



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:46 PM UTC

PDB ID : 2Z0V / pdb_00002z0v
Title : Crystal structure of BAR domain of Endophilin-III
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Deposited on : 2007-05-07
Resolution : 2.49 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
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

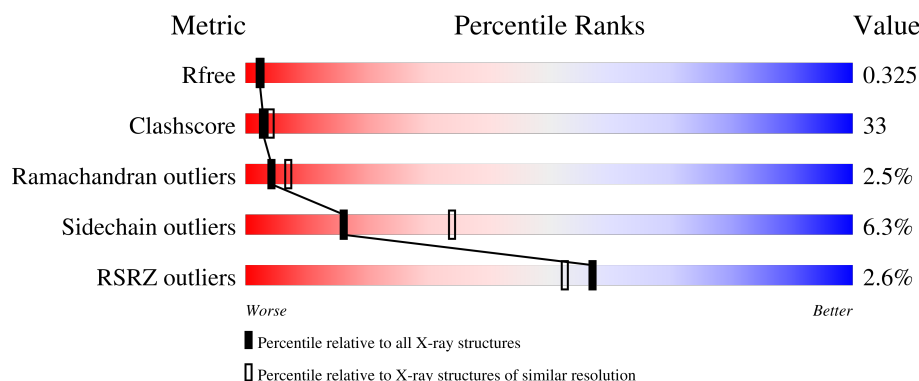
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	240	
1	B	240	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3657 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 SH3-containing GRB2-like protein 3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	Se	0	0	0
			1788	1123	304	353	1	7			
1	B	221	Total	C	N	O	S	Se	0	0	0
			1776	1113	304	351	1	7			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	GLY	-	expression tag	UNP Q99963
A	-5	SER	-	expression tag	UNP Q99963
A	-4	SER	-	expression tag	UNP Q99963
A	-3	GLY	-	expression tag	UNP Q99963
A	-2	SER	-	expression tag	UNP Q99963
A	-1	SER	-	expression tag	UNP Q99963
A	0	GLY	-	expression tag	UNP Q99963
B	-6	GLY	-	expression tag	UNP Q99963
B	-5	SER	-	expression tag	UNP Q99963
B	-4	SER	-	expression tag	UNP Q99963
B	-3	GLY	-	expression tag	UNP Q99963
B	-2	SER	-	expression tag	UNP Q99963
B	-1	SER	-	expression tag	UNP Q99963
B	0	GLY	-	expression tag	UNP Q99963

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	45	Total	O	0	0
			45	45		
2	B	48	Total	O	0	0
			48	48		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	56.82Å 194.17Å 43.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.04 – 2.49 49.04 – 2.49	Depositor EDS
% Data completeness (in resolution range)	94.2 (49.04-2.49) 94.2 (49.04-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.242 , 0.322 0.246 , 0.325	Depositor DCC
R_{free} test set	1676 reflections (10.03%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.958	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3657	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% 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	A	0.45	0/1803	1.04	10/2407 (0.4%)
1	B	0.45	0/1789	0.95	8/2385 (0.3%)
All	All	0.45	0/3592	1.00	18/4792 (0.4%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	227	VAL	CA-C-N	11.58	134.31	119.84
1	A	227	VAL	C-N-CA	11.58	134.31	119.84
1	B	58	VAL	N-CA-C	-7.90	105.49	113.47
1	B	183	GLU	N-CA-C	-7.04	103.26	112.41
1	A	189	VAL	N-CA-C	-6.26	103.22	111.05
1	A	65	GLN	N-CA-C	-6.18	99.82	109.72
1	A	186	VAL	N-CA-C	-6.17	103.59	110.62
1	A	228	PRO	N-CA-C	5.70	124.22	112.47
1	B	17	ILE	N-CA-C	-5.70	103.97	111.09
1	B	200	LEU	N-CA-C	5.65	117.44	111.28
1	A	127	ASP	N-CA-C	5.43	117.62	111.11
1	B	63	TYR	CA-C-N	5.33	126.03	120.12
1	B	63	TYR	C-N-CA	5.33	126.03	120.12
1	A	121	GLN	N-CA-C	-5.20	105.79	111.82
1	B	181	PHE	N-CA-C	-5.19	107.02	113.55
1	A	114	GLN	N-CA-C	5.11	116.66	111.14
1	B	137	LYS	N-CA-C	-5.08	105.65	111.14
1	A	52	SER	N-CA-C	-5.04	103.02	110.48

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	1788	0	1812	142	0
1	B	1776	0	1795	117	0
2	A	45	0	0	7	0
2	B	48	0	0	6	0
All	All	3657	0	3607	234	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 33.

All (234) 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:A:155:ILE:HD12	1:A:159:GLU:HG2	1.39	1.02
1:A:69:LEU:HB3	1:B:30:LYS:HZ3	1.38	0.88
1:B:158:GLU:HA	1:B:161:ARG:HD2	1.58	0.85
1:A:89:ASN:HB3	1:A:220:ARG:NH2	1.93	0.83
1:B:51:VAL:HG13	1:B:118:ASP:HB3	1.60	0.83
1:B:24:VAL:O	1:B:28:LEU:HG	1.80	0.81
1:B:70:LEU:HD23	1:B:98:MSE:HE2	1.62	0.80
1:A:13:MSE:HE1	1:A:175:GLU:HG3	1.65	0.78
1:A:23:VAL:HG22	1:B:77:TYR:HB3	1.65	0.77
1:B:124:GLN:HA	1:B:128:LEU:HD12	1.66	0.77
1:B:7:ASP:O	1:B:11:LEU:HD23	1.84	0.77
1:B:111:ASN:O	1:B:115:THR:HG22	1.86	0.76
1:A:174:ALA:O	1:A:178:MSE:HG3	1.87	0.75
1:B:85:SER:HA	2:B:252:HOH:O	1.87	0.74
1:A:188:GLN:HA	1:A:191:GLN:HE21	1.53	0.73
1:A:17:ILE:HD13	1:A:131:ILE:HG21	1.71	0.73
1:B:145:TYR:CE1	1:B:164:VAL:HG22	2.25	0.72
1:A:15:ARG:O	1:A:19:VAL:HG23	1.90	0.71
1:A:186:VAL:O	1:A:189:VAL:HG22	1.89	0.71
1:B:19:VAL:O	1:B:23:VAL:HG22	1.91	0.71
1:A:77:TYR:HB2	1:B:23:VAL:HG13	1.72	0.70
1:A:15:ARG:HH11	1:A:15:ARG:HB3	1.57	0.70
1:A:80:GLU:HG2	1:B:19:VAL:HG21	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:GLU:C	1:A:184:ASN:HD22	1.99	0.70
1:A:161:ARG:O	1:A:164:VAL:HG22	1.92	0.69
1:A:168:GLU:HA	1:A:171:LYS:HG2	1.75	0.69
1:B:29:SER:O	1:B:33:GLU:HG3	1.92	0.69
1:A:29:SER:O	1:A:33:GLU:HG3	1.94	0.68
1:A:55:ARG:HH11	1:A:55:ARG:HB2	1.58	0.68
1:A:161:ARG:NH1	1:A:162:GLN:HE22	1.91	0.68
1:A:156:PRO:HD2	1:A:159:GLU:OE1	1.93	0.68
1:A:188:GLN:HE21	1:A:191:GLN:NE2	1.92	0.67
1:A:89:ASN:HB3	1:A:220:ARG:HH21	1.59	0.67
1:B:54:ILE:HG13	1:B:55:ARG:NH1	2.09	0.67
1:A:15:ARG:HB3	1:A:15:ARG:NH1	2.10	0.67
1:B:21:ASN:ND2	1:B:128:LEU:HD13	2.10	0.67
1:A:184:ASN:HD22	1:A:184:ASN:N	1.92	0.66
1:A:69:LEU:HB3	1:B:30:LYS:NZ	2.09	0.66
1:A:30:LYS:NZ	1:B:73:CYS:SG	2.70	0.64
1:B:106:ASP:O	1:B:110:ILE:HG13	1.97	0.64
1:A:36:GLN:HG2	1:A:41:TYR:HB2	1.80	0.64
1:A:158:GLU:O	1:A:162:GLN:HG2	1.98	0.64
1:A:21:ASN:OD1	1:A:128:LEU:HD13	1.98	0.64
1:B:178:MSE:O	1:B:182:LEU:HG	1.98	0.63
1:B:17:ILE:HD13	1:B:131:ILE:HG21	1.81	0.63
1:B:55:ARG:HA	2:B:280:HOH:O	1.99	0.63
1:B:93:GLU:OE2	1:B:220:ARG:HD2	1.98	0.63
1:B:176:ARG:HG3	1:B:176:ARG:HH11	1.64	0.63
1:A:39:PRO:HA	1:A:42:ARG:HG3	1.80	0.62
1:B:75:LEU:HD23	1:B:92:ILE:HA	1.82	0.62
1:A:216:LYS:O	1:A:220:ARG:HG3	1.99	0.62
1:B:116:PHE:C	1:B:119:PRO:HD2	2.25	0.62
1:A:135:LEU:HD23	1:A:174:ALA:HB1	1.82	0.61
1:A:123:LEU:HA	2:A:251:HOH:O	1.99	0.61
1:A:31:THR:O	1:A:34:TYR:HB3	2.00	0.61
1:A:216:LYS:HA	1:A:219:MSE:HE3	1.84	0.60
1:B:184:ASN:ND2	1:B:187:GLU:OE2	2.35	0.60
1:A:17:ILE:CD1	1:A:131:ILE:HG21	2.32	0.59
1:A:55:ARG:HB2	1:A:55:ARG:NH1	2.16	0.59
1:A:31:THR:HG22	1:A:117:ILE:HD11	1.83	0.59
1:A:228:PRO:O	1:A:229:ARG:HB3	2.02	0.59
1:B:127:ASP:O	1:B:131:ILE:HG12	2.01	0.59
1:A:53:LYS:HD3	1:A:57:GLN:HG3	1.83	0.59
1:A:201:ASP:CG	1:A:205:GLN:HE21	2.11	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:VAL:HG11	1:B:86:THR:HG22	1.84	0.58
1:B:111:ASN:O	1:B:115:THR:CG2	2.51	0.58
1:B:184:ASN:O	1:B:186:VAL:N	2.36	0.58
1:B:64:PRO:O	1:B:65:GLN:HB2	2.04	0.58
1:A:17:ILE:HD12	1:A:178:MSE:HG2	1.86	0.58
1:B:130:GLU:HG3	2:B:279:HOH:O	2.03	0.58
1:B:137:LYS:O	1:B:141:ARG:HG2	2.04	0.58
1:A:37:PRO:HB2	1:B:37:PRO:HB2	1.84	0.58
1:B:48:LEU:HD21	1:B:58:VAL:HG13	1.84	0.58
1:B:114:GLN:HA	1:B:118:ASP:OD2	2.04	0.57
1:A:188:GLN:HG3	1:A:191:GLN:NE2	2.20	0.57
1:A:220:ARG:HG2	1:A:220:ARG:HH11	1.70	0.57
1:B:21:ASN:HD21	1:B:128:LEU:HD13	1.70	0.57
1:A:137:LYS:HG2	2:A:268:HOH:O	2.04	0.57
1:B:13:MSE:O	1:B:17:ILE:HG23	2.05	0.57
1:B:216:LYS:O	1:B:220:ARG:HG3	2.05	0.57
1:A:23:VAL:HG22	1:B:77:TYR:CB	2.33	0.56
1:A:185:ASP:HB2	1:B:87:PHE:HB2	1.87	0.56
1:B:36:GLN:OE1	1:B:59:LYS:HB3	2.04	0.56
1:A:200:LEU:HG	1:B:210:LEU:HB3	1.87	0.56
1:B:75:LEU:O	1:B:79:LYS:HG3	2.06	0.56
1:A:200:LEU:HD22	1:A:204:ARG:NH1	2.21	0.55
1:A:161:ARG:NH1	1:A:162:GLN:NE2	2.53	0.55
1:A:187:GLU:O	1:A:190:SER:HB3	2.07	0.55
1:A:78:GLY:N	2:A:237:HOH:O	2.40	0.55
1:A:155:ILE:CD1	1:A:159:GLU:HG2	2.23	0.55
1:A:225:SER:C	1:A:227:VAL:H	2.14	0.55
1:A:214:GLN:HA	1:A:214:GLN:OE1	2.06	0.55
1:B:176:ARG:HG3	1:B:176:ARG:NH1	2.22	0.55
1:A:108:LEU:O	1:A:112:VAL:HG23	2.06	0.55
1:A:116:PHE:O	1:A:120:LEU:HB2	2.06	0.54
1:A:164:VAL:HG23	1:A:165:GLU:N	2.21	0.54
1:A:209:ILE:HG23	2:A:270:HOH:O	2.06	0.54
1:B:95:GLY:O	1:B:99:LYS:HG2	2.08	0.54
1:B:20:THR:O	1:B:23:VAL:HG23	2.07	0.54
1:A:145:TYR:HA	1:A:163:ALA:HB1	1.89	0.54
1:A:201:ASP:OD1	1:A:205:GLN:NE2	2.40	0.54
1:B:17:ILE:HG13	1:B:18:ASP:N	2.19	0.54
1:A:77:TYR:CB	1:B:23:VAL:HG13	2.38	0.53
1:A:141:ARG:NH1	1:A:166:LYS:HD2	2.23	0.53
1:A:161:ARG:HH11	1:A:162:GLN:NE2	2.05	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:LYS:C	1:B:150:LYS:H	2.15	0.53
1:A:19:VAL:O	1:A:23:VAL:HG23	2.08	0.53
1:B:22:LYS:O	1:B:26:GLU:HG2	2.09	0.53
1:B:32:THR:HG21	1:B:48:LEU:HD13	1.90	0.53
1:B:137:LYS:HD2	1:B:170:SER:OG	2.08	0.53
1:B:108:LEU:HD22	1:B:202:TYR:CD2	2.44	0.53
1:A:36:GLN:CG	1:A:41:TYR:HB2	2.39	0.52
1:A:168:GLU:O	1:A:171:LYS:HG2	2.11	0.51
1:B:212:GLU:O	1:B:216:LYS:HG2	2.09	0.51
1:A:128:LEU:HD21	1:A:181:PHE:CZ	2.46	0.51
1:B:54:ILE:HG13	1:B:55:ARG:HH12	1.75	0.51
1:A:164:VAL:CG2	1:A:165:GLU:N	2.74	0.50
1:B:70:LEU:O	1:B:74:MSE:HG3	2.12	0.50
1:B:118:ASP:N	1:B:119:PRO:CD	2.75	0.50
1:A:115:THR:HB	1:A:194:VAL:HG11	1.93	0.50
1:B:108:LEU:HD22	1:B:202:TYR:CG	2.46	0.50
1:B:161:ARG:C	1:B:163:ALA:H	2.20	0.50
1:A:19:VAL:HA	1:A:22:LYS:HD3	1.93	0.50
1:B:26:GLU:OE1	1:B:26:GLU:HA	2.11	0.50
1:B:165:GLU:O	1:B:169:GLU:HG3	2.10	0.50
1:A:196:ILE:O	1:A:200:LEU:HB2	2.11	0.50
1:B:184:ASN:O	1:B:186:VAL:HG12	2.12	0.50
1:A:23:VAL:CG2	1:B:77:TYR:HB3	2.38	0.50
1:B:158:GLU:O	1:B:162:GLN:HG2	2.12	0.49
1:A:210:LEU:HB3	1:B:200:LEU:HD21	1.94	0.49
1:B:51:VAL:HG13	1:B:118:ASP:CB	2.37	0.49
1:A:43:ALA:C	1:A:45:LEU:H	2.20	0.49
1:A:164:VAL:O	1:A:167:PHE:HB3	2.11	0.49
1:B:172:GLU:O	1:B:175:GLU:N	2.45	0.49
1:A:192:LEU:O	1:A:196:ILE:HG12	2.13	0.49
1:A:184:ASN:N	1:A:184:ASN:ND2	2.57	0.49
1:A:220:ARG:HG2	1:A:220:ARG:NH1	2.27	0.49
1:A:155:ILE:HD12	1:A:159:GLU:CG	2.27	0.49
1:A:12:ASP:O	1:A:16:LYS:HG2	2.13	0.48
1:A:69:LEU:HD13	1:B:30:LYS:HE2	1.96	0.48
1:B:181:PHE:O	1:B:184:ASN:HB2	2.14	0.48
1:A:8:ASP:C	1:A:10:PHE:N	2.68	0.48
1:A:80:GLU:HG2	1:B:19:VAL:CG2	2.42	0.48
1:A:18:ASP:OD1	1:A:22:LYS:HD2	2.14	0.48
1:A:37:PRO:O	1:A:42:ARG:NH2	2.38	0.47
1:B:70:LEU:HD11	1:B:74:MSE:HE3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:ASN:O	1:B:185:ASP:C	2.57	0.47
1:A:168:GLU:HA	1:A:171:LYS:CG	2.44	0.47
1:B:146:ASP:O	1:B:149:LYS:HG2	2.14	0.47
1:A:187:GLU:O	1:A:191:GLN:HG3	2.15	0.47
1:B:148:LYS:O	1:B:150:LYS:N	2.48	0.47
1:B:227:VAL:C	2:B:268:HOH:O	2.57	0.47
1:A:69:LEU:HB3	1:B:30:LYS:CE	2.45	0.46
1:B:148:LYS:C	1:B:150:LYS:N	2.73	0.46
1:A:145:TYR:CZ	1:A:149:LYS:HD2	2.51	0.46
1:A:77:TYR:HB2	1:B:23:VAL:CG1	2.43	0.46
1:B:94:VAL:HG22	1:B:217:LEU:HD11	1.98	0.46
1:A:161:ARG:HH12	1:A:162:GLN:HE22	1.62	0.46
1:A:31:THR:CG2	1:A:117:ILE:HD11	2.45	0.46
1:B:196:ILE:N	1:B:196:ILE:HD12	2.31	0.45
1:B:58:VAL:O	1:B:58:VAL:HG12	2.17	0.45
1:B:55:ARG:NH1	1:B:55:ARG:HG3	2.31	0.45
1:B:67:GLU:OE2	1:B:105:LYS:HG3	2.17	0.45
1:A:100:LEU:O	1:A:104:VAL:HG23	2.17	0.45
1:A:92:ILE:O	1:A:96:GLU:HG3	2.17	0.45
1:A:118:ASP:HB2	1:A:119:PRO:HD3	1.98	0.45
1:A:138:LEU:HD13	1:A:170:SER:HB3	1.99	0.45
1:A:138:LEU:HD22	1:A:171:LYS:HA	1.99	0.45
1:B:115:THR:HG23	1:B:194:VAL:HG12	1.99	0.44
1:B:92:ILE:O	1:B:96:GLU:HG3	2.17	0.44
1:A:17:ILE:HD11	1:A:178:MSE:HA	1.99	0.44
1:B:51:VAL:CG1	1:B:118:ASP:HB3	2.41	0.44
1:A:8:ASP:HA	1:A:11:LEU:HB2	2.00	0.44
1:A:65:GLN:CD	1:A:105:LYS:HG3	2.43	0.44
1:A:77:TYR:HB2	2:A:237:HOH:O	2.15	0.44
1:A:86:THR:HB	1:B:185:ASP:HB3	1.99	0.44
1:A:144:ASP:O	1:A:147:TYR:HB3	2.18	0.44
1:B:29:SER:HB2	2:B:251:HOH:O	2.17	0.44
1:A:38:ASN:ND2	1:A:40:ALA:HB3	2.32	0.44
1:A:87:PHE:O	1:A:91:LEU:HB2	2.17	0.44
1:A:113:LYS:HD3	1:A:113:LYS:C	2.42	0.44
1:A:127:ASP:O	1:A:131:ILE:HG13	2.18	0.44
1:A:184:ASN:C	1:A:186:VAL:N	2.75	0.44
1:B:192:LEU:O	1:B:195:PHE:HB3	2.16	0.44
1:A:26:GLU:O	1:A:30:LYS:HG2	2.18	0.44
1:A:43:ALA:C	1:A:45:LEU:N	2.75	0.44
1:A:55:ARG:NH1	1:A:55:ARG:CB	2.79	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:GLN:NE2	1:A:105:LYS:HG3	2.33	0.44
1:A:227:VAL:HA	1:A:228:PRO:HD3	1.81	0.43
1:B:158:GLU:N	1:B:161:ARG:NH1	2.66	0.43
1:B:162:GLN:O	1:B:166:LYS:HG3	2.19	0.43
1:A:70:LEU:HD21	1:A:74:MSE:HE3	2.01	0.43
1:A:17:ILE:HD12	1:A:178:MSE:CG	2.49	0.43
1:B:151:ARG:HA	1:B:151:ARG:HD3	1.81	0.43
1:A:17:ILE:CD1	1:A:178:MSE:HG2	2.48	0.43
1:A:70:LEU:HG	1:B:30:LYS:HB3	2.00	0.43
1:A:189:VAL:CG2	1:B:221:ILE:HG12	2.49	0.43
1:A:151:ARG:O	1:A:155:ILE:HG12	2.18	0.43
1:A:168:GLU:CA	1:A:171:LYS:HG2	2.47	0.43
1:B:8:ASP:O	1:B:9:GLU:C	2.62	0.43
1:B:107:SER:O	1:B:108:LEU:C	2.62	0.43
1:B:181:PHE:C	1:B:183:GLU:H	2.26	0.43
1:A:17:ILE:HD13	1:A:131:ILE:HD13	2.01	0.42
1:B:48:LEU:CD2	1:B:58:VAL:HG13	2.49	0.42
1:A:53:LYS:NZ	1:A:59:LYS:HE2	2.35	0.42
1:B:97:SER:CB	1:B:213:LEU:HD22	2.49	0.42
1:A:91:LEU:HD12	1:A:91:LEU:HA	1.84	0.42
1:A:30:LYS:HB2	1:B:70:LEU:HD13	2.00	0.42
1:A:168:GLU:HA	1:A:171:LYS:CD	2.50	0.42
1:A:210:LEU:HB3	1:B:200:LEU:CD2	2.48	0.42
1:B:113:LYS:HG3	2:B:266:HOH:O	2.19	0.42
1:A:55:ARG:HA	2:A:247:HOH:O	2.19	0.42
1:B:9:GLU:O	1:B:10:PHE:C	2.63	0.41
1:A:89:ASN:HB3	1:A:220:ARG:HH22	1.79	0.41
1:A:138:LEU:HD13	1:A:171:LYS:N	2.34	0.41
1:B:141:ARG:HG3	1:B:141:ARG:HH11	1.84	0.41
1:B:134:HIS:O	1:B:138:LEU:HB2	2.20	0.41
1:B:138:LEU:C	1:B:138:LEU:HD22	2.45	0.41
1:A:8:ASP:O	1:A:9:GLU:C	2.61	0.41
1:B:55:ARG:HG3	1:B:55:ARG:HH11	1.86	0.41
1:B:216:LYS:N	1:B:216:LYS:HD3	2.34	0.41
1:A:160:VAL:O	1:A:164:VAL:HG13	2.20	0.41
1:A:225:SER:C	1:A:227:VAL:N	2.79	0.41
1:B:87:PHE:O	1:B:91:LEU:HG	2.21	0.41
1:B:118:ASP:HB2	1:B:119:PRO:HD3	2.03	0.41
1:B:141:ARG:HG3	1:B:141:ARG:NH1	2.36	0.41
1:A:200:LEU:HA	1:B:210:LEU:HD23	2.02	0.41
1:A:212:GLU:HG2	1:A:216:LYS:NZ	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:GLN:O	1:A:212:GLU:C	2.64	0.40
1:A:118:ASP:N	1:A:119:PRO:CD	2.85	0.40
1:A:189:VAL:CG1	1:B:86:THR:HG22	2.51	0.40
1:A:103:GLU:HB3	2:A:244:HOH:O	2.21	0.40
1:B:28:LEU:HD22	1:B:117:ILE:HG23	2.04	0.40
1:A:17:ILE:CD1	1:A:131:ILE:HD13	2.51	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	221/240 (92%)	208 (94%)	11 (5%)	2 (1%)	14	27
1	B	217/240 (90%)	190 (88%)	18 (8%)	9 (4%)	2	3
All	All	438/480 (91%)	398 (91%)	29 (7%)	11 (2%)	4	7

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	228	PRO
1	B	185	ASP
1	B	151	ARG
1	B	161	ARG
1	B	162	GLN
1	A	125	ASP
1	B	149	LYS
1	B	176	ARG
1	B	186	VAL
1	B	10	PHE
1	B	152	VAL

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	199/204 (98%)	187 (94%)	12 (6%)	17	36
1	B	197/204 (97%)	184 (93%)	13 (7%)	15	32
All	All	396/408 (97%)	371 (94%)	25 (6%)	16	34

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	SER
1	A	36	GLN
1	A	70	LEU
1	A	91	LEU
1	A	103	GLU
1	A	107	SER
1	A	130	GLU
1	A	184	ASN
1	A	200	LEU
1	A	210	LEU
1	A	217	LEU
1	A	228	PRO
1	B	17	ILE
1	B	21	ASN
1	B	23	VAL
1	B	49	ASN
1	B	105	LYS
1	B	115	THR
1	B	138	LEU
1	B	139	GLU
1	B	164	VAL
1	B	172	GLU
1	B	176	ARG
1	B	210	LEU
1	B	214	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	162	GLN
1	A	184	ASN
1	A	191	GLN
1	A	205	GLN
1	A	218	GLN
1	B	21	ASN
1	B	49	ASN
1	B	121	GLN
1	B	188	GLN
1	B	218	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	A	216/240 (90%)	0.26	5 (2%) 61 57	14, 27, 60, 72	0
1	B	214/240 (89%)	0.17	6 (2%) 55 50	10, 26, 59, 80	0
All	All	430/480 (89%)	0.22	11 (2%) 57 52	10, 27, 60, 80	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	228	PRO	4.0
1	A	58	VAL	2.8
1	B	151	ARG	2.7
1	A	9	GLU	2.7
1	B	160	VAL	2.6
1	B	153	GLY	2.5
1	A	227	VAL	2.4
1	B	152	VAL	2.1
1	A	160	VAL	2.1
1	B	11	LEU	2.1
1	B	231	GLU	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.