



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 26, 2021 – 10:14 AM EST

PDB ID : 6XF2
Title : Nesprin-1G (aa2070-2200)-FHOD1(aa1-339) complex, H. sapiens
Authors : Lim, S.M.; Schwartz, T.U.
Deposited on : 2020-06-15
Resolution : 7.11 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.16
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.16

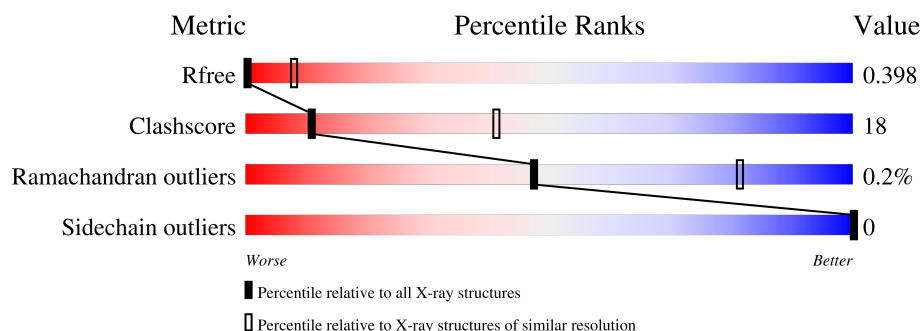
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 7.11 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}	130704	1004 (10.00-3.90)
Clashscore	141614	1069 (10.00-3.90)
Ramachandran outliers	138981	1002 (10.00-3.90)
Sidechain outliers	138945	1002 (10.00-3.86)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	131	
1	C	131	
2	B	321	
2	D	321	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6075 atoms, of which 0 are hydrogens and 0 are deuteriums.

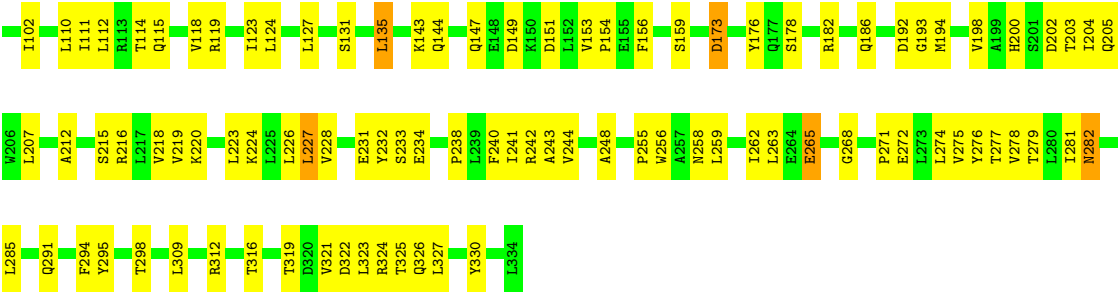
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Nesprin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	118	Total	C	N	O		0	0	0
			591	354	118	119				
1	C	109	Total	C	N	O	S	0	0	0
			554	331	109	111	3			

- Molecule 2 is a protein called FH1/FH2 domain-containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	321	Total	C	N	O	S	0	0	0
			2465	1571	414	472	8			
2	D	321	Total	C	N	O	S	0	0	0
			2465	1571	414	472	8			



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	209.37Å 142.05Å 58.05Å 90.00° 106.31° 90.00°	Depositor
Resolution (Å)	60.58 – 7.11 60.58 – 7.11	Depositor EDS
% Data completeness (in resolution range)	94.5 (60.58-7.11) 94.5 (60.58-7.11)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 7.40Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.340 , 0.397 0.339 , 0.398	Depositor DCC
R_{free} test set	244 reflections (10.26%)	wwPDB-VP
Wilson B-factor (Å ²)	464.1	Xtriage
Anisotropy	0.724	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 500.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.367 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.63	EDS
Total number of atoms	6075	wwPDB-VP
Average B, all atoms (Å ²)	320.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.69% 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	0.40	0/588	0.62	0/817
1	C	0.40	0/551	0.54	0/762
2	B	0.44	0/2510	0.82	1/3424 (0.0%)
2	D	0.46	0/2510	0.99	8/3424 (0.2%)
All	All	0.44	0/6159	0.86	9/8427 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1
2	D	0	1
All	All	0	2

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	282	ASN	CB-CG-OD1	21.41	164.41	121.60
2	D	282	ASN	CB-CG-ND2	-18.99	71.11	116.70
2	D	282	ASN	OD1-CG-ND2	-9.41	100.25	121.90
2	D	227	LEU	CA-CB-CG	5.83	128.71	115.30
2	D	135	LEU	CA-CB-CG	5.83	128.70	115.30
2	D	51	LEU	CA-CB-CG	-5.56	102.52	115.30
2	B	51	LEU	CA-CB-CG	-5.53	102.58	115.30
2	D	228	VAL	CG1-CB-CG2	-5.53	102.05	110.90
2	D	173	ASP	CB-CG-OD1	5.43	123.19	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	24	GLU	Sidechain
2	D	24	GLU	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	591	0	260	2	0
1	C	554	0	256	4	0
2	B	2465	0	2469	102	0
2	D	2465	0	2469	102	0
All	All	6075	0	5454	210	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 18.

All (210) 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:D:220:LYS:HE2	2:D:224:LYS:HE2	1.46	0.98
2:B:319:THR:HB	2:B:324:ARG:HH21	1.43	0.81
2:D:47:GLY:HA2	2:D:88:LEU:HD12	1.65	0.77
2:B:22:TYR:HA	2:B:112:LEU:HB3	1.68	0.75
2:B:58:VAL:HG23	2:B:61:LEU:HD23	1.69	0.74
2:B:19:ARG:HD2	2:B:40:ALA:HB3	1.67	0.74
2:D:319:THR:HB	2:D:324:ARG:HH21	1.53	0.73
2:B:216:ARG:NH2	2:B:270:ASP:OD1	2.18	0.72
2:B:72:ALA:HB2	2:B:115:GLN:HG2	1.71	0.71
2:D:69:GLU:OE1	2:D:69:GLU:N	2.27	0.68
2:B:124:LEU:HD21	2:B:161:GLY:HA2	1.76	0.67
2:B:47:GLY:HA2	2:B:88:LEU:HD12	1.78	0.66
2:D:131:SER:HA	2:D:135:LEU:HD22	1.78	0.65
2:B:69:GLU:OE1	2:B:69:GLU:N	2.29	0.64
1:C:2131:ALA:O	1:C:2135:MET:HG2	1.98	0.64
2:B:200:HIS:NE2	2:B:202:ASP:HB2	2.12	0.64
2:D:45:LEU:HD22	2:D:54:GLN:HB3	1.79	0.64
2:D:319:THR:HB	2:D:324:ARG:NH2	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:149:ASP:OD2	2:D:151:ASP:HB2	1.98	0.63
2:D:279:THR:HG22	2:D:326:GLN:HG2	1.79	0.63
2:D:241:ILE:HG23	2:D:256:TRP:HZ2	1.62	0.63
2:D:200:HIS:NE2	2:D:202:ASP:HB2	2.14	0.62
2:D:47:GLY:O	2:D:88:LEU:HB2	1.99	0.62
2:B:241:ILE:HG23	2:B:256:TRP:HZ2	1.64	0.62
2:B:216:ARG:NH2	2:B:272:GLU:OE2	2.33	0.62
2:D:263:LEU:HD13	2:D:274:LEU:HD12	1.82	0.62
2:D:58:VAL:HG23	2:D:61:LEU:HD23	1.82	0.62
2:D:23:LEU:N	2:D:112:LEU:O	2.31	0.61
2:D:200:HIS:CD2	2:D:202:ASP:HB2	2.36	0.61
2:B:73:LEU:HB3	2:B:82:LEU:HD12	1.83	0.59
2:D:154:PRO:HG3	2:D:192:ASP:CG	2.23	0.59
2:B:149:ASP:OD2	2:B:151:ASP:HB2	2.03	0.59
2:B:23:LEU:HD22	2:B:111:ILE:HG21	1.83	0.59
2:B:23:LEU:N	2:B:112:LEU:O	2.35	0.58
2:B:325:THR:OG1	2:B:326:GLN:OE1	2.21	0.58
2:B:277:THR:O	2:B:281:ILE:HG12	2.02	0.58
1:C:2132:LYS:HA	1:C:2135:MET:CG	2.34	0.58
2:D:325:THR:OG1	2:D:326:GLN:OE1	2.22	0.58
2:D:72:ALA:HB2	2:D:115:GLN:HG2	1.84	0.58
2:D:275:VAL:HA	2:D:278:VAL:HG12	1.86	0.58
2:D:321:VAL:O	2:D:325:THR:HG23	2.03	0.58
2:D:144:GLN:OE1	2:D:147:GLN:NE2	2.37	0.58
2:B:279:THR:HG22	2:B:326:GLN:HG2	1.85	0.57
2:D:204:ILE:HA	2:D:207:LEU:HD13	1.85	0.57
2:B:295:TYR:HA	2:B:298:THR:HG22	1.88	0.56
2:B:29:PHE:CE1	2:B:126:LYS:HD2	2.41	0.56
2:B:80:TYR:CE2	2:B:82:LEU:HD23	2.41	0.56
2:D:316:THR:O	2:D:324:ARG:NH2	2.38	0.56
2:D:212:ALA:HB2	2:D:258:ASN:OD1	2.05	0.56
2:D:277:THR:O	2:D:281:ILE:HG12	2.06	0.55
2:B:271:PRO:O	2:B:275:VAL:HG12	2.07	0.55
2:D:24:GLU:OE1	2:D:65:PRO:HG2	2.07	0.55
2:B:131:SER:HA	2:B:135:LEU:HD22	1.89	0.55
2:D:309:LEU:HG	2:D:312:ARG:NH2	2.21	0.55
2:B:203:THR:O	2:B:207:LEU:HD12	2.06	0.55
2:B:309:LEU:HG	2:B:312:ARG:NH2	2.22	0.55
2:B:50:PRO:HA	2:B:87:SER:HA	1.88	0.55
2:D:271:PRO:O	2:D:275:VAL:HG12	2.07	0.54
1:C:2132:LYS:HA	1:C:2135:MET:HG2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:240:PHE:O	2:D:244:VAL:HG23	2.07	0.54
2:B:275:VAL:HA	2:B:323:LEU:HD13	1.88	0.54
2:B:219:VAL:O	2:B:223:LEU:HD23	2.06	0.54
2:D:259:LEU:HA	2:D:262:ILE:HD12	1.89	0.54
2:D:327:LEU:O	2:D:330:TYR:HB3	2.09	0.53
2:D:76:SER:HB3	2:D:77:PRO:HD3	1.90	0.53
2:B:200:HIS:CD2	2:B:202:ASP:HB2	2.42	0.53
2:B:47:GLY:O	2:B:88:LEU:HB2	2.08	0.53
2:D:83:ASP:HB3	2:D:86:LEU:CD2	2.39	0.53
2:D:23:LEU:HD22	2:D:111:ILE:HG21	1.91	0.53
2:D:295:TYR:HA	2:D:298:THR:HG22	1.91	0.52
2:D:75:VAL:HA	2:D:110:LEU:CD2	2.40	0.52
2:D:80:TYR:CE2	2:D:82:LEU:HD23	2.45	0.52
2:B:73:LEU:CB	2:B:82:LEU:HD12	2.39	0.52
2:B:319:THR:HB	2:B:324:ARG:NH2	2.19	0.52
2:D:50:PRO:HA	2:D:87:SER:HA	1.92	0.52
2:B:265:GLU:O	2:B:265:GLU:HG2	2.10	0.52
2:D:265:GLU:O	2:D:265:GLU:HG2	2.10	0.52
2:D:19:ARG:HD2	2:D:40:ALA:HB3	1.91	0.52
2:B:205:GLN:HG3	2:B:243:ALA:HB1	1.92	0.51
2:B:76:SER:HB3	2:B:77:PRO:HD3	1.93	0.51
2:D:194:MET:HG2	2:D:232:TYR:CE1	2.46	0.51
2:B:51:LEU:O	2:B:55:ILE:HG13	2.10	0.51
2:B:173:ASP:O	2:B:177:GLN:HG3	2.11	0.51
2:B:321:VAL:O	2:B:325:THR:HG23	2.11	0.51
2:B:52:GLY:N	2:B:84:THR:O	2.44	0.51
2:D:240:PHE:HE2	2:D:256:TRP:CZ3	2.28	0.51
2:D:194:MET:O	2:D:198:VAL:HG23	2.11	0.51
2:B:58:VAL:HA	2:B:61:LEU:HB3	1.94	0.50
2:B:116:LEU:HD22	2:B:152:LEU:HD23	1.93	0.50
2:D:272:GLU:H	2:D:272:GLU:CD	2.15	0.49
2:D:233:SER:OG	2:D:234:GLU:N	2.45	0.49
2:D:291:GLN:O	2:D:294:PHE:HB3	2.12	0.49
2:D:51:LEU:O	2:D:55:ILE:HG13	2.12	0.49
2:D:83:ASP:HB3	2:D:86:LEU:HD22	1.94	0.49
2:B:173:ASP:OD1	2:B:176:TYR:HD2	1.94	0.49
2:D:153:VAL:HG11	2:D:193:GLY:H	1.78	0.49
2:D:205:GLN:HG3	2:D:243:ALA:HB1	1.95	0.49
2:B:220:LYS:O	2:B:224:LYS:HG3	2.13	0.48
2:B:233:SER:OG	2:B:234:GLU:N	2.46	0.48
2:D:70:ASP:HB3	2:D:118:VAL:HG21	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:178:SER:OG	2:D:182:ARG:NH2	2.46	0.48
2:B:220:LYS:HE2	2:B:224:LYS:HE2	1.96	0.48
2:D:278:VAL:HG22	2:D:282:ASN:OD1	2.13	0.48
2:B:45:LEU:HD22	2:B:54:GLN:HB3	1.96	0.48
2:B:240:PHE:CE2	2:B:244:VAL:HG21	2.49	0.47
2:B:288:LEU:O	2:B:288:LEU:HG	2.13	0.47
2:B:80:TYR:HE2	2:B:82:LEU:HD23	1.80	0.47
2:B:238:PRO:O	2:B:242:ARG:HG3	2.14	0.47
2:B:240:PHE:O	2:B:244:VAL:HG23	2.14	0.47
2:B:16:VAL:HG23	2:B:45:LEU:HB2	1.97	0.47
2:B:58:VAL:O	2:B:61:LEU:HB3	2.15	0.47
2:D:153:VAL:HG12	2:D:192:ASP:HB2	1.97	0.47
2:B:86:LEU:HB2	2:B:90:GLU:OE1	2.15	0.47
2:B:83:ASP:HB3	2:B:86:LEU:CD2	2.45	0.46
2:D:95:LEU:HB3	2:D:98:PHE:HB3	1.97	0.46
2:B:278:VAL:HG11	2:B:323:LEU:HD12	1.96	0.46
2:D:219:VAL:O	2:D:223:LEU:HD23	2.14	0.46
2:D:322:ASP:O	2:D:326:GLN:OE1	2.33	0.46
2:D:203:THR:O	2:D:207:LEU:HD12	2.15	0.46
2:D:227:LEU:O	2:D:231:GLU:HG3	2.15	0.46
2:D:248:ALA:CB	2:D:255:PRO:HA	2.45	0.46
1:A:2107:SER:C	1:A:2109:ILE:H	2.18	0.46
2:D:215:SER:O	2:D:219:VAL:HG23	2.15	0.46
2:D:259:LEU:O	2:D:263:LEU:HD23	2.15	0.46
2:D:47:GLY:HA2	2:D:88:LEU:CD1	2.38	0.46
2:B:327:LEU:HD12	2:B:330:TYR:HB3	1.98	0.46
2:B:93:GLU:HG2	2:B:94:MET:HG2	1.97	0.46
2:D:38:ARG:NH2	2:D:39:ARG:HD2	2.31	0.46
2:D:258:ASN:O	2:D:262:ILE:HG13	2.15	0.46
2:D:18:VAL:O	2:D:42:THR:HA	2.16	0.46
2:B:124:LEU:HD11	2:B:159:SER:O	2.16	0.45
2:D:238:PRO:O	2:D:242:ARG:HG3	2.15	0.45
2:D:80:TYR:HE2	2:D:82:LEU:HD23	1.82	0.45
2:B:154:PRO:HG3	2:B:192:ASP:CG	2.37	0.45
2:D:265:GLU:HB2	2:D:268:GLY:H	1.82	0.45
2:D:275:VAL:O	2:D:279:THR:HG23	2.16	0.45
2:B:19:ARG:HB2	2:B:109:THR:HG22	1.97	0.45
2:B:275:VAL:HA	2:B:278:VAL:HG12	1.99	0.45
2:B:205:GLN:HG2	2:B:247:VAL:HG23	1.97	0.45
2:D:75:VAL:HA	2:D:110:LEU:HD22	1.98	0.45
2:D:309:LEU:HG	2:D:312:ARG:HH22	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:248:ALA:CB	2:B:255:PRO:HA	2.47	0.45
2:B:309:LEU:HG	2:B:312:ARG:HH22	1.82	0.45
2:B:322:ASP:O	2:B:326:GLN:OE1	2.35	0.45
2:B:327:LEU:O	2:B:330:TYR:HB3	2.16	0.45
2:B:83:ASP:HB3	2:B:86:LEU:HD22	1.99	0.45
2:B:128:TYR:CE1	2:B:164:CYS:HA	2.51	0.45
2:B:197:VAL:HG12	2:B:229:PHE:CE1	2.52	0.45
2:B:99:TYR:HA	2:B:102:ILE:HD12	2.00	0.44
1:C:2132:LYS:HA	1:C:2135:MET:HG3	1.98	0.44
2:D:173:ASP:OD1	2:D:176:TYR:HD2	2.00	0.44
1:A:2095:LYS:HA	1:A:2098:VAL:CG2	2.47	0.44
2:B:212:ALA:HB2	2:B:258:ASN:OD1	2.18	0.44
2:B:275:VAL:O	2:B:279:THR:HG23	2.17	0.44
2:B:285:LEU:HD23	2:B:285:LEU:HA	1.83	0.44
2:D:173:ASP:C	2:D:173:ASP:OD1	2.55	0.44
2:D:274:LEU:HD13	2:D:274:LEU:HA	1.85	0.44
2:D:224:LYS:O	2:D:227:LEU:HB2	2.18	0.44
2:D:143:LYS:O	2:D:186:GLN:HG2	2.18	0.44
2:D:240:PHE:HE2	2:D:256:TRP:CH2	2.36	0.44
2:B:24:GLU:OE1	2:B:65:PRO:HG2	2.18	0.43
2:B:251:THR:HG23	2:B:253:ALA:H	1.82	0.43
2:D:153:VAL:O	2:D:156:PHE:HB3	2.18	0.43
2:B:272:GLU:H	2:B:272:GLU:CD	2.17	0.43
2:D:98:PHE:CZ	2:D:102:ILE:HD11	2.53	0.43
2:B:308:ALA:O	2:B:312:ARG:HG2	2.19	0.43
2:B:325:THR:O	2:B:328:VAL:HG22	2.19	0.43
2:D:215:SER:HB3	2:D:218:VAL:HB	2.00	0.43
2:B:151:ASP:O	2:B:154:PRO:HD2	2.17	0.43
2:B:173:ASP:C	2:B:173:ASP:OD1	2.56	0.43
2:B:272:GLU:OE1	2:B:272:GLU:N	2.42	0.43
2:D:114:THR:O	2:D:119:ARG:NH2	2.38	0.43
2:B:263:LEU:HD13	2:B:274:LEU:HD12	2.01	0.43
2:B:47:GLY:HA2	2:B:88:LEU:HB2	2.01	0.43
2:D:224:LYS:HD3	2:D:276:TYR:CE2	2.54	0.43
2:D:226:LEU:HD23	2:D:226:LEU:HA	1.87	0.42
2:B:240:PHE:HE2	2:B:256:TRP:CZ3	2.37	0.42
2:D:216:ARG:NH2	2:D:272:GLU:OE2	2.52	0.42
2:D:86:LEU:HD12	2:D:90:GLU:HB2	2.00	0.42
2:B:319:THR:HG23	2:B:323:LEU:HD23	2.01	0.42
2:D:58:VAL:HA	2:D:61:LEU:HB3	2.01	0.42
2:D:124:LEU:HD11	2:D:159:SER:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:275:VAL:CA	2:D:323:LEU:HD13	2.50	0.42
2:D:282:ASN:O	2:D:285:LEU:HB2	2.20	0.42
2:B:98:PHE:CZ	2:B:102:ILE:HD11	2.54	0.42
2:B:194:MET:O	2:B:198:VAL:HG23	2.20	0.42
2:B:47:GLY:HA2	2:B:88:LEU:CD1	2.46	0.42
2:B:259:LEU:O	2:B:263:LEU:HD23	2.20	0.42
2:B:75:VAL:HA	2:B:110:LEU:CD2	2.50	0.41
2:D:49:LEU:HB2	2:D:54:GLN:NE2	2.35	0.41
2:B:226:LEU:HD23	2:B:226:LEU:HA	1.86	0.41
2:D:204:ILE:H	2:D:204:ILE:HD12	1.85	0.41
2:B:275:VAL:CA	2:B:323:LEU:HD13	2.50	0.41
2:D:58:VAL:O	2:D:61:LEU:HB3	2.20	0.41
2:D:275:VAL:HA	2:D:323:LEU:HD13	2.02	0.41
2:D:92:ARG:HG3	2:D:99:TYR:CE2	2.56	0.41
2:B:19:ARG:O	2:B:109:THR:HA	2.20	0.41
2:B:282:ASN:ND2	2:B:326:GLN:HB3	2.36	0.41
2:B:29:PHE:HE2	2:B:138:SER:HB3	1.85	0.41
2:D:25:ASP:CG	2:D:114:THR:H	2.24	0.41
2:B:123:ILE:O	2:B:127:LEU:HG	2.21	0.41
2:B:147:GLN:HG2	2:B:186:GLN:CD	2.41	0.41
2:B:178:SER:OG	2:B:182:ARG:NH2	2.53	0.41
2:B:288:LEU:HD21	2:B:294:PHE:HD1	1.86	0.41
2:D:200:HIS:HE2	2:D:202:ASP:HB2	1.82	0.41
2:D:22:TYR:HA	2:D:112:LEU:HB3	2.03	0.41
2:D:73:LEU:CB	2:D:82:LEU:HD12	2.51	0.41
2:B:204:ILE:HD12	2:B:204:ILE:H	1.85	0.41
2:D:123:ILE:O	2:D:127:LEU:HG	2.20	0.41
2:D:73:LEU:HB3	2:D:82:LEU:HD12	2.02	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	112/131 (86%)	108 (96%)	4 (4%)	0	100	100
1	C	103/131 (79%)	99 (96%)	4 (4%)	0	100	100
2	B	319/321 (99%)	309 (97%)	9 (3%)	1 (0%)	41	77
2	D	319/321 (99%)	309 (97%)	9 (3%)	1 (0%)	41	77
All	All	853/904 (94%)	825 (97%)	26 (3%)	2 (0%)	47	81

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	265	GLU
2	D	265	GLU

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	2/124 (2%)	2 (100%)	0	100	100
1	C	5/124 (4%)	5 (100%)	0	100	100
2	B	269/276 (98%)	269 (100%)	0	100	100
2	D	269/276 (98%)	269 (100%)	0	100	100
All	All	545/800 (68%)	545 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	282	ASN

5.3.3 RNA ⓘ

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 monosaccharides 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](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.