



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:17 AM UTC

PDB ID : 5CWS / pdb_00005cws
Title : Crystal structure of the intact Chaetomium thermophilum Nsp1-Nup49-Nup57 channel nucleoporin heterotrimer bound to its Nic96 nuclear pore complex attachment site
Authors : Bley, C.J.; Petrovic, S.; Paduch, M.; Lu, V.; Kossiakoff, A.A.; Hoelz, A.
Deposited on : 2015-07-28
Resolution : 3.77 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

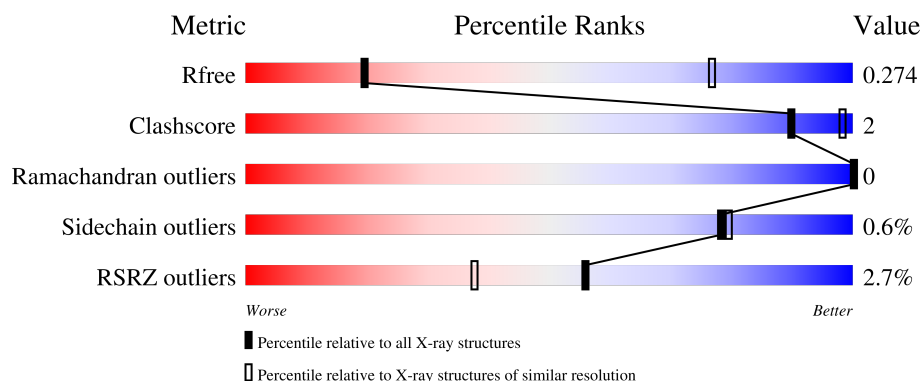
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.77 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	1085 (3.94-3.62)
Clashscore	190562	1029 (3.92-3.64)
Ramachandran outliers	187476	1064 (3.94-3.62)
Sidechain outliers	187428	1059 (3.94-3.62)
RSRZ outliers	180081	1084 (3.94-3.62)

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	237	86% 12%
1	G	237	3% 84% 12%
2	B	266	2% 81% 16%
2	H	266	3% 80% 16%
3	C	208	0% 78% 19%

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Mol	Chain	Length	Quality of chain
3	I	208	% 79% 19% .
4	D	227	3% 76% 21% .
4	J	227	3% 75% 21% .
5	E	247	3% 87% 9% ..
5	K	247	5% 88% 9% .
6	F	74	50% 46% .
6	L	74	3% 51% 46% .

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 32898 atoms, of which 16334 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 sAB-158 Fab Light Chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	208	Total	C	H	N	O	S	0	4	0
			3077	990	1503	260	319	5			
1	G	208	Total	C	H	N	O	S	0	4	0
			3077	990	1503	260	319	5			

- Molecule 2 is a protein called sAB-158 Fab Heavy Chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	223	Total	C	H	N	O	S	0	0	0
			3240	1043	1588	278	326	5			
2	H	223	Total	C	H	N	O	S	0	0	0
			3240	1043	1588	278	326	5			

- Molecule 3 is a protein called Nucleoporin NSP1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	169	Total	C	H	N	O	S	0	0	0
			2736	855	1364	235	276	6			
3	I	169	Total	C	H	N	O	S	0	0	0
			2736	855	1364	235	276	6			

- Molecule 4 is a protein called Nucleoporin NUP49.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	D	180	Total	C	H	N	O	S	0	0	0
			2803	876	1402	236	282	7			
4	J	180	Total	C	H	N	O	S	0	0	0
			2803	876	1402	236	282	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	244	MET	-	initiating methionine	UNP G0S4X2
D	245	SER	-	expression tag	UNP G0S4X2
J	244	MET	-	initiating methionine	UNP G0S4X2
J	245	SER	-	expression tag	UNP G0S4X2

- Molecule 5 is a protein called Nucleoporin NUP57.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	241	Total	C	H	N	O	0	0	0
			3947	1239	1976	360	368			
5	K	241	Total	C	H	N	O	0	0	0
			3947	1239	1976	360	368			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	70	MET	-	initiating methionine	UNP G0S0R2
K	70	MET	-	initiating methionine	UNP G0S0R2

- Molecule 6 is a protein called Nucleoporin NIC96.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	40	Total	C	H	N	O	0	0	0
			645	195	334	59	57			
6	L	40	Total	C	H	N	O	0	0	0
			645	195	334	59	57			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	140	SER	-	expression tag	UNP G0S024
L	140	SER	-	expression tag	UNP G0S024

- Molecule 7 is OSMIUM ION (CCD ID: OS) (formula: Os).

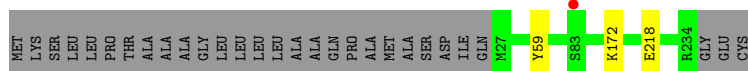
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total	Os	0	0
			1	1		
7	I	1	Total	Os	0	0
			1	1		

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.

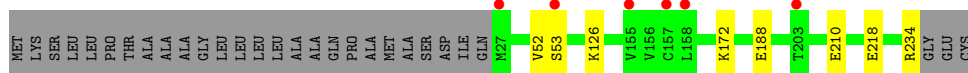
- Molecule 1: sAB-158 Fab Light Chain

Chain A:  86% 12%



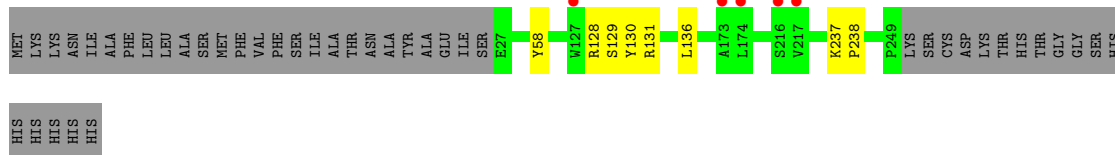
- Molecule 1: sAB-158 Fab Light Chain

Chain G:  84% 12% 3%



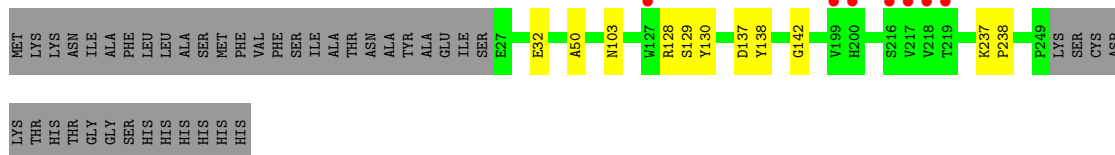
- Molecule 2: sAB-158 Fab Heavy Chain

Chain B:  81% 16% 2%



- Molecule 2: sAB-158 Fab Heavy Chain

Chain H:  80% 16% 3%



- Molecule 3: Nucleoporin NSP1

Chain C:  78% 19% 3%



- Molecule 3: Nucleoporin NSP1



- Molecule 4: Nucleoporin NUP49



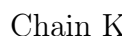
- Molecule 4: Nucleoporin NUP49



- Molecule 5: Nucleoporin NUP57



- Molecule 5: Nucleoporin NUP57

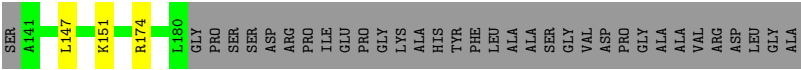


0314

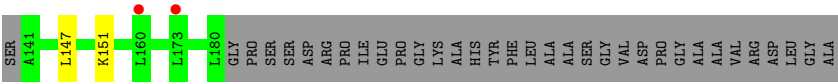
HIS

ASN

● Molecule 6: Nucleoporin NIC96



● Molecule 6: Nucleoporin NIC96



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	123.18Å 162.82Å 212.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.67 – 3.77 49.67 – 3.77	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.67-3.77) 99.9 (49.67-3.77)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 3.77Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.229 , 0.265 0.236 , 0.274	Depositor DCC
R_{free} test set	2000 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å ²)	179.4	Xtriage
Anisotropy	0.344	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 202.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	32898	wwPDB-VP
Average B, all atoms (Å ²)	264.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
OS

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	0.28	0/1621	0.59	0/2207
1	G	0.29	0/1621	0.60	0/2207
2	B	0.29	0/1694	0.63	0/2317
2	H	0.28	0/1694	0.59	0/2317
3	C	0.30	0/1388	0.66	0/1866
3	I	0.28	0/1388	0.68	1/1866 (0.1%)
4	D	0.30	0/1413	0.67	0/1898
4	J	0.30	0/1413	0.67	0/1898
5	E	0.34	0/2011	0.68	1/2715 (0.0%)
5	K	0.31	0/2011	0.69	1/2715 (0.0%)
6	F	0.29	0/312	0.62	0/416
6	L	0.29	0/312	0.62	0/416
All	All	0.30	0/16878	0.65	3/22838 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	535	GLU	N-CA-C	-5.77	105.50	113.18
5	K	108	VAL	CB-CA-C	-5.30	108.73	114.35
5	E	108	VAL	CB-CA-C	-5.11	108.93	114.35

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	1574	1503	1488	2	0
1	G	1574	1503	1488	4	0
2	B	1652	1588	1584	7	0
2	H	1652	1588	1584	7	0
3	C	1372	1364	1363	5	0
3	I	1372	1364	1363	1	3
4	D	1401	1402	1398	7	0
4	J	1401	1402	1398	8	3
5	E	1971	1976	1968	14	0
5	K	1971	1976	1968	14	0
6	F	311	334	334	2	0
6	L	311	334	334	2	0
7	C	1	0	0	0	0
7	I	1	0	0	0	0
All	All	16564	16334	16270	58	3

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 2.

All (58) 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:172:LYS:NZ	1:A:218:GLU:OE1	2.28	0.63
2:B:58:TYR:OH	5:E:171:ASP:OD2	2.10	0.58
1:G:172:LYS:NZ	1:G:218:GLU:OE1	2.33	0.58
3:C:598:GLU:OE1	5:E:254:ARG:NH1	2.40	0.55
4:J:368:ASN:OD1	5:K:210:ARG:NH2	2.40	0.54
1:G:210:GLU:O	1:G:234:ARG:NH1	2.41	0.54
2:H:50:ALA:O	2:H:103:ASN:ND2	2.41	0.54
5:K:81:LEU:O	5:K:85:LYS:HG2	2.08	0.53
4:J:304:ASP:OD2	5:K:177:HIS:ND1	2.41	0.53
4:D:468:MET:HE1	5:K:143:PRO:HA	1.91	0.52
5:E:81:LEU:O	5:E:85:LYS:HG2	2.10	0.52
4:J:280:LEU:HG	5:K:156:ILE:HD11	1.92	0.52
3:C:494:GLU:OE2	3:C:497:ARG:NH2	2.42	0.51
3:C:605:THR:O	5:E:261:LYS:NZ	2.43	0.51
5:E:142:PHE:HA	5:E:145:ILE:HG22	1.94	0.49
2:B:128:ARG:HG3	4:D:289:PHE:HZ	1.77	0.48
4:J:463:GLN:O	4:J:467:PHE:CD2	2.66	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:85:LYS:HD2	5:E:94:ALA:HB2	1.96	0.48
3:C:495:TRP:CG	4:D:280:LEU:HD13	2.48	0.47
6:L:147:LEU:O	6:L:151:LYS:N	2.47	0.47
5:K:108:VAL:N	5:K:109:PRO:CD	2.78	0.47
5:E:108:VAL:N	5:E:109:PRO:CD	2.78	0.46
4:J:284:PRO:HA	4:J:287:VAL:HG12	1.96	0.46
3:I:618:GLN:HG2	6:L:147:LEU:HD13	1.97	0.46
5:K:85:LYS:HD3	5:K:92:ASN:O	2.17	0.45
5:K:99:LEU:O	5:K:137:VAL:N	2.50	0.45
5:E:90:HIS:N	5:E:91:PRO:CD	2.79	0.45
4:D:284:PRO:HA	4:D:287:VAL:HG12	1.98	0.45
2:H:137:ASP:OD2	2:H:138:TYR:N	2.50	0.44
2:B:129:SER:N	2:B:130:TYR:HA	2.33	0.44
5:E:156:ILE:CG2	5:E:157:THR:N	2.81	0.44
4:J:463:GLN:HB3	4:J:467:PHE:CE2	2.53	0.44
5:K:90:HIS:N	5:K:91:PRO:CD	2.81	0.44
5:E:98:TYR:HD2	5:E:125:ALA:HB2	1.83	0.44
4:D:433:GLU:OE2	6:F:174:ARG:NH2	2.51	0.44
2:B:131:ARG:NH2	4:D:286:ASP:OD1	2.49	0.43
1:G:126:LYS:NZ	1:G:188:GLU:OE1	2.51	0.43
5:E:99:LEU:O	5:E:137:VAL:N	2.51	0.43
2:B:129:SER:H	2:B:130:TYR:HA	1.84	0.43
1:A:59:TYR:HH	2:B:136:LEU:H	1.67	0.43
6:F:147:LEU:O	6:F:151:LYS:N	2.52	0.43
5:E:156:ILE:HG22	5:E:157:THR:N	2.34	0.43
2:H:128:ARG:HG3	4:J:289:PHE:HZ	1.84	0.42
5:K:142:PHE:N	5:K:143:PRO:HD2	2.34	0.42
4:D:429:GLU:N	4:D:429:GLU:OE2	2.52	0.42
2:H:129:SER:N	2:H:130:TYR:HA	2.34	0.42
2:H:237:LYS:N	2:H:238:PRO:CD	2.83	0.42
2:B:237:LYS:N	2:B:238:PRO:CD	2.84	0.41
4:J:276:HIS:NE2	5:K:152:GLN:HB3	2.35	0.41
5:K:142:PHE:HA	5:K:145:ILE:HG22	2.02	0.41
5:E:75:ILE:N	5:E:76:PRO:CD	2.83	0.41
5:K:100:TYR:OH	5:K:128:ARG:O	2.39	0.41
2:H:129:SER:H	2:H:130:TYR:HA	1.86	0.41
3:C:614:ASP:HB2	3:C:615:PRO:HD2	2.03	0.41
5:E:98:TYR:CD1	5:E:138:LEU:HA	2.56	0.41
1:G:52:VAL:HG22	1:G:53:SER:N	2.36	0.41
5:K:152:GLN:HA	5:K:155:VAL:HG12	2.04	0.40
2:H:32:GLU:OE2	2:H:142:GLY:N	2.47	0.40

All (3) 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)
3:I:472:THR:O	4:J:432:TYR:OH[4_455]	2.07	0.13
3:I:476:THR:HG1	4:J:432:TYR:OH[4_455]	1.54	0.06
3:I:472:THR:O	4:J:432:TYR:HH[4_455]	1.59	0.01

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	210/237 (89%)	197 (94%)	13 (6%)	0	100	100
1	G	210/237 (89%)	197 (94%)	13 (6%)	0	100	100
2	B	221/266 (83%)	204 (92%)	17 (8%)	0	100	100
2	H	221/266 (83%)	206 (93%)	15 (7%)	0	100	100
3	C	165/208 (79%)	158 (96%)	7 (4%)	0	100	100
3	I	165/208 (79%)	158 (96%)	7 (4%)	0	100	100
4	D	174/227 (77%)	172 (99%)	2 (1%)	0	100	100
4	J	174/227 (77%)	171 (98%)	3 (2%)	0	100	100
5	E	239/247 (97%)	212 (89%)	27 (11%)	0	100	100
5	K	239/247 (97%)	215 (90%)	24 (10%)	0	100	100
6	F	38/74 (51%)	31 (82%)	7 (18%)	0	100	100
6	L	38/74 (51%)	30 (79%)	8 (21%)	0	100	100
All	All	2094/2518 (83%)	1951 (93%)	143 (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	174/204 (85%)	174 (100%)	0	100	100
1	G	174/204 (85%)	174 (100%)	0	100	100
2	B	180/221 (81%)	180 (100%)	0	100	100
2	H	180/221 (81%)	180 (100%)	0	100	100
3	C	146/171 (85%)	146 (100%)	0	100	100
3	I	146/171 (85%)	145 (99%)	1 (1%)	76	77
4	D	150/176 (85%)	150 (100%)	0	100	100
4	J	150/176 (85%)	150 (100%)	0	100	100
5	E	213/219 (97%)	208 (98%)	5 (2%)	44	62
5	K	213/219 (97%)	209 (98%)	4 (2%)	50	65
6	F	32/55 (58%)	32 (100%)	0	100	100
6	L	32/55 (58%)	32 (100%)	0	100	100
All	All	1790/2092 (86%)	1780 (99%)	10 (1%)	78	79

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	E	107	THR
5	E	156	ILE
5	E	163	LEU
5	E	246	LEU
5	E	287	LEU
3	I	641	GLN
5	K	96	LYS
5	K	163	LEU
5	K	246	LEU
5	K	287	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	161	ASN
2	B	103	ASN
3	C	502	ASN
4	D	276	HIS
4	D	298	HIS
5	E	133	ASN
5	E	152	GLN
5	E	296	HIS
1	G	112	GLN
1	G	123	GLN
2	H	103	ASN
3	I	507	GLN
3	I	618	GLN
5	K	133	ASN
5	K	296	HIS
5	K	300	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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	208/237 (87%)	-0.19	1 (0%) 87 70	144, 280, 330, 376	3 (1%)
1	G	208/237 (87%)	-0.03	6 (2%) 53 36	132, 264, 319, 335	3 (1%)
2	B	223/266 (83%)	-0.14	5 (2%) 62 43	159, 252, 382, 453	0
2	H	223/266 (83%)	-0.21	7 (3%) 51 34	158, 257, 368, 426	0
3	C	169/208 (81%)	-0.28	2 (1%) 76 54	143, 245, 351, 397	0
3	I	169/208 (81%)	-0.25	3 (1%) 67 46	159, 267, 383, 427	0
4	D	180/227 (79%)	-0.14	6 (3%) 49 33	145, 245, 343, 358	0
4	J	180/227 (79%)	-0.03	7 (3%) 43 30	151, 276, 363, 402	0
5	E	241/247 (97%)	-0.22	7 (2%) 53 36	132, 255, 370, 428	0
5	K	241/247 (97%)	-0.12	12 (4%) 34 25	136, 279, 371, 416	0
6	F	40/74 (54%)	-0.12	0 100 100	201, 254, 349, 360	0
6	L	40/74 (54%)	0.01	2 (5%) 34 25	198, 281, 369, 391	0
All	All	2122/2518 (84%)	-0.15	58 (2%) 56 38	132, 262, 365, 453	6 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	E	79	ILE	6.0
2	H	217	VAL	4.9
5	E	83	VAL	4.8
2	B	216	SER	4.4
2	H	127	TRP	4.1
1	G	158	LEU	4.1
2	H	216	SER	4.1
2	B	217	VAL	4.0
1	G	157	CYS	3.8
1	G	155	VAL	3.6
5	K	163	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
5	K	156	ILE	3.3
5	K	279	ILE	3.3
1	G	53	SER	3.1
5	K	82	ILE	3.0
5	K	283	ALA	2.9
4	D	252	ILE	2.9
5	E	86	TRP	2.9
4	J	290	VAL	2.9
5	K	79	ILE	2.8
4	D	406	HIS	2.8
2	B	174	LEU	2.7
6	L	160	LEU	2.7
5	E	163	LEU	2.7
4	J	273	LEU	2.7
3	I	506	ILE	2.6
5	E	82	ILE	2.6
5	E	280	GLU	2.6
3	I	627	LEU	2.5
5	K	159	PHE	2.5
1	G	203[A]	THR	2.5
1	G	27	MET	2.4
5	K	86	TRP	2.4
2	H	218	VAL	2.4
5	K	170	LEU	2.4
4	J	367	SER	2.4
4	J	272	PHE	2.4
4	J	311	LEU	2.4
3	C	499	LEU	2.3
4	D	248	LEU	2.3
2	H	219	THR	2.3
3	C	506	ILE	2.2
4	D	311	LEU	2.2
4	D	396	ILE	2.2
2	B	127	TRP	2.2
2	H	200	HIS	2.1
5	K	280	GLU	2.1
1	A	83	SER	2.1
2	H	199	VAL	2.1
4	J	414	ALA	2.1
4	J	431	VAL	2.1
3	I	618	GLN	2.1
6	L	173	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
4	D	287	VAL	2.0
5	K	162	LYS	2.0
5	E	166	ILE	2.0
5	K	287	LEU	2.0
2	B	173	ALA	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	OS	I	701	1/1	0.97	0.06	301,301,301,301	0
7	OS	C	701	1/1	0.99	0.08	296,296,296,296	0

6.5 Other polymers [i](#)

There are no such residues in this entry.