



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:59 PM UTC

PDB ID : 4NYL / pdb_00004nyl
Title : Crystal structure of adalimumab FAB fragment
Authors : Fan, L.J.; Lv, L.L.; Zhang, Q.J.; Chen, C.L.
Deposited on : 2013-12-10
Resolution : 2.80 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
Xtrriage (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

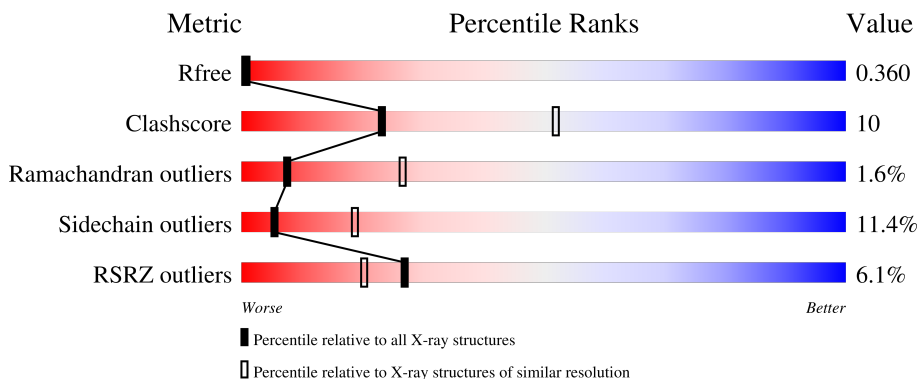
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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.



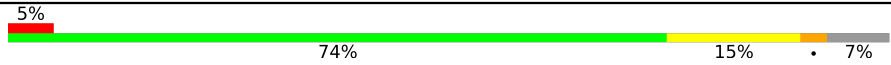

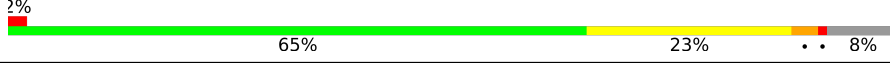
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	B	214	
1	D	214	
1	F	214	
1	L	214	
2	A	230	

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Mol	Chain	Length	Quality of chain
2	C	230	
2	E	230	
2	H	230	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10495 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 Adalimumab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	L	211	Total 1546	C 966	N 262	O 313	S 5	0	0	0
1	B	117	Total 829	C 518	N 144	O 164	S 3	0	0	0
1	D	211	Total 1557	C 972	N 261	O 319	S 5	0	0	0
1	F	124	Total 876	C 543	N 149	O 181	S 3	0	0	0

- Molecule 2 is a protein called Adalimumab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	212	Total 1564	C 991	N 263	O 304	S 6	0	0	0
2	A	199	Total 1405	C 879	N 239	O 281	S 6	0	0	0
2	C	213	Total 1561	C 987	N 263	O 305	S 6	0	0	0
2	E	162	Total 1155	C 723	N 200	O 228	S 4	0	1	0

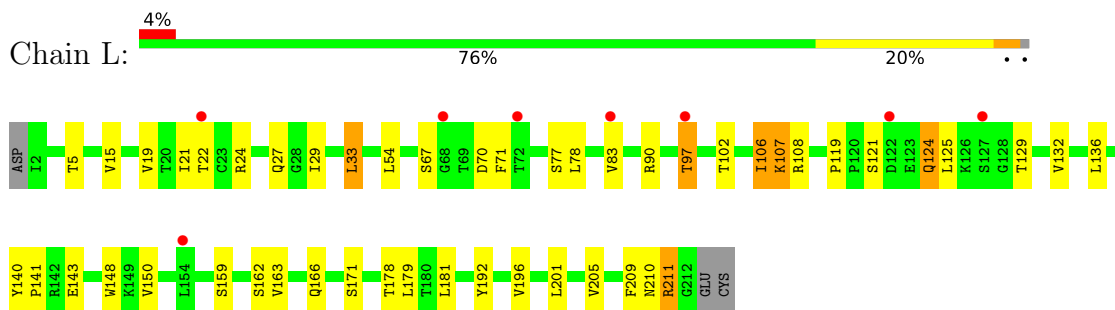
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total O 1 1	0	0
3	C	1	Total O 1 1	0	0

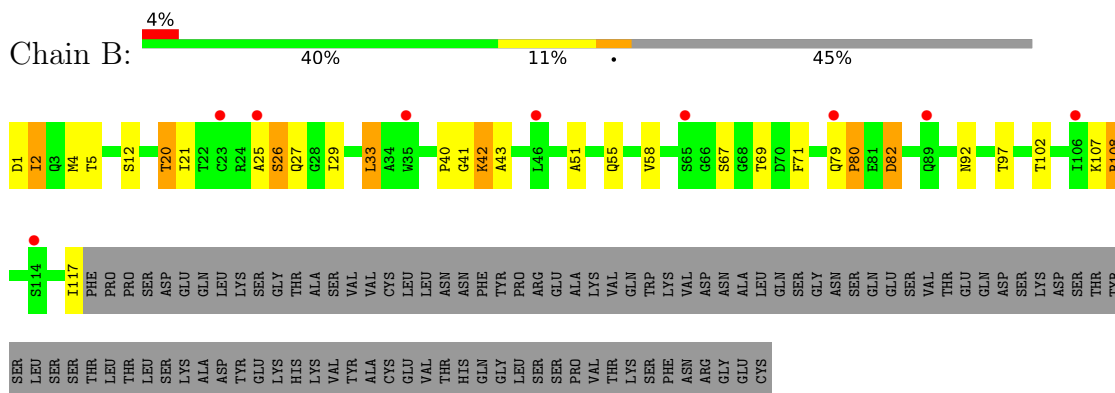
3 Residue-property plots [i](#)

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.

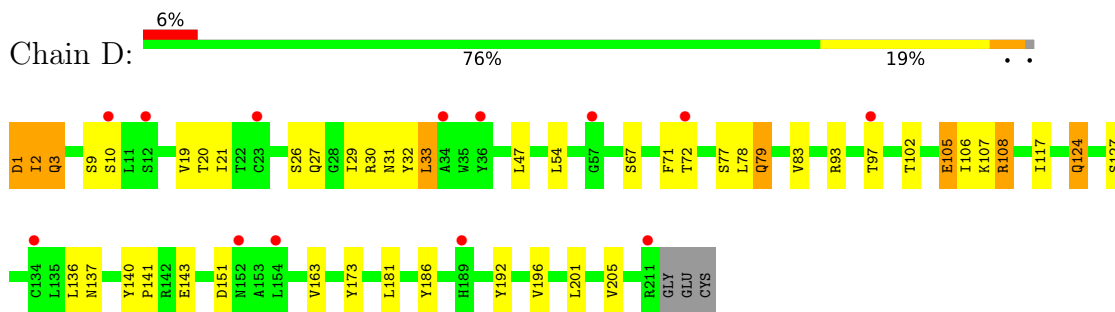
- Molecule 1: Adalimumab Light Chain



- Molecule 1: Adalimumab Light Chain

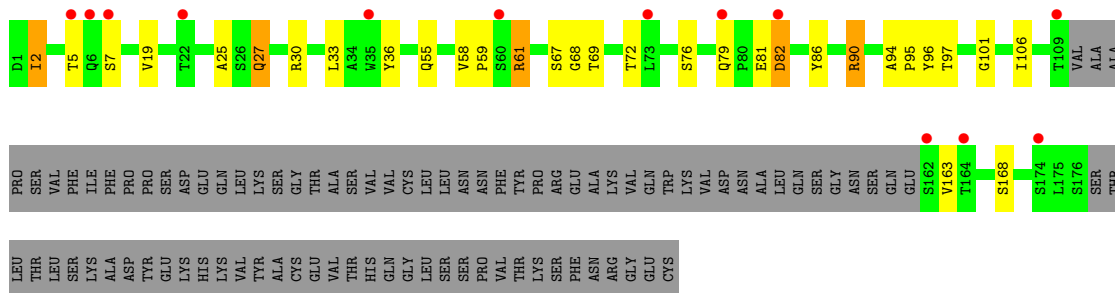


- Molecule 1: Adalimumab Light Chain

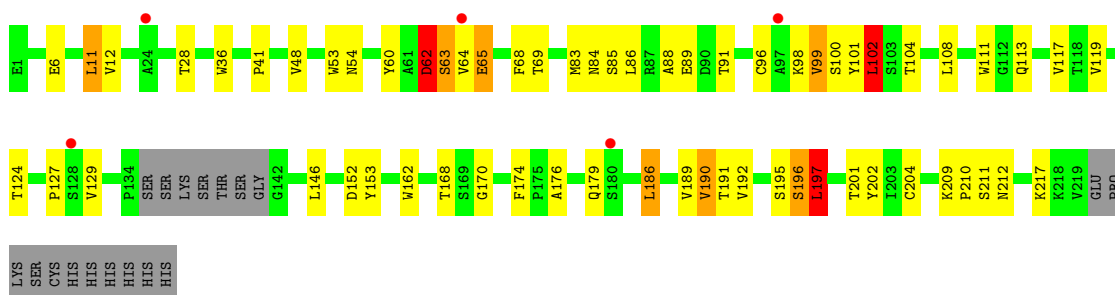


- Molecule 1: Adalimumab Light Chain

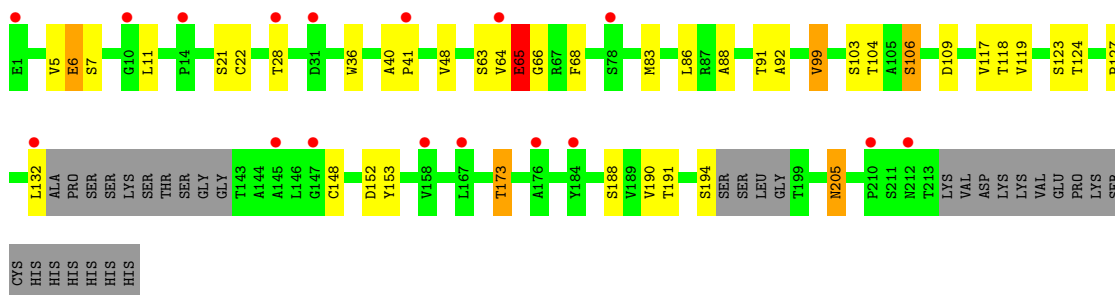




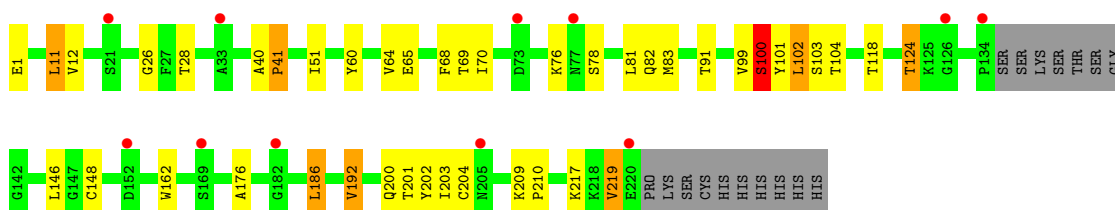
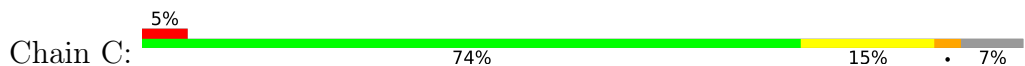
● Molecule 2: Adalimumab Heavy Chain



● Molecule 2: Adalimumab Heavy Chain

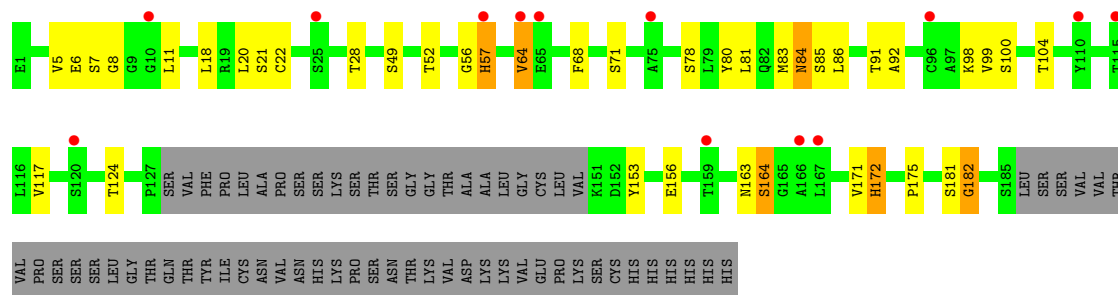


● Molecule 2: Adalimumab Heavy Chain



● Molecule 2: Adalimumab Heavy Chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	131.42Å 47.05Å 168.51Å 90.00° 101.33° 90.00°	Depositor
Resolution (Å)	42.03 – 2.80 42.03 – 2.80	Depositor EDS
% Data completeness (in resolution range)	87.1 (42.03-2.80) 87.0 (42.03-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.81Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.275 , 0.358 0.273 , 0.360	Depositor DCC
R_{free} test set	2249 reflections (4.42%)	wwPDB-VP
Wilson B-factor (Å ²)	58.0	Xtrriage
Anisotropy	0.157	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	10495	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.60% 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

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	B	0.79	0/848	0.91	0/1161
1	D	0.78	0/1591	0.95	1/2177 (0.0%)
1	F	0.71	0/892	0.95	3/1221 (0.2%)
1	L	0.73	0/1580	0.94	1/2162 (0.0%)
2	A	0.69	0/1438	0.90	2/1974 (0.1%)
2	C	0.80	1/1599 (0.1%)	0.97	2/2189 (0.1%)
2	E	0.72	0/1186	0.91	0/1623
2	H	0.86	1/1602 (0.1%)	0.99	2/2189 (0.1%)
All	All	0.76	2/10736 (0.0%)	0.94	11/14696 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	41	PRO	CA-C	-7.25	1.44	1.52
2	C	82	GLN	N-CA	-5.82	1.38	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	103	SER	N-CA-C	-8.59	95.49	108.46
1	L	83	VAL	CB-CA-C	-6.54	104.17	111.23
2	A	103	SER	N-CA-C	6.21	119.05	110.35
2	A	173	THR	N-CA-C	-6.16	100.54	109.59
2	H	62	ASP	N-CA-C	-5.91	101.43	110.30
1	F	68	GLY	N-CA-C	-5.79	106.34	111.67
1	F	101	GLY	N-CA-C	5.67	119.34	112.48
1	F	79	GLN	N-CA-C	5.25	112.75	108.07
1	D	137	ASN	N-CA-C	5.19	117.43	109.07
2	H	41	PRO	CA-C-O	-5.06	115.05	120.97
2	C	100	SER	N-CA-C	5.02	116.57	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	B	829	0	747	21	0
1	D	1557	0	1433	26	0
1	F	876	0	773	16	0
1	L	1546	0	1422	30	0
2	A	1405	0	1243	21	0
2	C	1561	0	1472	25	0
2	E	1155	0	1019	22	0
2	H	1564	0	1489	46	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
All	All	10495	0	9598	207	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 10.

All (207) 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:H:196:SER:OG	2:H:197:LEU:HD23	1.51	1.10
2:A:40:ALA:HB1	2:A:41:PRO:HD2	1.42	1.01
1:L:107:LYS:CB	1:L:108:ARG:HA	1.92	1.00
2:H:88:ALA:O	2:H:91:THR:HG23	1.66	0.94
2:H:197:LEU:HD23	2:H:197:LEU:N	1.87	0.88
2:H:64:VAL:HG11	2:H:68:PHE:HB2	1.55	0.88
1:F:81:GLU:C	1:F:82:ASP:OD1	2.18	0.87
1:L:107:LYS:CB	1:L:108:ARG:CA	2.54	0.84
1:B:20:THR:CG2	1:B:20:THR:O	2.25	0.84
2:A:64:VAL:O	2:A:65:GLU:C	2.21	0.83
2:H:88:ALA:O	2:H:91:THR:CG2	2.28	0.81
1:D:1:ASP:O	1:D:2:ILE:C	2.27	0.76
2:C:81:LEU:HD23	2:C:83:MET:HE3	1.66	0.76
1:B:20:THR:O	1:B:20:THR:HG22	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:64:VAL:HG11	2:A:68:PHE:HB2	1.70	0.73
2:H:196:SER:O	2:H:197:LEU:HB2	1.89	0.72
1:D:107:LYS:O	1:D:140:TYR:CZ	2.43	0.72
1:L:106:ILE:CG2	1:L:171:SER:HB3	2.19	0.72
2:C:91:THR:HG23	2:C:118:THR:HA	1.72	0.71
2:H:197:LEU:N	2:H:197:LEU:CD2	2.54	0.69
2:H:209:LYS:O	2:H:210:PRO:C	2.35	0.69
2:H:209:LYS:O	2:H:212:ASN:N	2.26	0.69
1:F:81:GLU:O	1:F:82:ASP:OD1	2.10	0.69
1:L:210:ASN:O	1:L:211:ARG:CG	2.41	0.68
2:H:91:THR:HG22	2:H:119:VAL:H	1.57	0.68
1:L:106:ILE:HG21	1:L:171:SER:HB3	1.76	0.67
2:H:89:GLU:N	2:H:89:GLU:OE1	2.26	0.67
2:A:64:VAL:O	2:A:66:GLY:N	2.27	0.67
2:A:91:THR:HG23	2:A:118:THR:HA	1.76	0.66
2:E:64:VAL:HG21	2:E:68:PHE:HB2	1.76	0.66
2:H:11:LEU:HD12	2:H:124:THR:HG22	1.78	0.66
2:C:102:LEU:HD12	2:C:102:LEU:O	1.96	0.65
1:L:210:ASN:O	1:L:211:ARG:CB	2.43	0.65
2:H:152:ASP:OD1	2:H:179:GLN:NE2	2.32	0.62
2:E:83:MET:HE1	2:E:117:VAL:HG11	1.82	0.62
1:D:30:ARG:HG2	1:D:31:ASN:N	2.16	0.61
1:L:106:ILE:HG22	1:L:166:GLN:NE2	2.17	0.60
1:D:1:ASP:O	1:D:3:GLN:N	2.35	0.60
2:H:102:LEU:C	2:H:102:LEU:HD13	2.27	0.59
2:H:196:SER:OG	2:H:197:LEU:CD2	2.39	0.59
1:B:107:LYS:O	1:B:108:ARG:HB3	2.02	0.59
2:H:195:SER:C	2:H:196:SER:O	2.44	0.59
2:H:209:LYS:O	2:H:211:SER:N	2.36	0.58
1:L:106:ILE:HG21	1:L:171:SER:CB	2.34	0.58
2:A:40:ALA:HB1	2:A:41:PRO:CD	2.28	0.58
1:D:136:LEU:HD21	1:D:196:VAL:HG13	1.85	0.58
2:C:1:GLU:O	2:C:26:GLY:HA3	2.03	0.58
1:F:94:ALA:HA	1:F:95:PRO:C	2.29	0.58
2:A:5:VAL:O	2:A:22:CYS:HA	2.04	0.57
1:L:136:LEU:HD21	1:L:196:VAL:HG13	1.86	0.56
2:C:64:VAL:CG1	2:C:68:PHE:CG	2.89	0.56
2:H:36:TRP:O	2:H:48:VAL:HB	2.06	0.56
2:H:127:PRO:HB3	2:H:153:TYR:HB3	1.88	0.56
2:C:100:SER:HB3	2:C:101:TYR:CD2	2.41	0.55
2:C:64:VAL:HG11	2:C:68:PHE:CD2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:201:LEU:HD13	1:L:205:VAL:HG23	1.89	0.55
2:H:64:VAL:HG12	2:H:65:GLU:N	2.20	0.55
2:E:64:VAL:HG21	2:E:68:PHE:CB	2.37	0.54
2:H:162:TRP:CH2	2:H:204:CYS:HB3	2.43	0.54
2:C:219:VAL:O	2:C:219:VAL:CG2	2.54	0.54
2:C:40:ALA:O	2:C:41:PRO:C	2.50	0.54
2:A:64:VAL:HG11	2:A:68:PHE:CB	2.38	0.54
1:B:20:THR:O	1:B:20:THR:HG23	2.04	0.54
1:L:148:TRP:CD1	1:L:159:SER:HG	2.26	0.54
2:C:203:ILE:HG23	2:C:217:LYS:O	2.08	0.53
2:H:196:SER:OG	2:H:197:LEU:N	2.35	0.53
2:E:5:VAL:O	2:E:22:CYS:HA	2.08	0.53
1:B:21:ILE:HD13	1:B:102:THR:HG21	1.91	0.53
1:F:2:ILE:HG21	1:F:90:ARG:CZ	2.39	0.53
2:C:11:LEU:HD13	2:C:124:THR:HG22	1.90	0.52
2:H:62:ASP:O	2:H:63:SER:HB2	2.09	0.52
2:H:192:VAL:HG11	2:H:202:TYR:CE2	2.44	0.52
1:B:40:PRO:CG	1:B:40:PRO:O	2.56	0.52
2:C:162:TRP:CH2	2:C:204:CYS:HB3	2.45	0.52
1:D:77:SER:O	1:D:79:GLN:OE1	2.27	0.52
1:L:24:ARG:CD	1:L:70:ASP:OD1	2.58	0.51
2:E:56:GLY:O	2:E:57[A]:HIS:CG	2.63	0.51
1:B:40:PRO:O	1:B:40:PRO:HG2	2.10	0.51
1:L:27:GLN:O	1:L:29:ILE:HG23	2.10	0.51
2:E:18:LEU:CD2	2:E:20:LEU:HD23	2.41	0.51
2:A:83:MET:HE2	2:A:86:LEU:HD21	1.92	0.50
2:A:83:MET:HB3	2:A:86:LEU:HD21	1.93	0.50
2:A:173:THR:HA	2:A:188:SER:HA	1.92	0.50
2:C:1:GLU:O	2:C:26:GLY:CA	2.60	0.50
1:L:90:ARG:CD	1:L:97:THR:HG23	2.42	0.50
1:F:82:ASP:OD1	1:F:82:ASP:N	2.45	0.50
1:L:15:VAL:HA	1:L:78:LEU:O	2.12	0.50
1:B:102:THR:HG23	1:B:102:THR:O	2.11	0.50
2:C:146:LEU:C	2:C:146:LEU:HD12	2.37	0.50
1:B:92:ASN:N	1:B:92:ASN:OD1	2.45	0.49
2:H:6:GLU:HG3	2:H:96:CYS:SG	2.51	0.49
1:D:107:LYS:O	1:D:140:TYR:CE2	2.65	0.49
1:D:136:LEU:HD21	1:D:196:VAL:CG1	2.43	0.49
2:E:6:GLU:HA	2:E:21:SER:O	2.12	0.49
1:D:1:ASP:O	1:D:3:GLN:CB	2.61	0.49
2:C:64:VAL:HG12	2:C:68:PHE:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:MET:HE1	1:B:25:ALA:HB2	1.94	0.49
2:E:83:MET:HB3	2:E:86:LEU:HD21	1.94	0.49
1:D:83:VAL:O	1:D:83:VAL:HG13	2.13	0.48
2:H:195:SER:O	2:H:196:SER:O	2.32	0.48
2:E:163:ASN:O	2:E:164:SER:CB	2.61	0.48
2:H:108:LEU:HD12	2:H:111:TRP:CZ2	2.49	0.48
1:D:27:GLN:O	1:D:29:ILE:HG23	2.12	0.48
1:D:30:ARG:HG2	1:D:31:ASN:H	1.75	0.48
2:E:8:GLY:O	2:E:18:LEU:HD21	2.12	0.48
2:H:62:ASP:O	2:H:63:SER:CB	2.61	0.48
2:A:127:PRO:HB3	2:A:153:TYR:HB3	1.94	0.48
2:C:219:VAL:O	2:C:219:VAL:HG23	2.12	0.48
1:B:27:GLN:O	1:B:29:ILE:HG23	2.12	0.48
2:E:52:THR:CG2	2:E:57[B]:HIS:HB2	2.44	0.48
2:C:64:VAL:CG1	2:C:68:PHE:HB2	2.43	0.48
1:B:107:LYS:O	1:B:108:ARG:CB	2.62	0.48
2:C:64:VAL:O	2:C:65:GLU:C	2.56	0.48
1:L:210:ASN:O	1:L:211:ARG:HB2	2.13	0.48
2:H:174:PHE:O	2:H:186:LEU:HD22	2.14	0.48
1:L:19:VAL:HG21	1:L:78:LEU:HD22	1.96	0.47
2:C:209:LYS:O	2:C:210:PRO:C	2.57	0.47
2:E:52:THR:HG23	2:E:57[B]:HIS:HB2	1.95	0.47
2:E:64:VAL:O	2:E:64:VAL:HG22	2.13	0.47
1:L:148:TRP:CE3	1:L:179:LEU:HD22	2.49	0.47
2:H:196:SER:O	2:H:197:LEU:CB	2.51	0.47
1:B:40:PRO:HA	1:B:41:GLY:HA2	1.58	0.47
1:D:105:GLU:OE1	1:D:173:TYR:OH	2.27	0.47
2:E:18:LEU:HD21	2:E:20:LEU:HD23	1.97	0.47
2:E:84:ASN:O	2:E:85:SER:C	2.58	0.47
1:L:107:LYS:CB	1:L:108:ARG:CB	2.92	0.47
1:D:201:LEU:HD13	1:D:205:VAL:HG23	1.95	0.47
1:L:90:ARG:HD3	1:L:97:THR:HG23	1.96	0.47
2:H:91:THR:HG22	2:H:119:VAL:N	2.29	0.46
2:H:129:VAL:O	2:H:217:LYS:HE3	2.15	0.46
1:D:1:ASP:OD1	1:D:26:SER:OG	2.33	0.46
1:B:33:LEU:HD22	1:B:71:PHE:CG	2.51	0.46
1:D:124:GLN:O	1:D:127:SER:CB	2.64	0.46
1:L:107:LYS:CB	1:L:108:ARG:HB3	2.45	0.46
2:C:192:VAL:HG11	2:C:202:TYR:CE1	2.50	0.46
2:H:196:SER:O	2:H:197:LEU:O	2.33	0.46
2:H:53:TRP:CZ2	2:H:54:ASN:ND2	2.83	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:ASP:O	1:B:2:ILE:HG22	2.16	0.46
1:D:141:PRO:HB2	1:D:143:GLU:OE1	2.16	0.45
2:E:20:LEU:HD12	2:E:81:LEU:HD23	1.98	0.45
1:L:121:SER:O	1:L:125:LEU:HD12	2.17	0.45
2:A:36:TRP:O	2:A:48:VAL:HB	2.16	0.45
2:C:176:ALA:HA	2:C:186:LEU:HD23	1.98	0.44
1:F:94:ALA:HB2	1:F:96:TYR:CE1	2.52	0.44
2:E:64:VAL:HG21	2:E:68:PHE:CG	2.52	0.44
1:B:55:GLN:O	1:B:58:VAL:HB	2.17	0.44
2:A:6:GLU:HA	2:A:21:SER:O	2.16	0.44
1:D:29:ILE:O	1:D:32:TYR:HD1	2.00	0.44
2:H:170:GLY:O	2:H:190:VAL:HA	2.17	0.44
1:B:25:ALA:O	1:B:69:THR:HB	2.18	0.44
2:A:88:ALA:HA	2:A:119:VAL:HB	1.99	0.44
2:E:153:TYR:CD1	2:E:153:TYR:C	2.96	0.44
1:F:25:ALA:O	1:F:69:THR:HB	2.17	0.44
1:B:33:LEU:HD22	1:B:71:PHE:CD1	2.52	0.44
1:B:42:LYS:HG3	1:B:43:ALA:H	1.83	0.44
1:F:2:ILE:HD13	1:F:2:ILE:C	2.42	0.44
2:H:83:MET:HE1	2:H:117:VAL:HG11	1.99	0.44
2:H:101:TYR:O	2:H:102:LEU:C	2.61	0.44
1:D:1:ASP:O	1:D:3:GLN:HB2	2.18	0.44
1:L:140:TYR:CG	1:L:141:PRO:HA	2.53	0.43
2:A:83:MET:HE1	2:A:117:VAL:HG11	1.99	0.43
2:H:176:ALA:HA	2:H:186:LEU:HD23	2.00	0.43
1:D:107:LYS:O	1:D:108:ARG:O	2.36	0.43
1:D:33:LEU:HD22	1:D:71:PHE:CD2	2.53	0.43
1:L:119:PRO:HB3	1:L:209:PHE:CE2	2.53	0.43
1:F:90:ARG:NH1	1:F:95:PRO:O	2.52	0.43
1:L:150:VAL:HG13	1:L:192:TYR:CE2	2.54	0.42
2:C:101:TYR:O	2:C:102:LEU:HG	2.18	0.42
1:F:2:ILE:HG23	1:F:2:ILE:O	2.18	0.42
2:H:83:MET:HB3	2:H:86:LEU:HD21	2.01	0.42
1:D:30:ARG:CG	1:D:31:ASN:N	2.75	0.42
1:D:186:TYR:HA	1:D:192:TYR:OH	2.19	0.42
1:F:36:TYR:O	1:F:86:TYR:HA	2.19	0.42
1:F:2:ILE:HG13	1:F:27:GLN:HG3	2.01	0.42
1:F:59:PRO:HB2	1:F:61:ARG:HD2	2.01	0.42
2:H:196:SER:HG	2:H:197:LEU:HD23	1.76	0.42
2:C:60:TYR:OH	2:C:69:THR:HA	2.19	0.42
2:E:91:THR:O	2:E:92:ALA:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:159:SER:HA	1:L:178:THR:O	2.20	0.42
1:L:132:VAL:HB	1:L:179:LEU:HB3	2.00	0.42
1:D:21:ILE:HG12	1:D:102:THR:HG21	2.02	0.42
1:F:2:ILE:CG1	1:F:27:GLN:HG3	2.50	0.42
2:A:40:ALA:CB	2:A:41:PRO:CD	2.95	0.41
1:L:33:LEU:HD22	1:L:71:PHE:CG	2.55	0.41
2:H:98:LYS:HG2	2:H:99:VAL:N	2.33	0.41
2:H:68:PHE:CE2	2:H:83:MET:HG2	2.56	0.41
2:E:71:SER:O	2:E:80:TYR:N	2.49	0.41
2:H:217:LYS:HD3	2:H:217:LYS:O	2.21	0.41
1:B:2:ILE:HA	1:B:26:SER:OG	2.19	0.41
1:D:3:GLN:HB3	1:D:26:SER:HB3	2.02	0.41
2:E:181:SER:N	2:E:182:GLY:HA2	2.36	0.41
1:L:136:LEU:HD21	1:L:196:VAL:CG1	2.50	0.41
2:H:60:TYR:OH	2:H:69:THR:HA	2.21	0.41
2:A:91:THR:O	2:A:92:ALA:HB2	2.21	0.41
2:A:205:ASN:OD1	2:A:205:ASN:N	2.54	0.41
1:F:55:GLN:O	1:F:58:VAL:HB	2.21	0.41
1:F:61:ARG:HB3	1:F:76:SER:OG	2.21	0.41
2:C:64:VAL:CG1	2:C:68:PHE:CD2	3.04	0.41
2:E:171:VAL:O	2:E:172:HIS:ND1	2.54	0.40
1:L:21:ILE:HG12	1:L:102:THR:HG21	2.03	0.40
2:C:51:ILE:HG22	2:C:70:ILE:HD11	2.03	0.40
2:A:99:VAL:HG13	2:A:106:SER:HB3	2.03	0.40
2:H:84:ASN:O	2:H:85:SER:C	2.65	0.40
2:H:102:LEU:HD22	2:H:102:LEU:HA	1.85	0.40
1:B:79:GLN:CB	1:B:80:PRO:CD	2.99	0.40
2:A:190:VAL:HG12	2:A:191:THR:N	2.37	0.40
1:D:19:VAL:HG21	1:D:78:LEU:HD22	2.04	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	B	115/214 (54%)	97 (84%)	13 (11%)	5 (4%)	2	7
1	D	209/214 (98%)	183 (88%)	24 (12%)	2 (1%)	12	38
1	F	120/214 (56%)	108 (90%)	12 (10%)	0	100	100
1	L	209/214 (98%)	195 (93%)	10 (5%)	4 (2%)	6	22
2	A	193/230 (84%)	168 (87%)	22 (11%)	3 (2%)	7	27
2	C	209/230 (91%)	196 (94%)	12 (6%)	1 (0%)	24	55
2	E	159/230 (69%)	139 (87%)	17 (11%)	3 (2%)	6	22
2	H	208/230 (90%)	189 (91%)	14 (7%)	5 (2%)	4	17
All	All	1422/1776 (80%)	1275 (90%)	124 (9%)	23 (2%)	7	27

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	107	LYS
1	L	124	GLN
1	L	211	ARG
2	H	102	LEU
2	H	196	SER
1	B	82	ASP
2	A	65	GLU
1	D	108	ARG
1	L	77	SER
2	H	63	SER
2	H	65	GLU
1	B	2	ILE
1	B	108	ARG
2	A	152	ASP
2	E	164	SER
1	B	80	PRO
1	D	2	ILE
2	E	175	PRO
2	E	182	GLY
1	B	51	ALA
2	H	197	LEU
2	A	63	SER
2	C	41	PRO

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	B	80/186 (43%)	70 (88%)	10 (12%)	4	15
1	D	165/186 (89%)	145 (88%)	20 (12%)	5	16
1	F	85/186 (46%)	69 (81%)	16 (19%)	1	5
1	L	162/186 (87%)	149 (92%)	13 (8%)	11	34
2	A	138/194 (71%)	123 (89%)	15 (11%)	6	21
2	C	165/194 (85%)	149 (90%)	16 (10%)	8	25
2	E	110/194 (57%)	94 (86%)	16 (14%)	3	11
2	H	166/194 (86%)	149 (90%)	17 (10%)	7	23
All	All	1071/1520 (70%)	948 (88%)	123 (12%)	5	18

All (123) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	5	THR
1	L	22	THR
1	L	33	LEU
1	L	54	LEU
1	L	67	SER
1	L	97	THR
1	L	106	ILE
1	L	124	GLN
1	L	129	THR
1	L	143	GLU
1	L	162	SER
1	L	163	VAL
1	L	181	LEU
2	H	11	LEU
2	H	12	VAL
2	H	28	THR
2	H	62	ASP
2	H	99	VAL
2	H	100	SER

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Mol	Chain	Res	Type
2	H	102	LEU
2	H	104	THR
2	H	113	GLN
2	H	146	LEU
2	H	168	THR
2	H	186	LEU
2	H	189	VAL
2	H	190	VAL
2	H	191	THR
2	H	197	LEU
2	H	201	THR
1	B	5	THR
1	B	12	SER
1	B	20	THR
1	B	26	SER
1	B	33	LEU
1	B	42	LYS
1	B	67	SER
1	B	82	ASP
1	B	97	THR
1	B	117	ILE
2	A	6	GLU
2	A	7	SER
2	A	11	LEU
2	A	28	THR
2	A	65	GLU
2	A	99	VAL
2	A	104	THR
2	A	106	SER
2	A	109	ASP
2	A	123	SER
2	A	124	THR
2	A	132	LEU
2	A	148	CYS
2	A	194	SER
2	A	205	ASN
1	D	1	ASP
1	D	3	GLN
1	D	9	SER
1	D	10	SER
1	D	20	THR
1	D	33	LEU

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Mol	Chain	Res	Type
1	D	47	LEU
1	D	54	LEU
1	D	67	SER
1	D	72	THR
1	D	79	GLN
1	D	93	ARG
1	D	97	THR
1	D	105	GLU
1	D	106	ILE
1	D	117	ILE
1	D	124	GLN
1	D	151	ASP
1	D	163	VAL
1	D	181	LEU
2	C	11	LEU
2	C	12	VAL
2	C	28	THR
2	C	76	LYS
2	C	78	SER
2	C	99	VAL
2	C	100	SER
2	C	102	LEU
2	C	104	THR
2	C	124	THR
2	C	148	CYS
2	C	186	LEU
2	C	192	VAL
2	C	200	GLN
2	C	201	THR
2	C	219	VAL
1	F	2	ILE
1	F	5	THR
1	F	7	SER
1	F	19	VAL
1	F	27	GLN
1	F	30	ARG
1	F	33	LEU
1	F	61	ARG
1	F	67	SER
1	F	72	THR
1	F	82	ASP
1	F	90	ARG

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Mol	Chain	Res	Type
1	F	97	THR
1	F	106	ILE
1	F	163	VAL
1	F	168	SER
2	E	7	SER
2	E	11	LEU
2	E	28	THR
2	E	49	SER
2	E	57[A]	HIS
2	E	57[B]	HIS
2	E	64	VAL
2	E	78	SER
2	E	84	ASN
2	E	98	LYS
2	E	99	VAL
2	E	100	SER
2	E	104	THR
2	E	124	THR
2	E	156	GLU
2	E	172	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	79	GLN
1	L	137	ASN
1	L	166	GLN
2	H	113	GLN
2	H	163	ASN
2	H	200	GLN
2	H	207	ASN
1	B	3	GLN
2	A	57	HIS
2	A	84	ASN
1	D	3	GLN
1	D	38	GLN
2	C	39	GLN
2	C	200	GLN
1	F	3	GLN
1	F	38	GLN
2	E	82	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](#)

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 [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	B	117/214 (54%)	0.80	9 (7%) 19 14	49, 83, 104, 140	0
1	D	211/214 (98%)	0.42	13 (6%) 26 20	44, 63, 82, 93	0
1	F	124/214 (57%)	0.68	13 (10%) 11 8	54, 83, 103, 121	0
1	L	211/214 (98%)	0.43	8 (3%) 44 35	46, 64, 85, 99	0
2	A	199/230 (86%)	0.75	17 (8%) 16 12	47, 75, 116, 126	0
2	C	213/230 (92%)	0.50	11 (5%) 33 25	37, 64, 79, 98	0
2	E	162/230 (70%)	0.65	13 (8%) 18 13	27, 69, 122, 140	1 (0%)
2	H	212/230 (92%)	0.45	5 (2%) 59 49	45, 61, 76, 101	0
All	All	1449/1776 (81%)	0.56	89 (6%) 27 20	27, 66, 102, 140	1 (0%)

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	120	SER	6.0
2	C	220	GLU	5.0
1	F	79	GLN	4.7
2	A	147	GLY	4.7
2	C	126	GLY	4.6
2	E	25	SER	4.2
2	E	167	LEU	3.5
2	A	158	VAL	3.5
2	E	159	THR	3.3
2	C	33	ALA	3.3
1	B	89	GLN	3.0
2	A	145	ALA	3.0
1	D	10	SER	3.0
2	A	10	GLY	3.0
2	E	57[A]	HIS	3.0
1	D	36	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	L	68	GLY	2.9
1	L	122	ASP	2.9
2	E	166	ALA	2.9
1	B	46	LEU	2.8
1	D	12	SER	2.8
1	F	162	SER	2.8
1	L	83	VAL	2.7
2	H	24	ALA	2.7
1	D	211	ARG	2.7
1	L	97	THR	2.7
2	E	10	GLY	2.7
1	L	72	THR	2.6
2	A	212	ASN	2.6
2	H	128	SER	2.6
1	F	82	ASP	2.6
2	C	21	SER	2.6
2	A	167	LEU	2.6
2	A	1	GLU	2.6
1	L	127	SER	2.5
1	L	154	LEU	2.5
1	F	6	GLN	2.5
1	F	22	THR	2.5
1	D	134	CYS	2.5
1	B	106	ILE	2.5
2	A	64	VAL	2.5
1	D	72	THR	2.5
2	C	152	ASP	2.4
1	F	174	SER	2.4
1	B	25	ALA	2.4
1	B	35	TRP	2.4
2	E	115	THR	2.4
2	A	210	PRO	2.4
2	C	205	ASN	2.3
1	B	79	GLN	2.3
2	H	180	SER	2.3
2	A	31	ASP	2.3
2	H	97	ALA	2.3
1	B	65	SER	2.3
1	D	154	LEU	2.2
1	F	60	SER	2.2
1	B	23	CYS	2.2
2	C	169	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	73	LEU	2.2
2	E	65	GLU	2.2
2	A	176	ALA	2.2
2	C	77	ASN	2.2
2	H	64	VAL	2.2
2	C	134	PRO	2.2
1	D	23	CYS	2.2
2	A	132	LEU	2.2
1	D	152	ASN	2.1
1	L	22	THR	2.1
1	D	57	GLY	2.1
1	F	5	THR	2.1
1	F	109	THR	2.1
2	A	78	SER	2.1
2	E	75	ALA	2.1
2	A	14	PRO	2.1
1	B	114	SER	2.1
1	D	189	HIS	2.1
2	A	41	PRO	2.1
2	C	73	ASP	2.1
2	C	182	GLY	2.1
2	E	96	CYS	2.1
1	D	97	THR	2.1
1	F	35	TRP	2.1
2	E	64	VAL	2.0
2	A	28	THR	2.0
1	D	34	ALA	2.0
1	F	7	SER	2.0
1	F	164	THR	2.0
2	A	184	TYR	2.0
2	E	110	TYR	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.