



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:30 pm GMT

PDB ID : 1ARZ
Title : ESCHERICHIA COLI DIHYDRODIPICOLINATE REDUCTASE IN COM-
PLEX WITH NADH AND 2,6 PYRIDINE DICARBOXYLATE
Authors : Scapin, G.; Reddy, S.G.; Zheng, R.; Blanchard, J.S.
Deposited on : 1997-08-08
Resolution : 2.60 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

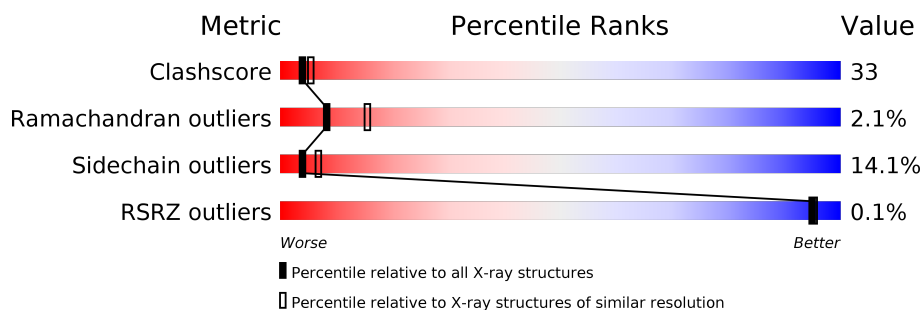
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

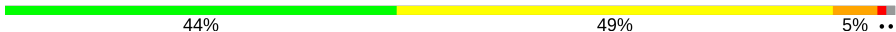



The reported resolution of this entry is 2.60 Å.

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	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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	273	 44% 49% 5% ..
1	B	273	 52% 37% 10% .
1	C	273	 50% 42% 7% .
1	D	273	 54% 36% 8% ..

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
5	PDC	D	303	-	-	X	-

2 Entry composition [i](#)

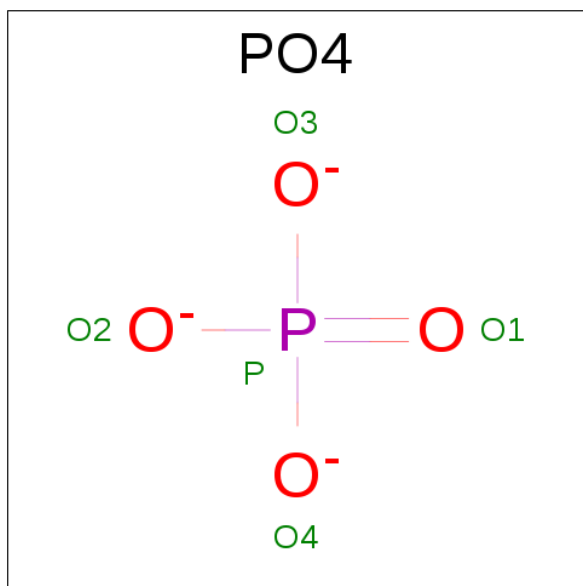
There are 6 unique types of molecules in this entry. The entry contains 8232 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DIHYDRODIPICOLINATE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			1939	1207	344	378	10			
1	B	269	Total	C	N	O	S	0	0	0
			1975	1225	356	384	10			
1	C	271	Total	C	N	O	S	0	0	0
			1980	1229	357	384	10			
1	D	271	Total	C	N	O	S	0	0	0
			1984	1232	357	385	10			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).

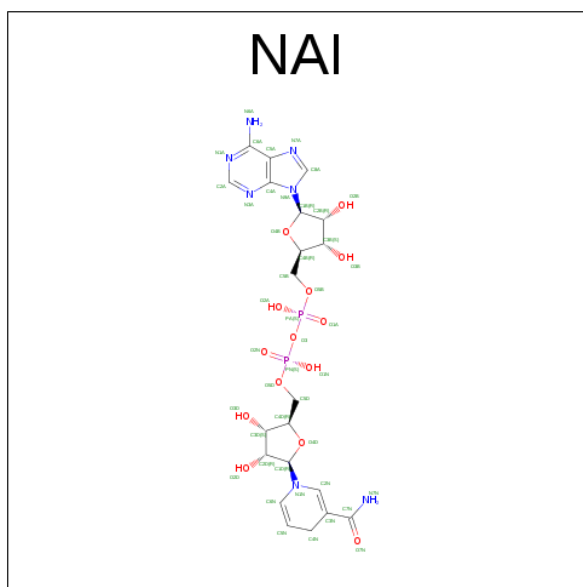


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	O P	0	0
			5 4 1			

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

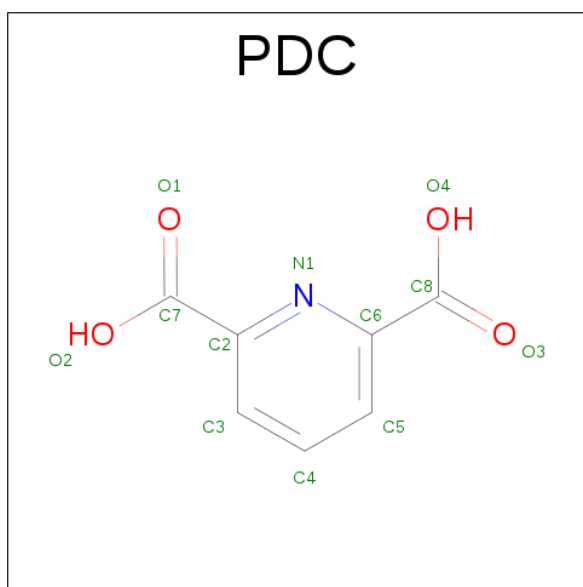
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total K 1 1	0	0
3	D	1	Total K 1 1	0	0

- Molecule 4 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: $C_{21}H_{29}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C N O P 44 21 7 14 2	0	0
4	C	1	Total C N O P 44 21 7 14 2	0	0
4	D	1	Total C N O P 44 21 7 14 2	0	0

- Molecule 5 is PYRIDINE-2,6-DICARBOXYLIC ACID (three-letter code: PDC) (formula: $C_7H_5NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			12	7	1	4		
5	C	1	Total	C	N	O	0	0
			12	7	1	4		
5	D	1	Total	C	N	O	0	0
			12	7	1	4		

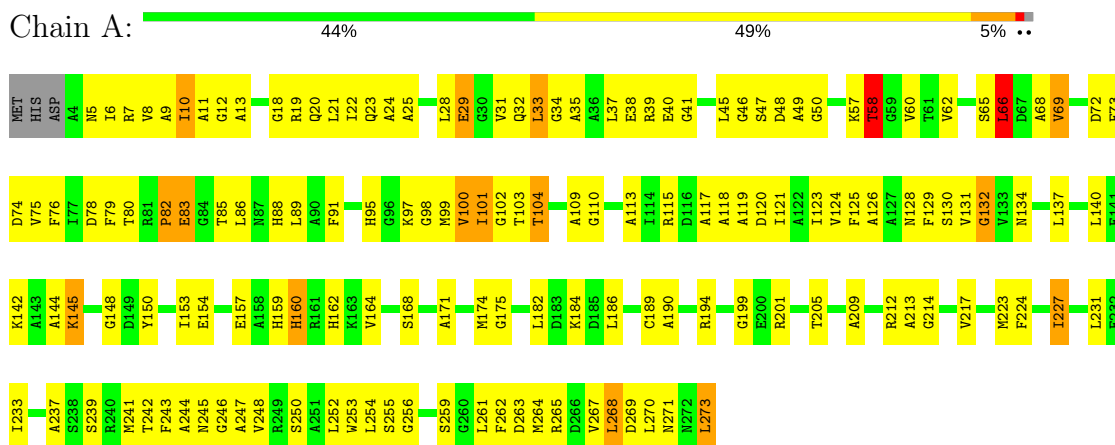
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	41	Total	O	0	0
			41	41		
6	B	47	Total	O	0	0
			47	47		
6	C	45	Total	O	0	0
			45	45		
6	D	46	Total	O	0	0
			46	46		

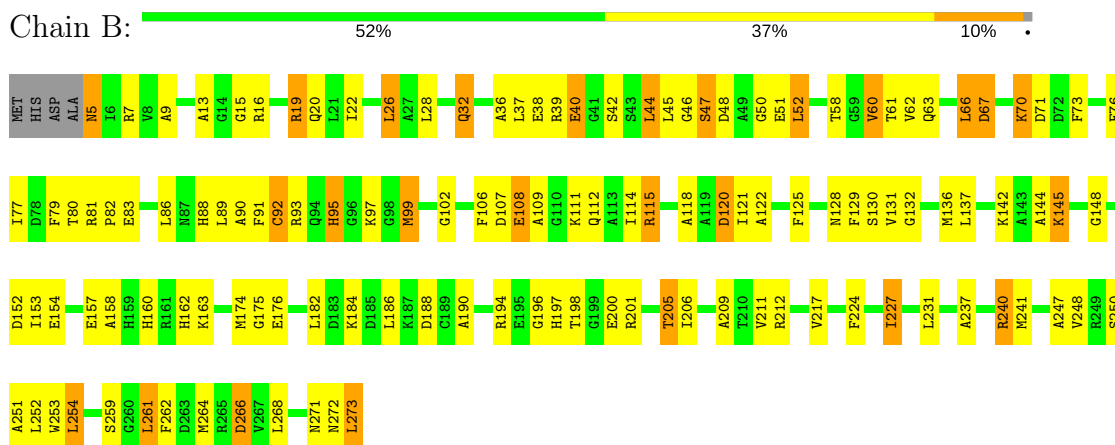
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

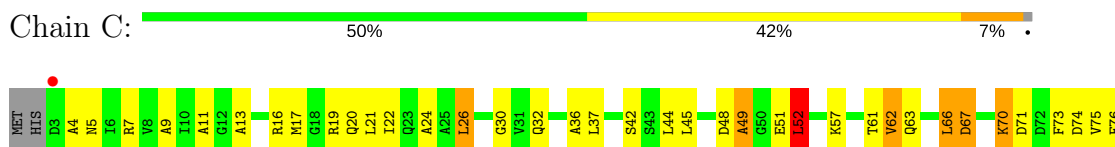
• Molecule 1: DIHYDRODIPICOLINATE REDUCTASE

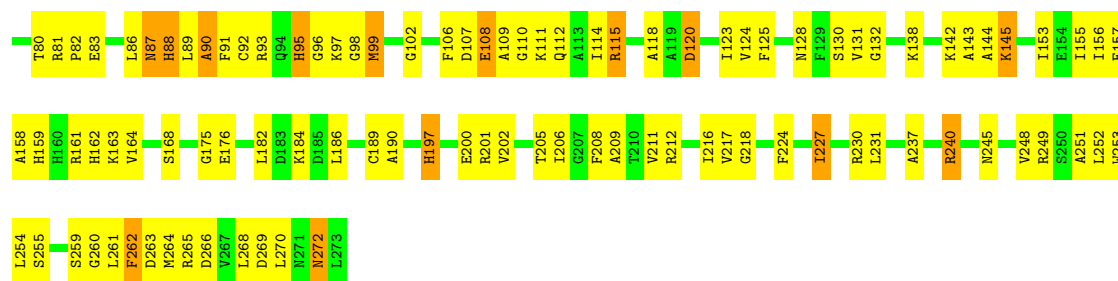


• Molecule 1: DIHYDRODIPICOLINATE REDUCTASE



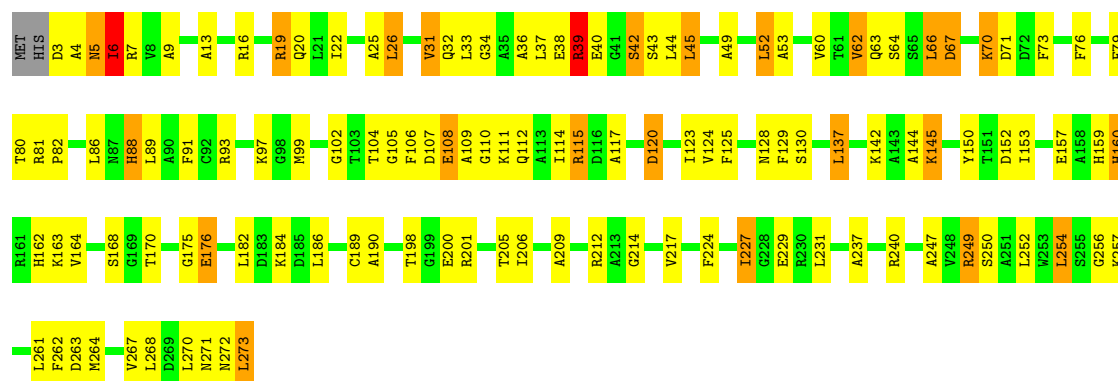
• Molecule 1: DIHYDRODIPICOLINATE REDUCTASE





● Molecule 1: DIHYDRODIPICOLINATE REDUCTASE

Chain D: 54% 36% 8% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	137.60Å 123.80Å 66.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60 23.21 – 2.52	Depositor EDS
% Data completeness (in resolution range)	89.9 (20.00-2.60) 85.3 (23.21-2.52)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 2.53Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.214 , 0.297 0.221 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.437	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.20 , 72.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8232	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K, NAI, PDC, PO4

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	0.73	0/1965	0.93	1/2658 (0.0%)
1	B	0.71	0/2001	0.92	2/2697 (0.1%)
1	C	0.72	0/2006	0.94	1/2705 (0.0%)
1	D	0.73	0/2010	0.93	4/2710 (0.1%)
All	All	0.73	0/7982	0.93	8/10770 (0.1%)

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

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	52	LEU	CA-CB-CG	6.70	130.71	115.30
1	B	52	LEU	CA-CB-CG	6.03	129.17	115.30
1	D	6	ILE	N-CA-C	5.59	126.09	111.00
1	D	45	LEU	CA-CB-CG	5.56	128.10	115.30
1	D	254	LEU	CA-CB-CG	5.34	127.58	115.30
1	A	72	ASP	N-CA-C	5.27	125.22	111.00
1	D	39	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	205	THR	N-CA-C	5.19	125.02	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	150	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	1939	0	1871	146	0
1	B	1975	0	1943	137	0
1	C	1980	0	1941	140	0
1	D	1984	0	1953	119	0
2	A	5	0	0	0	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
4	B	44	0	27	6	0
4	C	44	0	27	3	0
4	D	44	0	27	5	0
5	B	12	0	4	1	0
5	C	12	0	3	2	0
5	D	12	0	3	4	0
6	A	41	0	0	1	0
6	B	47	0	0	5	0
6	C	45	0	0	5	0
6	D	46	0	0	7	0
All	All	8232	0	7799	526	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 33.

All (526) 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:C:89:LEU:HA	1:C:99:MET:HE1	1.35	1.05
1:C:89:LEU:HA	1:C:99:MET:CE	1.93	0.99
1:C:45:LEU:HA	1:C:62:VAL:HG22	1.49	0.95
1:C:245:ASN:O	1:C:249:ARG:HD2	1.71	0.89
1:C:102:GLY:O	4:C:301:NAI:H2N	1.72	0.88
1:C:86:LEU:O	1:C:89:LEU:HB3	1.74	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:GLU:HA	1:C:111:LYS:HD3	1.55	0.88
1:C:153:ILE:HG22	1:C:206:ILE:HG23	1.56	0.87
1:B:108:GLU:HA	1:B:111:LYS:HD3	1.60	0.83
1:D:108:GLU:HA	1:D:111:LYS:HD3	1.60	0.82
1:A:101:ILE:N	1:A:101:ILE:HD12	1.95	0.81
1:C:218:GLY:HA3	6:C:412:HOH:O	1.77	0.81
1:D:102:GLY:O	4:D:302:NAI:H2N	1.82	0.79
1:C:48:ASP:HB3	1:C:51:GLU:HG2	1.62	0.79
1:C:138:LYS:HE2	1:C:272:ASN:HA	1.65	0.79
1:B:86:LEU:O	1:B:89:LEU:HG	1.84	0.78
1:C:67:ASP:HA	1:C:70:LYS:HB3	1.65	0.78
1:D:186:LEU:HD23	1:D:190:ALA:HB2	1.65	0.77
1:D:39:ARG:HG2	1:D:39:ARG:HH11	1.49	0.77
1:B:67:ASP:HA	1:B:70:LYS:HB3	1.64	0.77
1:D:67:ASP:HA	1:D:70:LYS:HB3	1.65	0.77
1:A:201:ARG:HD3	1:C:212:ARG:O	1.83	0.77
1:A:73:PHE:CE1	1:A:97:LYS:HE3	2.19	0.77
1:C:73:PHE:O	1:C:97:LYS:HE2	1.85	0.77
1:A:186:LEU:HD23	1:A:190:ALA:HB2	1.67	0.77
1:B:186:LEU:HD23	1:B:190:ALA:HB2	1.66	0.76
1:B:38:GLU:HG3	1:B:42:SER:OG	1.84	0.76
1:A:45:LEU:HD12	1:A:62:VAL:O	1.85	0.76
1:B:13:ALA:O	1:B:52:LEU:HD11	1.86	0.76
1:B:88:HIS:O	1:B:92:CYS:HB2	1.85	0.76
1:D:7:ARG:HB2	1:D:73:PHE:HA	1.69	0.75
1:D:66:LEU:HB3	1:D:91:PHE:CE2	2.22	0.75
1:C:186:LEU:HD23	1:C:190:ALA:HB2	1.70	0.74
1:C:49:ALA:O	1:C:52:LEU:HB3	1.87	0.74
1:D:249:ARG:HH11	1:D:249:ARG:HG2	1.53	0.74
1:D:49:ALA:O	1:D:52:LEU:HG	1.88	0.72
1:B:273:LEU:HA	6:B:438:HOH:O	1.89	0.72
1:A:129:PHE:O	1:A:246:GLY:HA3	1.90	0.72
1:C:216:ILE:HA	6:C:438:HOH:O	1.88	0.72
1:A:101:ILE:CD1	1:A:123:ILE:HG23	2.20	0.71
1:B:142:LYS:HE3	1:B:273:LEU:HG	1.71	0.70
1:B:76:PHE:HB2	1:B:92:CYS:SG	2.31	0.70
1:A:80:THR:O	1:A:103:THR:HG23	1.92	0.69
1:C:251:ALA:HA	1:C:254:LEU:HD12	1.74	0.68
1:C:48:ASP:OD2	1:C:57:LYS:HA	1.93	0.68
1:A:73:PHE:HE1	1:A:97:LYS:HB3	1.58	0.68
1:B:93:ARG:HA	1:B:121:ILE:HG21	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:ARG:HB2	1:C:73:PHE:HA	1.76	0.67
1:A:11:ALA:HB2	1:A:76:PHE:CZ	2.29	0.67
1:B:114:ILE:HG22	1:B:261:LEU:HD11	1.76	0.67
1:B:273:LEU:HD12	1:C:142:LYS:HD2	1.76	0.67
1:A:89:LEU:O	1:A:89:LEU:HD23	1.95	0.67
1:B:19:ARG:HD3	1:B:52:LEU:HD13	1.76	0.67
1:B:7:ARG:HB2	1:B:73:PHE:HA	1.76	0.67
1:A:124:VAL:HG23	1:A:254:LEU:HD21	1.78	0.66
4:D:302:NAI:H4B	6:D:426:HOH:O	1.94	0.66
1:B:76:PHE:CB	1:B:92:CYS:SG	2.83	0.66
1:A:79:PHE:CE1	1:A:102:GLY:HA3	2.30	0.66
1:A:46:GLY:H	1:A:62:VAL:HG13	1.60	0.66
1:D:19:ARG:HB3	1:D:53:ALA:HB2	1.78	0.66
1:C:17:MET:HB2	6:C:407:HOH:O	1.96	0.66
1:C:240:ARG:HD3	6:C:444:HOH:O	1.96	0.66
1:A:82:PRO:O	1:A:85:THR:N	2.29	0.66
1:C:262:PHE:HD1	1:C:262:PHE:N	1.94	0.65
1:C:264:MET:O	1:C:268:LEU:HD12	1.96	0.65
1:B:60:VAL:HG12	1:B:61:THR:H	1.60	0.65
1:D:93:ARG:HD2	1:D:117:ALA:HA	1.79	0.65
1:C:227:ILE:HD13	1:C:227:ILE:H	1.62	0.65
1:C:80:THR:O	4:C:301:NAI:H4D	1.95	0.65
1:C:262:PHE:CD1	1:C:262:PHE:N	2.64	0.65
1:D:99:MET:O	1:D:254:LEU:HD11	1.96	0.65
1:B:247:ALA:O	1:B:250:SER:HB2	1.97	0.64
1:B:44:LEU:HD13	1:B:44:LEU:O	1.96	0.64
1:D:3:ASP:N	1:D:31:VAL:H	1.96	0.64
1:A:79:PHE:CZ	1:A:102:GLY:HA3	2.32	0.64
1:B:114:ILE:HG22	1:B:261:LEU:CD1	2.28	0.64
1:C:128:ASN:HD22	1:C:130:SER:H	1.45	0.64
1:A:47:SER:O	1:A:62:VAL:HG12	1.98	0.63
1:C:93:ARG:O	1:C:96:GLY:N	2.31	0.63
1:B:120:ASP:O	1:B:259:SER:HB3	1.98	0.63
1:A:89:LEU:HD21	1:A:117:ALA:CB	2.27	0.63
1:A:264:MET:O	1:A:268:LEU:HB2	1.99	0.62
1:A:91:PHE:HE1	1:A:95:HIS:CD2	2.17	0.62
1:C:138:LYS:HE2	1:C:272:ASN:H	1.64	0.62
1:C:265:ARG:HG2	1:C:270:LEU:HD12	1.81	0.62
1:D:128:ASN:HD22	1:D:130:SER:H	1.48	0.62
1:A:131:VAL:HG13	1:A:268:LEU:HD21	1.81	0.62
1:B:152:ASP:HA	6:B:443:HOH:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:ILE:HD11	1:A:123:ILE:HG23	1.80	0.61
1:A:189:CYS:HB2	1:A:205:THR:HA	1.82	0.61
1:B:227:ILE:O	1:B:227:ILE:HG12	1.99	0.61
1:A:126:ALA:HB3	1:A:129:PHE:CE1	2.36	0.61
1:B:153:ILE:CG2	1:B:206:ILE:HD13	2.30	0.61
1:D:36:ALA:C	1:D:37:LEU:HD12	2.21	0.61
1:A:46:GLY:N	1:A:62:VAL:HG13	2.15	0.61
1:B:112:GLN:OE1	1:B:112:GLN:HA	1.99	0.61
1:A:227:ILE:H	1:A:227:ILE:HD13	1.65	0.60
1:A:265:ARG:NH1	1:A:271:ASN:HD21	2.00	0.60
1:B:48:ASP:HA	1:B:60:VAL:O	2.01	0.60
1:B:227:ILE:HD13	1:B:227:ILE:H	1.67	0.60
1:B:194:ARG:HH22	1:B:201:ARG:NH1	2.00	0.60
1:D:79:PHE:CZ	1:D:102:GLY:HA3	2.37	0.60
1:D:227:ILE:HD13	1:D:227:ILE:H	1.66	0.60
1:D:89:LEU:HD11	1:D:117:ALA:CB	2.31	0.60
1:B:112:GLN:O	1:B:115:ARG:HB3	2.02	0.60
1:B:28:LEU:HD12	1:B:248:VAL:HG11	1.84	0.59
1:B:36:ALA:C	1:B:37:LEU:HD12	2.22	0.59
1:C:227:ILE:HG12	1:C:227:ILE:O	2.02	0.59
1:B:118:ALA:HB1	1:B:261:LEU:N	2.17	0.59
1:B:142:LYS:HE3	1:B:273:LEU:CG	2.32	0.59
1:B:60:VAL:HG12	1:B:61:THR:N	2.18	0.59
1:A:153:ILE:N	1:A:201:ARG:HH21	2.01	0.58
1:C:138:LYS:HE2	1:C:272:ASN:CA	2.32	0.58
1:A:273:LEU:HD22	1:D:142:LYS:HD2	1.86	0.58
1:A:101:ILE:N	1:A:101:ILE:CD1	2.65	0.58
1:D:112:GLN:HA	1:D:112:GLN:OE1	2.02	0.58
1:A:91:PHE:CE1	1:A:95:HIS:CD2	2.91	0.58
1:A:78:ASP:OD2	1:A:85:THR:HG23	2.03	0.58
1:B:66:LEU:HB3	1:B:91:PHE:CZ	2.39	0.58
1:D:3:ASP:N	1:D:31:VAL:N	2.51	0.58
1:A:89:LEU:HD21	1:A:117:ALA:HB2	1.86	0.58
1:A:76:PHE:CZ	1:A:88:HIS:CD2	2.91	0.58
1:B:162:HIS:HA	1:D:200:GLU:OE1	2.04	0.58
1:D:227:ILE:HG12	1:D:227:ILE:O	2.04	0.58
1:C:112:GLN:HA	1:C:112:GLN:OE1	2.03	0.57
1:A:101:ILE:HD13	1:A:123:ILE:HG23	1.84	0.57
1:D:89:LEU:HD12	1:D:99:MET:HE1	1.85	0.57
1:A:89:LEU:HD23	1:A:89:LEU:C	2.25	0.57
1:B:120:ASP:C	1:B:259:SER:HB3	2.25	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:251:ALA:CA	1:C:254:LEU:HD12	2.35	0.57
1:C:74:ASP:O	1:C:97:LYS:HD3	2.04	0.57
1:C:66:LEU:HB3	1:C:91:PHE:CE1	2.40	0.57
1:C:261:LEU:HD23	1:C:262:PHE:N	2.19	0.57
1:A:153:ILE:H	1:A:201:ARG:NH2	2.03	0.57
1:A:19:ARG:O	1:A:22:ILE:HB	2.05	0.57
1:B:76:PHE:HB3	1:B:99:MET:HG3	1.87	0.57
1:D:39:ARG:CG	1:D:39:ARG:HH11	2.16	0.57
1:C:66:LEU:HB3	1:C:91:PHE:CZ	2.40	0.56
1:A:9:ALA:HB3	1:A:76:PHE:HA	1.87	0.56
1:A:8:VAL:HG22	1:A:33:LEU:HA	1.88	0.56
1:D:26:LEU:CD1	1:D:33:LEU:HB2	2.35	0.56
1:A:22:ILE:O	1:A:25:ALA:HB3	2.05	0.56
1:B:142:LYS:O	1:B:145:LYS:HB3	2.05	0.56
1:C:88:HIS:O	1:C:92:CYS:SG	2.63	0.56
1:D:102:GLY:O	4:D:302:NAI:C2N	2.51	0.56
1:A:83:GLU:HA	1:A:86:LEU:HD23	1.86	0.56
1:B:106:PHE:CD2	1:B:125:PHE:HZ	2.24	0.56
1:B:128:ASN:HD22	1:B:130:SER:H	1.53	0.56
1:A:261:LEU:HD23	1:A:262:PHE:N	2.21	0.56
1:D:76:PHE:HB3	1:D:99:MET:HG3	1.88	0.55
1:D:5:ASN:N	1:D:5:ASN:HD22	2.05	0.55
1:A:78:ASP:OD2	1:A:85:THR:HA	2.06	0.55
1:C:89:LEU:HA	1:C:99:MET:HE2	1.87	0.55
1:B:89:LEU:HA	1:B:99:MET:HE1	1.88	0.55
1:C:261:LEU:C	1:C:262:PHE:CD1	2.80	0.55
1:D:160:HIS:CE1	6:D:413:HOH:O	2.60	0.55
1:B:102:GLY:O	4:B:302:NAI:H2N	2.07	0.55
1:D:88:HIS:O	1:D:91:PHE:HB3	2.05	0.55
1:B:201:ARG:HG3	1:B:201:ARG:HH11	1.71	0.55
1:C:162:HIS:HE1	6:C:415:HOH:O	1.90	0.55
1:C:261:LEU:C	1:C:262:PHE:HD1	2.09	0.55
1:D:66:LEU:HB3	1:D:91:PHE:CZ	2.42	0.55
1:C:112:GLN:O	1:C:115:ARG:HB3	2.06	0.54
1:B:153:ILE:HG21	1:B:206:ILE:HD13	1.89	0.54
1:B:36:ALA:C	1:B:62:VAL:HG13	2.28	0.54
1:A:241:MET:CE	1:A:245:ASN:HD21	2.21	0.54
1:A:65:SER:HB3	1:A:68:ALA:HB2	1.90	0.54
1:A:79:PHE:CD1	1:A:102:GLY:HA3	2.43	0.54
1:D:249:ARG:NH1	1:D:249:ARG:HG2	2.19	0.54
1:A:268:LEU:HB3	1:A:270:LEU:HG	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:PHE:CE1	1:A:97:LYS:HB3	2.42	0.54
1:D:186:LEU:CD2	1:D:190:ALA:HB2	2.38	0.54
1:C:142:LYS:O	1:C:145:LYS:HB3	2.07	0.53
1:C:265:ARG:NH1	1:C:265:ARG:HB2	2.22	0.53
1:D:13:ALA:HB1	1:D:22:ILE:HD11	1.91	0.53
1:A:175:GLY:HA3	1:A:186:LEU:HD11	1.90	0.53
1:C:128:ASN:ND2	1:C:130:SER:H	2.06	0.53
1:C:36:ALA:C	1:C:37:LEU:HD12	2.29	0.53
1:C:13:ALA:HB1	1:C:22:ILE:HD11	1.91	0.53
1:D:104:THR:HG21	5:D:303:PDC:C7	2.39	0.53
1:B:252:LEU:HD23	1:B:253:TRP:N	2.23	0.53
1:C:45:LEU:HA	1:C:62:VAL:CG2	2.32	0.53
1:D:81:ARG:HB3	1:D:82:PRO:HD2	1.90	0.53
1:B:88:HIS:N	1:B:88:HIS:CD2	2.73	0.53
1:C:264:MET:CE	1:C:268:LEU:HD11	2.39	0.53
1:D:129:PHE:O	1:D:264:MET:HE2	2.08	0.53
1:B:194:ARG:NH2	1:B:201:ARG:NH1	2.55	0.53
1:C:163:LYS:NZ	5:C:302:PDC:N1	2.56	0.53
1:A:245:ASN:N	1:A:245:ASN:HD22	2.06	0.53
1:B:186:LEU:CD2	1:B:190:ALA:HB2	2.38	0.53
1:B:76:PHE:CB	1:B:92:CYS:HG	2.21	0.53
1:A:75:VAL:HG23	1:A:98:GLY:O	2.08	0.53
1:C:251:ALA:HA	1:C:254:LEU:CD1	2.38	0.53
1:C:253:TRP:HZ3	1:C:262:PHE:CD2	2.27	0.53
1:D:142:LYS:O	1:D:145:LYS:HB3	2.09	0.53
1:D:272:ASN:HB2	6:D:423:HOH:O	2.09	0.53
1:A:125:PHE:C	1:A:125:PHE:CD1	2.80	0.52
1:C:253:TRP:CZ3	1:C:262:PHE:CD2	2.97	0.52
1:D:170:THR:N	5:D:303:PDC:O1	2.42	0.52
1:D:19:ARG:CB	1:D:53:ALA:HB2	2.39	0.52
1:A:227:ILE:O	1:A:227:ILE:HG12	2.09	0.52
1:B:81:ARG:HB3	1:B:82:PRO:HD2	1.91	0.52
1:C:189:CYS:HB2	1:C:205:THR:HA	1.91	0.52
1:C:201:ARG:HH11	1:C:201:ARG:HG2	1.73	0.52
1:A:100:VAL:HG12	1:A:100:VAL:O	2.08	0.52
1:A:243:PHE:N	1:A:243:PHE:CD1	2.78	0.52
1:A:65:SER:HB3	1:A:68:ALA:CB	2.39	0.52
1:B:115:ARG:HA	1:B:261:LEU:HD12	1.92	0.52
1:A:209:ALA:HB1	1:C:209:ALA:HB1	1.92	0.52
1:C:106:PHE:CD2	1:C:125:PHE:HZ	2.28	0.52
1:C:186:LEU:CD2	1:C:190:ALA:HB2	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:138:LYS:HE2	1:C:272:ASN:N	2.24	0.52
1:A:273:LEU:HD22	1:D:142:LYS:CD	2.39	0.52
1:B:194:ARG:NH2	1:B:201:ARG:HH11	2.08	0.52
1:B:26:LEU:HD21	1:B:58:THR:HG21	1.91	0.52
1:B:38:GLU:OE1	4:B:302:NAI:H1B	2.10	0.52
1:A:142:LYS:O	1:A:145:LYS:HB3	2.10	0.52
1:A:186:LEU:CD2	1:A:190:ALA:HB2	2.38	0.52
1:A:79:PHE:CE2	1:A:102:GLY:HA3	2.45	0.52
1:D:112:GLN:O	1:D:115:ARG:HB3	2.10	0.52
1:C:95:HIS:N	1:C:95:HIS:ND1	2.56	0.51
1:B:67:ASP:O	1:B:70:LYS:HG2	2.11	0.51
1:D:19:ARG:HB3	1:D:53:ALA:CB	2.41	0.51
1:A:118:ALA:O	1:A:121:ILE:O	2.28	0.51
1:B:212:ARG:O	1:D:201:ARG:HD2	2.11	0.51
1:B:50:GLY:HA3	1:B:58:THR:HG23	1.93	0.51
1:A:128:ASN:O	1:A:134:ASN:ND2	2.44	0.51
1:B:88:HIS:H	1:B:88:HIS:CD2	2.27	0.51
1:C:265:ARG:O	1:C:269:ASP:N	2.43	0.51
1:D:106:PHE:CD2	1:D:125:PHE:HZ	2.28	0.51
1:A:119:ALA:O	1:A:259:SER:HB2	2.10	0.51
1:A:78:ASP:CG	1:A:85:THR:HG23	2.31	0.51
1:B:209:ALA:HB1	1:D:209:ALA:HB1	1.93	0.50
1:A:124:VAL:HG22	1:A:262:PHE:HB2	1.93	0.50
1:C:89:LEU:HD22	1:C:99:MET:CE	2.42	0.50
1:C:153:ILE:O	1:C:206:ILE:HA	2.11	0.50
1:C:21:LEU:O	1:C:24:ALA:HB3	2.12	0.50
1:D:7:ARG:CB	1:D:73:PHE:HA	2.40	0.50
1:A:144:ALA:O	1:A:182:LEU:HD21	2.12	0.50
1:A:160:HIS:HB2	1:A:214:GLY:O	2.11	0.50
1:C:73:PHE:CE1	1:C:97:LYS:HD2	2.47	0.50
1:B:106:PHE:HD2	1:B:125:PHE:HZ	1.58	0.50
1:B:46:GLY:O	1:B:61:THR:HB	2.11	0.50
1:C:5:ASN:HB2	1:C:30:GLY:O	2.12	0.50
1:D:152:ASP:HA	1:D:201:ARG:HH21	1.76	0.50
1:A:265:ARG:HH11	1:A:271:ASN:HD21	1.60	0.49
1:A:123:ILE:HG22	1:A:124:VAL:N	2.27	0.49
1:A:109:ALA:O	1:A:113:ALA:HB2	2.12	0.49
1:A:153:ILE:HG12	1:A:224:PHE:CD1	2.48	0.49
1:B:106:PHE:CD2	1:B:125:PHE:CZ	3.01	0.49
1:B:125:PHE:CB	1:B:261:LEU:HD21	2.41	0.49
1:D:9:ALA:HB2	1:D:73:PHE:CD2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:LEU:O	1:A:89:LEU:HB3	2.13	0.49
1:C:75:VAL:HG13	1:C:98:GLY:O	2.11	0.49
1:B:153:ILE:HG12	1:B:224:PHE:CD1	2.48	0.49
1:B:95:HIS:ND1	1:B:95:HIS:N	2.60	0.49
1:C:73:PHE:CD1	1:C:97:LYS:HD2	2.47	0.49
1:B:217:VAL:HB	1:B:237:ALA:O	2.12	0.49
1:B:252:LEU:C	1:B:252:LEU:HD23	2.32	0.49
1:D:263:ASP:HB2	6:D:416:HOH:O	2.12	0.49
1:D:86:LEU:O	1:D:89:LEU:HB3	2.13	0.49
1:A:131:VAL:O	1:A:132:GLY:C	2.51	0.49
1:B:175:GLY:C	1:B:186:LEU:HD12	2.33	0.49
1:A:102:GLY:O	1:A:104:THR:HG22	2.13	0.48
1:B:153:ILE:H	1:B:201:ARG:HH21	1.61	0.48
1:B:16:ARG:O	1:B:20:GLN:HG2	2.12	0.48
1:B:47:SER:O	1:B:61:THR:HA	2.13	0.48
1:B:83:GLU:O	1:B:86:LEU:HB2	2.13	0.48
6:B:409:HOH:O	1:C:230:ARG:HG2	2.13	0.48
1:D:150:TYR:HE2	6:D:432:HOH:O	1.96	0.48
1:D:39:ARG:O	1:D:42:SER:OG	2.32	0.48
1:A:162:HIS:HD2	1:C:200:GLU:OE1	1.97	0.48
1:D:249:ARG:HH11	1:D:249:ARG:CG	2.25	0.48
1:D:36:ALA:CB	1:D:62:VAL:HG23	2.44	0.48
1:C:82:PRO:HA	1:C:106:PHE:CE1	2.49	0.48
1:B:201:ARG:HD2	1:D:212:ARG:O	2.13	0.48
1:A:101:ILE:H	1:A:101:ILE:HD12	1.74	0.48
1:A:242:THR:HG23	1:D:229:GLU:CD	2.34	0.48
1:B:188:ASP:HA	6:B:428:HOH:O	2.13	0.48
1:B:201:ARG:NH2	1:B:205:THR:O	2.46	0.48
1:A:153:ILE:H	1:A:201:ARG:HH21	1.58	0.48
1:A:48:ASP:OD1	1:A:58:THR:HB	2.14	0.48
1:C:87:ASN:O	1:C:88:HIS:C	2.52	0.48
1:A:175:GLY:C	1:A:186:LEU:HD12	2.34	0.47
1:A:48:ASP:O	1:A:50:GLY:N	2.46	0.47
1:B:125:PHE:N	1:B:262:PHE:O	2.47	0.47
1:A:10:ILE:HD13	1:A:22:ILE:HA	1.96	0.47
1:C:264:MET:HE2	1:C:268:LEU:HD11	1.97	0.47
1:C:89:LEU:CD2	1:C:99:MET:HE1	2.45	0.47
1:B:7:ARG:CB	1:B:73:PHE:HA	2.42	0.47
1:C:175:GLY:C	1:C:186:LEU:HD12	2.35	0.47
1:D:128:ASN:ND2	1:D:130:SER:H	2.10	0.47
1:C:76:PHE:CZ	1:C:88:HIS:CD2	3.02	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:ASP:OD1	1:B:109:ALA:HB3	2.14	0.47
1:C:44:LEU:HD23	1:C:44:LEU:H	1.79	0.47
1:A:123:ILE:CG2	1:A:124:VAL:N	2.77	0.47
1:B:266:ASP:HA	1:B:271:ASN:ND2	2.30	0.47
1:B:273:LEU:HD22	6:B:438:HOH:O	2.14	0.47
1:C:153:ILE:HG12	1:C:224:PHE:CD1	2.49	0.47
1:C:90:ALA:O	1:C:93:ARG:HG3	2.15	0.47
1:A:140:LEU:HD12	1:A:233:ILE:HD12	1.97	0.47
1:A:243:PHE:H	1:A:243:PHE:HD1	1.61	0.47
1:B:20:GLN:NE2	1:B:241:MET:CE	2.78	0.47
1:C:270:LEU:O	1:C:272:ASN:N	2.47	0.47
1:A:38:GLU:HG3	1:A:39:ARG:N	2.30	0.47
1:B:125:PHE:HB3	1:B:261:LEU:HD21	1.96	0.47
1:B:144:ALA:O	1:B:182:LEU:HD21	2.14	0.47
1:D:152:ASP:CG	1:D:201:ARG:HE	2.17	0.47
1:A:100:VAL:HG11	1:A:247:ALA:HB1	1.97	0.47
1:A:65:SER:O	1:A:69:VAL:HG22	2.15	0.47
1:B:15:GLY:HA2	4:B:302:NAI:O1A	2.15	0.47
1:D:105:GLY:HA3	6:D:441:HOH:O	2.15	0.47
1:C:107:ASP:OD1	1:C:109:ALA:HB3	2.14	0.46
1:C:118:ALA:O	1:C:260:GLY:N	2.48	0.46
1:C:83:GLU:O	1:C:86:LEU:HB2	2.16	0.46
1:A:18:GLY:O	1:A:22:ILE:HD12	2.14	0.46
1:A:6:ILE:O	1:A:32:GLN:N	2.47	0.46
1:B:89:LEU:O	1:B:92:CYS:N	2.48	0.46
1:D:175:GLY:HA3	1:D:186:LEU:HD11	1.96	0.46
1:D:7:ARG:HH11	1:D:7:ARG:HG3	1.80	0.46
1:A:9:ALA:HB2	1:A:73:PHE:CD2	2.50	0.46
1:C:42:SER:C	1:C:44:LEU:H	2.18	0.46
1:C:90:ALA:O	1:C:93:ARG:HB2	2.15	0.46
1:A:7:ARG:HB3	1:A:34:GLY:HA3	1.98	0.46
1:A:160:HIS:HA	1:A:213:ALA:O	2.15	0.46
1:B:39:ARG:O	1:B:40:GLU:C	2.54	0.46
1:C:16:ARG:O	1:C:20:GLN:HG2	2.13	0.46
1:D:107:ASP:OD1	1:D:109:ALA:HB3	2.16	0.46
1:A:115:ARG:O	1:A:118:ALA:HB3	2.16	0.46
1:A:250:SER:O	1:A:253:TRP:HB3	2.16	0.46
1:C:144:ALA:O	1:C:182:LEU:HD21	2.16	0.46
1:C:175:GLY:HA3	1:C:186:LEU:HD11	1.98	0.46
1:C:44:LEU:HD23	1:C:44:LEU:N	2.31	0.46
1:D:93:ARG:HG3	1:D:120:ASP:OD1	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:ARG:HB2	1:A:73:PHE:HA	1.97	0.46
1:B:196:GLY:O	1:B:198:THR:HG23	2.16	0.46
1:D:175:GLY:C	1:D:186:LEU:HD12	2.36	0.46
1:A:241:MET:HE2	1:A:245:ASN:HD21	1.81	0.46
1:C:9:ALA:HB2	1:C:73:PHE:CG	2.51	0.46
1:A:33:LEU:O	1:A:60:VAL:HG21	2.16	0.45
1:B:118:ALA:CB	1:B:261:LEU:HB2	2.46	0.45
1:C:131:VAL:O	1:C:132:GLY:C	2.54	0.45
1:C:45:LEU:HD12	1:C:63:GLN:HA	1.98	0.45
1:D:36:ALA:O	1:D:37:LEU:HD12	2.16	0.45
1:A:85:THR:OG1	1:A:103:THR:HG21	2.17	0.45
1:A:217:VAL:HB	1:A:237:ALA:O	2.17	0.45
1:A:153:ILE:HA	1:A:223:MET:O	2.16	0.45
1:B:20:GLN:HG3	1:B:240:ARG:HB3	1.98	0.45
1:B:80:THR:O	4:B:302:NAI:H4D	2.17	0.45
1:D:163:LYS:NZ	5:D:303:PDC:O2	2.47	0.45
1:B:20:GLN:NE2	1:B:241:MET:HE2	2.31	0.45
1:B:9:ALA:HB3	1:B:76:PHE:HA	1.97	0.45
1:A:18:GLY:O	1:A:21:LEU:HB2	2.16	0.45
1:B:89:LEU:HB2	1:B:90:ALA:H	1.65	0.45
1:C:61:THR:O	1:C:63:GLN:HG2	2.17	0.45
1:C:92:CYS:HB2	1:C:99:MET:HE2	1.98	0.45
1:C:89:LEU:O	1:C:93:ARG:HG3	2.17	0.45
1:D:144:ALA:O	1:D:182:LEU:HD21	2.17	0.45
1:A:28:LEU:HG	1:A:248:VAL:HG12	1.99	0.45
1:B:13:ALA:HB1	1:B:22:ILE:HD11	1.98	0.45
1:C:201:ARG:NH1	1:C:205:THR:OG1	2.49	0.45
1:C:48:ASP:HB3	1:C:51:GLU:CG	2.42	0.45
1:C:87:ASN:C	1:C:89:LEU:N	2.68	0.45
1:D:176:GLU:N	1:D:186:LEU:HD12	2.32	0.45
1:D:206:ILE:N	1:D:206:ILE:HD12	2.31	0.45
1:A:24:ALA:HB2	1:A:241:MET:HE3	1.98	0.45
1:D:106:PHE:HD2	1:D:125:PHE:HZ	1.64	0.45
1:D:25:ALA:O	1:D:31:VAL:HG12	2.17	0.45
1:A:37:LEU:HD21	1:A:66:LEU:HD13	1.99	0.45
1:B:197:HIS:HA	1:D:164:VAL:O	2.17	0.45
1:B:261:LEU:O	1:B:262:PHE:CD1	2.70	0.45
1:B:76:PHE:HB3	1:B:92:CYS:SG	2.57	0.45
1:C:201:ARG:NH1	1:C:201:ARG:HG2	2.31	0.45
1:B:118:ALA:HB2	1:B:261:LEU:HB2	1.99	0.45
1:C:163:LYS:NZ	5:C:302:PDC:O2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:ILE:HG12	1:D:224:PHE:CD1	2.51	0.45
1:A:33:LEU:HD21	1:A:35:ALA:O	2.16	0.44
1:B:175:GLY:HA3	1:B:186:LEU:HD11	1.98	0.44
1:C:120:ASP:C	1:C:259:SER:HB2	2.37	0.44
1:D:106:PHE:CD2	1:D:125:PHE:CZ	3.05	0.44
1:A:97:LYS:HD2	1:A:97:LYS:HA	1.76	0.44
1:C:81:ARG:HB3	1:C:82:PRO:HD2	1.98	0.44
1:D:247:ALA:O	1:D:250:SER:HB2	2.17	0.44
1:D:264:MET:HE1	1:D:268:LEU:HD21	1.98	0.44
4:D:302:NAI:H2D	4:D:302:NAI:H6N	1.64	0.44
1:A:171:ALA:O	1:A:174:MET:HB2	2.17	0.44
1:C:248:VAL:O	1:C:251:ALA:HB3	2.16	0.44
1:D:45:LEU:HD12	1:D:63:GLN:HA	2.00	0.44
1:D:97:LYS:HA	6:D:428:HOH:O	2.16	0.44
1:A:110:GLY:O	1:A:113:ALA:HB3	2.18	0.44
4:B:302:NAI:H2B	4:B:302:NAI:H8A	1.64	0.44
1:A:118:ALA:HA	1:A:123:ILE:HD12	2.00	0.44
1:A:40:GLU:HG3	1:A:41:GLY:N	2.32	0.44
1:B:264:MET:HE3	1:B:268:LEU:HD11	1.99	0.44
1:C:89:LEU:HD22	1:C:99:MET:HE1	1.99	0.44
1:A:199:GLY:C	1:C:164:VAL:HG12	2.38	0.44
1:D:82:PRO:HA	1:D:106:PHE:CE1	2.53	0.44
1:A:201:ARG:HB3	1:C:161:ARG:HB2	1.99	0.44
1:C:158:ALA:HA	1:C:211:VAL:O	2.18	0.44
1:C:89:LEU:CA	1:C:99:MET:HE1	2.26	0.44
1:D:40:GLU:CD	1:D:40:GLU:N	2.70	0.44
1:A:18:GLY:O	1:A:19:ARG:C	2.56	0.44
1:B:44:LEU:HD11	1:B:47:SER:HB2	2.00	0.44
1:C:106:PHE:CD2	1:C:125:PHE:CZ	3.06	0.44
1:D:42:SER:C	1:D:44:LEU:H	2.22	0.44
1:C:17:MET:HB2	4:C:301:NAI:O2N	2.17	0.43
1:D:31:VAL:HG23	1:D:252:LEU:HD11	1.99	0.43
1:D:38:GLU:HG3	1:D:42:SER:OG	2.17	0.43
1:C:67:ASP:O	1:C:70:LYS:HG2	2.17	0.43
1:D:159:HIS:CE1	1:D:168:SER:HB3	2.53	0.43
1:B:38:GLU:HG3	1:B:42:SER:CB	2.48	0.43
1:C:202:VAL:O	1:C:205:THR:HG23	2.18	0.43
1:D:163:LYS:NZ	5:D:303:PDC:N1	2.65	0.43
1:D:4:ALA:HB1	1:D:5:ASN:HD22	1.83	0.43
1:A:144:ALA:O	1:A:148:GLY:HA3	2.18	0.43
1:B:137:LEU:HD12	1:B:137:LEU:HA	1.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ALA:HB1	1:A:248:VAL:HG21	2.01	0.43
1:A:252:LEU:O	1:A:252:LEU:HD12	2.18	0.43
1:B:60:VAL:CG1	1:B:61:THR:H	2.30	0.43
1:C:159:HIS:CE1	1:C:168:SER:HB3	2.53	0.43
1:A:10:ILE:HD11	1:A:25:ALA:HB2	2.00	0.43
1:B:48:ASP:O	1:B:50:GLY:N	2.52	0.43
1:C:7:ARG:CB	1:C:73:PHE:HA	2.47	0.43
1:C:86:LEU:O	1:C:89:LEU:CB	2.55	0.43
1:D:52:LEU:H	1:D:52:LEU:HD23	1.83	0.43
1:D:9:ALA:HB2	1:D:73:PHE:CG	2.53	0.43
1:B:122:ALA:O	1:B:254:LEU:HD11	2.18	0.43
1:B:7:ARG:HG3	1:B:7:ARG:HH11	1.83	0.43
1:D:153:ILE:H	1:D:201:ARG:NH2	2.16	0.43
1:D:264:MET:O	1:D:268:LEU:HD23	2.18	0.43
1:C:11:ALA:HB2	1:C:76:PHE:CZ	2.54	0.43
1:B:273:LEU:CD1	1:C:142:LYS:HD2	2.47	0.43
1:C:138:LYS:CE	1:C:272:ASN:H	2.32	0.43
1:D:80:THR:O	4:D:302:NAI:H4D	2.19	0.43
1:A:89:LEU:HD21	1:A:117:ALA:HB1	2.01	0.42
1:C:7:ARG:HA	1:C:32:GLN:O	2.19	0.42
1:B:129:PHE:HE2	1:B:250:SER:HG	1.64	0.42
1:D:16:ARG:H	1:D:16:ARG:HG2	1.61	0.42
1:D:270:LEU:O	1:D:272:ASN:N	2.52	0.42
1:B:120:ASP:C	1:B:259:SER:CB	2.87	0.42
1:D:42:SER:O	1:D:45:LEU:HD23	2.20	0.42
1:A:83:GLU:CA	1:A:86:LEU:HD23	2.48	0.42
1:B:163:LYS:NZ	5:B:303:PDC:N1	2.67	0.42
1:C:155:ILE:O	1:C:208:PHE:HA	2.19	0.42
1:C:16:ARG:H	1:C:16:ARG:HG2	1.67	0.42
1:B:131:VAL:O	1:B:132:GLY:C	2.57	0.42
1:B:77:ILE:HG22	1:B:77:ILE:O	2.17	0.42
1:A:82:PRO:O	1:A:83:GLU:C	2.57	0.42
1:C:114:ILE:H	1:C:114:ILE:HD12	1.85	0.42
1:D:42:SER:O	1:D:45:LEU:CD2	2.67	0.42
1:A:241:MET:O	1:A:244:ALA:N	2.53	0.42
1:D:123:ILE:HG22	1:D:124:VAL:N	2.34	0.42
1:A:160:HIS:CB	1:A:214:GLY:O	2.67	0.42
1:B:92:CYS:O	1:B:97:LYS:O	2.37	0.42
1:A:65:SER:CB	1:A:68:ALA:HB2	2.50	0.42
1:B:158:ALA:HA	1:B:211:VAL:O	2.20	0.42
1:B:48:ASP:C	1:B:50:GLY:N	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:160:HIS:HB2	1:D:214:GLY:O	2.20	0.42
1:D:33:LEU:O	1:D:60:VAL:HG21	2.20	0.41
1:D:110:GLY:O	1:D:114:ILE:HD12	2.21	0.41
1:D:45:LEU:HD11	1:D:64:SER:HB3	2.02	0.41
1:A:254:LEU:O	1:A:256:GLY:N	2.53	0.41
1:A:28:LEU:HD13	1:A:29:GLU:O	2.21	0.41
1:B:91:PHE:C	1:B:93:ARG:H	2.23	0.41
1:D:19:ARG:HA	1:D:53:ALA:CB	2.50	0.41
1:D:5:ASN:N	1:D:5:ASN:ND2	2.67	0.41
1:A:201:ARG:NH2	1:A:205:THR:O	2.52	0.41
1:A:124:VAL:CG2	1:A:254:LEU:HD21	2.47	0.41
1:A:57:LYS:O	1:A:58:THR:O	2.38	0.41
1:B:20:GLN:HE21	1:B:241:MET:HE1	1.85	0.41
1:D:189:CYS:HB2	1:D:205:THR:HA	2.03	0.41
1:D:67:ASP:O	1:D:70:LYS:HG2	2.20	0.41
1:A:212:ARG:HA	1:A:212:ARG:HD3	1.87	0.41
1:A:130:SER:OG	1:D:229:GLU:OE1	2.37	0.41
1:A:89:LEU:C	1:A:89:LEU:CD2	2.89	0.41
1:C:110:GLY:O	1:C:114:ILE:HD12	2.21	0.41
1:C:87:ASN:O	1:C:90:ALA:N	2.53	0.41
1:A:154:GLU:OE1	1:A:194:ARG:NH1	2.44	0.41
1:A:265:ARG:O	1:A:269:ASP:N	2.54	0.41
1:A:267:VAL:HG12	6:A:430:HOH:O	2.21	0.41
1:B:82:PRO:HA	1:B:106:PHE:CE1	2.56	0.41
1:B:136:MET:HE3	1:B:174:MET:CE	2.51	0.41
1:B:67:ASP:HB3	1:B:70:LYS:HD3	2.02	0.41
1:C:67:ASP:HB3	1:C:70:LYS:HD3	2.03	0.41
1:C:217:VAL:HB	1:C:237:ALA:O	2.20	0.41
1:C:26:LEU:HA	1:C:26:LEU:HD12	1.85	0.41
1:A:159:HIS:CE1	1:A:168:SER:HB3	2.56	0.41
1:B:5:ASN:HA	1:B:32:GLN:OE1	2.20	0.41
1:D:137:LEU:HD12	1:D:137:LEU:HA	1.72	0.41
1:B:200:GLU:CD	1:D:162:HIS:HD2	2.24	0.41
1:D:217:VAL:HB	1:D:237:ALA:O	2.21	0.41
1:D:32:GLN:O	1:D:34:GLY:N	2.54	0.41
1:B:79:PHE:CZ	1:B:102:GLY:HA3	2.56	0.41
1:B:20:GLN:HE21	1:B:241:MET:CE	2.34	0.41
1:B:114:ILE:CG2	1:B:261:LEU:HD11	2.47	0.41
1:B:45:LEU:HD11	1:B:63:GLN:C	2.41	0.41
1:C:263:ASP:N	1:C:266:ASP:OD1	2.54	0.41
1:D:76:PHE:HB3	1:D:99:MET:CG	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:PRO:O	1:C:86:LEU:HG	2.21	0.41
1:D:114:ILE:HD12	1:D:114:ILE:H	1.85	0.41
1:D:16:ARG:O	1:D:20:GLN:HG2	2.20	0.41
1:A:73:PHE:C	1:A:73:PHE:CD1	2.94	0.40
1:B:26:LEU:HA	1:B:26:LEU:HD12	1.80	0.40
1:D:273:LEU:HD22	1:D:273:LEU:OXT	2.21	0.40
1:D:88:HIS:N	1:D:88:HIS:ND1	2.60	0.40
1:A:23:GLN:HB2	1:A:241:MET:HE1	2.03	0.40
1:A:273:LEU:HD23	1:A:273:LEU:HA	1.37	0.40
1:B:154:GLU:OE2	1:B:201:ARG:NH1	2.55	0.40
1:B:60:VAL:CG1	1:B:61:THR:N	2.84	0.40
1:B:266:ASP:HA	1:B:271:ASN:HD21	1.86	0.40
1:C:123:ILE:HG22	1:C:124:VAL:N	2.36	0.40
1:A:164:VAL:O	1:C:197:HIS:HA	2.21	0.40
1:A:21:LEU:HD23	1:A:21:LEU:HA	1.84	0.40
1:B:15:GLY:HA3	4:B:302:NAI:H4B	2.02	0.40
1:B:251:ALA:O	1:B:252:LEU:C	2.60	0.40
1:B:89:LEU:HD22	1:B:99:MET:HE1	2.03	0.40
1:C:143:ALA:O	1:C:144:ALA:C	2.59	0.40
1:A:199:GLY:O	1:C:164:VAL:HA	2.21	0.40
1:C:156:ILE:HG23	1:C:209:ALA:HB3	2.02	0.40
1:C:91:PHE:CZ	1:C:95:HIS:CD2	3.10	0.40
1:D:25:ALA:HA	1:D:31:VAL:CG1	2.52	0.40
1:A:79:PHE:CD2	1:A:102:GLY:HA3	2.56	0.40
1:A:74:ASP:O	1:A:97:LYS:O	2.39	0.40
1:B:144:ALA:O	1:B:148:GLY:HA3	2.22	0.40
1:C:265:ARG:O	1:C:269:ASP:CA	2.70	0.40
1:D:256:GLY:O	1:D:257:LYS:C	2.60	0.40

There are no symmetry-related clashes.

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	268/273 (98%)	218 (81%)	40 (15%)	10 (4%)	4	5
1	B	267/273 (98%)	214 (80%)	50 (19%)	3 (1%)	17	35
1	C	269/273 (98%)	224 (83%)	40 (15%)	5 (2%)	9	18
1	D	269/273 (98%)	230 (86%)	34 (13%)	5 (2%)	9	18
All	All	1073/1092 (98%)	886 (83%)	164 (15%)	23 (2%)	8	15

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	ALA
1	A	49	ALA
1	A	58	THR
1	A	82	PRO
1	B	47	SER
1	C	4	ALA
1	D	6	ILE
1	A	255	SER
1	B	70	LYS
1	C	70	LYS
1	C	90	ALA
1	D	70	LYS
1	A	12	GLY
1	C	49	ALA
1	D	160	HIS
1	D	271	ASN
1	A	160	HIS
1	B	160	HIS
1	C	272	ASN
1	A	66	LEU
1	D	267	VAL
1	A	132	GLY
1	A	10	ILE

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	185/201 (92%)	161 (87%)	24 (13%)	5	8
1	B	195/201 (97%)	166 (85%)	29 (15%)	3	6
1	C	193/201 (96%)	168 (87%)	25 (13%)	5	8
1	D	195/201 (97%)	165 (85%)	30 (15%)	3	5
All	All	768/804 (96%)	660 (86%)	108 (14%)	4	7

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	20	GLN
1	A	29	GLU
1	A	31	VAL
1	A	33	LEU
1	A	58	THR
1	A	66	LEU
1	A	69	VAL
1	A	83	GLU
1	A	99	MET
1	A	100	VAL
1	A	101	ILE
1	A	104	THR
1	A	120	ASP
1	A	137	LEU
1	A	145	LYS
1	A	157	GLU
1	A	184	LYS
1	A	227	ILE
1	A	231	LEU
1	A	239	SER
1	A	263	ASP
1	A	268	LEU
1	A	273	LEU
1	B	5	ASN
1	B	19	ARG
1	B	26	LEU
1	B	32	GLN
1	B	40	GLU
1	B	44	LEU
1	B	51	GLU
1	B	60	VAL
1	B	66	LEU

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Mol	Chain	Res	Type
1	B	67	ASP
1	B	71	ASP
1	B	92	CYS
1	B	95	HIS
1	B	99	MET
1	B	108	GLU
1	B	115	ARG
1	B	120	ASP
1	B	145	LYS
1	B	157	GLU
1	B	176	GLU
1	B	184	LYS
1	B	227	ILE
1	B	231	LEU
1	B	240	ARG
1	B	254	LEU
1	B	261	LEU
1	B	266	ASP
1	B	272	ASN
1	B	273	LEU
1	C	19	ARG
1	C	26	LEU
1	C	52	LEU
1	C	62	VAL
1	C	66	LEU
1	C	67	ASP
1	C	71	ASP
1	C	87	ASN
1	C	88	HIS
1	C	95	HIS
1	C	99	MET
1	C	108	GLU
1	C	115	ARG
1	C	120	ASP
1	C	145	LYS
1	C	157	GLU
1	C	176	GLU
1	C	184	LYS
1	C	197	HIS
1	C	227	ILE
1	C	231	LEU
1	C	240	ARG

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Mol	Chain	Res	Type
1	C	252	LEU
1	C	255	SER
1	C	262	PHE
1	D	5	ASN
1	D	6	ILE
1	D	19	ARG
1	D	26	LEU
1	D	31	VAL
1	D	39	ARG
1	D	42	SER
1	D	43	SER
1	D	52	LEU
1	D	62	VAL
1	D	66	LEU
1	D	67	ASP
1	D	71	ASP
1	D	88	HIS
1	D	108	GLU
1	D	115	ARG
1	D	120	ASP
1	D	137	LEU
1	D	145	LYS
1	D	157	GLU
1	D	176	GLU
1	D	184	LYS
1	D	198	THR
1	D	227	ILE
1	D	231	LEU
1	D	240	ARG
1	D	249	ARG
1	D	261	LEU
1	D	262	PHE
1	D	273	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	20	GLN
1	A	162	HIS
1	A	245	ASN
1	B	20	GLN

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Mol	Chain	Res	Type
1	B	88	HIS
1	B	128	ASN
1	B	159	HIS
1	B	271	ASN
1	C	20	GLN
1	C	128	ASN
1	C	162	HIS
1	C	197	HIS
1	D	5	ASN
1	D	20	GLN
1	D	128	ASN
1	D	162	HIS
1	D	197	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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	$\# Z > 2$
2	PO4	A	301	-	4,4,4	1.18	1 (25%)	6,6,6	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAI	B	302	-	40,48,48	1.66	9 (22%)	41,73,73	2.08	5 (12%)
5	PDC	B	303	-	6,12,12	2.87	2 (33%)	8,16,16	0.58	0
4	NAI	C	301	-	40,48,48	1.73	8 (20%)	41,73,73	2.02	8 (19%)
5	PDC	C	302	-	6,12,12	3.08	2 (33%)	8,16,16	1.33	0
4	NAI	D	302	-	40,48,48	1.50	6 (15%)	41,73,73	2.08	5 (12%)
5	PDC	D	303	-	6,12,12	2.13	2 (33%)	8,16,16	1.60	3 (37%)

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
2	PO4	A	301	-	-	0/0/0/0	0/0/0/0
4	NAI	B	302	-	-	0/25/72/72	0/5/5/5
5	PDC	B	303	-	-	0/0/8/8	0/1/1/1
4	NAI	C	301	-	-	0/25/72/72	0/5/5/5
5	PDC	C	302	-	-	0/0/8/8	0/1/1/1
4	NAI	D	302	-	-	0/25/72/72	0/5/5/5
5	PDC	D	303	-	-	0/0/8/8	0/1/1/1

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	301	NAI	C2B-C1B	-4.94	1.45	1.53
4	C	301	NAI	C4N-C5N	-4.69	1.38	1.49
4	D	302	NAI	C4N-C5N	-3.89	1.40	1.49
4	C	301	NAI	C5A-C4A	-3.38	1.32	1.40
4	D	302	NAI	C5A-C4A	-3.30	1.33	1.40
4	D	302	NAI	C2B-C1B	-3.08	1.48	1.53
4	B	302	NAI	C5A-C4A	-2.53	1.34	1.40
4	B	302	NAI	C4N-C5N	-2.32	1.44	1.49
4	B	302	NAI	C2B-C1B	-2.08	1.50	1.53
4	C	301	NAI	C5B-C4B	2.07	1.58	1.51
4	D	302	NAI	O5B-C5B	2.16	1.53	1.44
4	D	302	NAI	C2A-N3A	2.17	1.35	1.32
4	C	301	NAI	C2N-C3N	2.18	1.41	1.34
4	C	301	NAI	O4B-C1B	2.27	1.44	1.41
4	B	302	NAI	O5B-C5B	2.29	1.53	1.44
2	A	301	PO4	P-O1	2.32	1.55	1.50
4	B	302	NAI	C2A-N1A	2.33	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	302	NAI	C1D-N1N	2.62	1.54	1.46
4	C	301	NAI	O5B-C5B	2.70	1.55	1.44
4	B	302	NAI	C2A-N3A	2.81	1.36	1.32
4	B	302	NAI	C2N-C3N	2.83	1.43	1.34
5	D	303	PDC	C2-N1	3.19	1.39	1.34
4	D	302	NAI	C6N-C5N	3.52	1.39	1.33
5	D	303	PDC	C6-N1	3.69	1.40	1.34
4	C	301	NAI	C6N-C5N	3.95	1.40	1.33
5	B	303	PDC	C2-N1	4.65	1.42	1.34
5	B	303	PDC	C6-N1	4.75	1.42	1.34
5	C	302	PDC	C2-N1	4.82	1.42	1.34
5	C	302	PDC	C6-N1	5.21	1.42	1.34
4	B	302	NAI	C6N-C5N	5.68	1.43	1.33

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	301	NAI	N3A-C2A-N1A	-3.95	125.41	128.86
4	C	301	NAI	O3B-C3B-C4B	-2.81	102.89	111.09
4	C	301	NAI	O4D-C1D-N1N	-2.44	103.15	108.07
5	D	303	PDC	C5-C6-N1	-2.15	119.06	121.97
4	D	302	NAI	O3B-C3B-C4B	-2.06	105.06	111.09
4	B	302	NAI	C2D-C3D-C4D	-2.01	98.71	102.62
5	D	303	PDC	C6-N1-C2	2.02	120.52	118.04
4	C	301	NAI	O4B-C4B-C5B	2.22	116.89	109.40
4	D	302	NAI	C4A-C5A-N7A	2.32	111.65	109.41
5	D	303	PDC	C5-C6-C8	2.56	123.14	120.16
4	B	302	NAI	C4A-C5A-N7A	2.60	111.93	109.41
4	D	302	NAI	O5D-C5D-C4D	2.69	118.54	109.00
4	B	302	NAI	O5B-C5B-C4B	3.43	121.15	109.00
4	C	301	NAI	C4A-C5A-N7A	3.47	112.77	109.41
4	B	302	NAI	O5D-C5D-C4D	3.60	121.75	109.00
4	C	301	NAI	O5D-C5D-C4D	4.08	123.47	109.00
4	C	301	NAI	O5B-C5B-C4B	4.20	123.91	109.00
4	D	302	NAI	O5B-C5B-C4B	4.26	124.12	109.00
4	C	301	NAI	C1B-N9A-C4A	7.80	140.11	126.64
4	D	302	NAI	C1B-N9A-C4A	10.04	143.99	126.64
4	B	302	NAI	C1B-N9A-C4A	10.16	144.19	126.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	302	NAI	6	0
5	B	303	PDC	1	0
4	C	301	NAI	3	0
5	C	302	PDC	2	0
4	D	302	NAI	5	0
5	D	303	PDC	4	0

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	270/273 (98%)	-0.73	0	100	100	5, 30, 64, 87	0
1	B	269/273 (98%)	-0.60	0	100	100	8, 31, 64, 77	0
1	C	271/273 (99%)	-0.61	1 (0%)	92	91	8, 28, 62, 91	0
1	D	271/273 (99%)	-0.67	0	100	100	9, 29, 62, 77	0
All	All	1081/1092 (98%)	-0.65	1 (0%)	95	95	5, 30, 63, 91	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	3	ASP	2.9

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
5	PDC	D	303	12/12	0.96	0.12	0.23	2,17,22,29	0
5	PDC	C	302	12/12	0.96	0.12	-0.02	2,11,15,15	0
4	NAI	D	302	44/44	0.96	0.11	-0.79	2,25,51,52	0
4	NAI	B	302	44/44	0.95	0.11	-0.95	4,32,49,51	0
4	NAI	C	301	44/44	0.97	0.10	-1.11	7,15,32,35	0
3	K	D	301	1/1	0.98	0.08	-1.63	24,24,24,24	0
5	PDC	B	303	12/12	0.98	0.08	-1.64	2,4,12,14	0
3	K	B	301	1/1	0.99	0.06	-3.09	29,29,29,29	0
2	PO4	A	301	5/5	0.91	0.18	-	37,41,50,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.