU:HD23	2:B:117:LYS:N	2.35	0.42
1:A:92:GLU:HB3	1:A:218:SER:CB	2.36	0.42
1:A:254:LEU:HD23	1:A:254:LEU:HA	1.78	0.42
1:A:130:TYR:HB3	1:A:133:ASN:HB2	2.01	0.42
1:A:69:VAL:C	2:B:127:PRO:HB3	2.45	0.41
1:A:70:CYS:HB2	2:B:127:PRO:CG	2.51	0.41
2:B:147:ILE:N	2:B:148:PRO:HD2	2.35	0.41
2:B:141:PHE:CD1	2:B:141:PHE:N	2.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:GLU:HG3	1:A:388:PRO:HD2	2.02	0.41
1:A:466:VAL:HG12	1:A:509:ALA:HB1	2.03	0.41
1:A:80:TRP:NE1	1:A:136:LYS:HE3	2.35	0.41
1:A:459:PRO:C	1:A:461:PRO:HD2	2.45	0.41
2:B:126:THR:HA	2:B:127:PRO:HD3	1.74	0.41
1:A:99:LYS:HA	1:A:164:ASN:O	2.21	0.40
1:A:154:SER:C	1:A:156:PHE:H	2.29	0.40
1:A:192:LEU:HD21	2:B:128:PHE:CE2	2.56	0.40
2:B:122:PHE:CZ	2:B:161:VAL:HG11	2.57	0.40
1:A:146:ALA:HA	1:A:147:PRO:HD3	1.92	0.40
1:A:403:GLU:HA	1:A:404:PRO:HD3	1.84	0.40
1:A:448:SER:HA	1:A:527:LEU:HD22	2.04	0.40
2:B:102:LYS:HB3	2:B:102:LYS:HZ3	1.84	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	476/545 (87%)	436 (92%)	37 (8%)	3 (1%)	21	58
2	B	137/177 (77%)	123 (90%)	6 (4%)	8 (6%)	1	14
All	All	613/722 (85%)	559 (91%)	43 (7%)	11 (2%)	6	33

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	112	PRO
2	B	154	ARG
2	B	113	ASN
1	A	480	ASN
2	B	33	ALA

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Mol	Chain	Res	Type
2	B	88	PHE
2	B	150	ASP
2	B	152	GLY
1	A	237	PRO
1	A	294	PRO
2	B	74	PRO

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	402/464 (87%)	379 (94%)	23 (6%)	18	41
2	B	127/159 (80%)	124 (98%)	3 (2%)	43	63
All	All	529/623 (85%)	503 (95%)	26 (5%)	22	44

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	128	LEU
1	A	150	ILE
1	A	157	GLU
1	A	174	THR
1	A	188	CYS
1	A	191	LEU
1	A	192	LEU
1	A	204	LEU
1	A	215	ILE
1	A	284	GLU
1	A	349	THR
1	A	379	CYS
1	A	387	GLU
1	A	392	LEU
1	A	397	VAL
1	A	411	THR

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Mol	Chain	Res	Type
1	A	441	LYS
1	A	442	VAL
1	A	458	ILE
1	A	462	GLN
1	A	484	VAL
1	A	494	LEU
2	B	101	PHE
2	B	102	LYS
2	B	154	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	77	GLN
1	A	79	ASN
1	A	259	GLN
1	A	274	GLN
1	A	333	HIS
2	B	48	GLN
2	B	68	HIS

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](#)

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

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	488/545 (89%)	0.99	78 (15%) 5 10	98, 250, 285, 314	2 (0%)
2	B	139/177 (78%)	1.30	29 (20%) 2 7	253, 269, 293, 298	0
All	All	627/722 (86%)	1.06	107 (17%) 4 9	98, 259, 287, 314	2 (0%)

All (107) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	412	VAL	7.2
2	B	139	GLU	6.7
1	A	470	GLU	6.3
2	B	104	TRP	5.6
2	B	64	VAL	5.6
1	A	516	GLY	5.2
1	A	182	PHE	5.0
1	A	344	VAL	4.5
2	B	159	LEU	4.5
1	A	484	VAL	4.4
1	A	471	VAL	4.4
1	A	134	PHE	4.4
2	B	54	ILE	4.3
2	B	51	ASP	4.3
1	A	364	VAL	4.2
1	A	505	VAL	4.1
1	A	472	THR	4.1
1	A	492	VAL	4.1
1	A	393	THR	4.0
1	A	482	TYR	4.0
2	B	62	LEU	4.0
2	B	50	GLY	3.9
1	A	119	PHE	3.9
1	A	463	GLN	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	481	SER	3.7
1	A	432	VAL	3.6
2	B	122	PHE	3.6
2	B	161	VAL	3.6
1	A	410	PHE	3.6
1	A	511	THR	3.6
1	A	81	LEU	3.6
1	A	223	LEU	3.5
1	A	180	LEU	3.4
1	A	517	ALA	3.3
1	A	486	ARG	3.2
1	A	348	TRP	3.2
1	A	241	GLU	3.2
2	B	48	GLN	3.1
2	B	142	TYR	3.1
1	A	221	PRO	3.1
1	A	83	THR	3.1
2	B	98	SER	3.1
1	A	220	ALA	3.1
1	A	96	ILE	3.0
2	B	102	LYS	2.9
1	A	189	VAL	2.9
2	B	141	PHE	2.9
1	A	334	TYR	2.8
1	A	439	PRO	2.8
1	A	191	LEU	2.8
2	B	53	HIS	2.8
1	A	462	GLN	2.8
1	A	456	TRP	2.7
1	A	84[A]	ASN	2.7
1	A	98	LEU	2.7
1	A	143	ASP	2.6
1	A	48	TYR	2.6
1	A	238	PRO	2.6
1	A	490	PHE	2.6
1	A	485	ARG	2.6
2	B	113	ASN	2.6
1	A	375	GLU	2.6
1	A	176	LYS	2.6
2	B	34	ASP	2.6
2	B	171	LYS	2.5
2	B	100	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	479	SER	2.5
1	A	438	GLU	2.5
1	A	494	LEU	2.5
1	A	338	VAL	2.5
1	A	423	THR	2.5
2	B	36	TYR	2.4
1	A	335	LEU	2.4
1	A	257	ILE	2.4
1	A	362	TYR	2.4
1	A	82	ARG	2.4
1	A	489	GLY	2.4
1	A	63	PRO	2.3
1	A	518	GLY	2.3
1	A	528	SER	2.3
1	A	495	ASP	2.3
1	A	236	VAL	2.3
1	A	347[A]	ARG	2.3
1	A	367	GLU	2.3
2	B	41	ASN	2.3
1	A	424	SER	2.3
1	A	255	VAL	2.3
1	A	121	LEU	2.3
1	A	373	SER	2.2
1	A	76	ASP	2.2
2	B	112	PRO	2.2
1	A	67	TYR	2.2
1	A	254	LEU	2.2
1	A	512	GLN	2.2
1	A	132	THR	2.2
2	B	128	PHE	2.1
1	A	61	ASP	2.1
1	A	506	GLN	2.1
2	B	160	LYS	2.1
1	A	402	LEU	2.1
2	B	151	ASN	2.0
2	B	43	SER	2.0
2	B	56	VAL	2.0
1	A	394	ARG	2.0
1	A	487	THR	2.0
1	A	464	SER	2.0
2	B	118	PHE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.