



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:02 PM GMT

PDB ID : 4NKH
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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

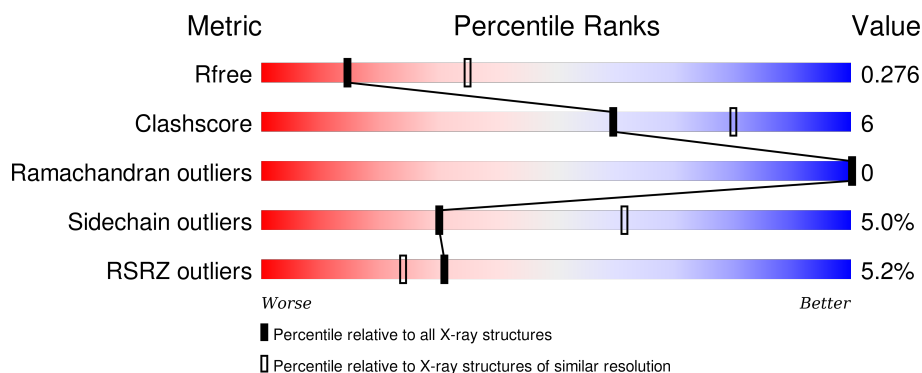
1 Overall quality at a glance

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}	91344	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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. 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	
1	B	239	
1	C	239	
1	D	239	
1	E	239	

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Mol	Chain	Length	Quality of chain
1	F	239	<div><div></div><div>15%</div><div></div><div>79%</div><div></div><div>18%</div><div></div><div>• •</div></div>

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			

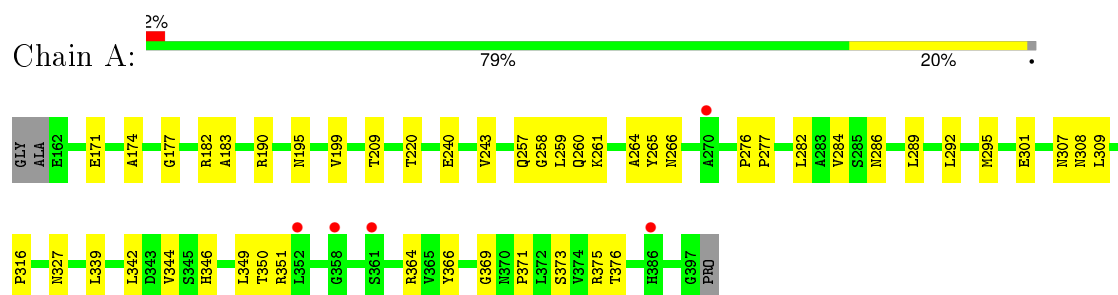
There are 6 discrepancies between the modelled and reference sequences:

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

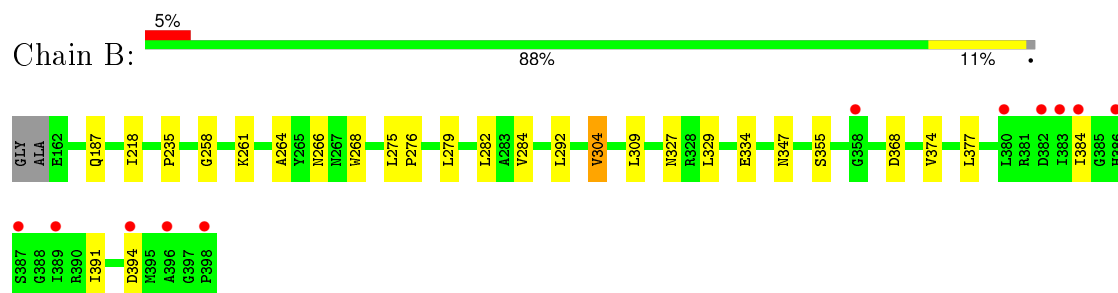
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA 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.

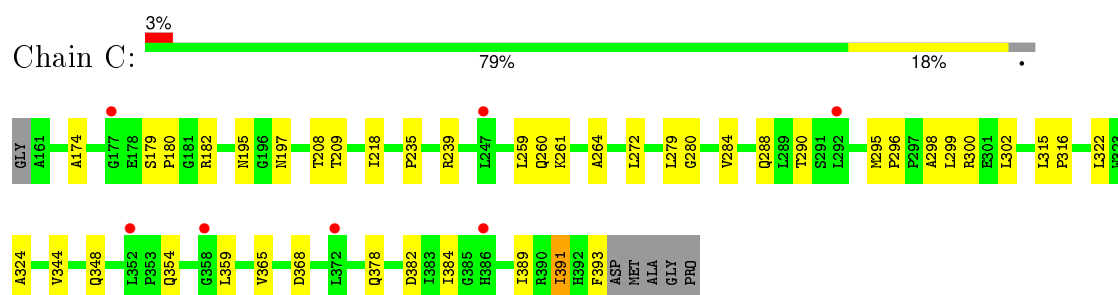
- Molecule 1: E3 ubiquitin-protein ligase sspH1



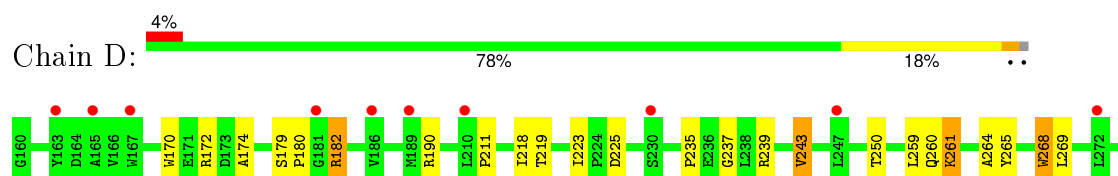
- Molecule 1: E3 ubiquitin-protein ligase sspH1



- Molecule 1: E3 ubiquitin-protein ligase sspH1

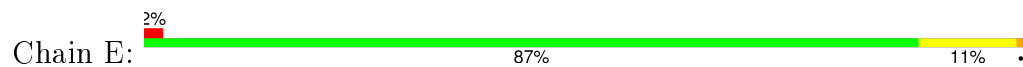


- Molecule 1: E3 ubiquitin-protein ligase sspH1

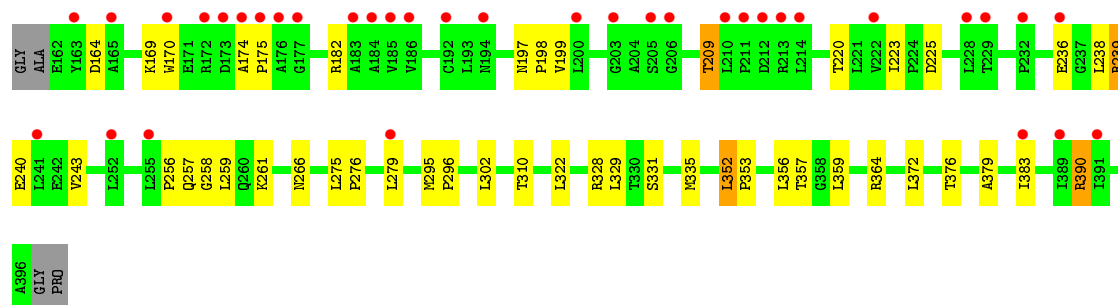
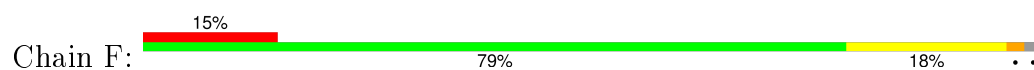




- Molecule 1: E3 ubiquitin-protein ligase sspH1



- 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 49.81 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.86-2.75) 98.9 (49.81-2.75)	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 0.239 , 0.276	Depositor DCC
R_{free} test set	2247 reflections (5.34%)	DCC
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 , 41.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	5 of 44291 reflections (0.011%)	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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.35	1/1793 (0.1%)	0.58	1/2464 (0.0%)
1	B	0.31	0/1820	0.56	0/2500
1	C	0.31	0/1763	0.53	0/2426
1	D	0.31	0/1761	0.53	0/2424
1	E	0.32	0/1829	0.55	0/2511
1	F	0.30	0/1734	0.55	0/2388
All	All	0.32	1/10700 (0.0%)	0.55	1/14713 (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	-7.73	1.23	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	ARG	NE-CZ-NH2	-6.42	117.09	120.30

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
1	B	1778	0	1769	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1722	0	1704	24	0
1	D	1720	0	1680	23	0
1	E	1787	0	1778	13	1
1	F	1696	0	1636	23	0
All	All	10455	0	10293	119	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 (119) 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: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: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:364:ARG:HD3	1:A:366:TYR:OH	2.09	0.53
1:A:351:ARG:HA	1:A:376:THR:HG21	1.91	0.53
1:E:174:ALA:HB2	1:E:182:ARG:HD2	1.89	0.53
1:C:365:VAL:HB	1:C:391:ILE:HG23	1.91	0.53
1:A:177:GLY:H	1:B:266:ASN:HD21	1.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:CE3	1:D:268:TRP:HA	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: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:C:354:GLN:CD	1:C:354:GLN:H	2.15	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:F:310:THR:HG22	1:F:328:ARG:HB2	1.93	0.49
1:C:280:GLY:HA2	1:C:299:LEU:HA	1.95	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
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:B:304:VAL:HG11	1:B:309:LEU:HD11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:282:LEU:HD23	1:A:295:MET:HE1	1.98	0.45
1:A:276:PRO:HA	1:A:277:PRO:HD2	1.74	0.45
1:D:287:ASN:HB3	1:D:288:GLN:H	1.64	0.45
1:C:302:LEU:HD23	1:C:322:LEU:HD13	2.00	0.45
1:B:187:GLN:HE22	1:C:261:LYS:NZ	2.15	0.45
1:E:392:HIS:N	1:E:392:HIS:ND1	2.65	0.44
1:E:354:GLN:H	1:E:354:GLN:HE21	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:D:170:TRP:CD2	1:D:211:PRO:HG3	2.54	0.43
1:C:368:ASP:HA	1:C:393:PHE:HD1	1.84	0.43
1:F:275:LEU:HA	1:F:276:PRO:HD3	1.94	0.43
1:B:368:ASP:HA	1:B:394:ASP:HB2	2.00	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:F:182:ARG:NH2	1:F:209:THR:HG22	2.34	0.42
1:A:171:GLU:OE2	1:A:183:ALA:HA	2.18	0.42
1:E:382:ASP:O	1:E:386:HIS:CD2	2.73	0.42
1:A:260:GLN:NE2	1:F:239:ARG:HH11	2.17	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:E:179:SER:HB2	1:E:180:PRO:HD3	2.02	0.41
1:C:324:ALA:HB3	1:C:344:VAL:HG12	2.03	0.41
1:A:266:ASN:ND2	1:C:180:PRO:HD3	2.34	0.41
1:F:379:ALA:O	1:F:383:ILE:HG13	2.21	0.41
1:E:354:GLN:H	1:E:354:GLN:NE2	2.18	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:D:268:TRP:HE3	1:D:268:TRP:HA	1.86	0.41
1:B:374:VAL:HA	1:B:377:LEU:HD12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:259:LEU:HD23	1:C:279:LEU:HD13	2.03	0.41
1:E:339:LEU:HD21	1:E:342:LEU:HB2	2.03	0.41
1:F:295:MET:HA	1:F:296:PRO:HD3	1.97	0.40
1:B:268:TRP:HA	1:B:268:TRP:CE3	2.56	0.40
1:E:380:LEU:HD22	1:E:391:ILE:HG21	2.03	0.40
1:E:268:TRP:HA	1:E:268:TRP:CE3	2.56	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

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	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
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%)	184 (96%)	7 (4%)	41	74
1	B	198/203 (98%)	196 (99%)	2 (1%)	82	95
1	C	188/203 (93%)	182 (97%)	6 (3%)	46	78
1	D	183/203 (90%)	165 (90%)	18 (10%)	10	26
1	E	196/203 (97%)	184 (94%)	12 (6%)	23	52
1	F	178/203 (88%)	166 (93%)	12 (7%)	20	46
All	All	1134/1218 (93%)	1077 (95%)	57 (5%)	30	61

All (57) 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	301	GLU
1	A	308	ASN
1	A	344	VAL
1	A	375	ARG
1	B	261	LYS
1	B	304	VAL
1	C	288	GLN
1	C	290	THR
1	C	300	ARG
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	225	ASP
1	D	243	VAL
1	D	261	LYS

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Mol	Chain	Res	Type
1	D	265	TYR
1	D	268	TRP
1	D	269	LEU
1	D	285	SER
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	E	162	GLU
1	E	188	GLU
1	E	192	CYS
1	E	225	ASP
1	E	254	SER
1	E	329	LEU
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	164	ASP
1	F	169	LYS
1	F	209	THR
1	F	225	ASP
1	F	236	GLU
1	F	239	ARG
1	F	266	ASN
1	F	329	LEU
1	F	331	SER
1	F	335	MET
1	F	352	LEU
1	F	390	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	260	GLN
1	A	266	ASN
1	B	260	GLN

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Mol	Chain	Res	Type
1	B	347	ASN
1	B	386	HIS
1	C	348	GLN
1	D	386	HIS
1	E	257	GLN
1	E	354	GLN
1	F	266	ASN

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 ⓘ

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 ⓘ

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%)	0.32	5 (2%) 67 61	53, 75, 97, 110	1 (0%)
1	B	237/239 (99%)	0.27	11 (4%) 36 29	42, 58, 99, 123	0
1	C	233/239 (97%)	0.38	7 (3%) 54 47	55, 79, 104, 124	0
1	D	236/239 (98%)	0.46	10 (4%) 40 33	56, 85, 110, 136	0
1	E	239/239 (100%)	0.22	5 (2%) 67 61	37, 56, 95, 108	0
1	F	235/239 (98%)	0.94	36 (15%) 3 2	55, 79, 130, 169	0
All	All	1416/1434 (98%)	0.43	74 (5%) 31 24	37, 73, 112, 169	1 (0%)

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	172	ARG	6.1
1	F	176	ALA	4.6
1	F	205	SER	4.2
1	F	170	TRP	4.2
1	F	214	LEU	4.0
1	F	236	GLU	4.0
1	F	185	VAL	3.9
1	B	386	HIS	3.8
1	F	184	ALA	3.7
1	F	213	ARG	3.7
1	A	361	SER	3.6
1	F	210	LEU	3.6
1	F	211	PRO	3.5
1	F	232	PRO	3.5
1	D	165	ALA	3.5
1	B	382	ASP	3.4
1	C	352	LEU	3.4
1	F	255	LEU	3.3
1	F	212	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	F	203	GLY	3.2
1	F	206	GLY	3.2
1	B	383	ILE	3.2
1	F	389	ILE	3.2
1	D	186	VAL	3.2
1	A	270	ALA	3.1
1	F	252	LEU	3.0
1	F	229	THR	3.0
1	E	161	ALA	3.0
1	C	372	LEU	3.0
1	F	186	VAL	3.0
1	C	292	LEU	2.9
1	F	192	CYS	2.9
1	F	200	LEU	2.8
1	E	180	PRO	2.8
1	B	358	GLY	2.7
1	B	380	LEU	2.6
1	F	177	GLY	2.6
1	F	194	ASN	2.6
1	E	181	GLY	2.6
1	A	386	HIS	2.6
1	B	389	ILE	2.6
1	F	163	TYR	2.5
1	E	179	SER	2.5
1	D	167	TRP	2.5
1	D	210	LEU	2.5
1	C	386	HIS	2.5
1	D	189	MET	2.4
1	F	165	ALA	2.4
1	F	183	ALA	2.4
1	B	384	ILE	2.3
1	B	396	ALA	2.3
1	F	391	ILE	2.3
1	D	181	GLY	2.3
1	B	398	PRO	2.3
1	E	176	ALA	2.3
1	F	228	LEU	2.3
1	F	222	VAL	2.3
1	F	174	ALA	2.3
1	D	163	TYR	2.3
1	B	387	SER	2.2
1	B	394	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	173	ASP	2.2
1	D	247	LEU	2.2
1	D	230	SER	2.2
1	F	241	LEU	2.2
1	A	352	LEU	2.2
1	F	175	PRO	2.2
1	A	358	GLY	2.1
1	C	358	GLY	2.1
1	F	279	LEU	2.1
1	F	383	ILE	2.1
1	C	177	GLY	2.1
1	D	272	LEU	2.0
1	C	247	LEU	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 carbohydrates 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.