



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:20 PM GMT

PDB ID : 1OID
Title : 5'-NUCLEOTIDASE (E. COLI) WITH AN ENGINEERED DISULFIDE BRIDGE (S228C, P513C)
Authors : Schultz-Heienbrok, R.; Maier, T.; Straeter, N.
Deposited on : 2003-06-13
Resolution : 2.10 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

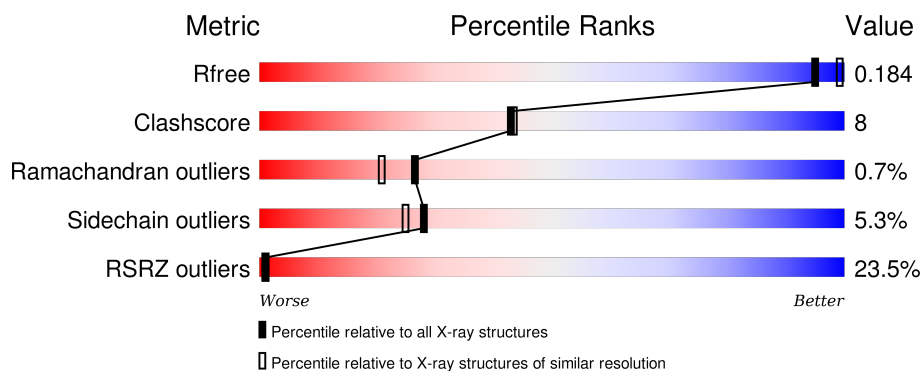
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	532	<div> <div>15%</div> <div>77%</div> <div>19%</div> <div>..</div> </div>
1	B	532	<div> <div>32%</div> <div>76%</div> <div>19%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8920 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 PROTEIN USHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	525	Total	C	N	O	S	0	0	0
			4098	2588	703	788	19			
1	B	523	Total	C	N	O	S	0	0	0
			4075	2572	699	785	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	CYS	SER	ENGINEERED MUTATION	UNP P07024
A	513	CYS	PRO	ENGINEERED MUTATION	UNP P07024
B	228	CYS	SER	ENGINEERED MUTATION	UNP P07024
B	513	CYS	PRO	ENGINEERED MUTATION	UNP P07024

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ni	0	0
			1	1		

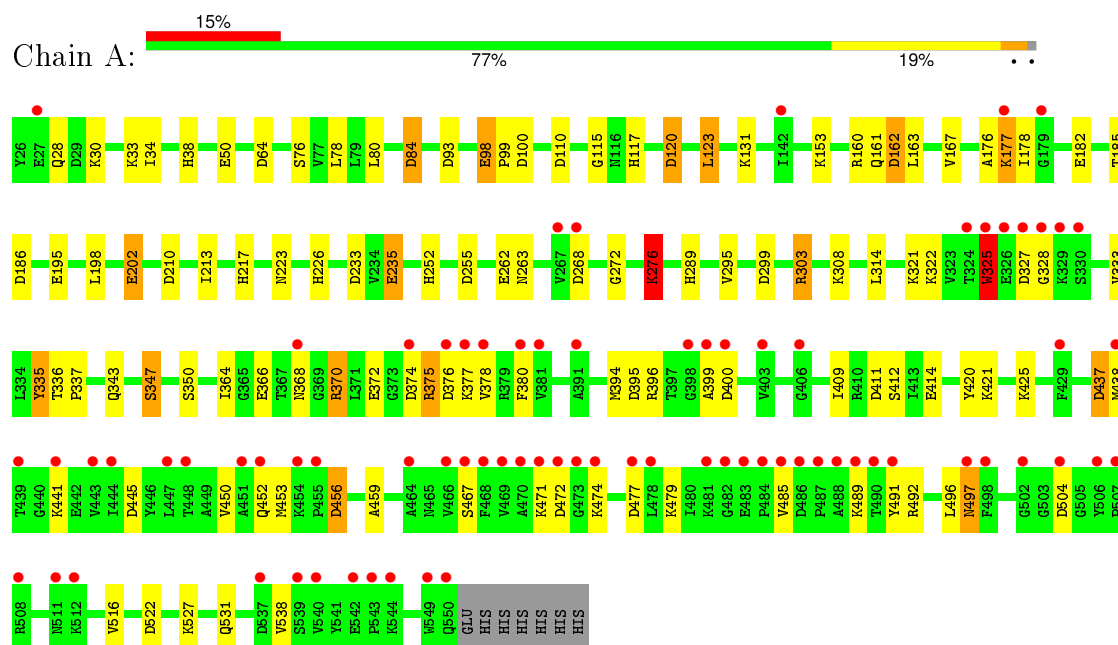
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	408	Total	O	0	0
			408	408		
3	B	338	Total	O	0	0
			338	338		

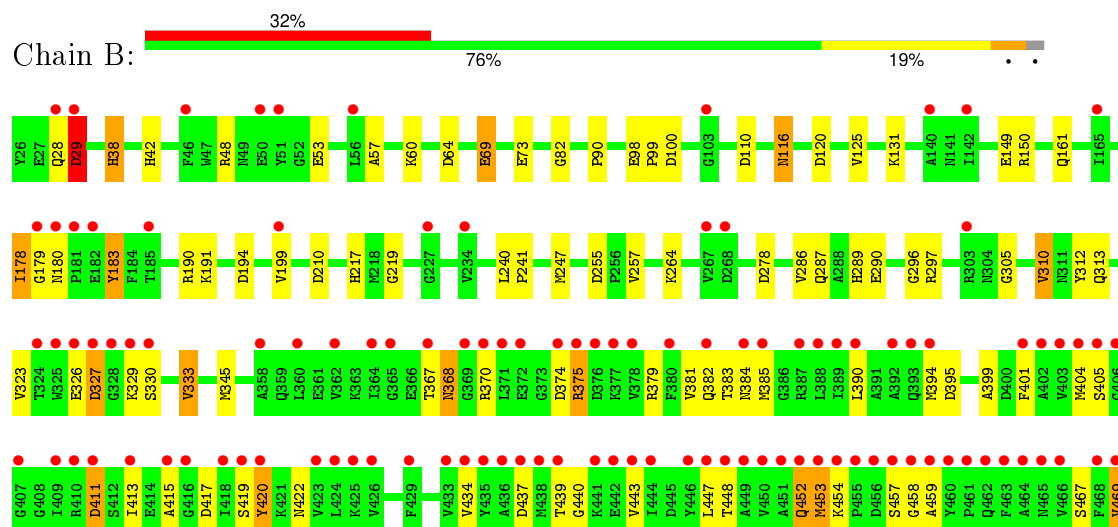
3 Residue-property plots

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 ($RSRZ > 2$). 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: PROTEIN USHA



• Molecule 1: PROTEIN USHA



L536	L537	D537	V538	S539	V540	V541	E542	P543	K544	G545	E546	V547	S548	TRP	GLN	GLU	HIS	HIS	HIS	HIS	HIS	HIS	L5470	K471	D472	K473	K474	L475	D476	D477	L478	K479	L480	K481	K482	E483	P484	V485	D486	P487	A488	K489	T490	T491	R492	M493	A494	T495	L496	N497	P498	N499	A500	T501	G502	G503	D504	G505	T506	P507	G515	V516	N517	L518	G519	F520	I521	D522	E523	E524	V525	L526	K527	E528	T529	I530	Q531	K532	S533	E534	E535
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.50Å 93.71Å 82.90Å 90.00° 97.71° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 19.94 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.3 (30.00-2.10) 98.4 (19.94-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.166 , 0.215 0.192 , 0.184	Depositor DCC
R_{free} test set	3529 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 66.4	EDS
Estimated twinning fraction	0.019 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 70046 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8920	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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

5.1 Standard geometry

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

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	1.44	15/4185 (0.4%)	1.24	39/5662 (0.7%)
1	B	1.28	17/4160 (0.4%)	1.16	27/5627 (0.5%)
All	All	1.36	32/8345 (0.4%)	1.20	66/11289 (0.6%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	333	VAL	CB-CG2	-10.24	1.31	1.52
1	B	333	VAL	CB-CG2	-9.18	1.33	1.52
1	A	347	SER	CB-OG	-8.95	1.30	1.42
1	A	235	GLU	CD-OE2	-8.45	1.16	1.25
1	B	69	GLU	CD-OE1	8.01	1.34	1.25
1	B	199	VAL	CB-CG1	-7.94	1.36	1.52
1	B	546	GLU	CD-OE1	7.42	1.33	1.25
1	A	98	GLU	CD-OE1	-7.21	1.17	1.25
1	B	69	GLU	CD-OE2	-6.82	1.18	1.25
1	A	343	GLN	CG-CD	6.77	1.66	1.51
1	A	347	SER	CA-CB	6.67	1.62	1.52
1	A	235	GLU	CG-CD	6.66	1.61	1.51
1	A	235	GLU	CD-OE1	-6.51	1.18	1.25
1	A	276	LYS	CD-CE	6.43	1.67	1.51
1	A	350	SER	CA-CB	6.21	1.62	1.52
1	A	202	GLU	CD-OE2	6.10	1.32	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	323	VAL	CB-CG2	-5.95	1.40	1.52
1	B	219	GLY	C-O	5.78	1.32	1.23
1	B	53	GLU	CD-OE1	-5.68	1.19	1.25
1	A	195	GLU	CD-OE2	-5.65	1.19	1.25
1	A	497	ASN	CB-CG	5.60	1.64	1.51
1	B	131	LYS	CD-CE	5.55	1.65	1.51
1	B	547	VAL	CA-CB	5.47	1.66	1.54
1	B	548	SER	CB-OG	5.45	1.49	1.42
1	B	247	MET	CG-SD	5.29	1.95	1.81
1	B	149	GLU	CD-OE1	-5.23	1.19	1.25
1	A	335	TYR	CE1-CZ	-5.17	1.31	1.38
1	A	272	GLY	C-O	-5.13	1.15	1.23
1	B	313	GLN	CB-CG	-5.09	1.38	1.52
1	B	497	ASN	CB-CG	5.09	1.62	1.51
1	B	125	VAL	CB-CG2	5.08	1.63	1.52
1	B	290	GLU	CG-CD	5.00	1.59	1.51

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	GLU	OE1-CD-OE2	-13.50	107.11	123.30
1	B	48	ARG	NE-CZ-NH1	13.28	126.94	120.30
1	B	64	ASP	CB-CG-OD1	11.16	128.34	118.30
1	B	210	ASP	CB-CG-OD2	10.76	127.99	118.30
1	A	162	ASP	CB-CG-OD2	10.71	127.94	118.30
1	B	48	ARG	NE-CZ-NH2	-9.66	115.47	120.30
1	B	255	ASP	CB-CG-OD2	8.69	126.12	118.30
1	A	347	SER	N-CA-CB	8.34	123.01	110.50
1	A	84	ASP	CB-CG-OD1	8.28	125.75	118.30
1	A	255	ASP	CB-CG-OD1	-8.14	110.97	118.30
1	A	303	ARG	NE-CZ-NH2	-8.13	116.24	120.30
1	B	255	ASP	CB-CG-OD1	-8.03	111.07	118.30
1	A	93	ASP	CB-CG-OD1	7.95	125.46	118.30
1	A	255	ASP	CB-CG-OD2	7.92	125.43	118.30
1	B	120	ASP	CB-CG-OD2	7.81	125.33	118.30
1	A	445	ASP	CB-CG-OD2	7.77	125.30	118.30
1	A	123	LEU	CB-CG-CD1	7.45	123.67	111.00
1	A	477	ASP	CB-CG-OD2	7.32	124.89	118.30
1	A	299	ASP	CB-CG-OD2	7.32	124.88	118.30
1	A	504	ASP	CB-CG-OD2	7.19	124.77	118.30
1	A	400	ASP	CB-CG-OD2	6.96	124.56	118.30
1	B	395	ASP	CB-CG-OD2	6.81	124.43	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	522	ASP	CB-CG-OD2	6.63	124.27	118.30
1	A	437	ASP	CB-CG-OD2	6.52	124.17	118.30
1	B	477	ASP	CB-CG-OD2	6.51	124.16	118.30
1	A	303	ARG	NE-CZ-NH1	6.46	123.53	120.30
1	A	347	SER	CA-CB-OG	-6.43	93.83	111.20
1	A	162	ASP	CB-CG-OD1	-6.26	112.67	118.30
1	A	110	ASP	CB-CG-OD2	6.21	123.89	118.30
1	B	374	ASP	CB-CG-OD2	6.19	123.87	118.30
1	A	78	LEU	CB-CG-CD1	6.12	121.41	111.00
1	A	235	GLU	CG-CD-OE1	6.12	130.54	118.30
1	B	194	ASP	CB-CG-OD2	6.04	123.73	118.30
1	A	370	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	B	297	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	A	210	ASP	CB-CG-OD2	5.84	123.55	118.30
1	A	100	ASP	CB-CG-OD2	5.82	123.54	118.30
1	B	110	ASP	CB-CG-OD2	5.80	123.52	118.30
1	B	504	ASP	CB-CG-OD2	5.79	123.51	118.30
1	A	268	ASP	CB-CG-OD2	5.72	123.45	118.30
1	A	325	TRP	CA-CB-CG	5.68	124.49	113.70
1	A	233	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	472	ASP	CB-CG-OD2	5.62	123.36	118.30
1	B	537	ASP	CB-CG-OD2	5.60	123.34	118.30
1	A	186	ASP	CB-CG-OD2	5.59	123.33	118.30
1	A	185	THR	OG1-CB-CG2	-5.56	97.22	110.00
1	A	327	ASP	CB-CG-OD2	5.53	123.28	118.30
1	B	29	ASP	CB-CG-OD2	5.51	123.26	118.30
1	A	370	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	278	ASP	CB-CG-OD2	5.49	123.24	118.30
1	A	347	SER	CB-CA-C	-5.38	99.89	110.10
1	B	437	ASP	CB-CG-OD2	5.36	123.12	118.30
1	A	522	ASP	CB-CG-OD1	5.32	123.08	118.30
1	B	327	ASP	CB-CG-OD2	5.30	123.07	118.30
1	B	69	GLU	CG-CD-OE2	-5.28	107.74	118.30
1	B	60	LYS	CD-CE-NZ	-5.18	99.79	111.70
1	B	150	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	B	310	VAL	CG1-CB-CG2	5.16	119.16	110.90
1	A	295	VAL	CG1-CB-CG2	-5.13	102.69	110.90
1	B	411	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	396	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	B	486	ASP	CB-CG-OD2	5.11	122.90	118.30
1	B	417	ASP	CB-CG-OD2	5.11	122.90	118.30
1	A	210	ASP	CB-CG-OD1	-5.07	113.73	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	ASP	CB-CG-OD2	5.06	122.86	118.30
1	A	64	ASP	CB-CG-OD1	5.05	122.84	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	437	ASP	Mainchain

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	4098	0	4015	58	1
1	B	4075	0	3997	74	0
2	A	1	0	0	0	0
3	A	408	0	0	13	1
3	B	338	0	0	9	0
All	All	8920	0	8012	132	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 8.

All (132) 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:38:HIS:CE1	1:B:296:GLY:HA3	1.80	1.15
1:B:401:PHE:CE2	1:B:493:MET:HG3	1.83	1.14
1:B:73:GLU:HG2	3:B:2058:HOH:O	1.59	1.02
1:A:84:ASP:CG	1:A:217:HIS:HE1	1.67	0.98
1:B:217:HIS:HE1	3:B:2099:HOH:O	1.51	0.94
1:B:401:PHE:CZ	1:B:493:MET:HB2	2.05	0.92
1:B:38:HIS:HE1	1:B:296:GLY:HA3	1.37	0.88
1:A:217:HIS:CD2	1:A:252:HIS:HB2	2.10	0.87
1:A:115:GLY:CA	1:A:217:HIS:ND1	2.40	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:401:PHE:CE2	1:B:493:MET:CG	2.60	0.84
1:A:115:GLY:HA2	1:A:217:HIS:ND1	1.92	0.83
1:A:30:LYS:NZ	3:A:2006:HOH:O	2.13	0.82
1:A:153:LYS:HD2	3:A:2066:HOH:O	1.80	0.80
1:A:84:ASP:CG	1:A:217:HIS:CE1	2.55	0.79
1:B:38:HIS:CD2	1:B:286:VAL:CG2	2.67	0.77
1:B:38:HIS:ND1	1:B:296:GLY:HA3	2.01	0.75
1:B:38:HIS:CD2	1:B:286:VAL:HG21	2.22	0.75
1:B:217:HIS:CE1	3:B:2099:HOH:O	2.34	0.72
1:B:401:PHE:CZ	1:B:493:MET:CB	2.72	0.71
1:A:117:HIS:HD2	1:A:120:ASP:OD2	1.73	0.71
1:A:115:GLY:HA3	1:A:217:HIS:ND1	2.05	0.70
1:B:28:GLN:HG3	1:B:29:ASP:OD1	1.91	0.69
1:B:116:ASN:H	1:B:116:ASN:HD22	1.41	0.69
1:B:401:PHE:CD2	1:B:493:MET:HG3	2.29	0.68
1:A:467:SER:OG	1:A:479:LYS:HB2	1.93	0.67
1:B:401:PHE:CZ	1:B:493:MET:HG3	2.29	0.67
1:B:401:PHE:CZ	1:B:493:MET:CG	2.79	0.66
1:B:38:HIS:CD2	1:B:286:VAL:HG23	2.31	0.65
1:A:153:LYS:CD	3:A:2066:HOH:O	2.42	0.65
1:B:367:THR:HG23	1:B:415:ALA:HA	1.80	0.63
1:A:527:LYS:HE2	1:A:531:GLN:NE2	2.17	0.60
1:B:394:MET:HG2	1:B:399:ALA:HB3	1.84	0.60
1:A:276:LYS:HE2	3:A:2264:HOH:O	2.01	0.60
1:B:178:ILE:HG22	1:B:179:GLY:H	1.65	0.60
1:B:180:ASN:O	1:B:183:TYR:HB3	2.03	0.59
1:A:84:ASP:OD1	1:A:217:HIS:HE1	1.86	0.58
1:B:116:ASN:ND2	1:B:217:HIS:HE1	2.02	0.58
1:A:177:LYS:C	1:A:178:ILE:HD12	2.24	0.58
1:B:439:THR:O	1:B:443:VAL:HG23	2.04	0.57
1:B:116:ASN:ND2	1:B:217:HIS:CE1	2.72	0.57
1:B:38:HIS:HD2	1:B:286:VAL:CG2	2.16	0.57
1:B:448:THR:O	1:B:452:GLN:NE2	2.38	0.56
1:B:385:MET:HE1	1:B:413:ILE:HD12	1.86	0.56
1:A:198:LEU:O	1:A:202:GLU:HG3	2.05	0.55
1:A:84:ASP:CB	1:A:217:HIS:HE1	2.19	0.55
1:A:226:HIS:CD2	1:A:235:GLU:OE2	2.60	0.55
1:A:325:TRP:NE1	3:A:2302:HOH:O	2.39	0.55
1:A:117:HIS:CD2	1:A:120:ASP:OD2	2.58	0.54
1:A:527:LYS:HE2	1:A:531:GLN:HE22	1.72	0.54
1:A:178:ILE:HG13	1:A:516:VAL:HG11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:ASN:OD1	1:A:538:VAL:HG22	2.08	0.54
1:B:384:ASN:ND2	3:B:2311:HOH:O	2.39	0.54
1:B:544:LYS:NZ	3:B:2338:HOH:O	2.41	0.54
1:B:82:GLY:O	1:B:217:HIS:HD2	1.91	0.53
1:A:226:HIS:HD2	1:A:235:GLU:OE2	1.91	0.53
1:A:380:PHE:CZ	1:A:456:ASP:HB2	2.44	0.53
1:B:440:GLY:HA3	1:B:486:ASP:O	2.10	0.52
1:B:240:LEU:HB3	1:B:241:PRO:HD2	1.93	0.51
1:A:376:ASP:O	1:A:377:LYS:C	2.49	0.51
1:B:191:LYS:HE2	3:B:2173:HOH:O	2.10	0.51
1:A:115:GLY:HA3	1:A:217:HIS:CE1	2.45	0.50
1:A:489:LYS:HB3	1:A:491:TYR:CE2	2.45	0.50
1:A:421:LYS:HD2	1:A:425:LYS:NZ	2.27	0.50
1:B:257:VAL:HG23	1:B:287:GLN:HB3	1.94	0.49
1:A:262:GLU:OE1	3:A:2244:HOH:O	2.20	0.49
1:A:372:GLU:OE2	1:A:374:ASP:HB3	2.12	0.49
1:B:116:ASN:HD21	1:B:217:HIS:HE1	1.59	0.49
1:A:153:LYS:CE	3:A:2066:HOH:O	2.60	0.49
1:B:405:SER:HB3	1:B:458:GLY:O	2.13	0.49
1:B:38:HIS:NE2	1:B:286:VAL:CG2	2.75	0.49
1:A:527:LYS:CE	3:A:2398:HOH:O	2.60	0.48
1:B:326:GLU:HG3	1:B:327:ASP:N	2.27	0.48
1:B:419:SER:H	1:B:422:ASN:ND2	2.11	0.48
1:B:401:PHE:CE1	1:B:493:MET:HB2	2.49	0.48
1:B:190:ARG:HD3	3:B:2170:HOH:O	2.14	0.48
1:B:38:HIS:HD2	1:B:286:VAL:HG23	1.78	0.48
1:A:84:ASP:OD1	1:A:217:HIS:CE1	2.66	0.48
1:B:486:ASP:OD1	1:B:489:LYS:HE3	2.13	0.48
1:A:176:ALA:HB3	3:A:2175:HOH:O	2.14	0.47
1:A:76:SER:HB2	1:A:160:ARG:HD2	1.97	0.47
1:B:401:PHE:CE2	1:B:493:MET:SD	3.08	0.47
1:B:183:TYR:C	1:B:183:TYR:CD1	2.88	0.47
1:A:399:ALA:HA	1:A:492:ARG:HB3	1.97	0.47
1:B:368:ASN:O	1:B:538:VAL:HG22	2.14	0.47
1:B:90:PRO:HB3	1:B:420:TYR:CE2	2.50	0.46
1:A:33:LYS:HE2	3:A:2007:HOH:O	2.15	0.46
1:A:34:ILE:HD11	1:A:163:LEU:HD12	1.97	0.46
1:B:379:ARG:NH1	1:B:458:GLY:N	2.64	0.46
1:B:467:SER:O	1:B:478:LEU:HA	2.15	0.46
1:B:98:GLU:N	1:B:99:PRO:HD2	2.30	0.46
1:A:84:ASP:CB	1:A:217:HIS:CE1	2.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:ALA:HA	1:B:345:MET:HG2	1.98	0.46
1:B:42:HIS:HE1	1:B:100:ASP:OD2	1.98	0.45
1:A:453:MET:HE2	1:A:459:ALA:HB1	1.98	0.45
1:B:453:MET:HG2	1:B:459:ALA:HB1	1.99	0.45
1:B:116:ASN:HD22	1:B:217:HIS:CE1	2.35	0.44
1:A:223:ASN:HB2	1:A:263:ASN:HB3	1.99	0.44
1:B:434:VAL:HG12	1:B:517:ASN:HA	1.99	0.44
1:B:38:HIS:CE1	1:B:312:TYR:OH	2.70	0.44
1:B:447:LEU:HD13	1:B:478:LEU:HD11	2.00	0.44
1:A:38:HIS:HA	1:A:80:LEU:O	2.17	0.44
1:B:448:THR:HG23	1:B:473:GLY:O	2.18	0.44
1:B:448:THR:OG1	1:B:475:LEU:CD1	2.66	0.44
1:B:161:GLN:NE2	3:B:2152:HOH:O	2.11	0.44
1:A:325:TRP:CE2	3:A:2302:HOH:O	2.57	0.43
1:B:116:ASN:H	1:B:116:ASN:ND2	2.12	0.43
1:B:38:HIS:CE1	1:B:296:GLY:CA	2.75	0.43
1:B:305:GLY:HA2	3:B:2193:HOH:O	2.18	0.43
1:A:364:ILE:HD13	1:A:527:LYS:HG2	2.01	0.42
1:A:497:ASN:HA	3:A:2390:HOH:O	2.19	0.42
1:A:375:ARG:CG	1:A:376:ASP:N	2.83	0.42
1:B:183:TYR:CD1	1:B:183:TYR:O	2.73	0.42
1:A:321:LYS:HB2	1:A:335:TYR:CZ	2.55	0.41
1:B:502:GLY:O	1:B:503:GLY:C	2.58	0.41
1:B:69:GLU:OE2	1:B:73:GLU:OE2	2.38	0.41
1:B:379:ARG:HD3	1:B:457:SER:C	2.41	0.41
1:A:98:GLU:N	1:A:99:PRO:HD2	2.35	0.41
1:B:390:LEU:HD21	1:B:404:MET:HG2	2.02	0.41
1:A:453:MET:HE3	3:A:2380:HOH:O	2.21	0.41
1:A:160:ARG:O	1:A:161:GLN:HB2	2.20	0.41
1:A:167:VAL:HA	1:A:213:ILE:O	2.21	0.41
1:B:447:LEU:CD1	1:B:478:LEU:HD11	2.51	0.41
1:A:378:VAL:HG11	1:A:409:ILE:HG21	2.03	0.41
1:B:469:VAL:HG13	1:B:548:SER:OG	2.20	0.41
1:A:394:MET:O	1:A:395:ASP:C	2.55	0.41
1:B:375:ARG:HG3	1:B:375:ARG:H	1.71	0.41
1:B:381:VAL:HG12	1:B:382:GLN:N	2.36	0.41
1:A:370:ARG:NH1	1:A:372:GLU:OE1	2.42	0.40
1:A:336:THR:HB	1:A:337:PRO:HD2	2.03	0.40
1:A:411:ASP:OD2	1:A:412:SER:N	2.45	0.40
1:A:314:LEU:HA	1:A:314:LEU:HD23	1.97	0.40
1:B:453:MET:CG	1:B:459:ALA:HB1	2.51	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:A:414:GLU:OE2	3:A:2286:HOH:O[2_945]	1.99	0.21

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	523/532 (98%)	495 (95%)	25 (5%)	3 (1%)	30	24
1	B	521/532 (98%)	496 (95%)	21 (4%)	4 (1%)	24	17
All	All	1044/1064 (98%)	991 (95%)	46 (4%)	7 (1%)	26	21

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	GLY
1	B	183	TYR
1	B	503	GLY
1	A	289	HIS
1	B	29	ASP
1	B	289	HIS
1	A	456	ASP

5.3.2 Protein sidechains [i](#)

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	433/440 (98%)	409 (94%)	24 (6%)	27	23
1	B	431/440 (98%)	409 (95%)	22 (5%)	29	26
All	All	864/880 (98%)	818 (95%)	46 (5%)	28	25

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	50	GLU
1	A	123	LEU
1	A	131	LYS
1	A	162	ASP
1	A	177	LYS
1	A	182	GLU
1	A	276	LYS
1	A	303	ARG
1	A	308	LYS
1	A	322	LYS
1	A	325	TRP
1	A	347	SER
1	A	366	GLU
1	A	375	ARG
1	A	420	TYR
1	A	438	MET
1	A	441	LYS
1	A	450	VAL
1	A	452	GLN
1	A	471	LYS
1	A	474	LYS
1	A	485	VAL
1	A	496	LEU
1	B	38	HIS
1	B	116	ASN
1	B	178	ILE
1	B	264	LYS
1	B	310	VAL
1	B	329	LYS
1	B	330	SER
1	B	333	VAL
1	B	368	ASN
1	B	370	ARG
1	B	375	ARG

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Mol	Chain	Res	Type
1	B	383	THR
1	B	411	ASP
1	B	420	TYR
1	B	452	GLN
1	B	453	MET
1	B	454	LYS
1	B	469	VAL
1	B	477	ASP
1	B	479	LYS
1	B	532	LYS
1	B	534	SER

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

Mol	Chain	Res	Type
1	A	43	HIS
1	A	117	HIS
1	A	204	GLN
1	A	217	HIS
1	A	226	HIS
1	A	266	GLN
1	A	289	HIS
1	A	531	GLN
1	B	38	HIS
1	B	42	HIS
1	B	116	ASN
1	B	217	HIS
1	B	279	GLN
1	B	353	GLN
1	B	422	ASN
1	B	431	ASN
1	B	452	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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	525/532 (98%)	0.61	78 (14%) 3 5	18, 32, 63, 82	0
1	B	523/532 (98%)	1.62	168 (32%) 1 1	22, 38, 80, 94	0
All	All	1048/1064 (98%)	1.11	246 (23%) 1 1	18, 35, 75, 94	0

All (246) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	401	PHE	10.2
1	B	491	TYR	9.8
1	B	463	PHE	9.8
1	B	454	LYS	9.5
1	B	403	VAL	9.2
1	B	389	ILE	9.1
1	B	402	ALA	8.5
1	B	490	THR	8.5
1	A	325	TRP	8.5
1	B	530	ILE	8.2
1	B	494	ALA	7.9
1	B	526	LEU	7.8
1	B	464	ALA	7.6
1	B	523	ALA	7.2
1	B	543	PRO	7.1
1	B	547	VAL	7.0
1	B	466	VAL	7.0
1	B	443	VAL	6.9
1	A	490	THR	6.8
1	A	484	PRO	6.7
1	B	478	LEU	6.6
1	A	473	GLY	6.6
1	B	404	MET	6.5
1	B	480	ILE	6.5

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Mol	Chain	Res	Type	RSRZ
1	B	493	MET	6.4
1	B	469	VAL	6.4
1	B	482	GLY	6.4
1	B	495	THR	6.3
1	B	473	GLY	6.3
1	B	520	PHE	6.3
1	B	502	GLY	6.2
1	A	488	ALA	6.2
1	B	423	VAL	6.2
1	B	434	VAL	6.2
1	A	482	GLY	6.1
1	B	436	ALA	6.1
1	B	324	THR	5.9
1	B	458	GLY	5.9
1	B	470	ALA	5.8
1	A	326	GLU	5.8
1	B	388	LEU	5.8
1	B	390	LEU	5.8
1	B	420	TYR	5.7
1	B	447	LEU	5.6
1	B	329	LYS	5.6
1	B	472	ASP	5.5
1	B	446	TYR	5.5
1	B	376	ASP	5.4
1	B	539	SER	5.4
1	B	468	PHE	5.4
1	B	527	LYS	5.4
1	B	409	ILE	5.3
1	B	327	ASP	5.3
1	B	180	ASN	5.3
1	B	485	VAL	5.2
1	B	459	ALA	5.2
1	A	328	GLY	5.1
1	B	433	VAL	5.1
1	A	485	VAL	5.0
1	B	378	VAL	5.0
1	B	393	GLN	4.9
1	A	448	THR	4.9
1	B	484	PRO	4.8
1	B	471	LYS	4.8
1	B	267	VAL	4.8
1	B	516	VAL	4.8

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Mol	Chain	Res	Type	RSRZ
1	B	525	VAL	4.8
1	B	416	GLY	4.8
1	B	367	THR	4.8
1	A	327	ASP	4.8
1	B	182	GLU	4.7
1	A	469	VAL	4.7
1	A	486	ASP	4.7
1	B	435	TYR	4.7
1	B	521	ILE	4.6
1	B	450	VAL	4.6
1	A	451	ALA	4.5
1	B	438	MET	4.5
1	B	456	ASP	4.5
1	A	454	LYS	4.4
1	B	475	LEU	4.4
1	B	536	LEU	4.3
1	B	418	ILE	4.3
1	B	462	GLN	4.2
1	B	56	LEU	4.1
1	A	439	THR	4.1
1	B	515	TYR	4.1
1	B	360	LEU	4.1
1	A	483	GLU	4.1
1	A	489	LYS	4.1
1	B	424	LEU	4.1
1	B	326	GLU	4.0
1	A	467	SER	4.0
1	A	470	ALA	4.0
1	B	519	GLY	4.0
1	B	364	ILE	4.0
1	A	477	ASP	3.9
1	B	489	LYS	3.9
1	B	518	THR	3.8
1	A	452	GLN	3.8
1	B	429	PHE	3.8
1	B	330	SER	3.7
1	B	380	PHE	3.7
1	B	375	ARG	3.7
1	A	487	PRO	3.7
1	B	268	ASP	3.7
1	B	451	ALA	3.7
1	B	524	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	460	TYR	3.7
1	B	453	MET	3.6
1	B	528	ALA	3.6
1	A	268	ASP	3.6
1	B	439	THR	3.6
1	B	544	LYS	3.6
1	B	465	ASN	3.6
1	A	504	ASP	3.5
1	A	329	LYS	3.5
1	B	457	SER	3.5
1	A	398	GLY	3.5
1	A	429	PHE	3.5
1	A	539	SER	3.5
1	B	548	SER	3.5
1	B	488	ALA	3.5
1	B	181	PRO	3.4
1	B	517	ASN	3.4
1	A	502	GLY	3.4
1	B	541	TYR	3.4
1	B	29	ASP	3.4
1	B	534	SER	3.4
1	A	399	ALA	3.4
1	A	471	LYS	3.3
1	B	474	LYS	3.3
1	A	466	VAL	3.3
1	A	540	VAL	3.3
1	B	405	SER	3.3
1	A	472	ASP	3.3
1	A	498	PHE	3.3
1	A	481	LYS	3.3
1	B	498	PHE	3.3
1	B	481	LYS	3.3
1	B	382	GLN	3.3
1	B	385	MET	3.3
1	B	452	GLN	3.3
1	A	324	THR	3.2
1	B	477	ASP	3.2
1	B	476	ASN	3.2
1	A	507	PRO	3.1
1	B	461	PRO	3.1
1	B	371	LEU	3.1
1	B	444	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	267	VAL	3.1
1	A	377	LYS	3.1
1	B	479	LYS	3.1
1	B	503	GLY	3.1
1	B	142	ILE	3.1
1	A	391	ALA	3.0
1	B	542	GLU	3.0
1	A	438	MET	3.0
1	A	455	PRO	3.0
1	B	411	ASP	3.0
1	B	413	ILE	3.0
1	B	51	TYR	3.0
1	A	550	GLN	3.0
1	A	537	ASP	2.9
1	A	378	VAL	2.9
1	A	512	LYS	2.9
1	A	508	ARG	2.9
1	A	544	LYS	2.9
1	B	545	GLY	2.9
1	B	496	LEU	2.9
1	A	368	ASN	2.9
1	B	392	ALA	2.9
1	A	474	LYS	2.8
1	A	443	VAL	2.8
1	B	46	PHE	2.8
1	B	384	ASN	2.8
1	A	376	ASP	2.8
1	B	504	ASP	2.8
1	A	403	VAL	2.8
1	B	426	VAL	2.8
1	B	328	GLY	2.8
1	B	500	ALA	2.8
1	B	325	TRP	2.7
1	B	415	ALA	2.7
1	B	441	LYS	2.7
1	B	50	GLU	2.7
1	A	542	GLU	2.6
1	A	380	PHE	2.6
1	B	425	LYS	2.6
1	B	394	MET	2.6
1	B	28	GLN	2.6
1	B	362	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	506	TYR	2.6
1	B	419	SER	2.6
1	B	372	GLU	2.6
1	B	227	GLY	2.5
1	B	448	THR	2.5
1	A	374	ASP	2.5
1	A	549	TRP	2.5
1	A	381	VAL	2.5
1	B	179	GLY	2.5
1	A	400	ASP	2.5
1	B	370	ARG	2.5
1	B	406	GLY	2.5
1	A	468	PHE	2.5
1	B	546	GLU	2.5
1	A	497	ASN	2.5
1	A	447	LEU	2.5
1	B	387	ARG	2.5
1	B	103	GLY	2.4
1	A	441	LYS	2.4
1	B	410	ARG	2.4
1	B	455	PRO	2.4
1	B	507	PRO	2.4
1	A	491	TYR	2.4
1	B	449	ALA	2.4
1	A	179	GLY	2.4
1	B	538	VAL	2.3
1	A	506	TYR	2.3
1	A	464	ALA	2.3
1	A	543	PRO	2.3
1	B	199	VAL	2.3
1	B	407	GLY	2.3
1	A	27	GLU	2.3
1	B	540	VAL	2.3
1	B	165	ILE	2.3
1	B	140	ALA	2.3
1	A	177	LYS	2.2
1	B	185	THR	2.2
1	A	330	SER	2.2
1	B	369	GLY	2.2
1	B	442	GLU	2.2
1	B	365	GLY	2.2
1	B	374	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	234	VAL	2.2
1	A	478	LEU	2.1
1	B	522	ASP	2.1
1	B	303	ARG	2.1
1	B	377	LYS	2.1
1	B	358	ALA	2.1
1	A	142	ILE	2.1
1	A	444	ILE	2.1
1	A	406	GLY	2.1
1	A	511	ASN	2.1
1	B	505	GLY	2.0
1	B	437	ASP	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](#)

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(\AA^2)	Q<0.9
2	NI	A	1551	1/1	0.97	0.12	0.47	66,66,66,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.