



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:49 PM GMT

PDB ID : 3ZC2
Title : STRUCTURE OF AN ACTIVITY SUPPRESSING FAB FRAGMENT TO
CYTOCHROME P450 AROMATASE
Authors : Sawicki, M.W.; Ng, P.C.; Burkhart, B.; Pletnev, V.; Higashiyama, T.; Osawa,
Y.; Ghosh, D.
Deposited on : 1999-04-21
Resolution : 3.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

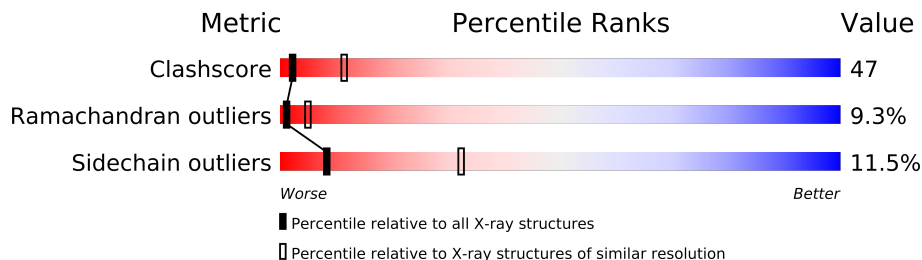
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	217	
2	B	218	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3326 atoms, of which 0 are hydrogen and 0 are deuterium.

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 IGG1 ANTIBODY 32C2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1673	1040	288	338	7			

- Molecule 2 is a protein called IGG1 ANTIBODY 32C2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1653	1048	266	333	6			

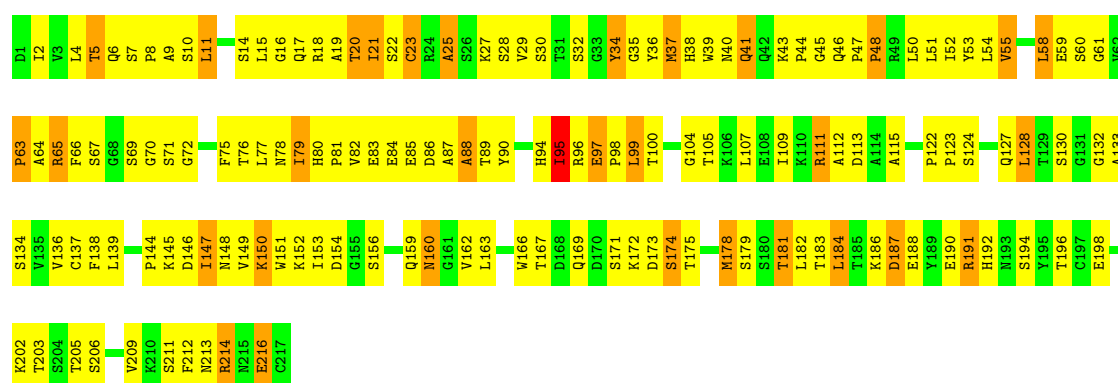
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

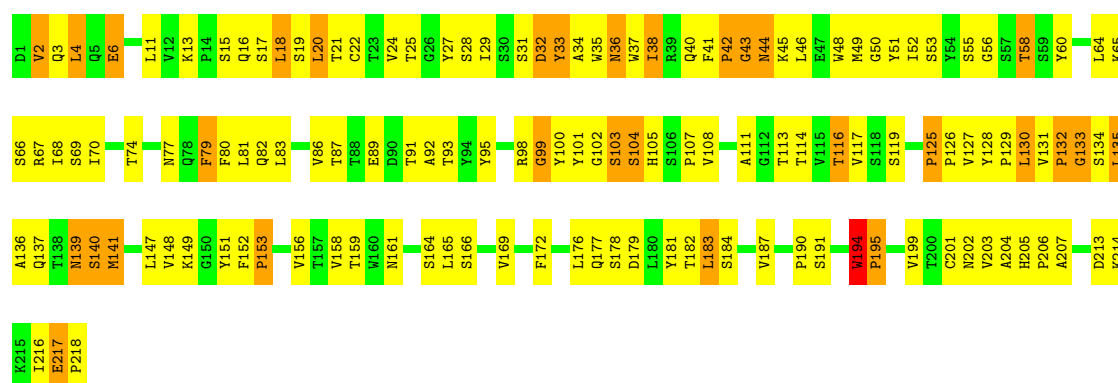
• Molecule 1: IGG1 ANTIBODY 32C2

Chain A:



• Molecule 2: IGG1 ANTIBODY 32C2

Chain B:



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	154.89Å 73.51Å 36.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.00 – 3.00	Depositor
% Data completeness (in resolution range)	88.3 (99.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.213 , 0.317	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3326	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

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	A	0.41	0/1711	0.67	0/2325
2	B	0.42	0/1701	0.71	0/2334
All	All	0.42	0/3412	0.69	0/4659

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1673	0	1610	166	0
2	B	1653	0	1603	156	0
All	All	3326	0	3213	310	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 47.

All (310) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:58:LEU:HD13	1:A:63:PRO:HB2	1.28	1.11
1:A:30:SER:HB3	1:A:72:GLY:H	1.20	1.03

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:163:LEU:HD11	2:B:177:GLN:HG3	1.43	0.98
1:A:80:HIS:HB2	1:A:81:PRO:HD3	1.45	0.96
1:A:136:VAL:HG22	1:A:181:THR:HG23	1.46	0.96
2:B:4:LEU:HB3	2:B:24:VAL:HG22	1.48	0.96
2:B:125:PRO:HB3	2:B:151:TYR:HB3	1.51	0.92
2:B:91:THR:HG23	2:B:116:THR:HA	1.50	0.92
1:A:30:SER:CB	1:A:72:GLY:H	1.85	0.90
2:B:158:VAL:HG22	2:B:203:VAL:HG22	1.55	0.88
2:B:33:TYR:HA	2:B:101:TYR:H	1.39	0.86
1:A:18:ARG:NH1	1:A:18:ARG:HB2	1.93	0.84
1:A:36:TYR:HB3	1:A:55:VAL:HG23	1.60	0.83
1:A:19:ALA:HB3	1:A:79:ILE:HG23	1.61	0.82
2:B:4:LEU:N	2:B:4:LEU:HD23	1.95	0.82
1:A:58:LEU:HD13	1:A:63:PRO:CB	2.09	0.81
2:B:31:SER:O	2:B:32:ASP:HB2	1.80	0.81
2:B:194:TRP:HB3	2:B:195:PRO:HD3	1.61	0.81
1:A:38:HIS:CE1	1:A:50:LEU:HD13	2.18	0.79
1:A:58:LEU:HB3	1:A:63:PRO:HB3	1.64	0.79
2:B:127:VAL:HG23	2:B:214:LYS:HD2	1.65	0.79
2:B:149:LYS:HG3	2:B:182:THR:OG1	1.85	0.77
2:B:166:SER:O	2:B:169:VAL:HG12	1.83	0.77
1:A:163:LEU:HD11	2:B:177:GLN:CG	2.15	0.77
1:A:28:SER:HA	1:A:72:GLY:O	1.84	0.77
1:A:136:VAL:HG22	1:A:181:THR:CG2	2.15	0.77
1:A:30:SER:HB3	1:A:72:GLY:N	1.97	0.76
2:B:20:LEU:HD23	2:B:20:LEU:H	1.49	0.76
2:B:158:VAL:HA	2:B:202:ASN:O	1.86	0.76
1:A:67:SER:HB2	1:A:78:ASN:ND2	2.01	0.75
1:A:150:LYS:HZ2	1:A:198:GLU:C	1.91	0.74
1:A:133:ALA:O	1:A:184:LEU:HD12	1.88	0.73
1:A:58:LEU:HD21	1:A:66:PHE:O	1.86	0.73
2:B:126:PRO:HD2	2:B:149:LYS:O	1.89	0.72
1:A:41:GLN:HG2	1:A:51:LEU:HD11	1.71	0.72
2:B:48:TRP:CZ2	2:B:50:GLY:HA2	2.25	0.72
2:B:44:ASN:HD22	2:B:45:LYS:NZ	1.87	0.72
2:B:32:ASP:HB3	2:B:33:TYR:HD1	1.54	0.72
2:B:16:GLN:HG2	2:B:17:SER:H	1.55	0.72
2:B:133:GLY:O	2:B:135:LEU:HD23	1.90	0.71
2:B:137:GLN:C	2:B:139:ASN:H	1.94	0.71
2:B:40:GLN:HG3	2:B:46:LEU:HD12	1.73	0.71
2:B:18:LEU:HD11	2:B:20:LEU:HD22	1.73	0.70
1:A:198:GLU:HG2	1:A:209:VAL:HG22	1.75	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:68:ILE:HD11	2:B:81:LEU:HD11	1.74	0.68
2:B:35:TRP:O	2:B:52:ILE:HG22	1.93	0.68
1:A:20:THR:OG1	1:A:78:ASN:HB3	1.94	0.67
1:A:95:ILE:HG21	2:B:103:SER:O	1.94	0.67
1:A:45:GLY:O	1:A:46:GLN:HG3	1.95	0.67
1:A:124:SER:OG	1:A:127:GLN:HB2	1.95	0.67
2:B:67:ARG:NH2	2:B:83:LEU:HD11	2.09	0.67
1:A:186:LYS:HE2	1:A:190:GLU:CD	2.15	0.67
1:A:55:VAL:HG11	1:A:71:SER:OG	1.94	0.67
2:B:40:GLN:OE1	2:B:46:LEU:HD11	1.96	0.66
1:A:163:LEU:HD21	2:B:177:GLN:HG2	1.77	0.66
1:A:52:ILE:HG22	1:A:53:TYR:N	2.11	0.66
2:B:15:SER:N	2:B:86:VAL:O	2.18	0.66
2:B:22:CYS:SG	2:B:79:PHE:CE1	2.89	0.66
2:B:129:PRO:HD3	2:B:214:LYS:HD3	1.78	0.65
1:A:10:SER:O	1:A:11:LEU:HB2	1.96	0.65
2:B:183:LEU:C	2:B:183:LEU:HD23	2.16	0.65
1:A:145:LYS:O	1:A:147:ILE:HG22	1.95	0.65
1:A:8:PRO:HD2	1:A:21:ILE:CG2	2.26	0.65
2:B:32:ASP:HB3	2:B:33:TYR:CD1	2.32	0.65
2:B:22:CYS:HB2	2:B:37:TRP:CH2	2.32	0.64
1:A:54:LEU:O	1:A:55:VAL:HB	1.96	0.64
2:B:176:LEU:HD23	2:B:176:LEU:N	2.12	0.64
1:A:16:GLY:HA2	1:A:81:PRO:HA	1.80	0.63
1:A:186:LYS:HE2	1:A:190:GLU:OE1	1.98	0.63
2:B:6:GLU:CD	2:B:6:GLU:H	2.02	0.63
1:A:96:ARG:HG2	1:A:96:ARG:HH11	1.64	0.63
1:A:18:ARG:HB2	1:A:18:ARG:HH11	1.62	0.63
1:A:4:LEU:HD11	1:A:94:HIS:HB3	1.81	0.63
1:A:94:HIS:ND1	1:A:95:ILE:N	2.48	0.62
2:B:38:ILE:HG23	2:B:48:TRP:HA	1.82	0.62
2:B:141:MET:HE2	2:B:190:PRO:HA	1.81	0.62
1:A:214:ARG:HH11	1:A:214:ARG:CB	2.11	0.62
2:B:204:ALA:O	2:B:206:PRO:HD3	2.00	0.62
2:B:131:VAL:HG11	2:B:217:GLU:O	2.00	0.62
2:B:202:ASN:ND2	2:B:213:ASP:OD2	2.33	0.61
1:A:94:HIS:C	1:A:95:ILE:HD13	2.20	0.61
1:A:41:GLN:O	1:A:48:PRO:HA	2.00	0.61
1:A:134:SER:HB3	1:A:183:THR:HG23	1.82	0.61
2:B:103:SER:O	2:B:104:SER:HB2	1.99	0.61
2:B:194:TRP:HH2	2:B:216:ILE:O	1.84	0.61
1:A:127:GLN:HG3	2:B:128:TYR:CE2	2.35	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:159:GLN:OE1	1:A:159:GLN:HA	2.01	0.60
1:A:83:GLU:HB3	1:A:86:ASP:CG	2.22	0.60
1:A:152:LYS:HA	1:A:156:SER:O	2.01	0.60
2:B:4:LEU:H	2:B:4:LEU:HD23	1.63	0.59
2:B:147:LEU:HD12	2:B:148:VAL:N	2.16	0.59
2:B:79:PHE:HD1	2:B:79:PHE:H	1.51	0.59
1:A:160:ASN:HD22	1:A:160:ASN:H	1.50	0.59
1:A:95:ILE:HG21	2:B:104:SER:HB2	1.83	0.59
1:A:35:GLY:HA3	1:A:96:ARG:HH12	1.67	0.59
2:B:20:LEU:N	2:B:20:LEU:HD23	2.16	0.59
2:B:44:ASN:C	2:B:45:LYS:HD2	2.23	0.59
1:A:97:GLU:HA	1:A:99:LEU:HD23	1.85	0.59
1:A:87:ALA:HA	1:A:107:LEU:HD12	1.85	0.59
1:A:169:GLN:OE1	1:A:174:SER:HB3	2.03	0.58
2:B:83:LEU:HD12	2:B:86:VAL:HG22	1.85	0.58
2:B:201:CYS:O	2:B:213:ASP:HA	2.03	0.58
1:A:160:ASN:HD22	1:A:160:ASN:N	2.01	0.58
1:A:95:ILE:HG22	1:A:99:LEU:HD13	1.86	0.58
2:B:194:TRP:CB	2:B:195:PRO:HD3	2.32	0.58
2:B:194:TRP:HB3	2:B:195:PRO:CD	2.34	0.58
2:B:44:ASN:HD22	2:B:45:LYS:HZ2	1.52	0.57
2:B:4:LEU:CB	2:B:24:VAL:HG22	2.31	0.57
2:B:45:LYS:O	2:B:46:LEU:HD12	2.05	0.57
1:A:167:THR:HG23	2:B:172:PHE:CD2	2.40	0.57
1:A:173:ASP:O	1:A:175:THR:N	2.38	0.57
1:A:99:LEU:HD23	1:A:99:LEU:N	2.20	0.57
1:A:65:ARG:NH2	1:A:83:GLU:OE1	2.37	0.56
1:A:144:PRO:HG3	1:A:202:LYS:HD3	1.86	0.56
2:B:98:ARG:HG2	2:B:98:ARG:O	2.05	0.56
1:A:98:PRO:O	1:A:100:THR:HG23	2.05	0.55
2:B:93:THR:HG23	2:B:114:THR:HG22	1.87	0.55
2:B:29:ILE:HD13	2:B:29:ILE:N	2.22	0.55
2:B:135:LEU:HD12	2:B:137:GLN:HB2	1.88	0.55
2:B:4:LEU:HD21	2:B:108:VAL:HG12	1.89	0.55
1:A:52:ILE:HG22	1:A:53:TYR:H	1.72	0.55
1:A:65:ARG:O	1:A:79:ILE:HA	2.07	0.54
1:A:41:GLN:HG2	1:A:51:LEU:HD21	1.89	0.54
2:B:34:ALA:N	2:B:101:TYR:HB2	2.22	0.54
2:B:52:ILE:HD13	2:B:58:THR:HG22	1.90	0.54
2:B:91:THR:CG2	2:B:116:THR:HA	2.31	0.54
1:A:178:MET:HG2	1:A:179:SER:N	2.23	0.54
1:A:113:ASP:OD2	1:A:202:LYS:HE2	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:22:CYS:SG	2:B:79:PHE:HE1	2.31	0.54
2:B:4:LEU:HD21	2:B:108:VAL:CG1	2.38	0.54
1:A:59:GLU:HG3	1:A:60:SER:N	2.23	0.53
2:B:41:PHE:HB2	2:B:45:LYS:HB2	1.90	0.53
1:A:196:THR:HA	1:A:211:SER:CB	2.38	0.53
2:B:86:VAL:HG12	2:B:117:VAL:HG11	1.90	0.53
1:A:137:CYS:HB2	1:A:151:TRP:CH2	2.43	0.53
1:A:109:ILE:HD11	1:A:174:SER:OG	2.09	0.53
1:A:150:LYS:HE3	1:A:150:LYS:N	2.24	0.53
1:A:39:TRP:O	1:A:51:LEU:HB2	2.09	0.53
2:B:40:GLN:HG3	2:B:46:LEU:CD1	2.39	0.53
1:A:53:TYR:HE2	1:A:59:GLU:OE1	1.92	0.53
1:A:196:THR:HA	1:A:211:SER:HB2	1.90	0.53
2:B:27:TYR:CD1	2:B:28:SER:N	2.77	0.53
2:B:27:TYR:CE1	2:B:31:SER:HB3	2.44	0.53
2:B:95:TYR:HE1	2:B:111:ALA:O	1.92	0.53
2:B:194:TRP:O	2:B:195:PRO:C	2.47	0.52
1:A:38:HIS:ND1	1:A:52:ILE:O	2.41	0.52
1:A:95:ILE:HA	1:A:99:LEU:HD22	1.91	0.52
2:B:194:TRP:CH2	2:B:216:ILE:O	2.61	0.52
1:A:80:HIS:CB	1:A:81:PRO:HD3	2.28	0.52
2:B:43:GLY:O	2:B:45:LYS:N	2.43	0.52
1:A:4:LEU:CD2	1:A:25:ALA:HB2	2.40	0.52
1:A:96:ARG:NH1	1:A:96:ARG:HG2	2.21	0.52
1:A:59:GLU:HG3	1:A:60:SER:H	1.75	0.52
1:A:190:GLU:HA	1:A:214:ARG:HH21	1.75	0.52
2:B:65:LYS:O	2:B:66:SER:HB2	2.10	0.52
1:A:50:LEU:HD11	2:B:107:PRO:HD2	1.91	0.52
1:A:166:TRP:CD1	1:A:178:MET:HB2	2.45	0.52
1:A:7:SER:O	1:A:21:ILE:HG22	2.10	0.51
2:B:33:TYR:CD2	2:B:100:TYR:HD1	2.28	0.51
2:B:139:ASN:O	2:B:140:SER:CB	2.58	0.51
1:A:69:SER:OG	1:A:70:GLY:N	2.44	0.51
2:B:60:TYR:HE1	2:B:70:ILE:HG13	1.76	0.51
2:B:151:TYR:CZ	2:B:156:VAL:HG21	2.45	0.51
2:B:65:LYS:O	2:B:66:SER:CB	2.59	0.51
1:A:83:GLU:HB3	1:A:86:ASP:OD1	2.11	0.51
2:B:13:LYS:HE2	2:B:16:GLN:OE1	2.11	0.51
1:A:47:PRO:HB3	2:B:95:TYR:CE1	2.46	0.51
1:A:128:LEU:HD12	1:A:133:ALA:HB2	1.93	0.50
1:A:6:GLN:HB2	1:A:23:CYS:SG	2.51	0.50
2:B:139:ASN:O	2:B:140:SER:HB3	2.11	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:36:ASN:HB3	2:B:51:TYR:HB3	1.94	0.50
2:B:41:PHE:HD2	2:B:45:LYS:HB3	1.76	0.50
2:B:95:TYR:CE1	2:B:111:ALA:O	2.64	0.50
1:A:84:GLU:H	1:A:84:GLU:CD	2.14	0.50
2:B:87:THR:O	2:B:117:VAL:HG11	2.11	0.50
1:A:81:PRO:O	1:A:83:GLU:N	2.45	0.50
1:A:52:ILE:CG2	1:A:53:TYR:N	2.75	0.50
2:B:141:MET:HA	2:B:191:SER:H	1.77	0.50
1:A:20:THR:HA	1:A:77:LEU:O	2.12	0.49
1:A:8:PRO:HD2	1:A:21:ILE:HG23	1.94	0.49
2:B:161:ASN:O	2:B:164:SER:HB3	2.12	0.49
1:A:83:GLU:OE1	1:A:86:ASP:OD1	2.29	0.49
2:B:137:GLN:C	2:B:139:ASN:N	2.64	0.49
2:B:187:VAL:HG13	2:B:187:VAL:O	2.12	0.49
1:A:83:GLU:HB3	1:A:86:ASP:OD2	2.13	0.49
1:A:128:LEU:O	1:A:186:LYS:HD3	2.12	0.49
2:B:176:LEU:HD22	2:B:181:TYR:CE1	2.47	0.49
2:B:194:TRP:CB	2:B:195:PRO:CD	2.90	0.49
2:B:51:TYR:HD1	2:B:52:ILE:N	2.10	0.49
1:A:75:PHE:CD1	1:A:75:PHE:N	2.80	0.49
1:A:196:THR:HG23	1:A:211:SER:HB3	1.95	0.49
2:B:129:PRO:O	2:B:130:LEU:HD12	2.13	0.49
1:A:115:ALA:HB2	1:A:203:THR:HG21	1.95	0.48
1:A:28:SER:O	1:A:96:ARG:HD2	2.13	0.48
1:A:149:VAL:C	1:A:150:LYS:HE3	2.33	0.48
2:B:21:THR:HG1	2:B:80:PHE:HE1	1.59	0.48
2:B:178:SER:OG	2:B:179:ASP:N	2.46	0.48
2:B:135:LEU:HD12	2:B:137:GLN:CB	2.44	0.48
2:B:194:TRP:CZ3	2:B:199:VAL:HG23	2.47	0.48
1:A:214:ARG:HH11	1:A:214:ARG:HB2	1.78	0.48
1:A:194:SER:HA	1:A:213:ASN:HA	1.94	0.48
1:A:58:LEU:CB	1:A:63:PRO:HB3	2.41	0.48
2:B:4:LEU:HD11	2:B:108:VAL:O	2.14	0.48
1:A:22:SER:HB3	1:A:76:THR:HG22	1.96	0.47
1:A:83:GLU:HG2	1:A:84:GLU:N	2.29	0.47
1:A:8:PRO:HD2	1:A:21:ILE:HG22	1.97	0.47
2:B:98:ARG:O	2:B:99:GLY:C	2.53	0.47
1:A:34:TYR:CD1	2:B:103:SER:OG	2.67	0.47
1:A:43:LYS:O	1:A:45:GLY:N	2.48	0.47
2:B:13:LYS:N	2:B:13:LYS:HD3	2.29	0.47
1:A:94:HIS:ND1	1:A:94:HIS:C	2.67	0.47
1:A:216:GLU:HG2	1:A:216:GLU:O	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:37:MET:HG3	1:A:38:HIS:H	1.80	0.47
1:A:94:HIS:O	1:A:95:ILE:HG23	2.14	0.47
1:A:107:LEU:O	1:A:107:LEU:HD13	2.15	0.47
1:A:136:VAL:HG12	1:A:137:CYS:N	2.29	0.47
2:B:20:LEU:N	2:B:20:LEU:CD2	2.78	0.46
1:A:34:TYR:HB3	2:B:103:SER:OG	2.15	0.46
1:A:14:SER:O	1:A:15:LEU:C	2.54	0.46
1:A:21:ILE:O	1:A:76:THR:HA	2.15	0.46
2:B:152:PHE:HA	2:B:153:PRO:HA	1.67	0.46
1:A:52:ILE:CG2	1:A:53:TYR:H	2.28	0.46
2:B:67:ARG:NH2	2:B:83:LEU:CD1	2.77	0.46
1:A:205:THR:OG1	1:A:206:SER:N	2.47	0.46
1:A:162:VAL:O	1:A:163:LEU:HD23	2.16	0.46
1:A:41:GLN:CG	1:A:51:LEU:HD11	2.44	0.45
2:B:44:ASN:HD22	2:B:45:LYS:HZ3	1.61	0.45
2:B:141:MET:HG3	2:B:190:PRO:HA	1.98	0.45
2:B:19:SER:HA	2:B:81:LEU:O	2.16	0.45
2:B:205:HIS:CE1	2:B:207:ALA:HB3	2.51	0.45
2:B:113:THR:O	2:B:113:THR:HG23	2.15	0.45
1:A:67:SER:O	1:A:78:ASN:OD1	2.35	0.45
1:A:18:ARG:HB2	1:A:18:ARG:CZ	2.45	0.45
2:B:35:TRP:HB2	2:B:52:ILE:HG23	1.97	0.45
1:A:111:ARG:HG2	1:A:174:SER:O	2.16	0.45
1:A:30:SER:HA	1:A:36:TYR:HA	1.99	0.45
1:A:186:LYS:O	1:A:190:GLU:HG3	2.16	0.45
1:A:85:GLU:HG3	1:A:171:SER:OG	2.16	0.45
2:B:4:LEU:CD2	2:B:4:LEU:N	2.67	0.45
1:A:127:GLN:O	1:A:128:LEU:C	2.55	0.45
1:A:139:LEU:HD12	1:A:139:LEU:N	2.32	0.45
1:A:80:HIS:HB2	1:A:81:PRO:CD	2.33	0.44
2:B:152:PHE:CE2	2:B:153:PRO:HB3	2.52	0.44
1:A:196:THR:CA	1:A:211:SER:HB2	2.48	0.44
2:B:40:GLN:O	2:B:92:ALA:HB1	2.17	0.44
2:B:183:LEU:HD23	2:B:184:SER:N	2.32	0.44
2:B:89:GLU:H	2:B:89:GLU:CD	2.21	0.44
1:A:2:ILE:HD11	1:A:25:ALA:HB1	1.99	0.44
2:B:164:SER:O	2:B:165:LEU:HB2	2.17	0.44
1:A:90:TYR:O	1:A:104:GLY:HA2	2.18	0.44
1:A:123:PRO:HB2	1:A:128:LEU:HD13	1.99	0.44
1:A:127:GLN:NE2	1:A:134:SER:OG	2.50	0.44
1:A:194:SER:HB3	1:A:213:ASN:OD1	2.17	0.44
1:A:54:LEU:O	1:A:55:VAL:CB	2.66	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:51:TYR:CD1	2:B:51:TYR:C	2.91	0.44
1:A:87:ALA:O	1:A:88:ALA:HB2	2.18	0.43
1:A:9:ALA:O	1:A:105:THR:HA	2.18	0.43
2:B:159:THR:OG1	2:B:202:ASN:HB2	2.18	0.43
2:B:34:ALA:HB2	2:B:101:TYR:HB2	1.98	0.43
2:B:136:ALA:HA	2:B:139:ASN:OD1	2.18	0.43
1:A:188:GLU:O	1:A:191:ARG:HB3	2.17	0.43
2:B:27:TYR:HE1	2:B:31:SER:HB3	1.83	0.43
2:B:131:VAL:HA	2:B:132:PRO:HD3	1.80	0.43
1:A:67:SER:HB2	1:A:78:ASN:HD21	1.78	0.43
2:B:53:SER:C	2:B:55:SER:N	2.72	0.43
1:A:39:TRP:N	1:A:39:TRP:CD1	2.87	0.43
1:A:4:LEU:O	1:A:5:THR:C	2.56	0.43
2:B:29:ILE:HG12	2:B:77:ASN:HA	2.00	0.43
2:B:53:SER:O	2:B:56:GLY:N	2.43	0.43
1:A:30:SER:HB2	1:A:75:PHE:CE2	2.54	0.42
2:B:68:ILE:HA	2:B:82:GLN:O	2.19	0.42
1:A:4:LEU:HD23	1:A:25:ALA:HB2	2.00	0.42
2:B:135:LEU:HB2	2:B:136:ALA:H	1.60	0.42
1:A:18:ARG:CB	1:A:18:ARG:HH11	2.29	0.42
2:B:51:TYR:CD1	2:B:52:ILE:N	2.87	0.42
2:B:176:LEU:CD2	2:B:176:LEU:N	2.81	0.42
1:A:138:PHE:C	1:A:139:LEU:HD12	2.40	0.42
2:B:3:GLN:HB2	2:B:25:THR:OG1	2.19	0.42
1:A:212:PHE:C	1:A:212:PHE:CD1	2.92	0.42
2:B:28:SER:O	2:B:31:SER:HB3	2.18	0.42
2:B:52:ILE:O	2:B:52:ILE:HG23	2.20	0.42
2:B:11:LEU:HD13	2:B:11:LEU:C	2.39	0.42
2:B:2:VAL:HG12	2:B:4:LEU:HD22	2.01	0.42
2:B:166:SER:O	2:B:169:VAL:CG1	2.62	0.42
1:A:154:ASP:HB2	1:A:192:HIS:HD2	1.84	0.42
1:A:130:SER:C	1:A:132:GLY:H	2.24	0.41
2:B:33:TYR:CD2	2:B:100:TYR:HA	2.55	0.41
1:A:41:GLN:HB3	1:A:51:LEU:HD11	2.01	0.41
2:B:16:GLN:HG2	2:B:17:SER:N	2.30	0.41
2:B:93:THR:HA	2:B:114:THR:HA	2.01	0.41
1:A:63:PRO:O	1:A:64:ALA:HB3	2.21	0.41
1:A:87:ALA:HA	1:A:107:LEU:CD1	2.49	0.41
1:A:111:ARG:HD3	1:A:112:ALA:O	2.20	0.41
2:B:49:MET:HE1	2:B:81:LEU:HD21	2.03	0.41
2:B:126:PRO:HG2	2:B:126:PRO:O	2.20	0.41
1:A:34:TYR:CG	2:B:103:SER:OG	2.74	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:45:GLY:C	1:A:46:GLN:HG3	2.42	0.41
2:B:172:PHE:O	2:B:183:LEU:HG	2.21	0.41
1:A:147:ILE:O	1:A:147:ILE:HG23	2.20	0.41
2:B:18:LEU:HD12	2:B:19:SER:N	2.36	0.41
1:A:83:GLU:N	1:A:86:ASP:OD2	2.54	0.40
1:A:20:THR:O	1:A:20:THR:HG22	2.20	0.40
1:A:122:PRO:HA	1:A:123:PRO:HD3	1.96	0.40
1:A:203:THR:O	1:A:203:THR:HG22	2.20	0.40
1:A:187:ASP:O	1:A:188:GLU:C	2.58	0.40
1:A:153:ILE:HG22	1:A:192:HIS:CD2	2.56	0.40
2:B:101:TYR:CG	2:B:102:GLY:N	2.88	0.40
2:B:68:ILE:HG12	2:B:69:SER:H	1.87	0.40
2:B:79:PHE:N	2:B:79:PHE:CD1	2.88	0.40
2:B:217:GLU:O	2:B:218:PRO:O	2.40	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	215/217 (99%)	153 (71%)	43 (20%)	19 (9%)	1	5
2	B	216/218 (99%)	161 (74%)	34 (16%)	21 (10%)	1	4
All	All	431/435 (99%)	314 (73%)	77 (18%)	40 (9%)	1	5

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	82	VAL
1	A	97	GLU
1	A	172	LYS
1	A	174	SER
1	A	216	GLU
2	B	44	ASN

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Mol	Chain	Res	Type
2	B	140	SER
2	B	194	TRP
1	A	61	GLY
1	A	187	ASP
2	B	32	ASP
2	B	103	SER
2	B	132	PRO
2	B	134	SER
2	B	139	ASN
1	A	11	LEU
1	A	32	SER
1	A	44	PRO
1	A	58	LEU
1	A	63	PRO
2	B	2	VAL
2	B	33	TYR
2	B	74	THR
2	B	125	PRO
2	B	141	MET
1	A	25	ALA
1	A	29	VAL
1	A	88	ALA
1	A	95	ILE
2	B	42	PRO
2	B	104	SER
2	B	105	HIS
2	B	153	PRO
1	A	48	PRO
1	A	55	VAL
2	B	43	GLY
2	B	133	GLY
1	A	147	ILE
2	B	99	GLY
2	B	195	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	190/190 (100%)	163 (86%)	27 (14%)	5	22
2	B	194/194 (100%)	177 (91%)	17 (9%)	14	48
All	All	384/384 (100%)	340 (88%)	44 (12%)	8	32

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	17	GLN
1	A	20	THR
1	A	21	ILE
1	A	23	CYS
1	A	27	LYS
1	A	34	TYR
1	A	37	MET
1	A	40	ASN
1	A	41	GLN
1	A	65	ARG
1	A	79	ILE
1	A	89	THR
1	A	95	ILE
1	A	99	LEU
1	A	111	ARG
1	A	128	LEU
1	A	146	ASP
1	A	148	ASN
1	A	150	LYS
1	A	160	ASN
1	A	178	MET
1	A	181	THR
1	A	182	LEU
1	A	184	LEU
1	A	191	ARG
1	A	214	ARG
2	B	4	LEU
2	B	6	GLU
2	B	18	LEU
2	B	20	LEU
2	B	36	ASN
2	B	38	ILE
2	B	42	PRO
2	B	58	THR

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Mol	Chain	Res	Type
2	B	64	LEU
2	B	79	PHE
2	B	116	THR
2	B	119	SER
2	B	130	LEU
2	B	135	LEU
2	B	183	LEU
2	B	194	TRP
2	B	217	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	127	GLN
1	A	160	ASN
1	A	164	ASN
1	A	192	HIS
1	A	193	ASN
1	A	201	HIS
2	B	44	ASN
2	B	177	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.