



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:35 PM UTC

PDB ID : 4GH4 / pdb_00004gh4
Title : Crystal Structure of Foot and Mouth Disease Virus A22 Serotype
Authors : Kotecha, A.; Jinshan, R.; Curry, S.; Fry, E.; Stuart, D.
Deposited on : 2012-08-07
Resolution : 3.00 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

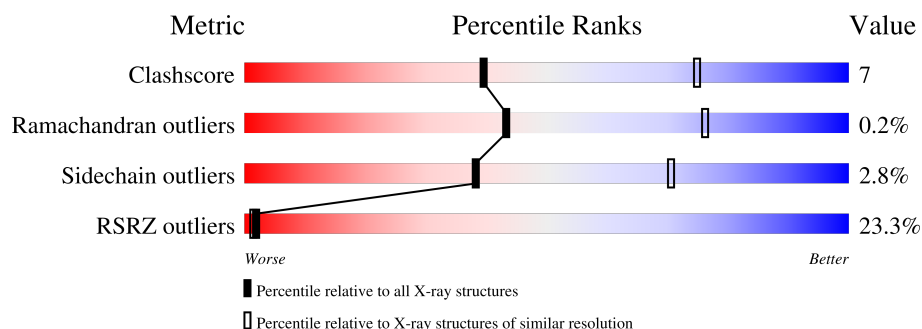
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	210	
2	B	207	
3	C	221	
4	D	45	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5457 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 capsid protein VP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1494	945	267	278	4			

- Molecule 2 is a protein called capsid protein VP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	207	Total	C	N	O	S	0	0	0
			1649	1053	286	305	5			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	13	ARG	ALA	conflict	UNP Q9Q2N8
B	14	LEU	ILE	conflict	UNP Q9Q2N8
B	135	ARG	ALA	conflict	UNP Q9Q2N8

- Molecule 3 is a protein called capsid protein VP3.

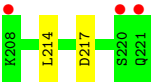
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	221	Total	C	N	O	S	0	0	0
			1707	1087	277	334	9			

- Molecule 4 is a protein called capsid protein VP4.

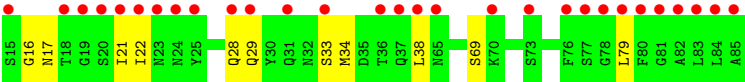
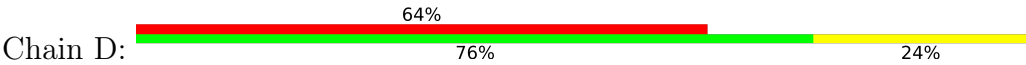
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	45	Total	C	N	O	S	0	0	0
			349	220	56	71	2			

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	84	Total 84	O 84	0	0
5	B	84	Total 84	O 84	0	0
5	C	75	Total 75	O 75	0	0
5	D	15	Total 15	O 15	0	0



● Molecule 4: capsid protein VP4



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	328.20Å 341.80Å 363.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 3.00 25.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	61.7 (25.00-3.00) 98.6 (25.00-3.00)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 3.01Å)	Xtriage
Refinement program	CNS, X-PLOR	Depositor
R, R_{free}	0.167 , (Not available) 0.177 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	22.5	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.16	EDS
Total number of atoms	5457	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.75% 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	0.66	2/1528 (0.1%)	1.05	9/2084 (0.4%)
2	B	0.60	0/1697	1.05	11/2312 (0.5%)
3	C	0.65	1/1755 (0.1%)	1.15	16/2400 (0.7%)
4	D	0.54	0/355	0.90	2/476 (0.4%)
All	All	0.63	3/5335 (0.1%)	1.08	38/7272 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	217	ASP	C-N	8.92	1.41	1.33
1	A	186	CYS	C-N	5.48	1.41	1.33
1	A	187	PRO	N-CD	5.41	1.55	1.47

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	54	PHE	N-CA-C	7.00	121.25	110.20
1	A	191	LEU	N-CA-C	6.98	121.04	109.95
3	C	199	THR	N-CA-C	6.74	120.59	110.14
2	B	107	VAL	N-CA-C	6.65	117.48	108.17
2	B	13	ARG	N-CA-C	-6.53	105.07	113.16
3	C	46	ASP	N-CA-C	-6.46	104.32	111.36
3	C	217	ASP	CA-C-N	-6.38	113.31	121.04
3	C	217	ASP	C-N-CA	-6.38	113.31	121.04
3	C	8	SER	N-CA-C	6.35	119.86	109.76
4	D	33	SER	N-CA-C	-6.29	102.05	110.55
1	A	32	VAL	N-CA-C	6.26	116.43	110.42
1	A	43	GLN	N-CA-C	6.19	120.95	112.04
1	A	100	ASN	N-CA-C	6.16	118.98	109.07
3	C	32	PRO	N-CA-C	-6.12	104.03	110.58
3	C	95	ALA	N-CA-C	6.10	118.73	111.71
2	B	61	PHE	N-CA-C	6.10	119.31	110.28
3	C	175	VAL	N-CA-C	-6.03	100.34	107.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	CYS	CA-C-N	-6.02	113.58	119.78
1	A	186	CYS	C-N-CA	-6.02	113.58	119.78
1	A	27	ARG	N-CA-C	6.01	120.20	112.26
3	C	38	PRO	N-CA-C	5.98	119.59	111.22
2	B	112	VAL	N-CA-C	5.93	117.32	108.54
2	B	126	VAL	N-CA-C	5.92	114.35	107.89
1	A	196	SER	N-CA-C	5.87	117.76	111.36
3	C	87	SER	N-CA-C	5.79	119.87	112.34
3	C	73	LEU	N-CA-C	-5.71	99.22	108.41
3	C	37	TYR	N-CA-C	-5.70	102.60	109.83
2	B	203	ILE	N-CA-C	5.69	116.67	108.48
2	B	113	GLY	N-CA-C	-5.54	100.05	113.18
2	B	194	ALA	N-CA-C	-5.34	103.64	110.53
1	A	6	GLU	N-CA-C	5.33	117.84	111.71
3	C	61	LYS	N-CA-C	-5.33	102.10	110.14
2	B	145	HIS	N-CA-C	5.22	115.95	108.74
3	C	23	ASP	CA-C-N	5.22	125.67	120.14
3	C	23	ASP	C-N-CA	5.22	125.67	120.14
2	B	175	LYS	CA-C-N	5.18	124.36	118.97
2	B	175	LYS	C-N-CA	5.18	124.36	118.97
4	D	69	SER	N-CA-C	-5.16	105.56	111.14

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	1494	0	1493	25	0
2	B	1649	0	1608	18	0
3	C	1707	0	1635	22	0
4	D	349	0	321	13	0
5	A	84	0	0	3	0
5	B	84	0	0	4	0
5	C	75	0	0	1	0
5	D	15	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5457	0	5057	69	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 7.

All (69) 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)
4:D:28:GLN:HG2	4:D:34:MET:HE1	1.48	0.96
2:B:81:LEU:HD11	2:B:85:THR:HG21	1.60	0.82
4:D:28:GLN:HG2	4:D:34:MET:CE	2.09	0.81
1:A:198:GLN:HG2	5:A:384:HOH:O	1.81	0.79
4:D:28:GLN:CG	4:D:34:MET:HE1	2.14	0.77
1:A:37:ASP:O	4:D:17:ASN:ND2	2.16	0.75
1:A:56:THR:O	1:A:67:ARG:NH2	2.20	0.73
1:A:195:VAL:HG22	1:A:201:HIS:HB2	1.74	0.69
2:B:172:LYS:HG3	5:B:329:HOH:O	1.96	0.66
1:A:197:SER:O	1:A:198:GLN:HB2	1.94	0.66
4:D:21:ILE:HG22	4:D:22:ILE:HD12	1.79	0.64
3:C:90:TYR:CE1	3:C:94:ILE:HD11	2.35	0.61
3:C:120:ARG:NH1	3:C:149:ASP:OD2	2.34	0.61
1:A:134:SER:O	3:C:177:GLU:HG3	2.01	0.60
4:D:22:ILE:HD12	4:D:22:ILE:N	2.16	0.59
1:A:37:ASP:OD2	4:D:16:GLY:O	2.20	0.59
1:A:58:GLN:HG2	5:A:332:HOH:O	2.03	0.58
2:B:101:MET:HG2	2:B:210:VAL:HG12	1.86	0.58
2:B:85:THR:HG22	5:B:371:HOH:O	2.05	0.57
1:A:101:THR:HG23	3:C:16:THR:HG21	1.87	0.56
1:A:7:SER:O	1:A:8:ALA:HB3	2.06	0.56
2:B:172:LYS:CG	5:B:329:HOH:O	2.51	0.55
1:A:73:PHE:C	1:A:73:PHE:CD1	2.85	0.54
2:B:13:ARG:O	2:B:29:SER:HB3	2.08	0.54
3:C:72:ARG:HD2	5:C:352:HOH:O	2.10	0.52
3:C:104:THR:CG2	3:C:159:SER:HB3	2.41	0.51
4:D:28:GLN:CD	4:D:34:MET:HE1	2.36	0.49
1:A:57:HIS:O	1:A:63:GLY:HA3	2.11	0.49
2:B:118:GLY:HA3	2:B:188:THR:OG1	2.12	0.48
2:B:116:PHE:CD2	2:B:192:VAL:HG12	2.48	0.48
4:D:22:ILE:N	4:D:22:ILE:CD1	2.76	0.48
3:C:133:PRO:HG3	3:C:184:TRP:CD2	2.49	0.47
2:B:214:LEU:HD23	3:C:127:PRO:HG2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:ASN:O	1:A:104:PRO:C	2.59	0.46
3:C:192:HIS:ND1	3:C:192:HIS:C	2.73	0.46
2:B:42:HIS:CG	4:D:38:LEU:CD1	3.00	0.45
2:B:40:GLU:HG3	2:B:207:HIS:CD2	2.52	0.45
3:C:194:LYS:HD2	3:C:194:LYS:HA	1.80	0.45
1:A:198:GLN:CG	5:A:384:HOH:O	2.53	0.44
1:A:105:THR:O	3:C:15:VAL:HA	2.16	0.44
1:A:79:VAL:HB	1:A:176:LEU:HB2	1.98	0.44
1:A:52:ASP:O	1:A:55:GLN:HB2	2.19	0.43
3:C:76:LYS:CD	3:C:133:PRO:HG2	2.49	0.43
4:D:38:LEU:HD12	4:D:38:LEU:C	2.43	0.43
3:C:181:VAL:O	3:C:181:VAL:HG13	2.19	0.43
1:A:199:ASP:OD1	1:A:200:ARG:N	2.51	0.43
2:B:133:THR:OG1	2:B:136:GLU:HG3	2.18	0.42
3:C:84:LYS:HD3	3:C:84:LYS:C	2.44	0.42
2:B:81:LEU:CD1	2:B:85:THR:HG21	2.40	0.42
2:B:216:SER:HB3	3:C:142:HIS:CD2	2.55	0.42
3:C:98:TYR:HA	3:C:214:LEU:O	2.20	0.42
1:A:39:PHE:CD1	1:A:178:ARG:HB2	2.54	0.42
2:B:39:GLN:HA	5:B:376:HOH:O	2.18	0.42
3:C:107:LEU:HG	3:C:160:ILE:HD11	2.01	0.42
4:D:29:GLN:H	4:D:29:GLN:CD	2.27	0.41
2:B:111:ALA:HB3	2:B:148:ILE:HG21	2.02	0.41
3:C:70:GLU:OE1	3:C:72:ARG:NH1	2.49	0.41
1:A:198:GLN:HE21	1:A:198:GLN:HB3	1.68	0.41
2:B:101:MET:HE2	2:B:101:MET:HB3	1.94	0.41
3:C:120:ARG:NH1	3:C:149:ASP:CG	2.78	0.41
1:A:54:MET:HG3	1:A:158:LEU:HD11	2.03	0.41
2:B:78:LEU:HD23	2:B:78:LEU:C	2.46	0.41
3:C:133:PRO:HA	3:C:134:PRO:HD3	1.94	0.41
1:A:37:ASP:HB3	4:D:17:ASN:HB3	2.03	0.41
3:C:76:LYS:HB3	3:C:76:LYS:HE2	1.65	0.41
1:A:157:GLN:N	1:A:157:GLN:CD	2.79	0.40
3:C:76:LYS:HD2	3:C:133:PRO:HG2	2.03	0.40
1:A:181:ARG:HH11	1:A:181:ARG:HD2	1.78	0.40
1:A:210:LYS:OXT	1:A:210:LYS:HG2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	187/210 (89%)	180 (96%)	7 (4%)	0	100	100
2	B	205/207 (99%)	196 (96%)	8 (4%)	1 (0%)	24	60
3	C	219/221 (99%)	204 (93%)	15 (7%)	0	100	100
4	D	41/45 (91%)	38 (93%)	3 (7%)	0	100	100
All	All	652/683 (96%)	618 (95%)	33 (5%)	1 (0%)	43	76

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	41	ASP

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	161/171 (94%)	155 (96%)	6 (4%)	30	64
2	B	182/182 (100%)	178 (98%)	4 (2%)	45	74
3	C	185/185 (100%)	180 (97%)	5 (3%)	39	71
4	D	37/37 (100%)	36 (97%)	1 (3%)	39	71
All	All	565/575 (98%)	549 (97%)	16 (3%)	38	70

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

Mol	Chain	Res	Type
1	A	25	GLN
1	A	30	THR
1	A	67	ARG
1	A	129	VAL
1	A	196	SER
1	A	198	GLN
2	B	63	LYS
2	B	108	GLU
2	B	140	LEU
2	B	191	THR
3	C	76	LYS
3	C	114	SER
3	C	147	GLU
3	C	179	THR
3	C	194	LYS
4	D	79	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	59	HIS
1	A	168	GLN
1	A	198	GLN
2	B	77	HIS
2	B	103	ASN
2	B	207	HIS
3	C	71	GLN
3	C	142	HIS
3	C	153	ASN
3	C	180	ASN
3	C	197	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
4	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	38:LEU	C	65:ASN	N	14.86

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	191/210 (90%)	1.57	41 (21%)	2 1	4, 11, 67, 101	0
2	B	207/207 (100%)	1.40	33 (15%)	5 3	5, 11, 39, 95	0
3	C	221/221 (100%)	1.49	52 (23%)	2 1	4, 11, 53, 96	0
4	D	45/45 (100%)	3.17	29 (64%)	0 0	10, 36, 75, 85	0
All	All	664/683 (97%)	1.60	155 (23%)	2 1	4, 12, 63, 101	0

All (155) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	218	GLU	11.9
1	A	198	GLN	10.4
1	A	156	ALA	9.7
4	D	23	ASN	9.3
1	A	210	LYS	8.9
2	B	12	ASP	8.6
2	B	39	GLN	8.2
2	B	42	HIS	8.1
4	D	81	GLY	7.7
1	A	197	SER	7.6
4	D	15	SER	7.5
1	A	135	LYS	7.5
3	C	70	GLU	7.4
3	C	177	GLU	7.4
1	A	199	ASP	7.1
3	C	176	ALA	7.1
4	D	85	ALA	7.0
1	A	157	GLN	7.0
2	B	41	ASP	6.9
4	D	65	ASN	6.8
2	B	217	LYS	6.7

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Mol	Chain	Res	Type	RSRZ
1	A	136	TYR	6.5
2	B	44	SER	6.4
3	C	131	GLU	6.2
1	A	209	ALA	5.9
3	C	181	VAL	5.7
4	D	22	ILE	5.5
3	C	178	THR	5.5
1	A	44	ASN	5.4
3	C	67	ARG	5.3
3	C	59	GLU	5.3
3	C	58	ASP	5.2
3	C	175	VAL	5.1
3	C	180	ASN	5.1
4	D	78	GLY	4.9
1	A	43	GLN	4.8
2	B	40	GLU	4.7
4	D	20	SER	4.5
3	C	68	THR	4.5
4	D	80	PHE	4.5
4	D	82	ALA	4.3
2	B	38	THR	4.3
4	D	37	GLN	4.3
4	D	21	ILE	4.3
2	B	29	SER	4.3
4	D	24	ASN	4.3
2	B	13	ARG	4.0
2	B	28	SER	4.0
4	D	19	GLY	3.9
3	C	174	ASP	3.9
3	C	220	SER	3.8
2	B	191	THR	3.8
4	D	18	THR	3.8
2	B	14	LEU	3.8
4	D	79	LEU	3.7
4	D	33	SER	3.6
2	B	130	LYS	3.6
3	C	130	VAL	3.5
2	B	93	HIS	3.4
4	D	38	LEU	3.4
2	B	190	ASN	3.4
1	A	25	GLN	3.4
3	C	208	LYS	3.4

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Mol	Chain	Res	Type	RSRZ
3	C	221	GLN	3.4
3	C	132	THR	3.3
3	C	21	THR	3.3
3	C	133	PRO	3.2
1	A	95	GLU	3.2
1	A	194	GLU	3.2
2	B	216	SER	3.2
3	C	197	GLN	3.2
4	D	28	GLN	3.1
3	C	114	SER	3.1
3	C	71	GLN	3.1
2	B	43	VAL	3.1
3	C	196	GLU	3.1
3	C	153	ASN	3.1
3	C	179	THR	3.0
1	A	202	LYS	3.0
1	A	204	LYS	3.0
4	D	36	THR	3.0
1	A	196	SER	2.9
2	B	119	GLY	2.9
4	D	76	PHE	2.9
2	B	86	ASP	2.9
1	A	206	ILE	2.9
1	A	13	THR	2.8
1	A	134	SER	2.8
1	A	46	ASN	2.8
4	D	70	LYS	2.8
1	A	28	GLN	2.8
3	C	6	ALA	2.8
1	A	207	ALA	2.7
1	A	1	THR	2.7
1	A	5	GLY	2.6
2	B	131	GLU	2.6
3	C	103	GLY	2.6
3	C	56	CYS	2.6
1	A	96	ALA	2.6
1	A	110	ALA	2.6
3	C	30	TYR	2.6
3	C	161	PRO	2.5
4	D	29	GLN	2.5
2	B	72	ASP	2.5
3	C	36	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
3	C	135	ASP	2.5
3	C	195	ALA	2.5
3	C	3	VAL	2.5
1	A	78	ILE	2.5
4	D	25	TYR	2.4
3	C	102	SER	2.4
1	A	24	VAL	2.4
3	C	101	TYR	2.4
4	D	77	SER	2.4
2	B	189	THR	2.4
1	A	205	ILE	2.4
3	C	189	GLN	2.4
2	B	192	VAL	2.4
1	A	85	ASN	2.3
2	B	36	TYR	2.3
1	A	159	PRO	2.3
3	C	35	THR	2.3
3	C	84	LYS	2.3
1	A	105	THR	2.3
2	B	71	PRO	2.3
4	D	31	GLN	2.3
3	C	173	SER	2.3
1	A	133	THR	2.2
3	C	44	LEU	2.2
2	B	46	PRO	2.1
3	C	72	ARG	2.1
3	C	129	GLY	2.1
3	C	18	ASP	2.1
1	A	22	THR	2.1
2	B	111	ALA	2.1
3	C	60	GLY	2.1
3	C	183	GLY	2.1
1	A	201	HIS	2.1
3	C	152	LEU	2.1
4	D	83	LEU	2.1
1	A	48	THR	2.1
3	C	17	THR	2.1
2	B	92	GLY	2.1
3	C	32	PRO	2.1
3	C	194	LYS	2.1
2	B	96	ASP	2.1
1	A	101	THR	2.1

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Mol	Chain	Res	Type	RSRZ
4	D	73	SER	2.0
1	A	3	THR	2.0
3	C	138	GLU	2.0
2	B	32	VAL	2.0
4	D	84	LEU	2.0
1	A	7	SER	2.0
2	B	79	GLU	2.0
1	A	176	LEU	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.