



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:35 PM GMT

PDB ID : 1BJM
Title : LOC NAKS, A LAMBDA 1 TYPE LIGHT-CHAIN DIMER (BENCE-JONES PROTEIN) CRYSTALLIZED IN NAKSO4
Authors : Schiffer, M.; Huang, D.B.
Deposited on : 1995-05-26
Resolution : 2.20 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

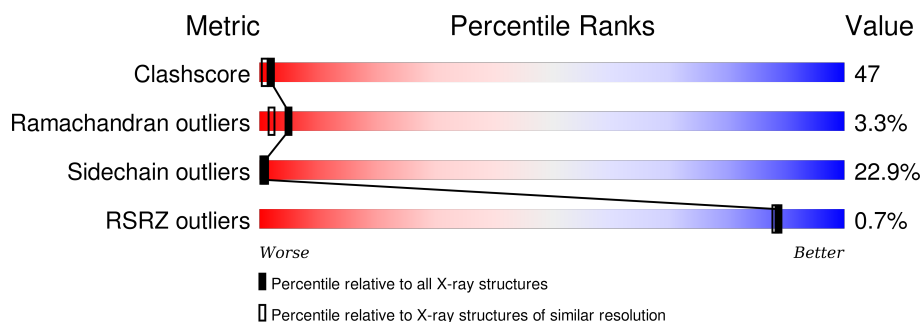
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	216	<div> <div></div> <div> <div></div> <div>30%</div> <div>47%</div> <div>17%</div> <div>6%</div> </div> </div>
1	B	216	<div> <div></div> <div> <div></div> <div>38%</div> <div>42%</div> <div>16%</div> <div>•</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3494 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 LOC - LAMBDA 1 TYPE LIGHT-CHAIN DIMER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1598	992	266	335	5			
1	B	216	Total	C	N	O	S	0	0	0
			1598	992	266	335	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	THR	ILE	CONFLICT	PIR S25754
A	31	GLU	GLY	CONFLICT	PIR S25754
A	33	SER	THR	CONFLICT	PIR S25754
A	35	THR	ASN	CONFLICT	PIR S25754
A	39	HIS	GLN	CONFLICT	PIR S25754
A	41	SER	PRO	CONFLICT	PIR S25754
A	43	THR	ARG	CONFLICT	PIR S25754
A	50	TYR	HIS	CONFLICT	PIR S25754
A	51	GLU	SER	CONFLICT	PIR S25754
A	52	ASP	ASN	CONFLICT	PIR S25754
A	54	SER	GLN	CONFLICT	PIR S25754
A	56	ALA	PRO	CONFLICT	PIR S25754
A	60	SER	PRO	CONFLICT	PIR S25754
A	65	ALA	GLY	CONFLICT	PIR S25754
A	81	PRO	SER	CONFLICT	PIR S25754
A	85	THR	ALA	CONFLICT	PIR S25754
A	?	-	ASN	DELETION	PIR S25754
A	97	ASP	GLY	CONFLICT	PIR S25754
A	98	VAL	ARG	CONFLICT	PIR S25754
A	99	ALA	TYR	CONFLICT	PIR S25754
B	19	THR	ILE	CONFLICT	PIR S25754
B	31	GLU	GLY	CONFLICT	PIR S25754
B	33	SER	THR	CONFLICT	PIR S25754
B	35	THR	ASN	CONFLICT	PIR S25754
B	39	HIS	GLN	CONFLICT	PIR S25754

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Chain	Residue	Modelled	Actual	Comment	Reference
B	41	SER	PRO	CONFLICT	PIR S25754
B	43	THR	ARG	CONFLICT	PIR S25754
B	50	TYR	HIS	CONFLICT	PIR S25754
B	51	GLU	SER	CONFLICT	PIR S25754
B	52	ASP	ASN	CONFLICT	PIR S25754
B	54	SER	GLN	CONFLICT	PIR S25754
B	56	ALA	PRO	CONFLICT	PIR S25754
B	60	SER	PRO	CONFLICT	PIR S25754
B	65	ALA	GLY	CONFLICT	PIR S25754
B	81	PRO	SER	CONFLICT	PIR S25754
B	85	THR	ALA	CONFLICT	PIR S25754
B	?	-	ASN	DELETION	PIR S25754
B	97	ASP	GLY	CONFLICT	PIR S25754
B	98	VAL	ARG	CONFLICT	PIR S25754
B	99	ALA	TYR	CONFLICT	PIR S25754

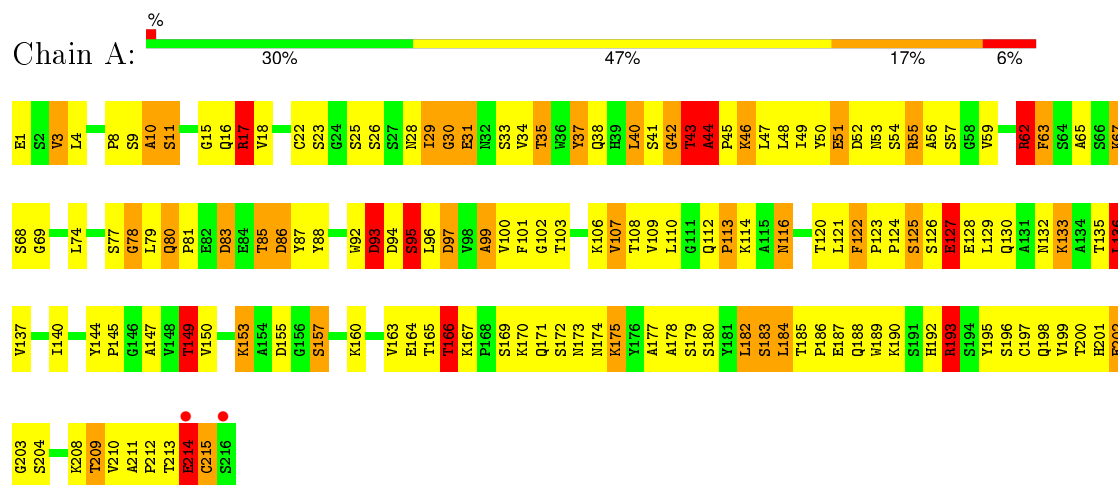
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	161	Total O 161 161	0	0
2	B	137	Total O 137 137	0	0

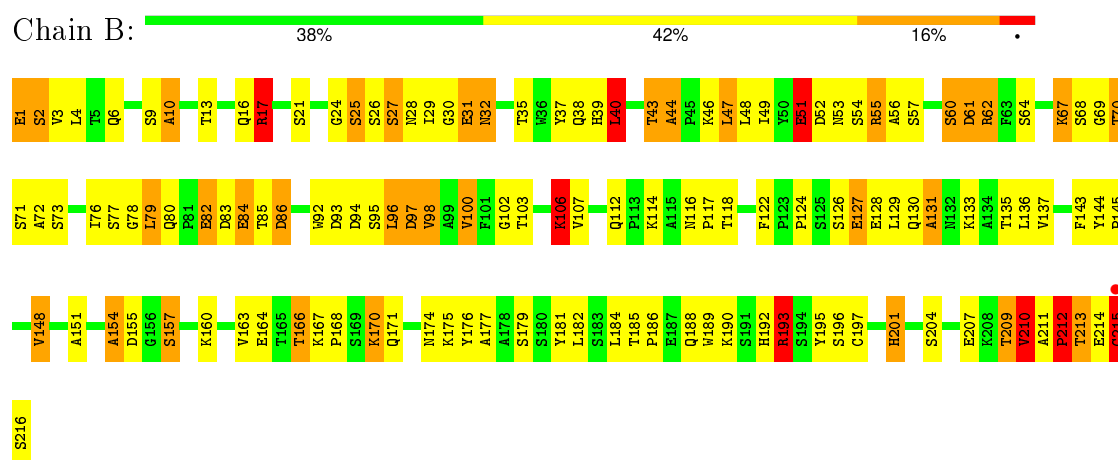
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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: LOC - LAMBDA 1 TYPE LIGHT-CHAIN DIMER



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.87Å 72.62Å 63.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 10.19 – 2.24	Depositor EDS
% Data completeness (in resolution range)	76.3 (10.00-2.20) 73.8 (10.19-2.24)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.158 , (Not available) 0.156 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.09 , 69.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ¹	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Outliers	0 of 14084 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3494	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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 ⓘ

5.1 Standard geometry ⓘ

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

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	1.04	1/1628 (0.1%)	1.93	41/2226 (1.8%)
1	B	0.99	1/1628 (0.1%)	1.99	45/2226 (2.0%)
All	All	1.02	2/3256 (0.1%)	1.96	86/4452 (1.9%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	GLY	C-N	-8.35	1.14	1.34
1	B	2	SER	N-CA	-5.22	1.35	1.46

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	PCA	C-N-CA	18.33	167.54	121.70
1	A	17	ARG	NE-CZ-NH2	9.99	125.29	120.30
1	A	17	ARG	NE-CZ-NH1	-9.64	115.48	120.30
1	A	95	SER	N-CA-CB	9.16	124.24	110.50
1	B	94	ASP	CB-CG-OD1	9.16	126.54	118.30
1	B	61	ASP	CB-CG-OD1	-9.05	110.15	118.30
1	B	106	LYS	CA-CB-CG	8.41	131.90	113.40
1	A	52	ASP	CB-CG-OD1	8.25	125.72	118.30
1	A	55	ARG	CD-NE-CZ	7.73	134.42	123.60
1	A	166	THR	CA-CB-CG2	7.68	123.16	112.40
1	A	127	GLU	CB-CG-CD	7.67	134.90	114.20
1	B	62	ARG	NE-CZ-NH2	7.61	124.10	120.30
1	A	193	ARG	CD-NE-CZ	7.38	133.93	123.60
1	B	100	VAL	CA-CB-CG1	7.16	121.63	110.90
1	B	84	GLU	OE1-CD-OE2	-7.13	114.75	123.30
1	A	127	GLU	CG-CD-OE1	6.93	132.16	118.30
1	B	154	ALA	CB-CA-C	6.93	120.49	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	ASP	N-CA-C	6.83	129.43	111.00
1	B	84	GLU	CG-CD-OE1	6.72	131.75	118.30
1	B	181	TYR	CB-CG-CD1	6.71	125.03	121.00
1	B	62	ARG	CA-CB-CG	6.63	128.00	113.40
1	A	116	ASN	CB-CA-C	6.63	123.66	110.40
1	B	86	ASP	CB-CG-OD1	6.63	124.27	118.30
1	B	197	CYS	CA-CB-SG	-6.54	102.23	114.00
1	B	148	VAL	O-C-N	6.53	133.14	122.70
1	B	193	ARG	NE-CZ-NH2	6.52	123.56	120.30
1	A	46	LYS	CA-CB-CG	6.45	127.58	113.40
1	A	37	TYR	CB-CG-CD2	6.36	124.82	121.00
1	B	166	THR	N-CA-CB	-6.36	98.22	110.30
1	B	10	ALA	CB-CA-C	6.30	119.55	110.10
1	B	69	GLY	C-N-CA	6.21	137.23	121.70
1	A	63	PHE	O-C-N	6.09	132.45	122.70
1	B	51	GLU	CG-CD-OE2	5.99	130.28	118.30
1	B	148	VAL	CA-CB-CG1	5.98	119.87	110.90
1	A	83	ASP	CB-CG-OD2	-5.91	112.98	118.30
1	B	155	ASP	CB-CG-OD2	5.88	123.59	118.30
1	A	133	LYS	N-CA-CB	5.86	121.15	110.60
1	B	62	ARG	NH1-CZ-NH2	-5.82	113.00	119.40
1	A	155	ASP	CB-CG-OD2	-5.79	113.09	118.30
1	B	40	LEU	N-CA-C	-5.78	95.40	111.00
1	A	129	LEU	CA-CB-CG	5.77	128.57	115.30
1	B	100	VAL	CB-CA-C	5.75	122.33	111.40
1	A	86	ASP	CB-CG-OD1	5.70	123.43	118.30
1	B	60	SER	N-CA-CB	-5.70	101.95	110.50
1	A	94	ASP	C-N-CA	5.68	135.90	121.70
1	A	127	GLU	OE1-CD-OE2	-5.66	116.50	123.30
1	B	164	GLU	CG-CD-OE1	5.66	129.61	118.30
1	A	147	ALA	N-CA-CB	-5.65	102.19	110.10
1	B	1	PCA	O-C-N	-5.65	113.66	122.70
1	B	213	THR	N-CA-CB	5.64	121.01	110.30
1	A	149	THR	CA-CB-CG2	5.59	120.22	112.40
1	A	95	SER	CA-C-N	5.58	129.49	117.20
1	B	166	THR	CA-CB-CG2	5.58	120.21	112.40
1	A	3	VAL	N-CA-CB	-5.58	99.23	111.50
1	A	86	ASP	CB-CA-C	5.58	121.55	110.40
1	B	57	SER	N-CA-CB	5.57	118.86	110.50
1	B	55	ARG	CD-NE-CZ	-5.55	115.82	123.60
1	B	86	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	B	135	THR	N-CA-CB	5.50	120.76	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	11	SER	N-CA-CB	-5.49	102.26	110.50
1	B	184	LEU	CA-CB-CG	5.46	127.85	115.30
1	B	17	ARG	CD-NE-CZ	5.43	131.20	123.60
1	B	137	VAL	O-C-N	5.39	131.32	122.70
1	A	52	ASP	CB-CG-OD2	-5.37	113.47	118.30
1	A	122	PHE	CA-CB-CG	5.37	126.78	113.90
1	A	62	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	A	54	SER	CB-CA-C	5.32	120.20	110.10
1	A	83	ASP	O-C-N	5.30	131.19	122.70
1	A	177	ALA	CB-CA-C	5.24	117.96	110.10
1	A	95	SER	N-CA-C	-5.23	96.87	111.00
1	B	151	ALA	N-CA-CB	5.21	117.40	110.10
1	A	214	GLU	CG-CD-OE2	5.21	128.72	118.30
1	B	201	HIS	CA-CB-CG	5.21	122.45	113.60
1	B	207	GLU	CG-CD-OE2	5.19	128.67	118.30
1	B	84	GLU	CB-CG-CD	5.18	128.18	114.20
1	B	62	ARG	NE-CZ-NH1	5.15	122.88	120.30
1	B	210	VAL	CA-CB-CG1	5.12	118.59	110.90
1	A	137	VAL	CA-CB-CG1	5.10	118.55	110.90
1	A	44	ALA	CB-CA-C	-5.10	102.45	110.10
1	B	70	THR	CA-CB-CG2	5.08	119.50	112.40
1	B	17	ARG	NE-CZ-NH2	5.07	122.84	120.30
1	B	212	PRO	C-N-CA	5.07	134.38	121.70
1	A	136	LEU	CA-CB-CG	5.05	126.91	115.30
1	A	99	ALA	N-CA-CB	5.04	117.15	110.10
1	A	10	ALA	CB-CA-C	5.02	117.63	110.10
1	A	29	ILE	CA-C-N	-5.02	106.16	116.20

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	1598	0	1535	161	0
1	B	1598	0	1536	137	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	161	0	0	49	0
2	B	137	0	0	24	1
All	All	3494	0	3071	292	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 47.

All (292) 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:30:GLY:O	1:A:31:GLU:HB2	1.49	1.12
1:A:95:SER:HB2	2:A:566:HOH:O	1.49	1.11
1:A:192:HIS:CE1	2:A:651:HOH:O	2.03	1.11
1:B:212:PRO:HG3	2:B:841:HOH:O	1.52	1.09
1:A:47:LEU:HD21	1:A:50:TYR:HB3	1.39	1.04
1:A:8:PRO:HG3	1:A:149:THR:HG22	1.42	1.02
1:B:192:HIS:O	1:B:212:PRO:HG2	1.59	1.00
2:A:688:HOH:O	1:B:166:THR:CG2	2.09	0.99
1:B:40:LEU:H	1:B:40:LEU:HD12	1.26	0.99
1:A:92:TRP:O	1:A:93:ASP:HB2	1.63	0.98
1:B:214:GLU:O	1:B:215:CYS:HB2	1.63	0.97
1:B:62:ARG:NH2	1:B:83:ASP:OD2	2.02	0.92
2:A:632:HOH:O	1:B:171:GLN:HA	1.70	0.92
1:A:59:VAL:HA	2:A:549:HOH:O	1.69	0.91
1:A:193:ARG:HA	2:A:653:HOH:O	1.70	0.91
1:A:166:THR:HG23	2:A:631:HOH:O	1.70	0.90
1:A:164:GLU:OE1	2:A:632:HOH:O	1.90	0.89
1:B:40:LEU:HD13	1:B:43:THR:OG1	1.73	0.89
1:A:126:SER:O	1:A:130:GLN:HG3	1.71	0.89
1:B:17:ARG:HB3	1:B:77:SER:HA	1.54	0.88
1:A:40:LEU:HD22	1:A:41:SER:H	1.38	0.87
1:B:25:SER:HB3	2:B:714:HOH:O	1.72	0.87
1:B:52:ASP:OD1	1:B:67:LYS:HD2	1.77	0.85
1:B:185:THR:OG1	1:B:188:GLN:OE1	1.95	0.84
1:B:28:ASN:HA	1:B:92:TRP:O	1.78	0.83
1:A:167:LYS:HB3	1:A:167:LYS:NZ	1.94	0.82
1:A:48:LEU:C	1:A:49:ILE:HD12	2.00	0.82
1:A:29:ILE:HG22	1:A:69:GLY:O	1.80	0.81
1:A:4:LEU:HB2	1:A:102:GLY:HA2	1.63	0.81
1:B:51:GLU:OE1	2:B:745:HOH:O	1.99	0.81
1:A:29:ILE:O	1:A:31:GLU:N	2.15	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:ASN:ND2	1:B:92:TRP:O	2.15	0.80
1:A:189:TRP:CZ2	1:A:212:PRO:HA	2.17	0.80
1:A:17:ARG:HG2	1:A:77:SER:HA	1.63	0.79
1:B:16:GLN:HG2	1:B:17:ARG:H	1.47	0.79
1:B:144:TYR:OH	2:B:818:HOH:O	2.00	0.79
1:B:95:SER:OG	1:B:96:LEU:HD22	1.83	0.79
1:B:9:SER:HB2	2:B:706:HOH:O	1.81	0.78
1:A:133:LYS:HE2	2:A:640:HOH:O	1.83	0.78
1:A:96:LEU:HD11	2:A:562:HOH:O	1.85	0.77
1:B:163:VAL:HG22	1:B:182:LEU:HD13	1.64	0.77
1:B:167:LYS:N	1:B:167:LYS:HD3	1.99	0.77
1:A:112:GLN:NE2	2:A:600:HOH:O	2.17	0.76
1:A:192:HIS:NE2	2:A:651:HOH:O	2.11	0.76
1:A:11:SER:HB2	1:A:108:THR:O	1.85	0.76
1:B:163:VAL:HB	2:B:882:HOH:O	1.87	0.75
1:A:93:ASP:O	2:A:574:HOH:O	2.04	0.74
1:B:157:SER:OG	2:B:826:HOH:O	2.02	0.74
1:B:16:GLN:HG2	1:B:17:ARG:N	2.04	0.73
1:B:133:LYS:HE2	2:B:606:HOH:O	1.87	0.72
1:A:86:ASP:OD1	1:A:106:LYS:HG3	1.90	0.72
1:A:190:LYS:HE3	2:A:646:HOH:O	1.88	0.72
1:A:167:LYS:HG3	2:A:837:HOH:O	1.89	0.72
1:A:18:VAL:HA	2:A:530:HOH:O	1.87	0.72
1:A:103:THR:HG22	2:A:525:HOH:O	1.89	0.72
1:A:175:LYS:HE3	2:A:601:HOH:O	1.90	0.71
1:A:48:LEU:HD22	1:A:59:VAL:HG11	1.73	0.71
1:A:35:THR:HG23	1:A:47:LEU:CD1	2.21	0.71
2:A:718:HOH:O	1:B:35:THR:HG23	1.90	0.71
1:B:40:LEU:HD13	1:B:43:THR:CB	2.22	0.70
1:B:31:GLU:O	2:B:716:HOH:O	2.11	0.69
1:A:173:ASN:O	1:A:174:ASN:HB2	1.93	0.68
1:B:86:ASP:OD2	2:B:719:HOH:O	2.09	0.68
1:A:150:VAL:HG22	1:A:199:VAL:HG22	1.76	0.68
1:B:62:ARG:HH21	1:B:83:ASP:CG	1.97	0.67
1:B:62:ARG:NH2	1:B:83:ASP:CG	2.47	0.67
1:B:189:TRP:HZ3	1:B:195:TYR:CD2	2.12	0.67
1:B:43:THR:O	1:B:44:ALA:CB	2.42	0.66
1:B:185:THR:CB	1:B:188:GLN:OE1	2.44	0.66
1:B:189:TRP:CZ3	1:B:195:TYR:HD2	2.14	0.66
1:A:45:PRO:HD2	2:A:540:HOH:O	1.94	0.66
1:B:4:LEU:HD13	1:B:100:VAL:HG12	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:LEU:CD1	2:A:562:HOH:O	2.41	0.65
1:A:38:GLN:HG3	1:A:85:THR:HG21	1.78	0.65
1:A:145:PRO:HD2	2:A:614:HOH:O	1.96	0.65
1:A:125:SER:HB2	1:A:127:GLU:CD	2.18	0.64
1:B:163:VAL:CG2	1:B:182:LEU:HD13	2.27	0.64
1:A:125:SER:HB2	1:A:127:GLU:HG2	1.79	0.64
1:A:40:LEU:C	1:A:42:GLY:H	2.00	0.64
1:B:195:TYR:N	1:B:210:VAL:O	2.27	0.64
1:B:212:PRO:CG	2:B:841:HOH:O	2.24	0.63
1:B:189:TRP:CZ3	1:B:195:TYR:CD2	2.85	0.63
1:A:4:LEU:HB2	1:A:102:GLY:CA	2.27	0.63
1:B:93:ASP:OD2	2:B:701:HOH:O	2.16	0.62
1:A:78:GLY:O	1:A:80:GLN:NE2	2.30	0.62
1:B:117:PRO:HD3	1:B:201:HIS:CD2	2.34	0.62
1:A:103:THR:CG2	2:A:525:HOH:O	2.46	0.62
1:A:153:LYS:HD3	1:A:196:SER:OG	1.99	0.62
1:A:1:PCA:HA	2:A:522:HOH:O	1.99	0.62
1:A:22:CYS:SG	1:A:29:ILE:HD11	2.39	0.62
1:B:190:LYS:HE2	1:B:190:LYS:HA	1.80	0.62
1:B:24:GLY:HA3	1:B:29:ILE:HD12	1.81	0.61
1:B:62:ARG:NH2	1:B:83:ASP:OD1	2.33	0.61
1:A:30:GLY:O	1:A:31:GLU:CB	2.37	0.61
1:B:40:LEU:CD1	1:B:43:THR:OG1	2.47	0.61
1:B:160:LYS:HE3	2:B:829:HOH:O	1.99	0.61
1:B:26:SER:HA	1:B:30:GLY:HA3	1.82	0.61
1:B:185:THR:HG23	1:B:188:GLN:OE1	2.01	0.61
1:A:29:ILE:HG23	1:A:67:LYS:HD3	1.83	0.61
1:B:157:SER:HB2	2:B:825:HOH:O	2.00	0.60
1:A:11:SER:CB	1:A:108:THR:O	2.48	0.60
1:B:98:VAL:HG12	2:B:743:HOH:O	2.01	0.60
1:A:40:LEU:CD2	1:A:41:SER:H	2.14	0.60
1:B:16:GLN:CG	1:B:17:ARG:H	2.15	0.60
1:A:167:LYS:NZ	1:A:167:LYS:CB	2.64	0.59
1:B:189:TRP:HZ3	1:B:195:TYR:HD2	1.45	0.59
1:A:35:THR:HG23	1:A:47:LEU:HD12	1.85	0.59
1:B:93:ASP:OD1	1:B:96:LEU:HB2	2.03	0.59
1:A:192:HIS:HE1	2:A:651:HOH:O	1.62	0.58
1:B:40:LEU:N	1:B:40:LEU:HD12	2.08	0.58
1:A:38:GLN:HE21	1:A:85:THR:HG21	1.67	0.58
1:B:28:ASN:O	1:B:32:ASN:ND2	2.37	0.58
1:A:15:GLY:N	1:A:79:LEU:O	2.26	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:SER:HB2	2:A:637:HOH:O	2.03	0.58
1:B:171:GLN:HG3	1:B:175:LYS:O	2.04	0.58
1:B:40:LEU:CD1	1:B:40:LEU:H	2.04	0.57
1:B:10:ALA:HB3	1:B:107:VAL:HG22	1.85	0.57
1:B:17:ARG:HA	1:B:76:ILE:O	2.04	0.56
1:A:16:GLN:HG2	1:A:17:ARG:H	1.69	0.56
1:A:30:GLY:HA2	1:A:69:GLY:O	2.06	0.56
1:A:140:ILE:CD1	1:A:150:VAL:HG21	2.36	0.56
1:A:92:TRP:O	1:A:93:ASP:CB	2.48	0.56
1:A:140:ILE:HD11	1:A:150:VAL:HG21	1.87	0.56
1:A:214:GLU:CA	1:A:214:GLU:OE1	2.53	0.56
1:A:167:LYS:HZ3	1:A:167:LYS:HB3	1.69	0.56
1:B:6:GLN:OE1	1:B:102:GLY:HA3	2.05	0.56
1:A:43:THR:O	1:A:44:ALA:CB	2.54	0.56
1:A:114:LYS:O	2:A:655:HOH:O	2.18	0.56
1:B:39:HIS:CD2	1:B:40:LEU:O	2.59	0.55
1:B:185:THR:CG2	1:B:188:GLN:OE1	2.53	0.55
1:A:62:ARG:HH22	1:A:83:ASP:CG	2.09	0.55
1:A:3:VAL:HG12	1:A:100:VAL:HG13	1.88	0.55
1:A:40:LEU:HD22	1:A:41:SER:N	2.15	0.55
1:A:182:LEU:CD1	1:A:184:LEU:HD11	2.36	0.55
1:B:168:PRO:HD3	2:B:834:HOH:O	2.06	0.55
1:B:133:LYS:CE	2:B:606:HOH:O	2.48	0.55
1:A:112:GLN:HB2	1:A:113:PRO:HD2	1.89	0.54
1:B:188:GLN:O	1:B:192:HIS:HD2	1.91	0.54
1:A:79:LEU:HD21	1:A:109:VAL:HG22	1.87	0.54
1:A:116:ASN:OD1	2:A:603:HOH:O	2.18	0.54
1:B:55:ARG:HD3	1:B:60:SER:O	2.08	0.54
1:B:49:ILE:HA	1:B:54:SER:O	2.07	0.54
1:A:215:CYS:HA	1:B:216:SER:C	2.28	0.54
1:B:116:ASN:HA	1:B:201:HIS:HD2	1.72	0.53
1:A:29:ILE:HG22	1:A:30:GLY:N	2.23	0.53
1:A:186:PRO:O	2:A:646:HOH:O	2.19	0.53
1:A:195:TYR:CD2	2:A:661:HOH:O	2.59	0.53
1:A:167:LYS:HE3	2:A:636:HOH:O	2.08	0.53
1:A:79:LEU:CD2	1:A:109:VAL:HG22	2.40	0.52
1:B:185:THR:HB	1:B:186:PRO:HD2	1.90	0.52
1:A:145:PRO:CD	2:A:614:HOH:O	2.55	0.52
1:B:143:PHE:CE2	1:B:148:VAL:HG13	2.43	0.52
1:A:35:THR:HG22	1:A:37:TYR:CE1	2.44	0.52
1:A:190:LYS:CE	2:A:646:HOH:O	2.54	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:ARG:HA	1:B:212:PRO:HD2	1.92	0.52
1:A:214:GLU:OE1	1:A:214:GLU:HA	2.10	0.52
1:A:130:GLN:HG2	2:A:609:HOH:O	2.09	0.51
1:B:13:THR:O	1:B:16:GLN:HB3	2.10	0.51
1:B:93:ASP:O	1:B:97:ASP:N	2.43	0.51
1:A:107:VAL:O	1:A:107:VAL:CG2	2.58	0.51
1:A:215:CYS:H	1:B:216:SER:C	2.13	0.51
2:A:612:HOH:O	1:B:118:THR:CG2	2.59	0.51
1:A:135:THR:O	1:B:122:PHE:CZ	2.64	0.51
1:B:40:LEU:CD1	1:B:40:LEU:N	2.72	0.50
1:B:16:GLN:CG	1:B:17:ARG:N	2.73	0.50
1:A:169:SER:HB3	2:A:831:HOH:O	2.11	0.50
1:A:3:VAL:HB	1:A:101:PHE:O	2.11	0.50
1:B:84:GLU:O	1:B:85:THR:HG22	2.11	0.50
1:B:126:SER:O	1:B:128:GLU:N	2.44	0.50
1:A:125:SER:HB2	1:A:127:GLU:CG	2.42	0.50
1:A:183:SER:O	1:A:184:LEU:HD12	2.12	0.50
1:A:49:ILE:N	1:A:49:ILE:HD12	2.27	0.50
1:A:33:SER:OG	1:A:51:GLU:HA	2.10	0.50
1:A:43:THR:HG22	2:B:741:HOH:O	2.11	0.50
1:A:166:THR:CG2	2:A:631:HOH:O	2.42	0.49
2:A:688:HOH:O	1:B:166:THR:HG23	1.91	0.49
1:A:87:TYR:CE1	1:A:107:VAL:HG13	2.47	0.49
1:A:38:GLN:HB2	1:A:48:LEU:HD11	1.93	0.49
1:A:43:THR:O	1:A:44:ALA:HB2	2.13	0.49
1:B:80:GLN:O	1:B:83:ASP:HB2	2.12	0.49
1:B:1:PCA:N	1:B:96:LEU:HD21	2.27	0.49
1:A:35:THR:HG23	1:A:47:LEU:HD11	1.91	0.49
1:A:196:SER:HA	1:A:208:LYS:O	2.13	0.49
1:A:112:GLN:HB2	1:A:113:PRO:CD	2.43	0.49
1:B:128:GLU:O	1:B:128:GLU:HG2	2.13	0.49
1:B:154:ALA:HB2	1:B:195:TYR:CE1	2.48	0.49
1:A:80:GLN:HB2	1:A:81:PRO:HD2	1.94	0.49
1:B:17:ARG:HH11	1:B:17:ARG:HG2	1.78	0.48
1:B:43:THR:O	1:B:44:ALA:HB3	2.12	0.48
1:A:157:SER:HB3	2:A:635:HOH:O	2.13	0.48
1:B:145:PRO:O	1:B:201:HIS:HE1	1.96	0.48
1:B:196:SER:HB3	1:B:209:THR:HB	1.96	0.47
1:B:96:LEU:HA	1:B:96:LEU:HD13	1.39	0.47
1:B:84:GLU:O	1:B:85:THR:CG2	2.62	0.47
1:A:124:PRO:HD3	1:A:136:LEU:HD22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:ALA:CB	1:B:195:TYR:CE1	2.97	0.47
1:A:9:SER:HA	1:A:106:LYS:O	2.14	0.47
1:B:47:LEU:HD13	1:B:56:ALA:HB2	1.95	0.47
1:A:85:THR:HG23	1:A:86:ASP:O	2.15	0.47
1:A:197:CYS:N	1:A:208:LYS:O	2.47	0.47
1:A:65:ALA:HB2	1:A:74:LEU:HD12	1.97	0.47
1:A:107:VAL:O	1:A:107:VAL:HG22	2.15	0.47
1:A:135:THR:O	1:B:122:PHE:HZ	1.98	0.47
1:A:166:THR:HB	1:A:179:SER:O	2.15	0.47
1:B:129:LEU:C	1:B:131:ALA:H	2.16	0.47
1:B:193:ARG:O	1:B:212:PRO:HD2	2.16	0.46
1:A:126:SER:O	1:A:130:GLN:CG	2.54	0.46
1:A:62:ARG:NH2	1:A:83:ASP:OD2	2.47	0.46
1:A:182:LEU:CD1	1:A:184:LEU:CD1	2.93	0.46
1:A:99:ALA:HB3	1:B:37:TYR:OH	2.16	0.46
1:B:193:ARG:C	1:B:212:PRO:HD2	2.37	0.46
1:B:171:GLN:OE1	1:B:177:ALA:HB2	2.15	0.46
1:B:93:ASP:OD1	1:B:96:LEU:HD23	2.15	0.46
1:B:163:VAL:CB	2:B:882:HOH:O	2.56	0.45
1:B:211:ALA:HA	1:B:212:PRO:HD3	1.85	0.45
1:A:112:GLN:HG3	1:A:144:TYR:CD2	2.51	0.45
1:B:78:GLY:O	1:B:79:LEU:C	2.54	0.45
1:B:27:SER:O	1:B:32:ASN:ND2	2.47	0.45
1:A:213:THR:O	1:B:216:SER:O	2.34	0.45
1:A:40:LEU:C	1:A:42:GLY:N	2.69	0.45
1:A:85:THR:HG23	1:A:86:ASP:N	2.32	0.45
1:A:130:GLN:NE2	2:A:609:HOH:O	2.32	0.45
1:A:201:HIS:O	1:A:202:GLU:C	2.53	0.45
1:A:166:THR:HG22	1:A:178:ALA:HA	1.98	0.45
1:A:211:ALA:HA	1:A:212:PRO:HD3	1.77	0.45
1:A:203:GLY:C	2:A:656:HOH:O	2.54	0.45
2:A:631:HOH:O	1:B:166:THR:HG23	2.17	0.45
1:B:93:ASP:CG	1:B:96:LEU:HB2	2.38	0.45
1:A:10:ALA:O	1:A:107:VAL:HA	2.16	0.44
1:A:132:ASN:HA	1:A:186:PRO:HG2	1.99	0.44
1:B:117:PRO:HB3	1:B:143:PHE:HB3	1.98	0.44
1:B:40:LEU:HB2	1:B:43:THR:OG1	2.16	0.44
1:B:43:THR:O	1:B:44:ALA:HB2	2.18	0.44
1:B:103:THR:HG22	2:B:705:HOH:O	2.18	0.44
1:B:126:SER:O	1:B:127:GLU:C	2.56	0.44
1:A:35:THR:CG2	1:A:37:TYR:CZ	3.01	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:SER:OG	1:B:96:LEU:CD2	2.58	0.44
1:B:196:SER:HB2	1:B:209:THR:HG22	1.98	0.44
1:A:171:GLN:NE2	1:A:171:GLN:HA	2.32	0.43
1:A:18:VAL:HG13	2:A:530:HOH:O	2.17	0.43
1:A:37:TYR:N	1:A:88:TYR:O	2.38	0.43
1:A:87:TYR:CE1	1:A:107:VAL:CG1	3.01	0.43
1:A:48:LEU:HB3	1:A:49:ILE:HD12	2.00	0.43
1:A:95:SER:C	2:A:566:HOH:O	2.56	0.43
1:A:170:LYS:HB2	1:A:170:LYS:HE3	1.76	0.43
1:A:87:TYR:CD1	1:A:107:VAL:HG13	2.53	0.43
1:A:128:GLU:HG2	1:A:133:LYS:HB2	2.00	0.42
1:A:116:ASN:ND2	1:A:204:SER:OG	2.52	0.42
1:A:93:ASP:OD1	1:A:97:ASP:HA	2.19	0.42
1:B:52:ASP:HB3	2:B:728:HOH:O	2.19	0.42
1:A:182:LEU:HD12	1:A:184:LEU:CD1	2.49	0.42
1:B:96:LEU:N	1:B:96:LEU:HD22	2.34	0.42
1:A:175:LYS:N	1:A:175:LYS:CD	2.83	0.42
1:A:165:THR:HG23	1:A:180:SER:HB2	2.00	0.42
1:A:120:THR:HG22	1:A:122:PHE:CE1	2.54	0.42
1:A:85:THR:HG22	1:A:87:TYR:CE1	2.54	0.42
1:B:24:GLY:CA	1:B:29:ILE:HD12	2.49	0.42
1:B:144:TYR:CD1	1:B:145:PRO:HA	2.55	0.42
1:B:128:GLU:O	1:B:133:LYS:O	2.38	0.42
1:A:122:PHE:HA	1:A:123:PRO:HD2	1.88	0.42
1:A:55:ARG:NH1	1:A:63:PHE:O	2.49	0.42
1:A:29:ILE:CG2	1:A:69:GLY:O	2.62	0.42
1:A:40:LEU:HG	1:A:85:THR:OG1	2.20	0.42
1:A:150:VAL:HA	1:A:198:GLN:O	2.20	0.42
1:A:183:SER:C	1:A:184:LEU:CD1	2.88	0.41
1:B:79:LEU:HD23	1:B:79:LEU:HA	1.88	0.41
1:B:124:PRO:HD3	1:B:136:LEU:HG	2.01	0.41
1:A:153:LYS:HZ2	1:A:153:LYS:HG2	1.48	0.41
2:A:688:HOH:O	1:B:166:THR:HG21	2.00	0.41
1:A:171:GLN:NE2	2:A:832:HOH:O	2.54	0.41
1:A:128:GLU:CG	1:A:133:LYS:HE3	2.50	0.41
1:B:68:SER:N	1:B:71:SER:O	2.40	0.41
1:B:9:SER:HB2	1:B:106:LYS:HD3	2.01	0.41
1:B:214:GLU:HB3	2:B:853:HOH:O	2.20	0.41
1:A:48:LEU:CB	1:A:49:ILE:HD12	2.51	0.41
1:B:67:LYS:HA	1:B:71:SER:O	2.21	0.41
1:B:196:SER:CB	1:B:209:THR:HG22	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:560:HOH:O	1:B:39:HIS:HE1	2.04	0.41
1:A:196:SER:CA	1:A:208:LYS:O	2.69	0.41
1:B:170:LYS:HD2	1:B:176:TYR:CZ	2.56	0.40
1:A:56:ALA:O	1:A:57:SER:C	2.58	0.40
1:A:35:THR:CG2	1:A:37:TYR:CE1	3.03	0.40
1:A:116:ASN:HD22	1:A:201:HIS:HD2	1.68	0.40
1:A:160:LYS:NZ	2:A:521:HOH:O	2.36	0.40
1:B:82:GLU:H	1:B:82:GLU:HG3	1.51	0.40
1:A:35:THR:HG21	1:A:37:TYR:OH	2.20	0.40
1:B:154:ALA:HB2	1:B:195:TYR:CD1	2.57	0.40
1:B:3:VAL:O	2:B:702:HOH:O	2.22	0.40
1:B:67:LYS:HA	1:B:72:ALA:HA	2.03	0.40
1:A:196:SER:HA	1:A:209:THR:HA	2.04	0.40
2:A:612:HOH:O	1:B:118:THR:HG23	2.21	0.40
1:B:64:SER:HA	2:B:729:HOH:O	2.21	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:B:127:GLU:OE1	2:B:803:HOH:O[3_545]	2.15	0.05

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	214/216 (99%)	188 (88%)	19 (9%)	7 (3%)	5	2
1	B	214/216 (99%)	191 (89%)	16 (8%)	7 (3%)	5	2
All	All	428/432 (99%)	379 (89%)	35 (8%)	14 (3%)	5	2

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	ALA
1	A	93	ASP
1	B	2	SER
1	B	44	ALA
1	B	213	THR
1	B	215	CYS
1	A	30	GLY
1	A	31	GLU
1	A	214	GLU
1	A	43	THR
1	B	127	GLU
1	B	212	PRO
1	B	131	ALA
1	A	78	GLY

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	181/181 (100%)	134 (74%)	47 (26%)	0	0
1	B	181/181 (100%)	145 (80%)	36 (20%)	1	1
All	All	362/362 (100%)	279 (77%)	83 (23%)	1	1

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	23	SER
1	A	25	SER
1	A	26	SER
1	A	28	ASN
1	A	34	VAL
1	A	35	THR
1	A	40	LEU
1	A	43	THR
1	A	46	LYS

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Mol	Chain	Res	Type
1	A	51	GLU
1	A	53	ASN
1	A	62	ARG
1	A	67	LYS
1	A	68	SER
1	A	80	GLN
1	A	85	THR
1	A	93	ASP
1	A	95	SER
1	A	97	ASP
1	A	107	VAL
1	A	110	LEU
1	A	113	PRO
1	A	121	LEU
1	A	125	SER
1	A	127	GLU
1	A	136	LEU
1	A	149	THR
1	A	153	LYS
1	A	157	SER
1	A	163	VAL
1	A	166	THR
1	A	172	SER
1	A	175	LYS
1	A	182	LEU
1	A	183	SER
1	A	184	LEU
1	A	185	THR
1	A	187	GLU
1	A	188	GLN
1	A	193	ARG
1	A	200	THR
1	A	202	GLU
1	A	209	THR
1	A	210	VAL
1	A	214	GLU
1	A	215	CYS
1	B	17	ARG
1	B	21	SER
1	B	25	SER
1	B	27	SER
1	B	31	GLU

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Mol	Chain	Res	Type
1	B	32	ASN
1	B	38	GLN
1	B	40	LEU
1	B	43	THR
1	B	46	LYS
1	B	47	LEU
1	B	48	LEU
1	B	51	GLU
1	B	53	ASN
1	B	61	ASP
1	B	67	LYS
1	B	70	THR
1	B	73	SER
1	B	79	LEU
1	B	82	GLU
1	B	96	LEU
1	B	97	ASP
1	B	98	VAL
1	B	106	LYS
1	B	112	GLN
1	B	114	LYS
1	B	130	GLN
1	B	157	SER
1	B	170	LYS
1	B	174	ASN
1	B	179	SER
1	B	193	ARG
1	B	204	SER
1	B	209	THR
1	B	210	VAL
1	B	215	CYS

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	38	GLN
1	A	53	ASN
1	A	116	ASN
1	A	201	HIS
1	B	32	ASN
1	B	39	HIS
1	B	53	ASN

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Mol	Chain	Res	Type
1	B	112	GLN
1	B	192	HIS
1	B	201	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	PCA	A	1	1	6,7,9	0.68	0	7,8,12	1.03	0
1	PCA	B	1	1	6,7,9	0.64	0	7,8,12	1.82	2 (28%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PCA	A	1	1	-	0/0/9/13	0/1/1/1
1	PCA	B	1	1	-	0/0/9/13	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	PCA	O-C-CA	2.37	131.70	125.44
1	B	1	PCA	CB-CA-C	2.84	116.64	112.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1	PCA	1	0
1	B	1	PCA	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	215/216 (99%)	-0.74	2 (0%) 85 85	8, 22, 45, 55	0
1	B	215/216 (99%)	-0.86	1 (0%) 91 91	5, 21, 35, 53	0
All	All	430/432 (99%)	-0.80	3 (0%) 89 88	5, 21, 42, 55	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	215	CYS	2.2
1	A	216	SER	2.1
1	A	214	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	PCA	B	1	7/9	0.77	0.21	-	45,47,48,48	0
1	PCA	A	1	7/9	0.84	0.17	-	44,44,45,45	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.