



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:16 PM UTC

PDB ID : 4NKH / pdb_00004nkh
Title : Crystal structure of SspH1 LRR domain
Authors : Keszei, A.F.A.; Xiaojing, T.; McCormick, C.; Zeqiraj, E.; Rohde, J.R.; Tyers, M.; Sicheri, F.
Deposited on : 2013-11-12
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

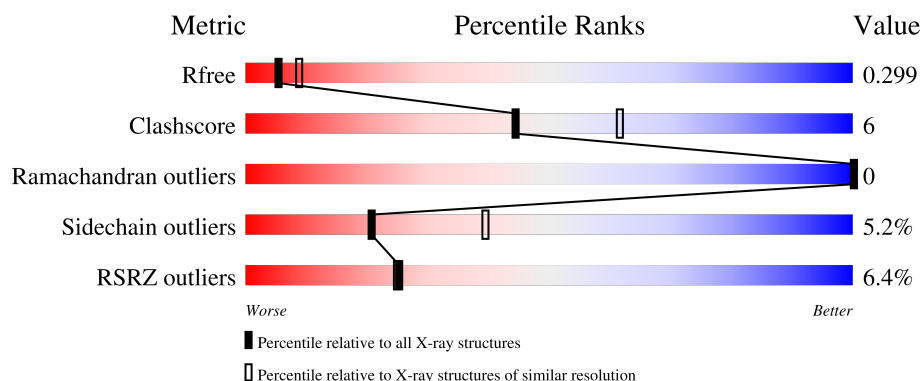
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	239	3% 78% 20% ..
1	B	239	3% 87% 11% ..
1	C	239	7% 79% 17% ..
1	D	239	10% 78% 18% ..
1	E	239	2% 86% 13% .

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Mol	Chain	Length	Quality of chain
1	F	239	13% 79% 18% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 10455 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 E3 ubiquitin-protein ligase sspH1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	236	Total	C	N	O	S	3	0	0
			1752	1103	309	335	5			
1	B	237	Total	C	N	O	S	0	0	0
			1778	1121	310	342	5			
1	C	233	Total	C	N	O	S	0	0	0
			1722	1089	302	327	4			
1	D	236	Total	C	N	O	S	0	0	0
			1720	1089	299	327	5			
1	E	239	Total	C	N	O	S	0	0	0
			1787	1124	319	339	5			
1	F	235	Total	C	N	O	S	0	0	0
			1696	1066	300	325	5			

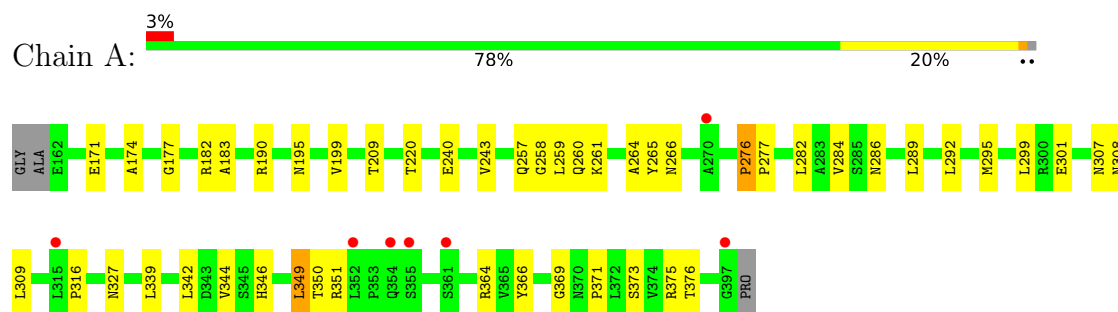
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	160	GLY	-	expression tag	UNP D0ZVG2
B	160	GLY	-	expression tag	UNP D0ZVG2
C	160	GLY	-	expression tag	UNP D0ZVG2
D	160	GLY	-	expression tag	UNP D0ZVG2
E	160	GLY	-	expression tag	UNP D0ZVG2
F	160	GLY	-	expression tag	UNP D0ZVG2

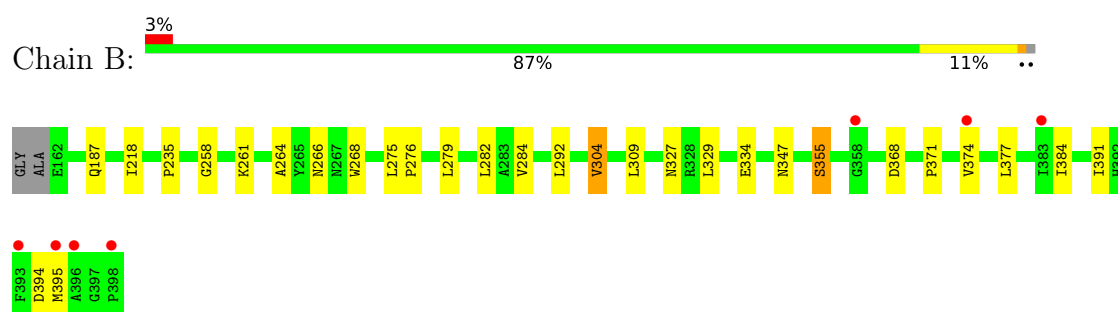
3 Residue-property plots [i](#)

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.

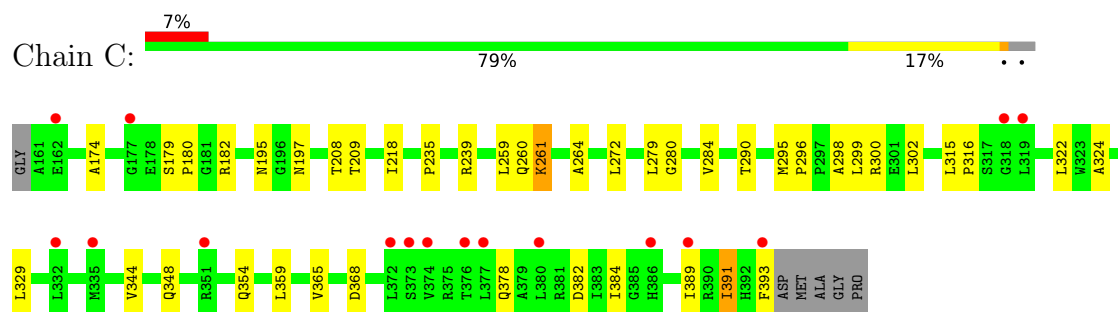
- Molecule 1: E3 ubiquitin-protein ligase sspH1



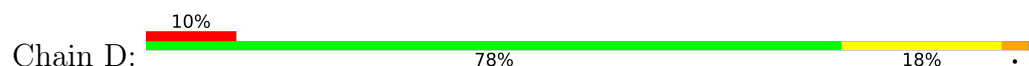
- Molecule 1: E3 ubiquitin-protein ligase sspH1

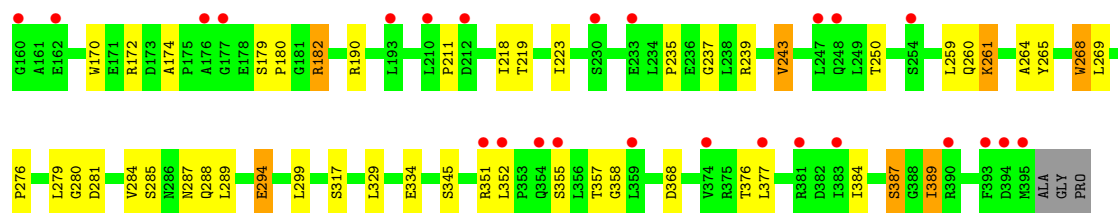


- Molecule 1: E3 ubiquitin-protein ligase sspH1

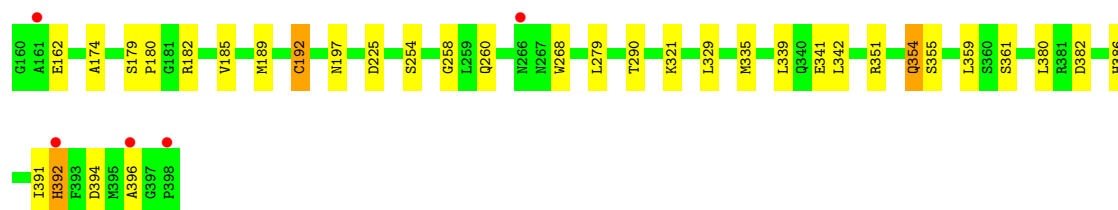
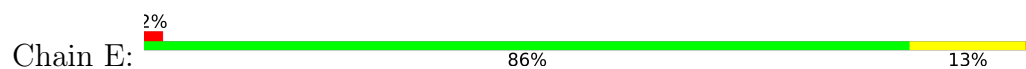


- Molecule 1: E3 ubiquitin-protein ligase sspH1

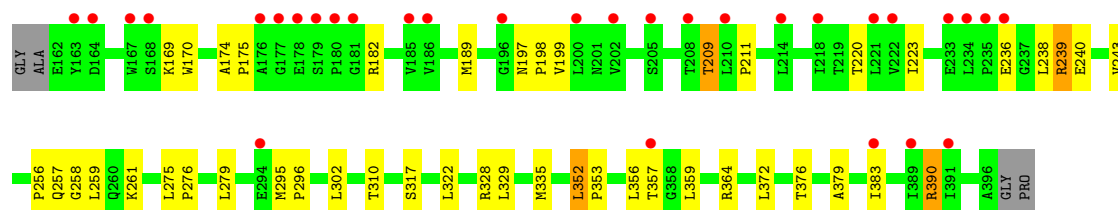
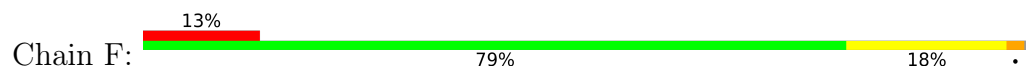




● Molecule 1: E3 ubiquitin-protein ligase sspH1



● Molecule 1: E3 ubiquitin-protein ligase sspH1



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	114.93Å 114.93Å 252.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.86 – 2.75 49.86 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.86-2.75) 98.8 (49.86-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.243 , 0.280 (Not available) , 0.299	Depositor DCC
R_{free} test set	2247 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	60.1	Xtriage
Anisotropy	0.437	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 60.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10455	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% 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	0.48	1/1793 (0.1%)	0.87	2/2464 (0.1%)
1	B	0.46	0/1820	0.85	0/2500
1	C	0.45	0/1763	0.87	0/2426
1	D	0.44	0/1761	0.84	0/2424
1	E	0.47	0/1829	0.80	0/2511
1	F	0.46	0/1734	0.86	0/2388
All	All	0.46	1/10700 (0.0%)	0.85	2/14713 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	190	ARG	NE-CZ	-9.14	1.23	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	276	PRO	CA-C-N	7.51	129.22	119.84
1	A	276	PRO	C-N-CA	7.51	129.22	119.84

There are no chirality outliers.

There are no planarity outliers.

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	1752	0	1726	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1778	0	1769	19	0
1	C	1722	0	1704	25	0
1	D	1720	0	1680	24	0
1	E	1787	0	1778	15	1
1	F	1696	0	1636	24	0
All	All	10455	0	10293	125	1

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 6.

All (125) 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:B:334:GLU:HG3	1:B:355:SER:OG	1.85	0.77
1:A:309:LEU:HB2	1:A:327:ASN:HD22	1.50	0.77
1:C:378:GLN:O	1:C:382:ASP:HB2	1.88	0.74
1:A:266:ASN:HD21	1:A:286:ASN:HD22	1.37	0.71
1:A:257:GLN:HB3	1:F:257:GLN:HB3	1.74	0.69
1:F:223:ILE:HB	1:F:243:VAL:HG12	1.77	0.66
1:A:265:TYR:HB3	1:A:266:ASN:HD22	1.61	0.66
1:F:364:ARG:HG2	1:F:390:ARG:HE	1.63	0.64
1:B:264:ALA:HB3	1:B:284:VAL:HG12	1.81	0.63
1:B:384:ILE:HD11	1:B:391:ILE:HB	1.80	0.62
1:B:279:LEU:HD21	1:B:282:LEU:HD13	1.83	0.60
1:B:371:PRO:HA	1:B:395:MET:HE3	1.83	0.59
1:D:179:SER:HB2	1:D:180:PRO:HD3	1.84	0.59
1:A:292:LEU:HG	1:A:309:LEU:HD21	1.84	0.58
1:C:264:ALA:HB3	1:C:284:VAL:HG12	1.86	0.57
1:B:258:GLY:HA2	1:E:258:GLY:HA2	1.87	0.57
1:A:309:LEU:HB2	1:A:327:ASN:ND2	2.21	0.56
1:F:197:ASN:OD1	1:F:198:PRO:HD2	2.05	0.56
1:E:260:GLN:HA	1:E:279:LEU:HA	1.89	0.55
1:D:264:ALA:HB3	1:D:284:VAL:HG12	1.88	0.55
1:A:266:ASN:ND2	1:A:286:ASN:HD22	2.05	0.55
1:D:260:GLN:HA	1:D:279:LEU:HA	1.88	0.54
1:F:372:LEU:HD22	1:F:376:THR:HG21	1.88	0.54
1:C:354:GLN:H	1:C:354:GLN:CD	2.15	0.54
1:D:261:LYS:HG3	1:D:281:ASP:HB3	1.90	0.54
1:C:195:ASN:OD1	1:C:197:ASN:HB2	2.08	0.53
1:A:351:ARG:HA	1:A:376:THR:HG21	1.91	0.53
1:A:364:ARG:HD3	1:A:366:TYR:OH	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:365:VAL:HB	1:C:391:ILE:HG23	1.91	0.53
1:E:174:ALA:HB2	1:E:182:ARG:HD2	1.89	0.53
1:A:177:GLY:H	1:B:266:ASN:HD21	1.56	0.53
1:A:349:LEU:O	1:A:371:PRO:O	2.27	0.53
1:A:199:VAL:HG22	1:A:220:THR:HB	1.91	0.53
1:F:352:LEU:HB2	1:F:353:PRO:HD2	1.91	0.53
1:F:357:THR:HG22	1:F:383:ILE:HG12	1.90	0.52
1:D:259:LEU:HB3	1:D:276:PRO:HG2	1.91	0.52
1:E:321:LYS:HG2	1:E:341:GLU:HB3	1.92	0.52
1:A:258:GLY:HA2	1:F:258:GLY:HA2	1.90	0.52
1:B:187:GLN:HE22	1:C:261:LYS:HZ2	1.58	0.52
1:A:174:ALA:HB2	1:A:182:ARG:HD2	1.92	0.52
1:A:350:THR:HG22	1:A:371:PRO:HB2	1.92	0.51
1:D:268:TRP:HA	1:D:268:TRP:CE3	2.45	0.51
1:F:302:LEU:HD23	1:F:322:LEU:HD13	1.92	0.51
1:D:174:ALA:HB2	1:D:182:ARG:CD	2.40	0.51
1:D:218:ILE:HD11	1:D:235:PRO:HG2	1.93	0.51
1:E:185:VAL:HG12	1:E:189:MET:HE2	1.92	0.51
1:E:192:CYS:HB2	1:E:197:ASN:HB3	1.93	0.50
1:C:218:ILE:HD11	1:C:235:PRO:HG2	1.93	0.50
1:B:327:ASN:HB2	1:B:347:ASN:HD21	1.77	0.49
1:B:292:LEU:HD21	1:B:304:VAL:HG11	1.95	0.49
1:B:329:LEU:H	1:B:347:ASN:HD22	1.60	0.49
1:C:280:GLY:HA2	1:C:299:LEU:HA	1.95	0.49
1:F:310:THR:HG22	1:F:328:ARG:HB2	1.93	0.49
1:D:280:GLY:HA2	1:D:299:LEU:HA	1.95	0.49
1:B:218:ILE:CG1	1:B:235:PRO:HG2	2.43	0.49
1:D:170:TRP:CE3	1:D:211:PRO:HG3	2.48	0.49
1:D:334:GLU:HB3	1:D:355:SER:OG	2.13	0.48
1:A:339:LEU:HD21	1:A:342:LEU:HB2	1.96	0.48
1:C:384:ILE:HG21	1:C:389:ILE:HG23	1.96	0.48
1:F:356:LEU:HA	1:F:359:LEU:HD13	1.96	0.48
1:D:345:SER:HB2	1:D:368:ASP:O	2.13	0.48
1:C:239:ARG:HG2	1:C:260:GLN:HG2	1.95	0.47
1:C:179:SER:N	1:C:180:PRO:HD2	2.29	0.47
1:B:284:VAL:HG23	1:B:304:VAL:HG13	1.97	0.47
1:F:279:LEU:HB3	1:F:296:PRO:HG2	1.95	0.47
1:D:387:SER:HB2	1:D:389:ILE:HG22	1.97	0.47
1:F:259:LEU:HB3	1:F:276:PRO:HG2	1.97	0.47
1:F:238:LEU:HB3	1:F:256:PRO:HG2	1.97	0.46
1:C:384:ILE:HG21	1:C:389:ILE:CG2	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:ALA:HB3	1:A:284:VAL:HG12	1.97	0.46
1:F:170:TRP:HE1	1:F:209:THR:HG23	1.81	0.46
1:D:174:ALA:CB	1:D:182:ARG:HD2	2.46	0.46
1:F:189:MET:HE3	1:F:211:PRO:HG2	1.98	0.46
1:B:304:VAL:HG11	1:B:309:LEU:HD11	1.98	0.45
1:A:282:LEU:HD23	1:A:295:MET:HE1	1.97	0.45
1:D:351:ARG:HA	1:D:376:THR:HG21	1.98	0.45
1:A:240:GLU:HG2	1:A:261:LYS:HB3	1.97	0.45
1:A:346:HIS:HA	1:A:369:GLY:O	2.16	0.45
1:D:223:ILE:HB	1:D:243:VAL:HG12	1.98	0.45
1:F:240:GLU:HG2	1:F:261:LYS:HB3	1.98	0.45
1:A:276:PRO:HA	1:A:277:PRO:HD2	1.74	0.45
1:B:187:GLN:HE22	1:C:261:LYS:NZ	2.15	0.45
1:C:302:LEU:HD23	1:C:322:LEU:HD13	2.00	0.45
1:D:287:ASN:HB3	1:D:288:GLN:H	1.64	0.45
1:E:392:HIS:N	1:E:392:HIS:ND1	2.65	0.44
1:E:354:GLN:HE21	1:E:354:GLN:H	1.66	0.44
1:A:373:SER:HB2	1:A:376:THR:HG23	2.00	0.43
1:D:250:THR:HA	1:D:268:TRP:O	2.17	0.43
1:C:368:ASP:HA	1:C:393:PHE:HD1	1.84	0.43
1:D:170:TRP:CD2	1:D:211:PRO:HG3	2.54	0.43
1:B:368:ASP:HA	1:B:394:ASP:HB2	2.00	0.43
1:F:275:LEU:HA	1:F:276:PRO:HD3	1.94	0.43
1:C:272:LEU:HD21	1:C:284:VAL:HG11	2.01	0.43
1:C:298:ALA:HB2	1:D:237:GLY:HA3	1.99	0.43
1:A:295:MET:HB2	1:A:316:PRO:HD3	2.02	0.42
1:A:171:GLU:OE2	1:A:183:ALA:HA	2.18	0.42
1:F:182:ARG:NH2	1:F:209:THR:HG22	2.34	0.42
1:A:260:GLN:NE2	1:F:239:ARG:HH11	2.17	0.42
1:E:382:ASP:O	1:E:386:HIS:CD2	2.73	0.42
1:B:275:LEU:HA	1:B:276:PRO:HD3	1.90	0.42
1:C:174:ALA:HB2	1:C:182:ARG:HD2	2.01	0.42
1:A:259:LEU:HB3	1:A:276:PRO:HG3	2.01	0.41
1:A:289:LEU:HD12	1:A:307:ASN:ND2	2.35	0.41
1:C:260:GLN:HB3	1:D:239:ARG:CZ	2.50	0.41
1:D:357:THR:HG23	1:D:358:GLY:N	2.35	0.41
1:C:295:MET:HA	1:C:296:PRO:HD2	1.80	0.41
1:C:324:ALA:HB3	1:C:344:VAL:HG12	2.03	0.41
1:E:179:SER:HB2	1:E:180:PRO:HD3	2.02	0.41
1:A:266:ASN:ND2	1:C:180:PRO:HD3	2.34	0.41
1:E:354:GLN:H	1:E:354:GLN:NE2	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:379:ALA:O	1:F:383:ILE:HG13	2.21	0.41
1:C:208:THR:HG22	1:C:209:THR:HG23	2.02	0.41
1:C:315:LEU:HA	1:C:316:PRO:HD3	1.94	0.41
1:B:374:VAL:HA	1:B:377:LEU:HD12	2.03	0.41
1:C:259:LEU:HD23	1:C:279:LEU:HD13	2.03	0.41
1:D:268:TRP:HA	1:D:268:TRP:HE3	1.86	0.41
1:E:339:LEU:HD21	1:E:342:LEU:HB2	2.03	0.41
1:B:268:TRP:CE3	1:B:268:TRP:HA	2.56	0.40
1:F:295:MET:HA	1:F:296:PRO:HD3	1.97	0.40
1:D:294:GLU:N	1:D:294:GLU:CD	2.80	0.40
1:E:268:TRP:CE3	1:E:268:TRP:HA	2.56	0.40
1:E:380:LEU:HD22	1:E:391:ILE:HG21	2.03	0.40
1:E:394:ASP:C	1:E:396:ALA:H	2.29	0.40
1:F:174:ALA:HA	1:F:175:PRO:HD2	1.97	0.40
1:F:199:VAL:HG23	1:F:220:THR:HB	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:290:THR:OG1	1:E:290:THR:OG1[2_555]	1.93	0.27

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	234/239 (98%)	217 (93%)	17 (7%)	0	100	100
1	B	235/239 (98%)	221 (94%)	14 (6%)	0	100	100
1	C	231/239 (97%)	210 (91%)	21 (9%)	0	100	100
1	D	234/239 (98%)	220 (94%)	14 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	237/239 (99%)	221 (93%)	16 (7%)	0	100	100
1	F	233/239 (98%)	212 (91%)	21 (9%)	0	100	100
All	All	1404/1434 (98%)	1301 (93%)	103 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	191/203 (94%)	182 (95%)	9 (5%)	23	45
1	B	198/203 (98%)	195 (98%)	3 (2%)	57	74
1	C	188/203 (93%)	181 (96%)	7 (4%)	30	54
1	D	183/203 (90%)	164 (90%)	19 (10%)	7	13
1	E	196/203 (97%)	184 (94%)	12 (6%)	17	31
1	F	178/203 (88%)	169 (95%)	9 (5%)	21	40
All	All	1134/1218 (93%)	1075 (95%)	59 (5%)	21	39

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	195	ASN
1	A	209	THR
1	A	243	VAL
1	A	299	LEU
1	A	301	GLU
1	A	308	ASN
1	A	344	VAL
1	A	349	LEU
1	A	375	ARG
1	B	261	LYS
1	B	304	VAL
1	B	355	SER

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Mol	Chain	Res	Type
1	C	261	LYS
1	C	290	THR
1	C	300	ARG
1	C	329	LEU
1	C	348	GLN
1	C	359	LEU
1	C	391	ILE
1	D	172	ARG
1	D	182	ARG
1	D	190	ARG
1	D	219	THR
1	D	243	VAL
1	D	261	LYS
1	D	265	TYR
1	D	268	TRP
1	D	269	LEU
1	D	285	SER
1	D	289	LEU
1	D	294	GLU
1	D	317	SER
1	D	329	LEU
1	D	352	LEU
1	D	377	LEU
1	D	384	ILE
1	D	387	SER
1	D	389	ILE
1	E	162	GLU
1	E	192	CYS
1	E	225	ASP
1	E	254	SER
1	E	329	LEU
1	E	335	MET
1	E	351	ARG
1	E	354	GLN
1	E	355	SER
1	E	359	LEU
1	E	361	SER
1	E	392	HIS
1	F	169	LYS
1	F	209	THR
1	F	236	GLU
1	F	239	ARG

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Mol	Chain	Res	Type
1	F	317	SER
1	F	329	LEU
1	F	335	MET
1	F	352	LEU
1	F	390	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	260	GLN
1	A	266	ASN
1	A	286	ASN
1	A	327	ASN
1	B	260	GLN
1	B	326	ASN
1	B	346	HIS
1	B	347	ASN
1	C	260	GLN
1	C	286	ASN
1	C	320	GLN
1	C	348	GLN
1	C	392	HIS
1	D	386	HIS
1	E	257	GLN
1	E	326	ASN
1	E	346	HIS
1	E	354	GLN
1	E	386	HIS
1	F	257	GLN
1	F	266	ASN
1	F	308	ASN
1	F	378	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	236/239 (98%)	0.46	7 (2%) 52 50	53, 75, 97, 110	1 (0%)
1	B	237/239 (99%)	0.18	7 (2%) 52 50	42, 58, 99, 123	0
1	C	233/239 (97%)	0.70	16 (6%) 23 22	55, 79, 104, 124	0
1	D	236/239 (98%)	0.93	25 (10%) 11 10	56, 85, 110, 136	0
1	E	239/239 (100%)	0.19	5 (2%) 63 61	37, 56, 95, 108	0
1	F	235/239 (98%)	1.00	31 (13%) 7 6	55, 79, 130, 169	0
All	All	1416/1434 (98%)	0.57	91 (6%) 25 25	37, 73, 112, 169	1 (0%)

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	178	GLU	4.2
1	F	176	ALA	4.0
1	F	179	SER	3.8
1	F	214	LEU	3.5
1	C	162	GLU	3.5
1	F	167	TRP	3.5
1	C	351	ARG	3.5
1	D	212	ASP	3.5
1	C	374	VAL	3.4
1	A	352	LEU	3.3
1	A	270	ALA	3.3
1	B	398	PRO	3.2
1	A	354	GLN	3.2
1	D	162	GLU	3.2
1	D	176	ALA	3.1
1	C	177	GLY	3.1
1	F	200	LEU	3.0
1	D	351	ARG	2.9
1	C	376	THR	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	391	ILE	2.9
1	F	222	VAL	2.9
1	B	396	ALA	2.7
1	A	361	SER	2.7
1	B	374	VAL	2.7
1	C	393	PHE	2.7
1	D	393	PHE	2.7
1	F	357	THR	2.7
1	D	354	GLN	2.7
1	C	377	LEU	2.7
1	F	185	VAL	2.7
1	F	202	VAL	2.7
1	D	352	LEU	2.6
1	F	205	SER	2.6
1	D	377	LEU	2.6
1	D	381	ARG	2.5
1	F	164	ASP	2.5
1	C	389	ILE	2.5
1	D	355	SER	2.4
1	F	177	GLY	2.4
1	F	163	TYR	2.4
1	F	186	VAL	2.4
1	B	393	PHE	2.4
1	D	383	ILE	2.4
1	D	374	VAL	2.4
1	B	358	GLY	2.4
1	F	233	GLU	2.3
1	F	236	GLU	2.3
1	F	196	GLY	2.3
1	C	332	LEU	2.3
1	F	208	THR	2.3
1	F	235	PRO	2.3
1	D	210	LEU	2.3
1	F	234	LEU	2.3
1	E	392	HIS	2.3
1	D	248	GLN	2.3
1	B	383	ILE	2.3
1	B	395	MET	2.3
1	D	230	SER	2.3
1	E	398	PRO	2.3
1	F	181	GLY	2.3
1	C	318	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	160	GLY	2.2
1	C	319	LEU	2.2
1	A	397	GLY	2.2
1	F	210	LEU	2.2
1	A	315	LEU	2.2
1	A	355	SER	2.2
1	C	380	LEU	2.2
1	D	254	SER	2.2
1	D	177	GLY	2.2
1	F	221	LEU	2.1
1	F	168	SER	2.1
1	D	193	LEU	2.1
1	D	359	LEU	2.1
1	F	383	ILE	2.1
1	D	233	GLU	2.1
1	C	386	HIS	2.1
1	D	394	ASP	2.1
1	C	373	SER	2.1
1	D	390	ARG	2.1
1	E	161	ALA	2.1
1	E	266	ASN	2.1
1	F	389	ILE	2.1
1	E	396	ALA	2.0
1	C	335	MET	2.0
1	C	372	LEU	2.0
1	F	294	GLU	2.0
1	D	395	MET	2.0
1	F	180	PRO	2.0
1	D	247	LEU	2.0
1	F	218	ILE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.