



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:19 PM GMT

PDB ID : 1T0L
Title : Crystal structure of human cytosolic NADP(+)-dependent isocitrate dehydrogenase in complex with NADP, isocitrate, and calcium(2+)
Authors : Xu, X.; Zhao, J.; Peng, B.; Huang, Q.; Arnold, E.; Ding, J.
Deposited on : 2004-04-10
Resolution : 2.41 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

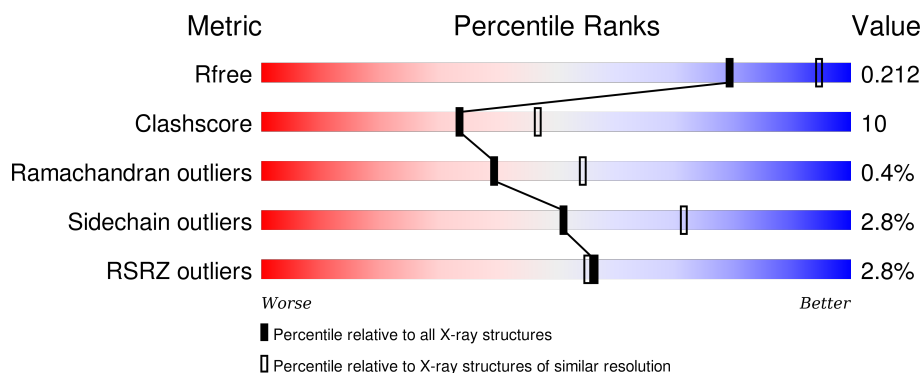
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (2.44-2.40)

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	414	<div> <div>2%</div> <div>84%</div> <div>15%</div> <div>.</div> </div>
1	B	414	<div> <div>3%</div> <div>83%</div> <div>15%</div> <div>.</div> </div>
1	C	414	<div> <div>2%</div> <div>84%</div> <div>15%</div> <div>.</div> </div>
1	D	414	<div> <div>3%</div> <div>83%</div> <div>14%</div> <div>.</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ICT	A	1616	-	-	-	X
4	ICT	B	1617	-	-	-	X
4	ICT	C	1618	-	-	-	X
4	ICT	D	1619	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14286 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.

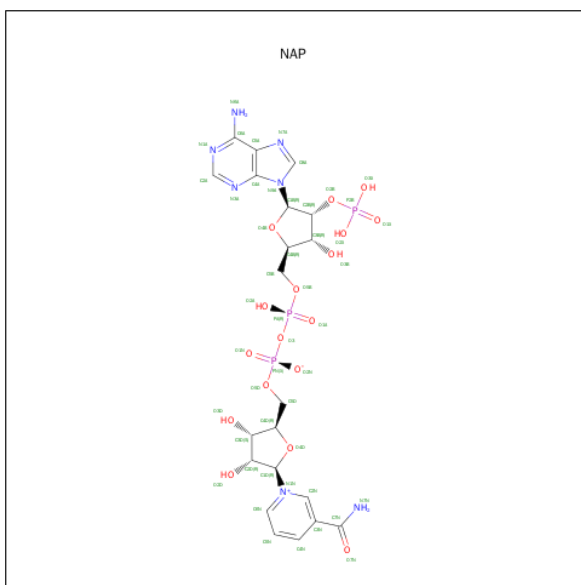
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP] cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	0	0
			3281	2085	556	621	19			
1	B	414	Total	C	N	O	S	0	0	0
			3281	2085	556	621	19			
1	C	414	Total	C	N	O	S	0	0	0
			3281	2085	556	621	19			
1	D	414	Total	C	N	O	S	0	0	0
			3281	2085	556	621	19			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

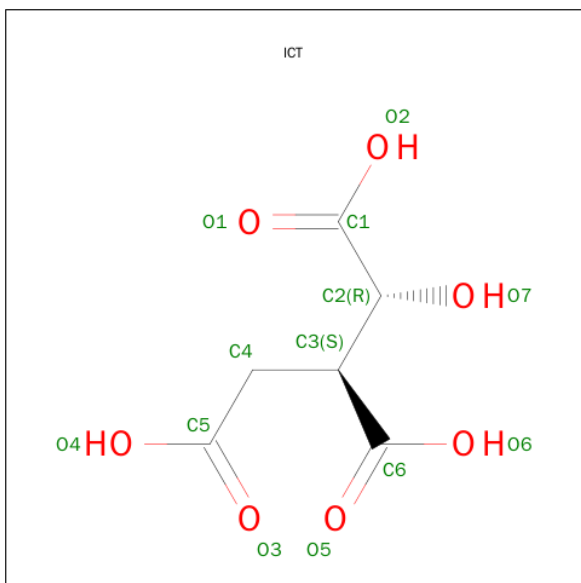
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
3	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
3	C	1	Total 48	C 21	N 7	O 17	P 3	0	0
3	D	1	Total 48	C 21	N 7	O 17	P 3	0	0

- Molecule 4 is ISOCITRIC ACID (three-letter code: ICT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	6	7		
4	B	1	Total	C	O	0	0
			13	6	7		
4	C	1	Total	C	O	0	0
			13	6	7		
4	D	1	Total	C	O	0	0
			13	6	7		

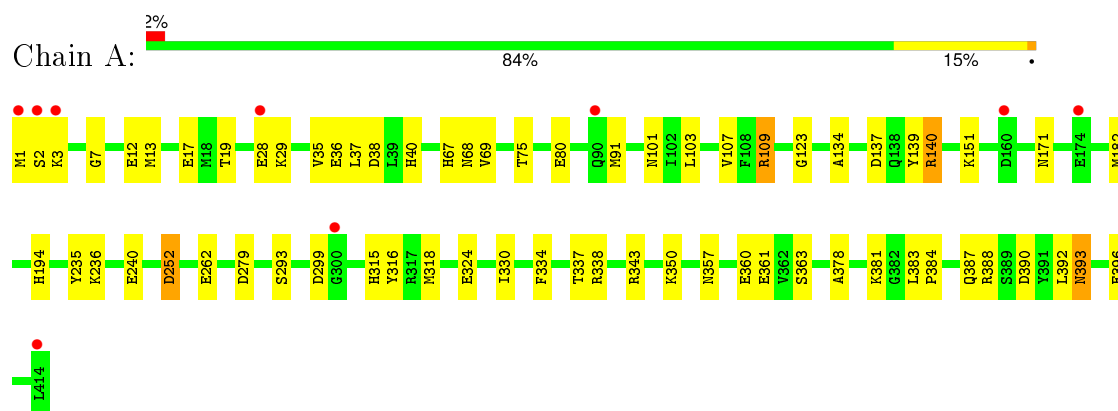
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	227	Total	O	0	0
			227	227		
5	B	221	Total	O	0	0
			221	221		
5	C	226	Total	O	0	0
			226	226		
5	D	240	Total	O	0	0
			240	240		

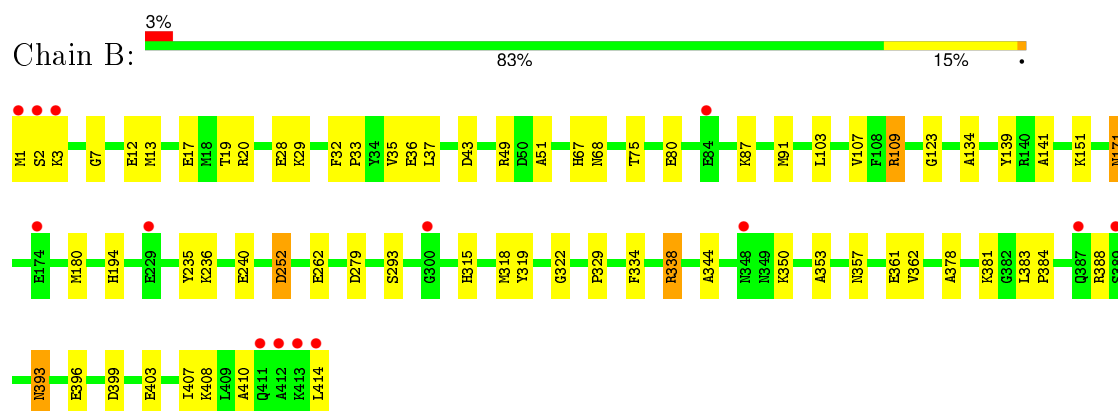
3 Residue-property plots [i](#)

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.

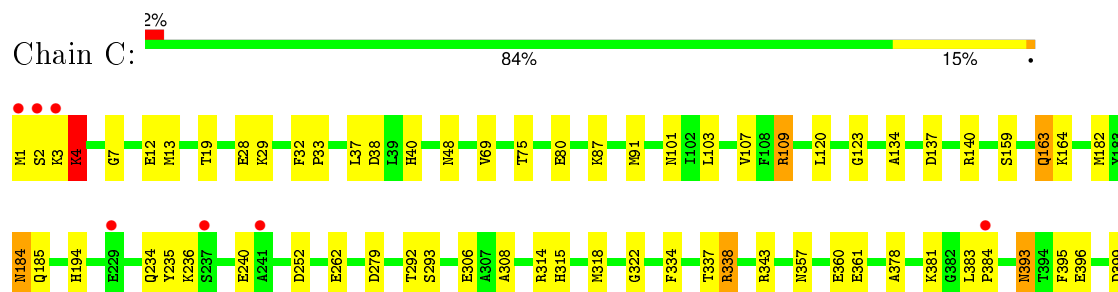
- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



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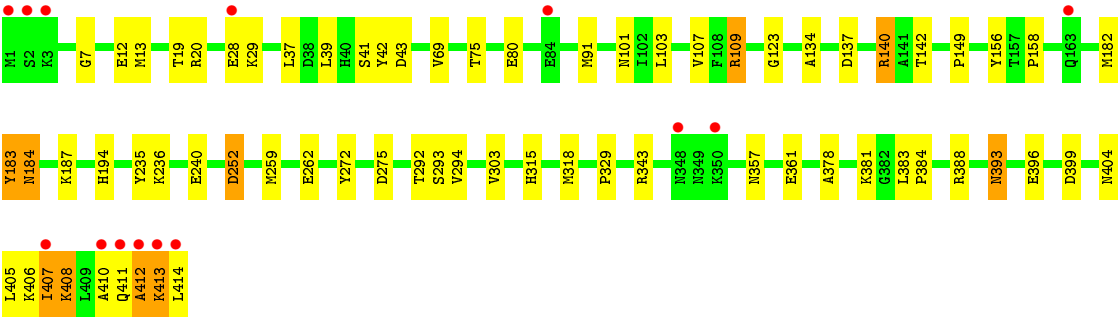
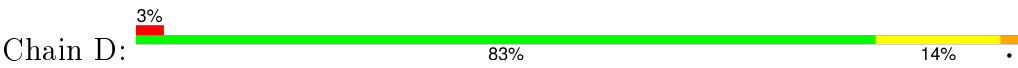


- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic





● Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	103.30Å 86.74Å 115.77Å 90.00° 107.15° 90.00°	Depositor
Resolution (Å)	12.00 – 2.41 11.99 – 2.41	Depositor EDS
% Data completeness (in resolution range)	94.2 (12.00-2.41) 94.3 (11.99-2.41)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.91 (at 2.40Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.214 , 0.253 0.216 , 0.212	Depositor DCC
R_{free} test set	3562 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.390	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 68.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 70684 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14286	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 8.22% 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: CA, NAP, ICT

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	0/3349	0.62	0/4513
1	B	0.34	0/3349	0.61	0/4513
1	C	0.35	0/3349	0.62	0/4513
1	D	0.35	0/3349	0.62	0/4513
All	All	0.35	0/13396	0.62	0/18052

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	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	183	TYR	Sidechain

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	3281	0	3256	59	0
1	B	3281	0	3256	57	0
1	C	3281	0	3256	63	0
1	D	3281	0	3256	94	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	48	0	25	3	0
3	B	48	0	25	4	0
3	C	48	0	25	5	0
3	D	48	0	25	4	0
4	A	13	0	5	3	0
4	B	13	0	5	2	0
4	C	13	0	5	3	0
4	D	13	0	5	2	0
5	A	227	0	0	3	0
5	B	221	0	0	4	0
5	C	226	0	0	4	0
5	D	240	0	0	4	0
All	All	14286	0	13144	270	0

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 10.

All (270) 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:D:407:ILE:HA	1:D:413:LYS:HB2	1.36	1.07
1:B:51:ALA:HB1	1:C:87:LYS:HG3	1.35	1.06
1:D:404:ASN:O	1:D:407:ILE:HB	1.55	1.05
1:C:393:ASN:ND2	1:C:396:GLU:H	1.66	0.94
1:D:412:ALA:O	1:D:413:LYS:HG3	1.68	0.93
1:A:3:LYS:HD3	1:A:36:GLU:HG3	1.50	0.91
1:D:407:ILE:HA	1:D:413:LYS:CB	2.01	0.89
1:D:407:ILE:CA	1:D:413:LYS:HB2	2.02	0.88
1:C:3:LYS:HB3	5:C:1875:HOH:O	1.72	0.88
1:C:1:MET:HA	1:C:32:PHE:HB3	1.56	0.88
1:D:407:ILE:HG23	1:D:413:LYS:N	1.88	0.87
1:A:393:ASN:C	1:A:393:ASN:HD22	1.78	0.87
1:A:393:ASN:ND2	1:A:396:GLU:H	1.73	0.86
1:D:137:ASP:HA	1:D:182:MET:HE3	1.64	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:393:ASN:C	1:D:393:ASN:HD22	1.85	0.79
1:D:407:ILE:HG12	1:D:413:LYS:HB2	1.62	0.79
1:A:2:SER:O	1:A:3:LYS:HG3	1.83	0.79
1:D:410:ALA:HB2	1:D:413:LYS:HZ1	1.48	0.79
1:A:3:LYS:HB3	1:A:36:GLU:HB2	1.65	0.77
1:D:404:ASN:HA	1:D:407:ILE:HG13	1.65	0.77
1:B:3:LYS:HG2	1:B:32:PHE:O	1.85	0.77
1:D:406:LYS:C	1:D:413:LYS:HD2	2.05	0.77
1:A:151:LYS:HE2	1:A:171:ASN:HD21	1.49	0.77
1:B:51:ALA:CB	1:C:87:LYS:HG3	2.14	0.76
1:B:315:HIS:HA	1:B:318:MET:HE3	1.68	0.75
1:A:137:ASP:HA	1:A:182:MET:HE3	1.69	0.75
1:C:393:ASN:HD21	1:C:396:GLU:H	1.33	0.73
1:D:407:ILE:HG12	1:D:413:LYS:CB	2.19	0.73
1:D:393:ASN:ND2	1:D:396:GLU:H	1.85	0.73
1:C:393:ASN:HD22	1:C:393:ASN:C	1.92	0.72
1:B:1:MET:H3	1:B:3:LYS:HE3	1.56	0.71
1:D:404:ASN:HA	1:D:407:ILE:CD1	2.21	0.70
1:C:292:THR:HG21	1:C:338:ARG:NH1	2.06	0.70
1:C:3:LYS:O	1:C:4:LYS:HB2	1.91	0.70
1:B:151:LYS:HG3	1:B:171:ASN:ND2	2.07	0.70
1:B:3:LYS:HD3	1:B:36:GLU:HA	1.73	0.70
1:B:7:GLY:HA3	1:B:37:LEU:HD23	1.74	0.70
1:D:404:ASN:O	1:D:407:ILE:CB	2.36	0.69
1:A:3:LYS:HE3	1:A:36:GLU:HA	1.73	0.69
1:B:393:ASN:C	1:B:393:ASN:HD22	1.94	0.69
1:C:101:ASN:HB2	5:C:1908:HOH:O	1.92	0.69
1:A:1:MET:HE2	1:A:3:LYS:HD2	1.75	0.69
1:D:406:LYS:HB3	1:D:413:LYS:HD2	1.74	0.68
1:D:404:ASN:HA	1:D:407:ILE:CG1	2.23	0.68
1:D:407:ILE:HG12	1:D:413:LYS:C	2.13	0.68
1:D:410:ALA:HB2	1:D:413:LYS:NZ	2.09	0.67
1:C:279:ASP:OD2	1:D:252:ASP:HB3	1.95	0.67
1:B:393:ASN:ND2	1:B:396:GLU:H	1.93	0.67
1:A:3:LYS:CD	1:A:36:GLU:HG3	2.24	0.66
1:D:414:LEU:HD23	5:D:1768:HOH:O	1.95	0.66
1:C:75:THR:H	3:C:1517:NAP:H71N	1.44	0.65
1:D:407:ILE:CG1	1:D:413:LYS:HB2	2.25	0.65
1:B:29:LYS:NZ	1:B:399:ASP:OD1	2.30	0.65
1:D:407:ILE:HG23	1:D:413:LYS:H	1.62	0.64
1:A:7:GLY:HA3	1:A:37:LEU:HD23	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:ASN:HD21	1:A:396:GLU:H	1.45	0.64
1:D:101:ASN:HB2	5:D:1734:HOH:O	1.97	0.64
1:D:412:ALA:O	1:D:413:LYS:CG	2.44	0.63
1:B:1:MET:N	1:B:3:LYS:HE3	2.11	0.63
1:A:393:ASN:C	1:A:393:ASN:ND2	2.52	0.62
1:B:151:LYS:HG3	1:B:171:ASN:HD21	1.65	0.62
1:B:3:LYS:CD	1:B:36:GLU:HA	2.30	0.62
1:C:75:THR:O	3:C:1517:NAP:H2N	2.00	0.62
1:C:2:SER:HB2	1:C:32:PHE:O	1.99	0.61
1:D:187:LYS:HE3	5:D:1954:HOH:O	2.00	0.61
1:C:38:ASP:OD2	1:C:40:HIS:HE1	1.82	0.61
1:B:75:THR:O	3:B:1516:NAP:H2N	2.00	0.61
1:D:156:TYR:CE2	1:D:158:PRO:HG3	2.36	0.61
1:D:39:LEU:HD12	1:D:41:SER:OG	2.01	0.60
3:B:1516:NAP:C4N	4:B:1617:ICT:H2	2.31	0.60
1:A:75:THR:H	3:A:1515:NAP:H71N	1.50	0.60
1:A:279:ASP:OD2	1:B:252:ASP:HB3	2.01	0.60
1:C:2:SER:HB2	1:C:33:PRO:HA	1.84	0.60
1:B:75:THR:H	3:B:1516:NAP:H71N	1.49	0.58
1:A:315:HIS:HA	1:A:318:MET:CE	2.33	0.58
1:C:2:SER:CB	1:C:33:PRO:HA	2.33	0.58
1:D:406:LYS:O	1:D:413:LYS:HD2	2.02	0.58
1:C:137:ASP:HA	1:C:182:MET:HE3	1.85	0.58
1:C:383:LEU:HB3	1:C:384:PRO:HD3	1.86	0.58
1:D:315:HIS:HA	1:D:318:MET:HE2	1.86	0.57
1:A:194:HIS:HD2	1:A:235:TYR:OH	1.87	0.57
1:B:383:LEU:HB3	1:B:384:PRO:HD3	1.86	0.57
1:C:194:HIS:HD2	1:C:235:TYR:OH	1.87	0.57
1:B:350:LYS:HG2	5:B:1743:HOH:O	2.04	0.57
1:D:393:ASN:ND2	1:D:396:GLU:HG3	2.19	0.57
1:D:194:HIS:HD2	1:D:235:TYR:OH	1.88	0.56
1:B:315:HIS:HA	1:B:318:MET:CE	2.34	0.56
1:B:3:LYS:HD3	1:B:35:VAL:O	2.05	0.56
1:B:151:LYS:NZ	1:B:171:ASN:HD21	2.04	0.56
1:D:383:LEU:HB3	1:D:384:PRO:HD3	1.86	0.56
1:B:87:LYS:HE2	1:C:48:ASN:OD1	2.06	0.56
1:A:383:LEU:HB3	1:A:384:PRO:HD3	1.87	0.56
1:C:393:ASN:ND2	1:C:396:GLU:HG3	2.20	0.56
3:D:1518:NAP:C4N	4:D:1619:ICT:H2	2.36	0.55
1:C:123:GLY:O	1:C:262:GLU:HA	2.06	0.55
1:B:3:LYS:HD2	1:B:36:GLU:CD	2.26	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:407:ILE:HD13	1:D:414:LEU:C	2.27	0.55
4:A:1616:ICT:H41	5:B:1886:HOH:O	2.07	0.55
1:C:69:VAL:HG11	1:C:343:ARG:HD2	1.88	0.55
1:D:406:LYS:CB	1:D:413:LYS:HD2	2.38	0.54
1:C:2:SER:O	1:C:3:LYS:HG3	2.06	0.54
1:B:3:LYS:HB2	1:B:36:GLU:OE1	2.07	0.54
1:B:194:HIS:HD2	1:B:235:TYR:OH	1.90	0.54
1:D:407:ILE:HA	1:D:413:LYS:CG	2.37	0.54
1:D:407:ILE:O	1:D:412:ALA:O	2.26	0.53
1:A:75:THR:O	3:A:1515:NAP:H2N	2.09	0.53
1:A:69:VAL:HG11	1:A:343:ARG:HD2	1.90	0.53
1:A:393:ASN:ND2	1:A:396:GLU:HG3	2.24	0.53
1:B:3:LYS:HD2	1:B:36:GLU:OE2	2.09	0.53
1:D:107:VAL:HG23	1:D:134:ALA:HB2	1.90	0.53
1:D:378:ALA:O	1:D:381:LYS:O	2.27	0.53
1:C:140:ARG:HG3	1:C:140:ARG:O	2.09	0.53
1:C:80:GLU:HG3	1:C:91:MET:CE	2.39	0.52
1:A:107:VAL:HG23	1:A:134:ALA:HB2	1.91	0.52
1:C:109:ARG:HD3	1:C:293:SER:OG	2.10	0.52
1:A:378:ALA:O	1:A:381:LYS:O	2.27	0.52
1:D:413:LYS:O	1:D:414:LEU:CB	2.57	0.52
3:C:1517:NAP:C4N	4:C:1618:ICT:H2	2.40	0.52
1:A:123:GLY:O	1:A:262:GLU:HA	2.09	0.52
1:D:407:ILE:HG22	1:D:412:ALA:HA	1.91	0.51
1:C:2:SER:OG	1:C:33:PRO:HA	2.10	0.51
1:B:378:ALA:O	1:B:381:LYS:O	2.28	0.51
1:B:80:GLU:HG3	1:B:91:MET:CE	2.40	0.51
1:A:3:LYS:HD3	1:A:36:GLU:CG	2.34	0.51
1:D:315:HIS:HA	1:D:318:MET:CE	2.40	0.51
1:C:378:ALA:O	1:C:381:LYS:O	2.28	0.51
1:B:141:ALA:HB1	1:B:180:MET:HE3	1.93	0.51
1:D:80:GLU:HG3	1:D:91:MET:CE	2.41	0.51
1:B:107:VAL:HG23	1:B:134:ALA:HB2	1.92	0.51
1:B:49:ARG:HD2	5:B:1736:HOH:O	2.11	0.51
1:D:39:LEU:CD1	1:D:41:SER:OG	2.58	0.51
1:D:404:ASN:CA	1:D:407:ILE:HG13	2.39	0.51
1:D:407:ILE:HG23	1:D:412:ALA:C	2.31	0.50
1:D:361:GLU:OE2	1:D:408:LYS:NZ	2.39	0.50
1:C:80:GLU:HG3	1:C:91:MET:HE1	1.92	0.50
1:D:142:THR:HG23	1:D:142:THR:O	2.12	0.50
1:B:109:ARG:HD3	1:B:293:SER:OG	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:315:HIS:HA	1:C:318:MET:CE	2.41	0.50
1:D:404:ASN:HD22	1:D:407:ILE:HD11	1.77	0.50
1:D:407:ILE:HG22	1:D:408:LYS:N	2.27	0.50
1:D:404:ASN:ND2	1:D:407:ILE:HD11	2.27	0.50
1:C:393:ASN:HD21	1:C:396:GLU:N	2.06	0.50
1:D:404:ASN:HA	1:D:407:ILE:HD12	1.92	0.50
3:A:1515:NAP:C4N	4:A:1616:ICT:H2	2.41	0.50
1:A:80:GLU:HG3	1:A:91:MET:CE	2.42	0.49
1:D:75:THR:O	3:D:1518:NAP:H2N	2.12	0.49
1:B:2:SER:O	1:B:33:PRO:HA	2.12	0.49
1:A:383:LEU:N	1:A:384:PRO:CD	2.75	0.49
1:D:29:LYS:NZ	1:D:399:ASP:OD1	2.28	0.49
1:D:383:LEU:N	1:D:384:PRO:CD	2.76	0.49
1:B:362:VAL:HG23	1:B:408:LYS:HD2	1.95	0.49
1:D:407:ILE:HG12	1:D:413:LYS:CA	2.43	0.49
1:D:407:ILE:O	1:D:408:LYS:C	2.51	0.49
1:A:315:HIS:HA	1:A:318:MET:HE2	1.93	0.49
1:C:29:LYS:NZ	1:C:399:ASP:OD1	2.30	0.49
1:D:407:ILE:HA	1:D:412:ALA:O	2.13	0.49
1:B:383:LEU:N	1:B:384:PRO:CD	2.77	0.48
1:B:410:ALA:O	1:B:414:LEU:HD13	2.13	0.48
1:C:334:PHE:CZ	1:C:360:GLU:HG2	2.47	0.48
1:D:393:ASN:C	1:D:393:ASN:ND2	2.59	0.48
1:D:407:ILE:HA	1:D:412:ALA:C	2.33	0.48
1:A:393:ASN:HD22	1:A:396:GLU:H	1.59	0.48
1:C:107:VAL:HG23	1:C:134:ALA:HB2	1.95	0.48
1:D:383:LEU:HD13	1:D:383:LEU:C	2.34	0.48
1:B:139:TYR:CD1	1:B:139:TYR:N	2.82	0.48
1:B:80:GLU:HG3	1:B:91:MET:HE1	1.95	0.48
1:A:324:GLU:OE1	1:A:388:ARG:NH1	2.46	0.48
1:C:383:LEU:N	1:C:384:PRO:CD	2.76	0.48
1:C:7:GLY:HA3	1:C:37:LEU:HD23	1.96	0.47
1:B:334:PHE:O	1:B:338:ARG:HB2	2.14	0.47
1:B:393:ASN:C	1:B:393:ASN:ND2	2.64	0.47
1:D:7:GLY:HA3	1:D:37:LEU:HD23	1.97	0.47
1:D:406:LYS:HG2	1:D:413:LYS:HE2	1.96	0.47
1:D:294:VAL:CG1	1:D:303:VAL:HG13	2.44	0.47
1:A:383:LEU:C	1:A:383:LEU:HD13	2.35	0.47
1:C:184:ASN:HD22	1:C:185:GLN:H	1.63	0.47
1:A:393:ASN:HD21	1:A:396:GLU:N	2.12	0.47
1:B:20:ARG:NH2	1:B:43:ASP:OD1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:406:LYS:O	1:D:413:LYS:CD	2.63	0.47
1:A:140:ARG:O	1:A:140:ARG:HG3	2.15	0.47
1:A:330:ILE:HD12	1:A:363:SER:HB3	1.96	0.47
1:D:405:LEU:C	1:D:407:ILE:N	2.66	0.47
1:A:109:ARG:HD3	1:A:293:SER:OG	2.15	0.47
1:A:3:LYS:HG2	1:A:35:VAL:O	2.15	0.46
1:A:388:ARG:HD2	1:A:388:ARG:O	2.15	0.46
1:A:252:ASP:HB3	1:B:279:ASP:OD2	2.14	0.46
1:C:383:LEU:HD13	1:C:383:LEU:C	2.35	0.46
1:A:324:GLU:CD	1:A:388:ARG:HH12	2.19	0.46
1:D:12:GLU:HG2	1:D:13:MET:N	2.30	0.46
1:D:407:ILE:CG2	1:D:412:ALA:HA	2.46	0.46
1:C:163:GLN:HG3	1:C:163:GLN:H	1.39	0.46
1:D:393:ASN:HD21	1:D:396:GLU:HG3	1.79	0.46
1:D:404:ASN:ND2	1:D:407:ILE:CD1	2.79	0.46
1:B:123:GLY:O	1:B:262:GLU:HA	2.16	0.46
1:B:236:LYS:O	1:B:240:GLU:HG3	2.17	0.45
1:C:109:ARG:NH2	4:C:1618:ICT:O2	2.39	0.45
1:C:314:ARG:HD2	3:C:1517:NAP:O3B	2.17	0.45
1:D:75:THR:H	3:D:1518:NAP:H71N	1.64	0.45
1:A:1:MET:HE2	1:A:3:LYS:CD	2.43	0.45
1:A:334:PHE:HA	1:A:337:THR:OG1	2.17	0.45
1:A:387:GLN:O	1:A:390:ASP:HB2	2.17	0.45
1:C:334:PHE:HA	1:C:337:THR:OG1	2.16	0.44
1:B:319:TYR:OH	1:C:322:GLY:HA2	2.17	0.44
1:C:315:HIS:HA	1:C:318:MET:HE3	2.00	0.44
1:D:292:THR:HA	5:D:1752:HOH:O	2.17	0.44
1:D:407:ILE:HA	1:D:413:LYS:HG3	1.99	0.44
1:B:383:LEU:HD13	1:B:383:LEU:C	2.37	0.44
1:A:29:LYS:HE3	5:A:1789:HOH:O	2.17	0.44
1:D:407:ILE:HG12	1:D:413:LYS:O	2.16	0.44
1:B:322:GLY:HA3	1:C:395:PHE:CD2	2.52	0.44
1:A:38:ASP:OD2	1:A:40:HIS:HE1	2.01	0.44
1:C:164:LYS:HE3	1:C:164:LYS:HB3	1.90	0.44
1:B:12:GLU:HG2	1:B:13:MET:N	2.31	0.44
1:A:236:LYS:O	1:A:240:GLU:HG3	2.18	0.44
1:C:308:ALA:HB3	5:C:1909:HOH:O	2.17	0.44
1:C:2:SER:HB2	1:C:33:PRO:CA	2.47	0.44
1:C:334:PHE:CE2	1:C:360:GLU:HG2	2.53	0.44
1:A:101:ASN:HB2	5:A:1772:HOH:O	2.18	0.44
1:D:357:ASN:O	1:D:361:GLU:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:403:GLU:O	1:B:407:ILE:HG13	2.18	0.44
1:C:236:LYS:O	1:C:240:GLU:HG3	2.18	0.44
1:D:13:MET:HA	1:D:42:TYR:O	2.17	0.43
1:A:12:GLU:HG2	1:A:13:MET:N	2.33	0.43
1:D:236:LYS:O	1:D:240:GLU:HG3	2.18	0.43
3:C:1517:NAP:H4N	4:C:1618:ICT:H2	2.01	0.43
1:A:357:ASN:O	1:A:361:GLU:HG2	2.18	0.43
1:D:80:GLU:HG3	1:D:91:MET:HE1	1.99	0.43
1:B:7:GLY:HA3	1:B:37:LEU:CD2	2.47	0.43
1:A:350:LYS:HD3	1:A:350:LYS:HA	1.81	0.43
1:A:3:LYS:CE	1:A:36:GLU:HG3	2.48	0.43
3:D:1518:NAP:H4N	4:D:1619:ICT:H2	2.01	0.43
1:A:252:ASP:HB2	5:B:1741:HOH:O	2.18	0.43
1:B:3:LYS:HB3	1:B:35:VAL:C	2.38	0.43
1:D:123:GLY:O	1:D:262:GLU:HA	2.19	0.43
1:C:357:ASN:O	1:C:361:GLU:HG2	2.19	0.43
1:A:315:HIS:HA	1:A:318:MET:HE3	2.01	0.42
1:A:17:GLU:HB3	1:A:316:TYR:CG	2.54	0.42
1:D:183:TYR:CD1	1:D:184:ASN:N	2.87	0.42
1:B:344:ALA:HB1	1:B:353:ALA:HB2	2.00	0.42
1:C:393:ASN:HD21	1:C:396:GLU:HG3	1.84	0.42
1:C:393:ASN:ND2	1:C:393:ASN:C	2.65	0.42
1:A:334:PHE:CE2	1:A:360:GLU:HG2	2.54	0.42
1:C:120:LEU:HD13	1:D:259:MET:HG3	2.02	0.42
1:D:20:ARG:NH2	1:D:43:ASP:OD1	2.51	0.42
1:D:109:ARG:HD3	1:D:293:SER:OG	2.19	0.42
1:B:357:ASN:O	1:B:361:GLU:HG2	2.20	0.42
1:D:393:ASN:HD21	1:D:396:GLU:H	1.60	0.42
1:C:120:LEU:CD1	1:D:259:MET:HG3	2.50	0.42
1:A:392:LEU:HB3	1:A:396:GLU:HB2	2.00	0.42
1:B:171:ASN:HA	1:B:171:ASN:HD22	1.62	0.41
1:A:80:GLU:HG3	1:A:91:MET:HE2	2.01	0.41
1:D:410:ALA:CB	1:D:413:LYS:HZ1	2.27	0.41
1:C:393:ASN:HD22	1:C:396:GLU:H	1.59	0.41
1:A:299:ASP:HB3	5:A:1893:HOH:O	2.21	0.41
1:D:406:LYS:HB3	1:D:413:LYS:CD	2.46	0.41
1:D:393:ASN:HD22	1:D:396:GLU:H	1.65	0.41
1:C:234:GLN:HG2	5:C:1924:HOH:O	2.20	0.41
1:B:393:ASN:HD21	1:B:396:GLU:H	1.64	0.41
1:C:1:MET:HG3	1:C:32:PHE:CG	2.55	0.41
1:D:101:ASN:ND2	1:D:140:ARG:HG3	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:80:GLU:HG3	1:D:91:MET:HE2	2.03	0.41
1:C:315:HIS:HA	1:C:318:MET:HE2	2.03	0.41
1:C:12:GLU:HG2	1:C:13:MET:N	2.35	0.41
1:A:139:TYR:HE1	4:A:1616:ICT:O5	2.04	0.41
1:D:69:VAL:HG11	1:D:343:ARG:HD2	2.02	0.41
1:B:67:HIS:O	1:B:68:ASN:HB2	2.21	0.40
1:C:159:SER:OG	1:D:149:PRO:HB2	2.21	0.40
3:B:1516:NAP:H4N	4:B:1617:ICT:H2	2.02	0.40
1:D:272:TYR:O	1:D:275:ASP:HB3	2.22	0.40
1:A:67:HIS:O	1:A:68:ASN:HB2	2.21	0.40

There are no symmetry-related clashes.

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	412/414 (100%)	392 (95%)	19 (5%)	1 (0%)	52	69
1	B	412/414 (100%)	396 (96%)	15 (4%)	1 (0%)	52	69
1	C	412/414 (100%)	390 (95%)	21 (5%)	1 (0%)	52	69
1	D	412/414 (100%)	388 (94%)	20 (5%)	4 (1%)	19	27
All	All	1648/1656 (100%)	1566 (95%)	75 (5%)	7 (0%)	39	54

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	411	GLN
1	D	413	LYS
1	C	4	LYS
1	D	407	ILE
1	D	412	ALA

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Mol	Chain	Res	Type
1	A	140	ARG
1	B	17	GLU

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	350/350 (100%)	343 (98%)	7 (2%)	63	80
1	B	350/350 (100%)	340 (97%)	10 (3%)	50	70
1	C	350/350 (100%)	339 (97%)	11 (3%)	47	68
1	D	350/350 (100%)	339 (97%)	11 (3%)	47	68
All	All	1400/1400 (100%)	1361 (97%)	39 (3%)	51	71

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	THR
1	A	28	GLU
1	A	103	LEU
1	A	109	ARG
1	A	252	ASP
1	A	338	ARG
1	A	393	ASN
1	B	19	THR
1	B	28	GLU
1	B	103	LEU
1	B	109	ARG
1	B	171	ASN
1	B	252	ASP
1	B	329	PRO
1	B	338	ARG
1	B	388	ARG
1	B	393	ASN
1	C	4	LYS
1	C	19	THR

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Mol	Chain	Res	Type
1	C	28	GLU
1	C	103	LEU
1	C	109	ARG
1	C	163	GLN
1	C	184	ASN
1	C	252	ASP
1	C	306	GLU
1	C	338	ARG
1	C	393	ASN
1	D	19	THR
1	D	28	GLU
1	D	103	LEU
1	D	109	ARG
1	D	140	ARG
1	D	184	ASN
1	D	252	ASP
1	D	329	PRO
1	D	388	ARG
1	D	393	ASN
1	D	408	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	A	90	GLN
1	A	163	GLN
1	A	171	ASN
1	A	185	GLN
1	A	194	HIS
1	A	198	GLN
1	A	228	GLN
1	A	257	GLN
1	A	277	GLN
1	A	323	GLN
1	A	393	ASN
1	A	404	ASN
1	B	40	HIS
1	B	90	GLN
1	B	171	ASN
1	B	194	HIS
1	B	198	GLN

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Mol	Chain	Res	Type
1	B	228	GLN
1	B	257	GLN
1	B	277	GLN
1	B	348	ASN
1	B	393	ASN
1	B	404	ASN
1	C	40	HIS
1	C	90	GLN
1	C	184	ASN
1	C	194	HIS
1	C	198	GLN
1	C	228	GLN
1	C	257	GLN
1	C	277	GLN
1	C	348	ASN
1	C	393	ASN
1	C	404	ASN
1	C	411	GLN
1	D	40	HIS
1	D	90	GLN
1	D	101	ASN
1	D	171	ASN
1	D	194	HIS
1	D	198	GLN
1	D	228	GLN
1	D	242	GLN
1	D	257	GLN
1	D	277	GLN
1	D	393	ASN
1	D	404	ASN
1	D	411	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

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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$
3	NAP	A	1515	-	42,52,52	2.16	10 (23%)	54,80,80	2.02	13 (24%)
4	ICT	A	1616	2	2,12,12	0.28	0	2,16,16	1.41	0
3	NAP	B	1516	-	42,52,52	2.20	13 (30%)	54,80,80	2.00	13 (24%)
4	ICT	B	1617	2	2,12,12	0.29	0	2,16,16	1.29	0
3	NAP	C	1517	-	42,52,52	2.26	12 (28%)	54,80,80	2.03	14 (25%)
4	ICT	C	1618	2	2,12,12	0.05	0	2,16,16	1.38	0
3	NAP	D	1518	-	42,52,52	2.19	11 (26%)	54,80,80	2.00	13 (24%)
4	ICT	D	1619	2	2,12,12	0.57	0	2,16,16	1.30	0

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
3	NAP	A	1515	-	-	0/27/67/67	0/5/5/5
4	ICT	A	1616	2	-	0/6/16/16	0/0/0/0
3	NAP	B	1516	-	-	0/27/67/67	0/5/5/5
4	ICT	B	1617	2	-	0/6/16/16	0/0/0/0
3	NAP	C	1517	-	-	0/27/67/67	0/5/5/5
4	ICT	C	1618	2	-	0/6/16/16	0/0/0/0
3	NAP	D	1518	-	-	0/27/67/67	0/5/5/5
4	ICT	D	1619	2	-	0/6/16/16	0/0/0/0

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1516	NAP	PA-O2A	-2.47	1.44	1.54
3	C	1517	NAP	PA-O2A	-2.41	1.44	1.54
3	D	1518	NAP	PA-O2A	-2.37	1.44	1.54
3	D	1518	NAP	P2B-O3X	-2.35	1.46	1.54
3	D	1518	NAP	P2B-O2X	-2.34	1.46	1.54
3	A	1515	NAP	PA-O2A	-2.28	1.45	1.54
3	A	1515	NAP	P2B-O2X	-2.26	1.46	1.54
3	C	1517	NAP	P2B-O2X	-2.22	1.46	1.54
3	B	1516	NAP	P2B-O2X	-2.22	1.46	1.54
3	C	1517	NAP	P2B-O3X	-2.19	1.46	1.54
3	B	1516	NAP	P2B-O3X	-2.17	1.46	1.54
3	B	1516	NAP	PN-O2N	-2.00	1.46	1.54
3	D	1518	NAP	O4B-C1B	2.02	1.43	1.41
3	C	1517	NAP	C2A-N3A	2.05	1.35	1.32
3	A	1515	NAP	O4B-C1B	2.16	1.43	1.41
3	B	1516	NAP	C3B-C2B	2.18	1.58	1.53
3	C	1517	NAP	O4B-C1B	2.29	1.44	1.41
3	B	1516	NAP	C2N-C3N	2.34	1.42	1.39
3	A	1515	NAP	C2A-N1A	2.70	1.39	1.33
3	B	1516	NAP	C2A-N1A	2.76	1.39	1.33
3	D	1518	NAP	C2A-N1A	2.85	1.39	1.33
3	C	1517	NAP	P2B-O2B	3.12	1.69	1.60
3	B	1516	NAP	P2B-O2B	3.16	1.69	1.60
3	D	1518	NAP	C6N-N1N	3.19	1.44	1.35
3	C	1517	NAP	C2A-N1A	3.27	1.40	1.33
3	A	1515	NAP	C4A-N3A	3.29	1.40	1.35
3	C	1517	NAP	C6N-N1N	3.33	1.44	1.35
3	D	1518	NAP	P2B-O2B	3.33	1.70	1.60
3	B	1516	NAP	C4A-N3A	3.36	1.40	1.35
3	A	1515	NAP	C6N-N1N	3.40	1.44	1.35
3	B	1516	NAP	C6N-N1N	3.45	1.44	1.35
3	A	1515	NAP	P2B-O2B	3.52	1.70	1.60
3	D	1518	NAP	C4A-N3A	3.87	1.41	1.35
3	C	1517	NAP	C4A-N3A	3.96	1.41	1.35
3	D	1518	NAP	C4N-C3N	4.66	1.47	1.39
3	B	1516	NAP	C4N-C3N	4.71	1.47	1.39
3	C	1517	NAP	C4N-C3N	4.94	1.47	1.39
3	A	1515	NAP	C4N-C3N	5.14	1.48	1.39
3	D	1518	NAP	C5N-C4N	5.20	1.49	1.38
3	B	1516	NAP	C5N-C4N	5.34	1.49	1.38
3	A	1515	NAP	C5N-C4N	5.42	1.50	1.38
3	C	1517	NAP	C5N-C4N	5.60	1.50	1.38
3	A	1515	NAP	C3N-C7N	6.67	1.61	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1518	NAP	C3N-C7N	7.00	1.61	1.50
3	B	1516	NAP	C3N-C7N	7.14	1.61	1.50
3	C	1517	NAP	C3N-C7N	7.39	1.62	1.50

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1515	NAP	C5N-C4N-C3N	-4.84	114.25	120.33
3	C	1517	NAP	C5N-C4N-C3N	-4.63	114.51	120.33
3	D	1518	NAP	C5N-C4N-C3N	-4.37	114.84	120.33
3	B	1516	NAP	C5N-C4N-C3N	-4.36	114.85	120.33
3	D	1518	NAP	O2B-P2B-O1X	-3.68	97.91	107.11
3	D	1518	NAP	C3N-C2N-N1N	-3.68	116.12	120.36
3	B	1516	NAP	C3N-C2N-N1N	-3.51	116.31	120.36
3	A	1515	NAP	O2B-P2B-O1X	-3.42	98.56	107.11
3	A	1515	NAP	C3N-C2N-N1N	-3.37	116.48	120.36
3	C	1517	NAP	C3N-C2N-N1N	-3.34	116.51	120.36
3	C	1517	NAP	O2B-P2B-O1X	-3.25	99.00	107.11
3	B	1516	NAP	O2B-P2B-O1X	-3.22	99.06	107.11
3	B	1516	NAP	N3A-C2A-N1A	-2.60	126.91	128.89
3	C	1517	NAP	C2B-C3B-C4B	-2.52	95.89	101.85
3	B	1516	NAP	C2B-C3B-C4B	-2.47	96.01	101.85
3	C	1517	NAP	N3A-C2A-N1A	-2.46	127.01	128.89
3	A	1515	NAP	C2B-C3B-C4B	-2.38	96.23	101.85
3	D	1518	NAP	N3A-C2A-N1A	-2.24	127.18	128.89
3	A	1515	NAP	N3A-C2A-N1A	-2.23	127.19	128.89
3	D	1518	NAP	C2B-C3B-C4B	-2.23	96.58	101.85
3	C	1517	NAP	O5B-PA-O1A	-2.05	101.67	109.62
3	D	1518	NAP	O3X-P2B-O2X	2.22	115.83	107.38
3	A	1515	NAP	O3X-P2B-O2X	2.27	116.04	107.38
3	C	1517	NAP	O3X-P2B-O2X	2.30	116.13	107.38
3	A	1515	NAP	N6A-C6A-N1A	2.32	124.19	119.20
3	B	1516	NAP	O3X-P2B-O2X	2.35	116.33	107.38
3	B	1516	NAP	O2X-P2B-O1X	2.57	118.85	110.58
3	B	1516	NAP	N6A-C6A-N1A	2.57	124.73	119.20
3	D	1518	NAP	O2X-P2B-O1X	2.60	118.94	110.58
3	A	1515	NAP	O2X-P2B-O1X	2.60	118.95	110.58
3	D	1518	NAP	N6A-C6A-N1A	2.65	124.89	119.20
3	C	1517	NAP	N6A-C6A-N1A	2.74	125.08	119.20
3	C	1517	NAP	O2X-P2B-O1X	2.77	119.50	110.58
3	C	1517	NAP	C6N-C5N-C4N	2.78	123.64	119.44
3	A	1515	NAP	C6N-C5N-C4N	2.79	123.65	119.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1516	NAP	C6N-C5N-C4N	2.90	123.83	119.44
3	D	1518	NAP	C6N-C5N-C4N	2.90	123.83	119.44
3	D	1518	NAP	O3B-C3B-C2B	3.12	120.18	111.16
3	C	1517	NAP	O3B-C3B-C2B	3.31	120.72	111.16
3	A	1515	NAP	O3B-C3B-C2B	3.38	120.91	111.16
3	B	1516	NAP	O3B-C3B-C2B	3.45	121.12	111.16
3	A	1515	NAP	C4A-C5A-N7A	4.64	113.75	109.48
3	B	1516	NAP	C4A-C5A-N7A	4.72	113.83	109.48
3	D	1518	NAP	C4A-C5A-N7A	4.75	113.84	109.48
3	C	1517	NAP	C4A-C5A-N7A	5.04	114.11	109.48
3	B	1516	NAP	C2N-C3N-C4N	5.07	123.94	118.29
3	D	1518	NAP	C2N-C3N-C4N	5.21	124.09	118.29
3	C	1517	NAP	C2N-C3N-C4N	5.29	124.19	118.29
3	A	1515	NAP	C2N-C3N-C4N	5.51	124.42	118.29
3	D	1518	NAP	PN-O3-PA	5.85	149.16	132.73
3	B	1516	NAP	PN-O3-PA	6.08	149.81	132.73
3	C	1517	NAP	PN-O3-PA	6.11	149.87	132.73
3	A	1515	NAP	PN-O3-PA	6.33	150.50	132.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1515	NAP	3	0
4	A	1616	ICT	3	0
3	B	1516	NAP	4	0
4	B	1617	ICT	2	0
3	C	1517	NAP	5	0
4	C	1618	ICT	3	0
3	D	1518	NAP	4	0
4	D	1619	ICT	2	0

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	414/414 (100%)	0.03	9 (2%) 65 64	9, 20, 41, 103	0
1	B	414/414 (100%)	0.07	14 (3%) 49 48	8, 22, 41, 97	0
1	C	414/414 (100%)	0.03	10 (2%) 62 61	8, 19, 42, 97	0
1	D	414/414 (100%)	0.11	14 (3%) 49 48	7, 21, 43, 114	0
All	All	1656/1656 (100%)	0.06	47 (2%) 56 55	7, 20, 42, 114	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	1	MET	17.5
1	C	414	LEU	12.4
1	A	414	LEU	11.3
1	D	414	LEU	10.3
1	C	2	SER	9.3
1	B	2	SER	8.7
1	A	1	MET	8.6
1	D	412	ALA	8.4
1	B	414	LEU	8.2
1	B	1	MET	7.2
1	D	413	LYS	6.8
1	D	2	SER	6.0
1	C	3	LYS	5.3
1	C	1	MET	5.3
1	A	2	SER	5.3
1	B	3	LYS	4.8
1	B	413	LYS	4.1
1	A	3	LYS	3.9
1	C	384	PRO	3.6
1	D	407	ILE	3.4
1	B	84	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	411	GLN	3.2
1	B	412	ALA	3.1
1	D	3	LYS	3.1
1	C	413	LYS	3.0
1	B	300	GLY	2.9
1	C	237	SER	2.8
1	A	174	GLU	2.8
1	D	28	GLU	2.8
1	B	387	GLN	2.6
1	C	412	ALA	2.6
1	D	163	GLN	2.6
1	B	348	ASN	2.6
1	D	410	ALA	2.4
1	A	28	GLU	2.4
1	D	348	ASN	2.4
1	B	174	GLU	2.4
1	B	411	GLN	2.4
1	B	229	GLU	2.4
1	A	300	GLY	2.3
1	A	160	ASP	2.2
1	C	229	GLU	2.2
1	D	84	GLU	2.2
1	A	90	GLN	2.2
1	B	389	SER	2.1
1	C	241	ALA	2.0
1	D	350	LYS	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
4	ICT	D	1619	13/13	0.78	0.32	7.07	30,36,39,41	0
4	ICT	A	1616	13/13	0.79	0.30	5.81	27,36,41,44	0
4	ICT	C	1618	13/13	0.82	0.26	5.51	28,31,36,38	0
4	ICT	B	1617	13/13	0.82	0.25	2.82	30,36,39,40	0
2	CA	A	1717	1/1	0.66	0.15	0.27	45,45,45,45	0
2	CA	B	1718	1/1	0.81	0.16	0.16	38,38,38,38	0
3	NAP	D	1518	48/48	0.95	0.14	-0.16	9,15,24,25	0
3	NAP	B	1516	48/48	0.95	0.13	-0.29	8,16,24,25	0
3	NAP	C	1517	48/48	0.95	0.12	-0.49	8,15,23,24	0
3	NAP	A	1515	48/48	0.96	0.12	-0.54	10,16,24,28	0
2	CA	D	1720	1/1	0.97	0.09	-1.31	26,26,26,26	0
2	CA	C	1719	1/1	0.97	0.07	-3.15	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.