



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:37 am GMT

PDB ID : 1Q3Q
Title : Crystal structure of the chaperonin from Thermococcus strain KS-1 (two-point mutant complexed with AMP-PNP)
Authors : Shomura, Y.; Yoshida, T.; Iizuka, R.; Maruyama, T.; Yohda, M.; Miki, K.
Deposited on : 2003-07-31
Resolution : 2.30 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

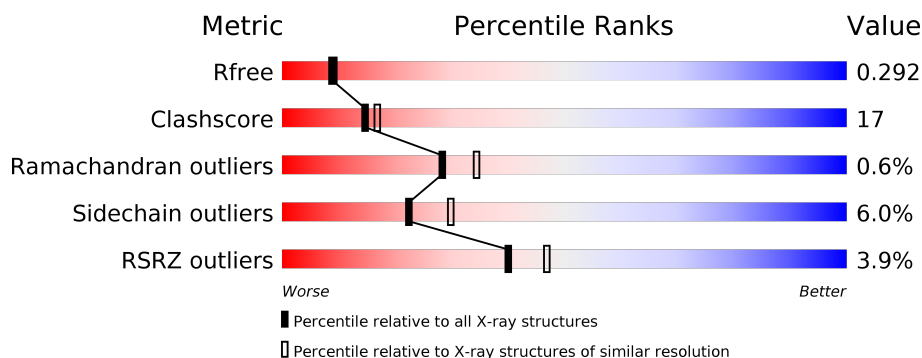
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	548	<div> <div>2%</div> <div> <div></div> <div>64%</div> <div>28%</div> <div>5%</div> </div> </div>
1	B	548	<div> <div>4%</div> <div> <div></div> <div>62%</div> <div>30%</div> <div>5%</div> </div> </div>
1	C	548	<div> <div>5%</div> <div> <div></div> <div>64%</div> <div>27%</div> <div>5%</div> </div> </div>
1	D	548	<div> <div>3%</div> <div> <div></div> <div>65%</div> <div>25%</div> <div>5%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ANP	B	2528	-	-	-	X
3	ANP	C	3528	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16610 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 Thermosome alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	518	Total	C	N	O	S	0	0	0
			3947	2485	673	772	17			
1	B	518	Total	C	N	O	S	0	0	0
			3947	2485	673	772	17			
1	C	518	Total	C	N	O	S	0	0	0
			3947	2485	673	772	17			
1	D	518	Total	C	N	O	S	0	0	0
			3947	2485	673	772	17			

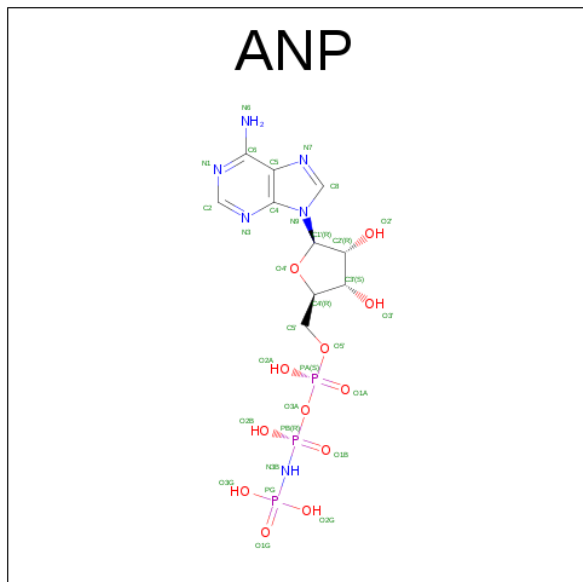
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	65	CYS	GLY	ENGINEERED	UNP O24729
A	125	THR	ILE	ENGINEERED	UNP O24729
B	65	CYS	GLY	ENGINEERED	UNP O24729
B	125	THR	ILE	ENGINEERED	UNP O24729
C	65	CYS	GLY	ENGINEERED	UNP O24729
C	125	THR	ILE	ENGINEERED	UNP O24729
D	65	CYS	GLY	ENGINEERED	UNP O24729
D	125	THR	ILE	ENGINEERED	UNP O24729

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
3	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
3	C	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
3	D	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

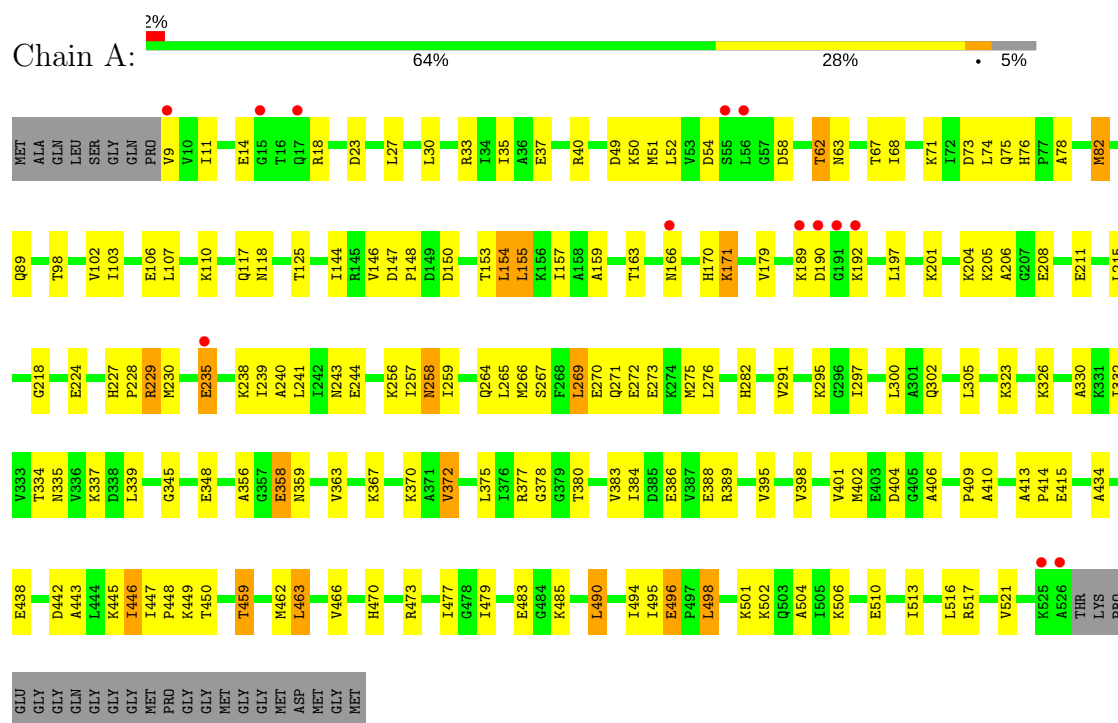
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	162	Total	O	0	0
			162	162		
4	B	174	Total	O	0	0
			174	174		
4	C	194	Total	O	0	0
			194	194		
4	D	164	Total	O	0	0
			164	164		

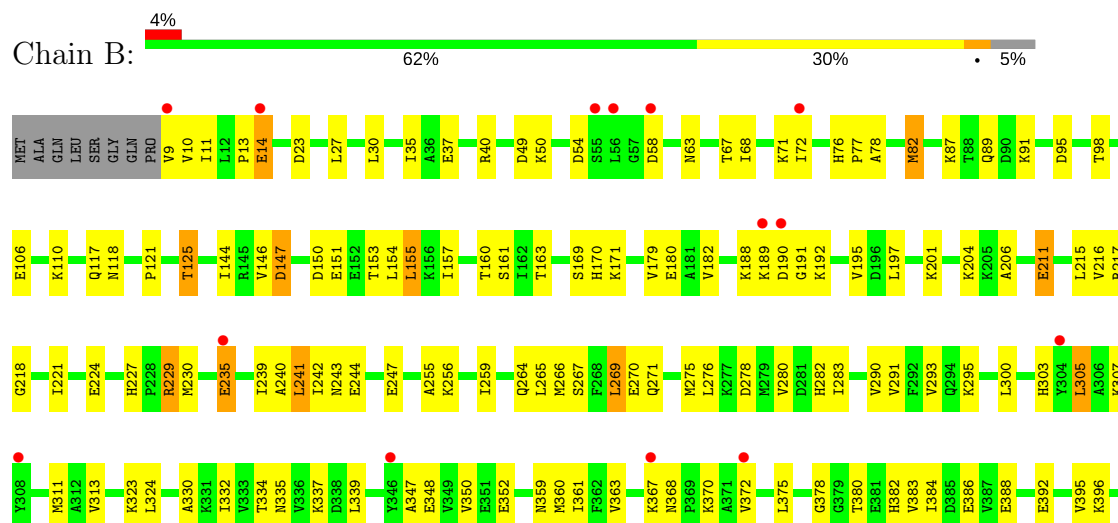
3 Residue-property plots [i](#)

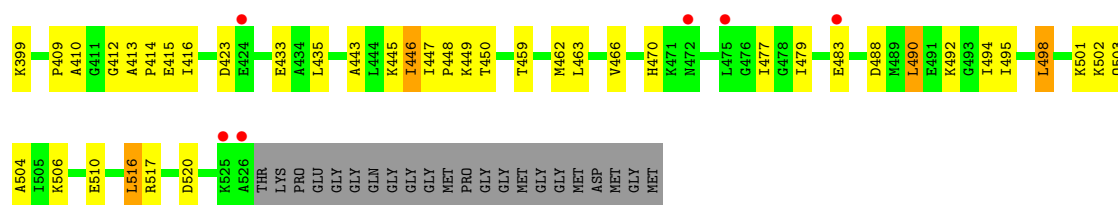
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Thermosome alpha subunit

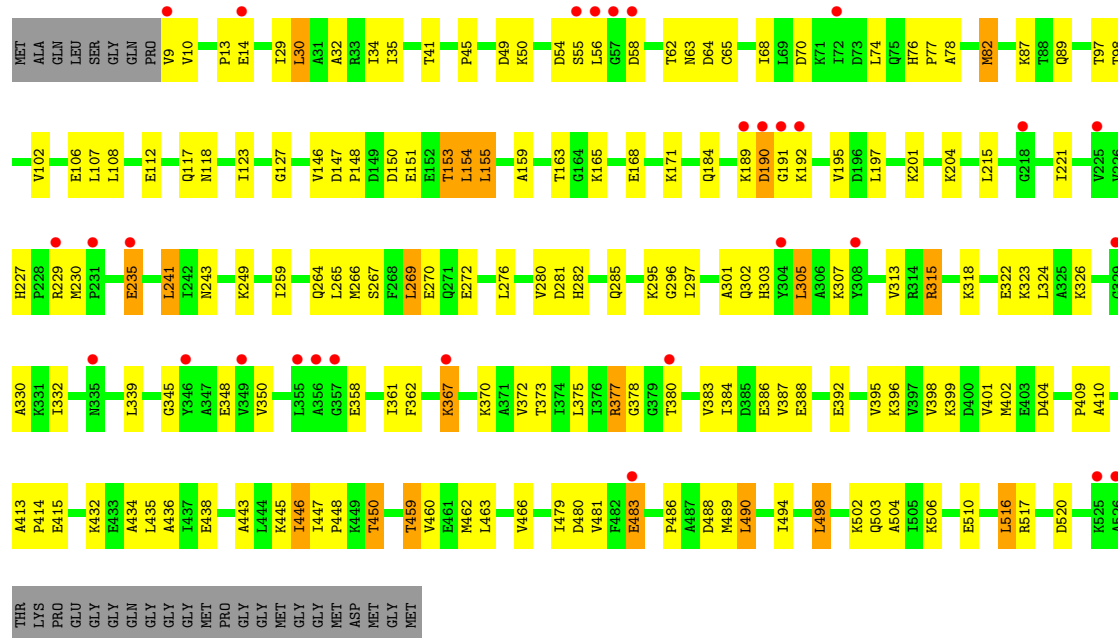


• Molecule 1: Thermosome alpha subunit

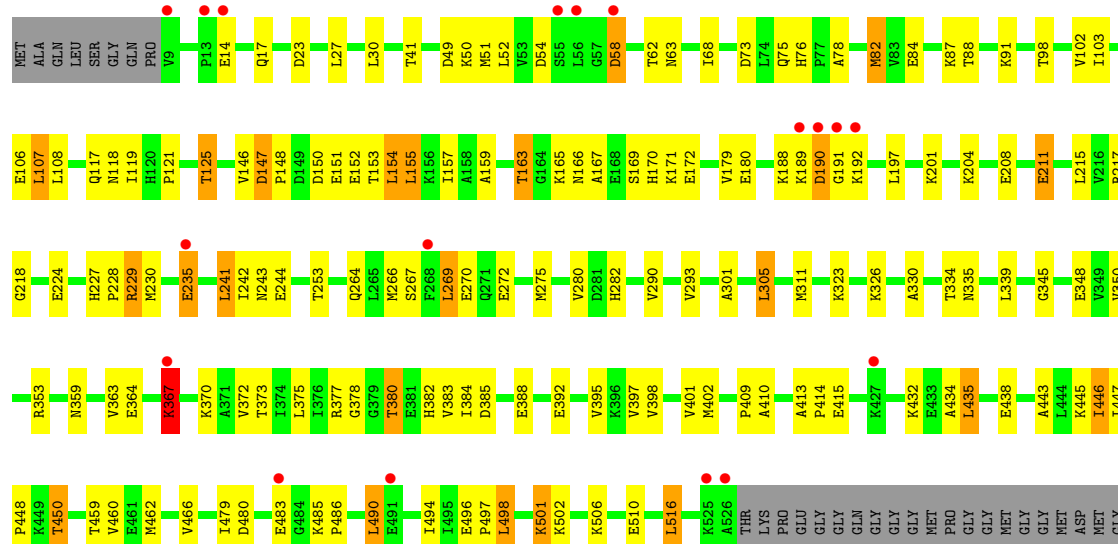




● Molecule 1: Thermosome alpha subunit



● Molecule 1: Thermosome alpha subunit



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4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	210.83Å 210.83Å 156.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.97 – 2.30 42.17 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.1 (41.97-2.30) 95.4 (42.17-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 2.29Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.213 , 0.250 0.268 , 0.292	Depositor DCC
R_{free} test set	7448 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	21.7	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.3	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	0.87	EDS
Total number of atoms	16610	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.32	0/3982	0.61	0/5367
1	B	0.32	0/3982	0.60	0/5367
1	C	0.33	0/3982	0.61	0/5367
1	D	0.32	0/3982	0.61	0/5367
All	All	0.32	0/15928	0.61	0/21468

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3947	0	4127	148	0
1	B	3947	0	4127	145	0
1	C	3947	0	4127	148	2
1	D	3947	0	4127	144	2
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	31	0	13	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	31	0	13	2	0
3	C	31	0	13	2	0
3	D	31	0	13	2	0
4	A	162	0	0	5	1
4	B	174	0	0	7	2
4	C	194	0	0	6	3
4	D	164	0	0	3	4
All	All	16610	0	16560	554	7

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 17.

All (554) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:VAL:HG23	4:A:1148:HOH:O	1.60	1.01
1:C:446:ILE:O	1:C:450:THR:HG23	1.65	0.97
1:D:446:ILE:O	1:D:450:THR:HG23	1.65	0.97
1:A:166:ASN:HB3	1:B:517:ARG:HD2	1.48	0.95
1:D:466:VAL:HG21	1:D:479:ILE:HG12	1.50	0.94
1:C:62:THR:HG22	1:C:64:ASP:H	1.31	0.93
1:C:503:GLN:HE22	1:D:208:GLU:H	1.07	0.93
1:A:462:MET:O	1:A:466:VAL:HG23	1.67	0.93
1:C:146:VAL:HG21	1:C:153:THR:HG21	1.51	0.90
1:A:208:GLU:H	1:B:503:GLN:HE22	1.13	0.90
1:A:410:ALA:HB3	1:A:494:ILE:HG22	1.54	0.90
1:D:380:THR:HG22	1:D:383:VAL:H	1.35	0.90
1:C:98:THR:H	3:C:3528:ANP:HNB1	1.20	0.89
1:A:163:THR:HG22	1:A:171:LYS:NZ	1.88	0.88
1:C:466:VAL:HG21	1:C:479:ILE:HG12	1.56	0.86
1:B:409:PRO:HB3	1:B:490:LEU:HD13	1.56	0.86
1:A:98:THR:H	3:A:1528:ANP:HNB1	1.23	0.85
1:A:459:THR:HG23	4:C:3189:HOH:O	1.75	0.85
1:B:466:VAL:HG21	1:B:479:ILE:HG12	1.60	0.83
1:D:410:ALA:HB3	1:D:494:ILE:HG22	1.58	0.83
1:C:410:ALA:HB3	1:C:494:ILE:HG22	1.61	0.83
1:B:462:MET:O	1:B:466:VAL:HG23	1.78	0.83
1:D:98:THR:H	3:D:4528:ANP:HNB1	1.27	0.82
1:A:163:THR:HG22	1:A:171:LYS:HZ2	1.44	0.81
1:C:503:GLN:HE22	1:D:208:GLU:N	1.78	0.81
1:B:450:THR:HG22	4:B:2054:HOH:O	1.80	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ASN:HB3	1:B:517:ARG:CD	2.12	0.80
1:B:98:THR:H	3:B:2528:ANP:HNB1	1.30	0.79
1:B:170:HIS:CD2	1:B:211:GLU:HG2	2.18	0.79
1:A:409:PRO:HB3	1:A:490:LEU:HD13	1.65	0.78
1:C:503:GLN:NE2	1:D:208:GLU:H	1.79	0.78
1:B:446:ILE:O	1:B:450:THR:HG23	1.83	0.78
1:C:447:ILE:HB	1:C:448:PRO:HD3	1.65	0.78
1:D:235:GLU:O	1:D:348:GLU:O	2.01	0.77
4:C:3074:HOH:O	1:D:380:THR:HG21	1.84	0.77
1:B:211:GLU:HB2	4:B:2030:HOH:O	1.86	0.76
1:B:160:THR:O	1:B:163:THR:HG22	1.85	0.76
1:B:23:ASP:O	1:B:27:LEU:HD23	1.85	0.76
1:C:409:PRO:HB3	1:C:490:LEU:HD13	1.68	0.76
1:A:380:THR:HG21	4:B:1067:HOH:O	1.85	0.75
1:A:466:VAL:HG21	1:A:479:ILE:HG12	1.66	0.75
1:C:462:MET:O	1:C:466:VAL:HG23	1.86	0.74
1:C:204:LYS:HB3	1:C:384:ILE:HG21	1.68	0.74
1:C:204:LYS:HD2	1:C:384:ILE:HG22	1.70	0.74
1:A:483:GLU:HG2	1:A:485:LYS:HE2	1.69	0.74
1:C:62:THR:HG23	1:C:386:GLU:OE2	1.88	0.74
1:D:445:LYS:O	1:D:448:PRO:HD2	1.88	0.74
1:D:170:HIS:CD2	1:D:211:GLU:HG2	2.23	0.74
1:B:163:THR:HB	1:B:171:LYS:NZ	2.02	0.73
1:D:434:ALA:O	1:D:438:GLU:HG3	1.89	0.73
1:D:462:MET:O	1:D:466:VAL:HG23	1.87	0.73
1:A:447:ILE:HB	1:A:448:PRO:HD3	1.69	0.73
1:A:150:ASP:OD2	1:A:153:THR:HG23	1.89	0.73
1:D:409:PRO:HB3	1:D:490:LEU:HD13	1.72	0.72
1:B:280:VAL:HG13	1:B:305:LEU:HD13	1.72	0.71
1:D:189:LYS:HG3	1:D:190:ASP:OD2	1.91	0.71
1:C:280:VAL:HG13	1:C:305:LEU:HD13	1.70	0.71
1:D:23:ASP:O	1:D:27:LEU:HD23	1.90	0.71
1:C:204:LYS:HB3	1:C:384:ILE:CG2	2.21	0.71
1:C:434:ALA:O	1:C:438:GLU:HG3	1.92	0.70
1:A:450:THR:HG22	4:A:1037:HOH:O	1.91	0.70
1:A:415:GLU:HG3	1:A:447:ILE:HB	1.73	0.70
1:C:380:THR:HG23	1:C:383:VAL:H	1.55	0.70
1:A:54:ASP:OD2	1:A:58:ASP:HB3	1.90	0.69
1:A:204:LYS:HB3	1:A:384:ILE:HG21	1.74	0.69
1:A:189:LYS:HG2	1:A:192:LYS:O	1.93	0.69
1:B:163:THR:HB	1:B:171:LYS:HZ3	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:415:GLU:HG3	1:C:447:ILE:HB	1.74	0.69
1:C:466:VAL:HG22	1:C:486:PRO:HG3	1.73	0.69
1:A:75:GLN:NE2	1:B:13:PRO:HG3	2.07	0.69
1:B:447:ILE:HB	1:B:448:PRO:HD3	1.75	0.69
1:C:54:ASP:OD2	1:C:58:ASP:HB2	1.92	0.69
1:A:443:ALA:O	1:A:446:ILE:HG13	1.93	0.68
1:B:201:LYS:HB2	1:B:323:LYS:HE2	1.76	0.68
1:B:215:LEU:HD21	1:B:372:VAL:HG21	1.76	0.68
1:B:443:ALA:O	1:B:446:ILE:HG13	1.93	0.68
1:A:208:GLU:N	1:B:503:GLN:HE22	1.90	0.68
1:B:227:HIS:CE1	1:B:229:ARG:HB2	2.29	0.67
1:A:23:ASP:O	1:A:27:LEU:HD23	1.93	0.67
1:A:50:LYS:HG2	1:B:520:ASP:HB3	1.77	0.67
1:C:380:THR:HG22	1:C:383:VAL:HG23	1.76	0.67
1:B:50:LYS:HD2	1:B:68:ILE:HD13	1.77	0.67
1:A:459:THR:CG2	4:C:3189:HOH:O	2.40	0.67
1:D:166:ASN:O	1:D:167:ALA:HB3	1.94	0.66
1:B:410:ALA:HB3	1:B:494:ILE:HG22	1.76	0.66
1:B:204:LYS:HB2	1:B:384:ILE:HG21	1.77	0.66
1:D:466:VAL:HG21	1:D:479:ILE:CG1	2.24	0.66
1:A:227:HIS:HD2	1:A:229:ARG:H	1.43	0.66
1:C:401:VAL:O	1:C:404:ASP:O	2.14	0.66
1:D:50:LYS:HD2	1:D:68:ILE:HD13	1.78	0.66
1:C:201:LYS:HB2	1:C:323:LYS:HE2	1.78	0.66
1:A:125:THR:HG21	4:A:1087:HOH:O	1.95	0.65
1:B:54:ASP:OD2	1:B:58:ASP:HB3	1.97	0.65
1:B:466:VAL:HG21	1:B:479:ILE:CG1	2.25	0.65
1:C:443:ALA:O	1:C:446:ILE:HG13	1.96	0.65
1:B:243:ASN:O	1:B:295:LYS:HG2	1.97	0.65
1:C:150:ASP:HB3	1:C:153:THR:HG23	1.79	0.65
1:B:266:MET:O	1:B:270:GLU:HG3	1.95	0.64
1:A:208:GLU:H	1:B:503:GLN:NE2	1.92	0.64
1:B:227:HIS:HB3	1:B:230:MET:HG3	1.79	0.64
1:A:82:MET:CE	1:A:82:MET:HA	2.28	0.64
1:C:445:LYS:O	1:C:448:PRO:HD2	1.97	0.64
1:D:146:VAL:HG21	1:D:153:THR:HG21	1.78	0.64
1:B:247:GLU:HA	1:B:276:LEU:HD21	1.79	0.64
1:D:498:LEU:HD22	1:D:502:LYS:HG3	1.79	0.64
1:A:335:ASN:OD1	1:A:337:LYS:HG2	1.97	0.64
1:D:447:ILE:HB	1:D:448:PRO:HD3	1.78	0.64
1:C:159:ALA:O	1:C:163:THR:HG23	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:LEU:HB3	1:B:332:ILE:HD13	1.80	0.63
1:B:204:LYS:HB2	1:B:384:ILE:CG2	2.28	0.63
1:C:215:LEU:HD11	1:C:372:VAL:HG13	1.80	0.63
1:A:243:ASN:OD1	1:A:334:THR:HA	1.97	0.63
1:A:410:ALA:HB1	3:A:1528:ANP:O2'	1.99	0.63
1:A:189:LYS:HG3	1:A:190:ASP:OD2	1.99	0.63
1:A:244:GLU:HB3	1:A:334:THR:O	1.99	0.63
1:C:197:LEU:HD22	1:C:395:VAL:CG1	2.28	0.63
1:A:50:LYS:HD2	1:A:68:ILE:HD13	1.81	0.62
1:D:49:ASP:OD1	1:D:63:ASN:HB2	1.99	0.62
1:A:155:LEU:HD12	1:A:179:VAL:HG21	1.81	0.62
1:C:410:ALA:HB1	3:C:3528:ANP:O2'	1.98	0.62
1:A:125:THR:HG23	1:A:513:ILE:HG23	1.80	0.62
1:B:241:LEU:HD22	1:B:330:ALA:HB3	1.81	0.62
1:D:190:ASP:C	1:D:192:LYS:H	2.00	0.62
1:C:32:ALA:HB1	1:C:82:MET:CE	2.29	0.62
1:D:150:ASP:OD1	1:D:153:THR:HG23	1.98	0.62
1:D:415:GLU:HG3	1:D:447:ILE:HB	1.80	0.62
1:A:276:LEU:HD23	1:A:300:LEU:HB3	1.81	0.62
1:C:215:LEU:HD11	1:C:372:VAL:CG1	2.29	0.62
1:A:197:LEU:HD22	1:A:395:VAL:CG1	2.30	0.61
1:A:446:ILE:O	1:A:450:THR:HG23	2.00	0.61
1:C:259:ILE:HD13	1:C:265:LEU:HD23	1.83	0.61
1:C:108:LEU:HD23	1:C:516:LEU:HD22	1.82	0.61
1:D:266:MET:HE2	1:D:269:LEU:HD23	1.82	0.61
1:A:146:VAL:CG2	4:A:1148:HOH:O	2.31	0.61
1:C:498:LEU:HD22	1:C:502:LYS:HG3	1.82	0.61
1:D:241:LEU:HD22	1:D:330:ALA:HB3	1.83	0.61
1:A:204:LYS:HB3	1:A:384:ILE:CG2	2.30	0.61
1:B:215:LEU:HD21	1:B:372:VAL:CG2	2.31	0.60
1:A:227:HIS:CD2	1:A:229:ARG:H	2.18	0.60
1:D:87:LYS:HA	1:D:87:LYS:HE2	1.82	0.60
1:D:76:HIS:HD2	1:D:78:ALA:H	1.48	0.60
1:B:215:LEU:HD23	1:B:216:VAL:N	2.16	0.60
1:D:443:ALA:O	1:D:446:ILE:HG13	2.01	0.60
1:B:278:ASP:OD2	4:B:2073:HOH:O	2.17	0.60
1:A:163:THR:HG22	1:A:171:LYS:HZ1	1.66	0.59
1:A:170:HIS:CD2	1:A:211:GLU:HG2	2.37	0.59
1:D:155:LEU:HD12	1:D:179:VAL:HG21	1.84	0.59
1:D:76:HIS:CD2	1:D:78:ALA:H	2.20	0.59
1:A:380:THR:HG23	1:A:383:VAL:H	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:VAL:HG22	1:A:147:ASP:N	2.17	0.59
1:C:56:LEU:HD12	1:C:56:LEU:N	2.17	0.59
1:D:378:GLY:N	1:D:384:ILE:HD11	2.17	0.59
1:D:410:ALA:HB1	3:D:4528:ANP:O2'	2.03	0.59
1:C:62:THR:HG22	1:C:63:ASN:H	1.67	0.59
1:D:201:LYS:HB2	1:D:323:LYS:HE2	1.85	0.58
1:B:206:ALA:HA	1:B:384:ILE:HD12	1.84	0.58
1:A:170:HIS:NE2	1:A:211:GLU:HG2	2.18	0.58
1:A:227:HIS:CD2	1:A:229:ARG:HB2	2.38	0.58
1:A:239:ILE:HD12	1:A:239:ILE:N	2.19	0.58
1:C:108:LEU:HD23	1:C:516:LEU:CD2	2.33	0.58
1:C:70:ASP:HB2	1:C:87:LYS:HE2	1.86	0.57
1:C:13:PRO:HG3	1:D:75:GLN:NE2	2.19	0.57
1:C:266:MET:O	1:C:270:GLU:HG3	2.04	0.57
1:A:204:LYS:HD2	1:A:384:ILE:HG22	1.86	0.57
1:B:466:VAL:CG2	1:B:479:ILE:HG12	2.34	0.57
1:C:108:LEU:CD2	1:C:516:LEU:HD22	2.35	0.57
1:D:244:GLU:HB2	1:D:334:THR:O	2.03	0.57
1:B:244:GLU:HB3	1:B:334:THR:O	2.04	0.57
1:C:191:GLY:C	1:C:192:LYS:HD2	2.25	0.57
1:C:62:THR:HG22	1:C:63:ASN:N	2.19	0.57
1:A:282:HIS:HE1	1:A:339:LEU:O	1.88	0.57
1:B:445:LYS:O	1:B:448:PRO:HD2	2.04	0.57
1:D:483:GLU:HG3	1:D:485:LYS:HE3	1.85	0.57
1:D:159:ALA:O	1:D:163:THR:HG22	2.05	0.56
1:B:87:LYS:O	1:B:91:LYS:HG2	2.05	0.56
1:A:148:PRO:O	1:A:402:MET:HG2	2.06	0.56
1:B:433:GLU:CD	1:B:433:GLU:H	2.09	0.56
1:A:380:THR:HG22	1:A:383:VAL:HB	1.87	0.56
1:B:151:GLU:HG2	1:B:155:LEU:HD22	1.87	0.56
1:A:117:GLN:O	1:A:118:ASN:HB2	2.06	0.56
1:C:380:THR:HG22	1:C:383:VAL:CG2	2.36	0.56
1:D:146:VAL:HG22	1:D:147:ASP:N	2.20	0.56
1:A:513:ILE:O	1:A:517:ARG:HG3	2.06	0.56
1:A:434:ALA:O	1:A:438:GLU:HG3	2.06	0.55
1:D:266:MET:O	1:D:270:GLU:HG3	2.05	0.55
1:D:413:ALA:HB3	1:D:414:PRO:HD3	1.88	0.55
1:B:224:GLU:HG2	1:B:359:ASN:HB2	1.88	0.55
1:A:146:VAL:HG21	1:A:153:THR:HG21	1.87	0.55
1:C:227:HIS:CD2	1:C:229:ARG:H	2.24	0.55
1:B:189:LYS:HG3	1:B:190:ASP:OD2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:LYS:HB2	1:C:323:LYS:CE	2.36	0.55
1:C:380:THR:CG2	1:C:383:VAL:HG23	2.37	0.55
1:B:498:LEU:HD22	1:B:502:LYS:HG3	1.89	0.55
1:C:30:LEU:HD22	1:C:34:ILE:HD11	1.89	0.55
1:A:264:GLN:HA	1:A:267:SER:OG	2.07	0.54
1:C:313:VAL:HG21	1:C:361:ILE:CD1	2.36	0.54
1:B:415:GLU:HG3	1:B:447:ILE:HB	1.89	0.54
1:C:227:HIS:HB3	1:C:230:MET:HG3	1.88	0.54
1:A:146:VAL:HG22	1:A:147:ASP:H	1.72	0.54
1:B:335:ASN:HB3	1:B:337:LYS:HG2	1.89	0.54
1:A:211:GLU:HB2	4:A:1110:HOH:O	2.07	0.54
1:B:37:GLU:HA	1:B:40:ARG:HG2	1.90	0.54
1:D:445:LYS:C	1:D:448:PRO:HD2	2.27	0.54
1:B:150:ASP:OD1	1:B:153:THR:HG23	2.08	0.54
1:C:281:ASP:O	1:C:285:GLN:HG3	2.08	0.54
1:B:410:ALA:HB1	3:B:2528:ANP:O2'	2.07	0.54
1:C:77:PRO:HB2	1:D:51:MET:CE	2.37	0.53
1:A:206:ALA:HA	1:A:384:ILE:HD12	1.90	0.53
1:C:517:ARG:NH1	1:D:165:LYS:HA	2.24	0.53
1:A:241:LEU:HD23	1:A:332:ILE:HD13	1.90	0.53
1:B:264:GLN:HA	1:B:267:SER:OG	2.08	0.53
1:B:388:GLU:O	1:B:392:GLU:HG3	2.09	0.53
1:B:155:LEU:HD13	1:B:179:VAL:HG21	1.89	0.53
1:C:76:HIS:CD2	1:C:78:ALA:H	2.27	0.53
1:C:89:GLN:HE22	1:C:504:ALA:N	2.07	0.53
1:B:449:LYS:HG3	1:B:459:THR:CG2	2.39	0.53
1:C:241:LEU:HD23	1:C:332:ILE:HD13	1.89	0.53
1:D:372:VAL:HG13	1:D:373:THR:N	2.23	0.53
1:D:91:LYS:HB2	1:D:91:LYS:HZ2	1.74	0.53
1:A:483:GLU:CG	1:A:485:LYS:HE2	2.39	0.52
1:C:282:HIS:HE1	1:C:339:LEU:O	1.91	0.52
1:D:280:VAL:HG13	1:D:305:LEU:HD13	1.91	0.52
1:A:201:LYS:HB2	1:A:323:LYS:HE2	1.92	0.52
1:C:517:ARG:NH1	1:D:166:ASN:H	2.07	0.52
1:B:216:VAL:O	1:B:372:VAL:HG22	2.09	0.52
1:B:380:THR:HG22	1:B:383:VAL:H	1.74	0.52
1:D:106:GLU:HB2	1:D:446:ILE:HG12	1.91	0.52
1:D:496:GLU:HG3	4:D:4061:HOH:O	2.09	0.52
1:D:82:MET:HA	1:D:82:MET:CE	2.39	0.52
1:A:82:MET:HE2	1:A:82:MET:HA	1.90	0.52
1:B:409:PRO:HB3	1:B:490:LEU:CD1	2.35	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:THR:HG22	1:A:383:VAL:CG2	2.40	0.52
1:B:380:THR:HG22	1:B:382:HIS:N	2.24	0.52
1:C:326:LYS:HB3	1:C:370:LYS:HB2	1.90	0.52
1:B:144:ILE:HD11	1:B:490:LEU:HD21	1.92	0.52
1:C:146:VAL:HG22	1:C:147:ASP:N	2.24	0.52
1:D:155:LEU:CD1	1:D:179:VAL:HG21	2.40	0.52
1:D:215:LEU:HD11	1:D:372:VAL:CG1	2.40	0.52
1:B:265:LEU:O	1:B:269:LEU:HD22	2.09	0.52
1:C:409:PRO:HB3	1:C:490:LEU:CD1	2.38	0.52
1:A:445:LYS:O	1:A:448:PRO:HD2	2.10	0.51
1:C:117:GLN:O	1:C:118:ASN:HB2	2.08	0.51
1:C:195:VAL:HB	1:C:399:LYS:HG3	1.91	0.51
1:D:388:GLU:O	1:D:392:GLU:HG3	2.10	0.51
1:A:239:ILE:HG22	1:A:241:LEU:HD13	1.92	0.51
1:A:380:THR:CG2	1:A:383:VAL:H	2.22	0.51
1:A:235:GLU:O	1:A:348:GLU:O	2.28	0.51
1:B:188:LYS:NZ	1:B:191:GLY:H	2.07	0.51
1:D:242:ILE:O	1:D:293:VAL:HA	2.11	0.51
1:D:41:THR:O	1:D:41:THR:HG22	2.09	0.51
1:B:82:MET:HA	1:B:82:MET:CE	2.41	0.51
1:A:227:HIS:HB3	1:A:230:MET:HG3	1.91	0.51
1:B:146:VAL:HG21	1:B:153:THR:HG21	1.91	0.51
1:D:154:LEU:HG	1:D:398:VAL:HG13	1.93	0.51
1:B:68:ILE:O	1:B:72:ILE:HG13	2.11	0.51
1:C:55:SER:OG	1:C:56:LEU:HD12	2.11	0.51
1:B:106:GLU:CB	1:B:446:ILE:HG12	2.41	0.51
1:D:204:LYS:HD2	1:D:384:ILE:HG23	1.92	0.51
1:A:258:ASN:HD22	1:A:258:ASN:N	2.08	0.50
1:A:144:ILE:HD11	1:A:490:LEU:HD21	1.93	0.50
1:D:272:GLU:HA	1:D:275:MET:CE	2.41	0.50
1:A:259:ILE:HD13	1:A:265:LEU:HD23	1.93	0.50
1:B:182:VAL:CG2	1:B:395:VAL:HG13	2.42	0.50
1:C:297:ILE:HG22	1:C:302:GLN:HG3	1.93	0.50
1:C:517:ARG:HD2	1:D:166:ASN:HB3	1.93	0.50
1:B:282:HIS:HE1	1:B:339:LEU:O	1.94	0.50
1:C:301:ALA:O	1:C:305:LEU:HD22	2.12	0.50
1:C:41:THR:HG22	1:C:41:THR:O	2.12	0.50
1:C:517:ARG:HH12	1:D:165:LYS:HA	1.76	0.50
1:D:204:LYS:HD2	1:D:384:ILE:CG2	2.41	0.50
1:A:297:ILE:HG22	1:A:302:GLN:HG3	1.92	0.50
1:D:334:THR:HG22	1:D:335:ASN:ND2	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:ILE:HD13	1:B:495:ILE:HG13	1.92	0.50
1:C:150:ASP:OD1	1:C:153:THR:HG22	2.12	0.50
1:D:146:VAL:HG22	1:D:147:ASP:H	1.77	0.50
1:B:117:GLN:O	1:B:118:ASN:HB2	2.11	0.50
1:C:150:ASP:HB3	1:C:153:THR:CG2	2.41	0.50
1:B:413:ALA:HB3	1:B:414:PRO:HD3	1.94	0.50
1:C:163:THR:HG22	1:C:171:LYS:HD3	1.94	0.50
1:C:378:GLY:N	1:C:384:ILE:HD11	2.27	0.50
1:C:410:ALA:O	1:C:489:MET:HG3	2.12	0.50
1:C:56:LEU:HD12	1:C:56:LEU:H	1.75	0.50
1:C:77:PRO:HB2	1:D:51:MET:HE1	1.94	0.50
1:C:480:ASP:OD1	1:C:483:GLU:HB2	2.11	0.49
1:A:218:GLY:HA3	1:A:363:VAL:O	2.12	0.49
1:B:190:ASP:C	1:B:192:LYS:H	2.16	0.49
1:B:76:HIS:HD2	1:B:78:ALA:H	1.58	0.49
1:A:466:VAL:HG21	1:A:479:ILE:CG1	2.39	0.49
1:C:313:VAL:HG21	1:C:361:ILE:HD12	1.93	0.49
1:C:13:PRO:HD3	1:D:73:ASP:O	2.12	0.49
1:B:488:ASP:O	1:B:492:LYS:HG2	2.12	0.49
1:A:35:ILE:HD11	1:A:74:LEU:HD22	1.94	0.49
1:B:9:VAL:O	1:B:11:ILE:HG12	2.12	0.49
1:B:242:ILE:O	1:B:293:VAL:HA	2.12	0.49
1:D:415:GLU:H	1:D:415:GLU:CD	2.16	0.49
1:B:259:ILE:HD13	1:B:265:LEU:HD23	1.93	0.49
1:D:188:LYS:HE3	1:D:191:GLY:HA2	1.94	0.49
1:D:119:ILE:HD11	1:D:435:LEU:HD23	1.95	0.49
1:A:206:ALA:HA	1:A:384:ILE:CD1	2.41	0.49
1:A:73:ASP:O	1:B:13:PRO:HD3	2.12	0.49
1:A:272:GLU:HA	1:A:275:MET:HE3	1.95	0.49
1:D:466:VAL:HG22	1:D:486:PRO:HG3	1.94	0.49
1:A:224:GLU:HG2	1:A:359:ASN:HB2	1.95	0.48
1:B:197:LEU:HD22	1:B:395:VAL:HG12	1.95	0.48
1:D:180:GLU:CD	1:D:215:LEU:HD23	2.33	0.48
1:D:197:LEU:HD22	1:D:395:VAL:CG1	2.42	0.48
1:A:330:ALA:HB2	1:A:345:GLY:HA3	1.94	0.48
1:D:506:LYS:O	1:D:510:GLU:HG3	2.14	0.48
1:C:259:ILE:CD1	1:C:265:LEU:HD23	2.43	0.48
1:C:330:ALA:HB2	1:C:345:GLY:HA3	1.95	0.48
1:D:224:GLU:HG2	1:D:359:ASN:HB2	1.96	0.48
1:B:412:GLY:O	1:B:415:GLU:HG2	2.13	0.48
1:C:215:LEU:CD1	1:C:372:VAL:HG13	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:388:GLU:O	1:C:392:GLU:HG3	2.13	0.48
1:D:218:GLY:HA3	1:D:363:VAL:O	2.13	0.48
1:D:119:ILE:CD1	1:D:435:LEU:HD23	2.43	0.48
1:D:157:ILE:HG13	1:D:401:VAL:HG21	1.96	0.48
1:A:9:VAL:O	1:A:9:VAL:HG13	2.12	0.48
1:C:517:ARG:CD	1:D:166:ASN:HB3	2.43	0.48
1:A:410:ALA:HB2	1:A:496:GLU:HG2	1.96	0.48
1:B:449:LYS:HG3	1:B:459:THR:HG23	1.96	0.48
1:C:76:HIS:HD2	1:C:78:ALA:H	1.60	0.48
1:C:9:VAL:O	1:C:9:VAL:HG13	2.14	0.48
1:A:272:GLU:HA	1:A:275:MET:CE	2.44	0.48
1:B:201:LYS:HB2	1:B:323:LYS:CE	2.43	0.48
1:C:106:GLU:HB2	1:C:446:ILE:HG12	1.95	0.48
1:A:62:THR:HG23	1:A:389:ARG:NE	2.28	0.48
1:C:32:ALA:HB1	1:C:82:MET:HE2	1.95	0.48
1:D:266:MET:CE	1:D:269:LEU:HD23	2.42	0.48
1:D:166:ASN:O	1:D:167:ALA:CB	2.60	0.47
1:A:413:ALA:HB3	1:A:414:PRO:HD3	1.96	0.47
1:D:103:ILE:HG22	1:D:107:LEU:HD22	1.95	0.47
1:C:30:LEU:HD22	1:C:34:ILE:CD1	2.44	0.47
1:A:157:ILE:HG13	1:A:401:VAL:HG21	1.95	0.47
1:A:380:THR:HG22	1:A:383:VAL:CB	2.44	0.47
1:B:224:GLU:HG2	1:B:359:ASN:CB	2.45	0.47
1:B:49:ASP:OD1	1:B:63:ASN:HB2	2.13	0.47
1:C:192:LYS:HD2	1:C:192:LYS:N	2.29	0.47
1:C:517:ARG:HH11	1:D:166:ASN:H	1.62	0.47
1:D:121:PRO:O	1:D:125:THR:HB	2.15	0.47
1:D:264:GLN:HA	1:D:267:SER:OG	2.14	0.47
1:B:89:GLN:HE22	1:B:504:ALA:N	2.13	0.47
1:D:432:LYS:O	1:D:435:LEU:HB2	2.15	0.47
1:A:76:HIS:CD2	1:A:78:ALA:H	2.33	0.47
1:B:95:ASP:OD1	1:B:161:SER:CB	2.63	0.47
1:B:271:GLN:HG3	1:B:275:MET:CE	2.45	0.47
1:B:303:HIS:CE1	1:B:307:LYS:HD2	2.50	0.47
1:C:190:ASP:C	1:C:192:LYS:H	2.18	0.47
1:C:148:PRO:O	1:C:402:MET:HG2	2.15	0.47
1:C:517:ARG:NH1	1:D:166:ASN:N	2.63	0.47
1:D:380:THR:HG23	1:D:382:HIS:H	1.80	0.47
1:A:380:THR:HG22	1:A:383:VAL:HG23	1.96	0.47
1:A:62:THR:CG2	1:A:389:ARG:NE	2.78	0.47
1:C:264:GLN:HA	1:C:267:SER:OG	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:204:LYS:HB3	1:D:384:ILE:CG2	2.45	0.47
1:C:350:VAL:HA	1:C:362:PHE:O	2.15	0.47
1:D:190:ASP:C	1:D:192:LYS:N	2.68	0.47
1:D:215:LEU:HD11	1:D:372:VAL:HG11	1.96	0.47
1:A:273:GLU:OE2	1:B:337:LYS:HE2	2.14	0.46
1:A:506:LYS:O	1:A:510:GLU:HG3	2.15	0.46
1:D:301:ALA:O	1:D:305:LEU:HD22	2.15	0.46
1:D:353:ARG:HH22	1:D:364:GLU:CD	2.17	0.46
1:A:89:GLN:HE22	1:A:504:ALA:N	2.13	0.46
1:C:506:LYS:O	1:C:510:GLU:HG3	2.16	0.46
1:D:151:GLU:HG2	1:D:155:LEU:HD22	1.98	0.46
1:B:188:LYS:HZ3	1:B:191:GLY:H	1.62	0.46
1:B:9:VAL:HG13	1:B:9:VAL:O	2.14	0.46
1:A:498:LEU:HD22	1:A:502:LYS:HG3	1.98	0.46
1:A:75:GLN:HE21	1:B:13:PRO:HG3	1.77	0.46
1:C:221:ILE:HD11	1:C:324:LEU:HD11	1.97	0.46
1:C:266:MET:HE2	1:C:269:LEU:HD23	1.98	0.46
1:D:117:GLN:O	1:D:118:ASN:HB2	2.15	0.46
1:B:206:ALA:HA	1:B:384:ILE:CD1	2.46	0.46
1:D:227:HIS:CD2	1:D:229:ARG:H	2.33	0.46
1:D:272:GLU:HG2	1:D:275:MET:HE3	1.96	0.46
1:D:326:LYS:HB3	1:D:370:LYS:HB2	1.98	0.46
1:B:239:ILE:HD11	1:B:347:ALA:CB	2.46	0.46
1:B:221:ILE:HD11	1:B:324:LEU:HD11	1.98	0.46
1:A:204:LYS:HD3	1:A:388:GLU:OE2	2.15	0.46
1:B:106:GLU:HB2	1:B:446:ILE:HG12	1.96	0.46
1:C:227:HIS:CD2	1:C:229:ARG:HB2	2.51	0.46
1:C:488:ASP:OD1	1:C:490:LEU:HB2	2.15	0.46
1:D:227:HIS:HD2	1:D:229:ARG:H	1.63	0.46
1:C:358:GLU:HB3	4:C:3188:HOH:O	2.15	0.45
1:C:35:ILE:HD13	1:C:35:ILE:HA	1.85	0.45
1:C:45:PRO:HB2	1:C:481:VAL:HG21	1.97	0.45
1:C:65:CYS:HB3	1:C:97:THR:OG1	2.16	0.45
1:C:303:HIS:O	1:C:307:LYS:HG3	2.16	0.45
1:D:466:VAL:CG2	1:D:479:ILE:HG12	2.35	0.45
1:D:76:HIS:HD2	1:D:78:ALA:N	2.13	0.45
1:B:215:LEU:HD23	1:B:215:LEU:C	2.37	0.45
1:B:311:MET:HE1	1:B:350:VAL:HG11	1.99	0.45
1:D:372:VAL:CG1	1:D:373:THR:N	2.79	0.45
1:A:51:MET:CE	1:B:77:PRO:HB2	2.46	0.45
1:B:76:HIS:CD2	1:B:78:ALA:H	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:MET:HE1	1:A:269:LEU:HD23	1.97	0.45
1:C:380:THR:O	1:C:384:ILE:HG12	2.15	0.45
1:D:54:ASP:OD2	1:D:58:ASP:HB2	2.16	0.45
1:C:189:LYS:HG3	1:C:190:ASP:OD2	2.16	0.45
1:B:235:GLU:O	1:B:348:GLU:O	2.35	0.45
1:C:318:LYS:O	1:C:322:GLU:HG3	2.16	0.45
1:D:98:THR:O	1:D:102:VAL:HG23	2.17	0.45
1:D:106:GLU:CB	1:D:446:ILE:HG12	2.47	0.45
1:D:243:ASN:OD1	1:D:334:THR:HA	2.17	0.45
1:C:249:LYS:HB2	1:D:269:LEU:HD21	1.99	0.45
1:A:215:LEU:HD11	1:A:372:VAL:CG1	2.47	0.45
1:A:378:GLY:N	1:A:384:ILE:HD11	2.33	0.45
1:D:290:VAL:HG13	1:D:311:MET:HE1	1.99	0.45
1:B:240:ALA:HB3	1:B:291:VAL:HG22	1.98	0.44
1:B:197:LEU:HD22	1:B:395:VAL:CG1	2.47	0.44
1:D:217:ARG:NH2	4:D:4002:HOH:O	2.45	0.44
1:A:243:ASN:O	1:A:295:LYS:HG2	2.17	0.44
1:B:180:GLU:HB3	1:B:215:LEU:HD12	1.99	0.44
1:D:165:LYS:O	1:D:166:ASN:C	2.53	0.44
1:A:159:ALA:O	1:A:163:THR:HG23	2.17	0.44
1:C:396:LYS:HE3	4:C:3119:HOH:O	2.17	0.44
1:C:520:ASP:HB3	1:D:50:LYS:HG2	1.98	0.44
1:B:204:LYS:HD2	1:B:384:ILE:HG22	1.98	0.44
1:C:154:LEU:HG	1:C:398:VAL:HG13	1.99	0.44
1:D:330:ALA:HB2	1:D:345:GLY:HA3	1.99	0.44
1:A:244:GLU:CB	1:A:334:THR:O	2.65	0.44
1:A:256:LYS:HE2	1:B:256:LYS:HB2	1.99	0.44
1:B:386:GLU:OE2	1:B:386:GLU:HA	2.17	0.44
1:C:235:GLU:O	1:C:348:GLU:O	2.36	0.44
1:C:459:THR:HG22	1:C:460:VAL:N	2.32	0.44
1:B:14:GLU:HG3	1:B:14:GLU:O	2.18	0.44
1:C:50:LYS:HD2	1:C:68:ILE:HD13	2.00	0.44
1:A:227:HIS:HA	1:A:228:PRO:HD3	1.86	0.44
1:A:52:LEU:HD11	1:A:68:ILE:HA	2.00	0.44
1:B:197:LEU:CD2	1:B:395:VAL:HG12	2.47	0.44
1:D:170:HIS:HD2	1:D:211:GLU:HG2	1.80	0.44
1:A:67:THR:O	1:A:71:LYS:HG2	2.18	0.44
1:C:35:ILE:HD11	1:C:74:LEU:HD22	1.99	0.44
1:B:488:ASP:OD1	1:B:490:LEU:HB2	2.18	0.43
1:A:9:VAL:O	1:A:11:ILE:HG12	2.18	0.43
1:C:383:VAL:O	1:C:387:VAL:HG23	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:480:ASP:HB3	1:D:485:LYS:O	2.18	0.43
1:D:49:ASP:HB3	1:D:62:THR:O	2.18	0.43
1:A:326:LYS:HB3	1:A:370:LYS:HB2	1.99	0.43
1:B:416:ILE:HD13	1:B:466:VAL:HG12	2.00	0.43
1:B:506:LYS:O	1:B:510:GLU:HG3	2.19	0.43
1:D:108:LEU:CD2	1:D:516:LEU:HD13	2.48	0.43
1:B:334:THR:HG22	1:B:335:ASN:ND2	2.34	0.43
1:C:243:ASN:O	1:C:295:LYS:HG2	2.19	0.43
1:D:501:LYS:HE3	1:D:501:LYS:HA	2.00	0.43
1:A:37:GLU:O	1:A:40:ARG:HG2	2.18	0.43
1:D:190:ASP:OD2	1:D:190:ASP:O	2.37	0.43
1:D:227:HIS:HB3	1:D:230:MET:HG3	2.01	0.43
1:D:204:LYS:HD3	1:D:388:GLU:OE2	2.18	0.43
1:A:266:MET:O	1:A:270:GLU:HG3	2.19	0.43
1:C:269:LEU:HD12	1:C:269:LEU:HA	1.92	0.43
1:B:370:LYS:HE3	4:B:2087:HOH:O	2.17	0.43
1:C:123:ILE:HG21	1:C:432:LYS:HB2	2.01	0.43
1:A:356:ALA:O	1:A:358:GLU:HG2	2.19	0.43
1:A:449:LYS:HG3	1:A:459:THR:OG1	2.18	0.43
1:A:157:ILE:HD13	1:A:495:ILE:HG13	2.00	0.43
1:C:413:ALA:HB3	1:C:414:PRO:HD3	1.99	0.43
1:D:450:THR:HG22	4:D:4013:HOH:O	2.18	0.43
1:A:103:ILE:HG12	1:A:446:ILE:HD11	2.00	0.42
1:B:121:PRO:O	1:B:125:THR:HB	2.19	0.42
1:B:170:HIS:HD2	1:B:211:GLU:HG2	1.78	0.42
1:C:49:ASP:OD1	1:C:63:ASN:HB2	2.18	0.42
1:D:334:THR:HG22	1:D:335:ASN:HD22	1.83	0.42
1:A:98:THR:O	1:A:102:VAL:HG23	2.19	0.42
1:C:106:GLU:CB	1:C:446:ILE:HG12	2.49	0.42
1:C:195:VAL:O	1:C:399:LYS:HE2	2.18	0.42
1:D:272:GLU:HA	1:D:275:MET:HE3	2.00	0.42
1:D:88:THR:HA	1:D:91:LYS:NZ	2.34	0.42
1:A:215:LEU:HD11	1:A:372:VAL:HG13	2.00	0.42
1:B:311:MET:CE	1:B:350:VAL:HG11	2.49	0.42
1:C:377:ARG:C	1:C:384:ILE:HD11	2.39	0.42
1:D:244:GLU:CB	1:D:334:THR:O	2.66	0.42
1:B:259:ILE:CD1	1:B:265:LEU:HD23	2.49	0.42
1:B:303:HIS:HD2	4:B:2090:HOH:O	2.02	0.42
1:C:184:GLN:NE2	1:C:184:GLN:HA	2.34	0.42
1:D:190:ASP:O	1:D:192:LYS:N	2.52	0.42
1:A:33:ARG:O	1:A:37:GLU:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:235:GLU:OE2	1:C:235:GLU:HA	2.20	0.42
1:C:32:ALA:HB1	1:C:82:MET:HE1	2.00	0.42
1:D:282:HIS:HE1	1:D:339:LEU:O	2.01	0.42
1:A:257:ILE:HB	1:B:255:ALA:CB	2.49	0.42
1:B:313:VAL:HG21	1:B:361:ILE:HD13	2.02	0.42
1:D:197:LEU:HD22	1:D:395:VAL:HG13	1.99	0.42
1:D:52:LEU:HD11	1:D:68:ILE:HA	2.02	0.42
1:A:205:LYS:HD3	1:A:205:LYS:HA	1.83	0.42
1:D:217:ARG:HA	1:D:372:VAL:HG22	2.02	0.42
1:D:88:THR:HA	1:D:91:LYS:HZ2	1.84	0.42
1:A:76:HIS:HD2	1:A:78:ALA:H	1.66	0.42
1:B:146:VAL:HG22	1:B:147:ASP:H	1.85	0.42
1:B:35:ILE:HA	1:B:35:ILE:HD13	1.89	0.42
1:C:98:THR:O	1:C:102:VAL:HG23	2.20	0.42
1:A:62:THR:CG2	1:A:386:GLU:OE2	2.68	0.42
1:A:166:ASN:CB	1:B:517:ARG:HD2	2.35	0.42
1:C:266:MET:HE2	1:C:266:MET:HA	2.02	0.42
1:A:227:HIS:HD2	1:A:229:ARG:HB2	1.84	0.41
1:B:378:GLY:N	1:B:384:ILE:HD11	2.35	0.41
1:B:470:HIS:HA	1:B:477:ILE:HB	2.02	0.41
1:A:110:LYS:HD2	1:A:442:ASP:OD1	2.20	0.41
1:B:195:VAL:HB	1:B:399:LYS:HG3	2.02	0.41
1:C:165:LYS:HE2	4:C:3120:HOH:O	2.18	0.41
1:C:76:HIS:HA	1:C:77:PRO:HD3	1.93	0.41
1:A:330:ALA:HB2	1:A:345:GLY:CA	2.51	0.41
1:B:276:LEU:HD13	1:B:300:LEU:HB3	2.03	0.41
1:A:490:LEU:HA	1:A:490:LEU:HD12	1.93	0.41
1:C:276:LEU:HA	1:C:276:LEU:HD12	1.95	0.41
1:C:372:VAL:HG13	1:C:373:THR:N	2.35	0.41
1:C:76:HIS:CD2	1:C:77:PRO:HD2	2.55	0.41
1:A:49:ASP:OD1	1:A:63:ASN:HB2	2.21	0.41
1:C:445:LYS:C	1:C:448:PRO:HD2	2.40	0.41
1:B:218:GLY:HA3	1:B:363:VAL:O	2.21	0.41
1:A:257:ILE:HB	1:B:255:ALA:HB2	2.01	0.41
1:B:283:ILE:HG21	1:B:291:VAL:HG21	2.02	0.41
1:B:352:GLU:HA	1:B:360:MET:O	2.20	0.41
1:B:516:LEU:HD12	1:B:516:LEU:HA	1.88	0.41
1:C:165:LYS:O	1:C:168:GLU:HG2	2.20	0.41
1:C:215:LEU:HD11	1:C:372:VAL:HG11	2.01	0.41
1:D:148:PRO:O	1:D:402:MET:HG2	2.21	0.41
1:D:459:THR:HG22	1:D:460:VAL:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:397:VAL:HB	1:D:497:PRO:HG2	2.03	0.41
1:A:240:ALA:HB3	1:A:291:VAL:HG22	2.03	0.41
1:B:271:GLN:O	1:B:275:MET:HE2	2.20	0.41
1:C:151:GLU:HG2	1:C:155:LEU:HD22	2.02	0.41
1:B:163:THR:HB	1:B:171:LYS:HZ1	1.80	0.41
1:B:290:VAL:HG13	1:B:311:MET:CE	2.51	0.41
1:B:67:THR:O	1:B:71:LYS:HG2	2.20	0.41
1:C:163:THR:HG22	1:C:171:LYS:NZ	2.36	0.41
1:C:29:ILE:HD13	1:C:112:GLU:HB2	2.02	0.41
1:D:171:LYS:HG3	1:D:172:GLU:N	2.36	0.41
1:A:146:VAL:CG2	1:A:147:ASP:N	2.84	0.41
1:A:266:MET:CE	1:A:269:LEU:HD23	2.51	0.41
1:A:271:GLN:HG3	1:A:275:MET:CE	2.50	0.41
1:A:470:HIS:HA	1:A:477:ILE:HB	2.03	0.41
1:B:217:ARG:HA	1:B:372:VAL:HG23	2.03	0.41
1:C:127:GLY:HA3	1:C:436:ALA:HB3	2.03	0.41
1:D:204:LYS:HB3	1:D:384:ILE:HG23	2.03	0.41
1:A:18:ARG:HA	1:A:521:VAL:O	2.20	0.40
1:B:239:ILE:HD11	1:B:347:ALA:HB2	2.03	0.40
1:B:396:LYS:HG3	4:B:2083:HOH:O	2.19	0.40
1:C:146:VAL:CG2	1:C:153:THR:HG21	2.37	0.40
1:D:384:ILE:HG22	1:D:385:ASP:N	2.36	0.40
1:D:76:HIS:CD2	1:D:78:ALA:HB3	2.56	0.40
1:A:404:ASP:O	1:A:406:ALA:N	2.50	0.40
1:A:106:GLU:HB2	1:A:446:ILE:HG12	2.03	0.40
1:A:473:ARG:HB2	1:A:477:ILE:HG13	2.03	0.40
1:A:154:LEU:HG	1:A:398:VAL:HG13	2.04	0.40
1:A:446:ILE:HD12	1:A:447:ILE:N	2.37	0.40
1:D:311:MET:HE1	1:D:350:VAL:HG11	2.02	0.40
1:A:271:GLN:HG3	1:A:275:MET:HE2	2.02	0.40
1:A:463:LEU:HD13	1:A:463:LEU:HA	1.90	0.40
1:B:110:LYS:HA	1:B:110:LYS:HE2	2.02	0.40
1:C:296:GLY:HA2	1:C:315:ARG:HD2	2.02	0.40
1:D:227:HIS:HA	1:D:228:PRO:HD3	1.86	0.40
1:D:367:LYS:NZ	1:D:367:LYS:HB3	2.37	0.40
1:D:446:ILE:O	1:D:450:THR:CG2	2.53	0.40

All (7) 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)
4:A:1162:HOH:O	4:B:2002:HOH:O[4_555]	0.56	1.64
4:C:3159:HOH:O	4:D:4143:HOH:O[4_555]	0.71	1.49
4:C:3137:HOH:O	4:D:4115:HOH:O[4_555]	0.74	1.46
1:C:272:GLU:OE2	1:D:253:THR:OG1[4_555]	2.01	0.19
4:C:3148:HOH:O	4:D:4139:HOH:O[4_555]	2.04	0.16
1:C:380:THR:OG1	1:D:84:GLU:OE1[4_555]	2.14	0.06
4:B:2067:HOH:O	4:D:4112:HOH:O[2_656]	2.15	0.05

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	516/548 (94%)	504 (98%)	10 (2%)	2 (0%)	38	47
1	B	516/548 (94%)	503 (98%)	10 (2%)	3 (1%)	28	34
1	C	516/548 (94%)	500 (97%)	12 (2%)	4 (1%)	22	26
1	D	516/548 (94%)	502 (97%)	11 (2%)	3 (1%)	28	34
All	All	2064/2192 (94%)	2009 (97%)	43 (2%)	12 (1%)	28	34

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	235	GLU
1	B	10	VAL
1	B	235	GLU
1	C	235	GLU
1	D	235	GLU
1	D	367	LYS
1	A	367	LYS
1	B	367	LYS
1	C	190	ASP
1	C	367	LYS
1	D	190	ASP
1	C	10	VAL

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	423/441 (96%)	398 (94%)	25 (6%)	23	30
1	B	423/441 (96%)	399 (94%)	24 (6%)	24	32
1	C	423/441 (96%)	400 (95%)	23 (5%)	26	35
1	D	423/441 (96%)	394 (93%)	29 (7%)	18	23
All	All	1692/1764 (96%)	1591 (94%)	101 (6%)	22	30

All (101) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLU
1	A	30	LEU
1	A	62	THR
1	A	82	MET
1	A	107	LEU
1	A	154	LEU
1	A	155	LEU
1	A	171	LYS
1	A	229	ARG
1	A	238	LYS
1	A	258	ASN
1	A	269	LEU
1	A	305	LEU
1	A	358	GLU
1	A	372	VAL
1	A	375	LEU
1	A	377	ARG
1	A	446	ILE
1	A	459	THR
1	A	463	LEU
1	A	490	LEU
1	A	496	GLU
1	A	498	LEU
1	A	501	LYS

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Mol	Chain	Res	Type
1	A	516	LEU
1	B	14	GLU
1	B	30	LEU
1	B	82	MET
1	B	125	THR
1	B	147	ASP
1	B	154	LEU
1	B	155	LEU
1	B	169	SER
1	B	211	GLU
1	B	229	ARG
1	B	241	LEU
1	B	269	LEU
1	B	305	LEU
1	B	368	ASN
1	B	375	LEU
1	B	423	ASP
1	B	435	LEU
1	B	446	ILE
1	B	463	LEU
1	B	483	GLU
1	B	490	LEU
1	B	498	LEU
1	B	501	LYS
1	B	516	LEU
1	C	14	GLU
1	C	30	LEU
1	C	82	MET
1	C	107	LEU
1	C	153	THR
1	C	154	LEU
1	C	155	LEU
1	C	241	LEU
1	C	269	LEU
1	C	305	LEU
1	C	315	ARG
1	C	367	LYS
1	C	375	LEU
1	C	377	ARG
1	C	435	LEU
1	C	446	ILE
1	C	450	THR

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Mol	Chain	Res	Type
1	C	459	THR
1	C	463	LEU
1	C	483	GLU
1	C	490	LEU
1	C	498	LEU
1	C	516	LEU
1	D	14	GLU
1	D	17	GLN
1	D	30	LEU
1	D	58	ASP
1	D	82	MET
1	D	107	LEU
1	D	125	THR
1	D	147	ASP
1	D	152	GLU
1	D	154	LEU
1	D	155	LEU
1	D	163	THR
1	D	169	SER
1	D	211	GLU
1	D	229	ARG
1	D	241	LEU
1	D	269	LEU
1	D	305	LEU
1	D	367	LYS
1	D	375	LEU
1	D	377	ARG
1	D	380	THR
1	D	435	LEU
1	D	446	ILE
1	D	450	THR
1	D	490	LEU
1	D	498	LEU
1	D	501	LYS
1	D	516	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	75	GLN
1	A	76	HIS

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Mol	Chain	Res	Type
1	A	89	GLN
1	A	117	GLN
1	A	118	ASN
1	A	184	GLN
1	A	227	HIS
1	A	258	ASN
1	A	282	HIS
1	A	285	GLN
1	A	368	ASN
1	B	28	ASN
1	B	76	HIS
1	B	89	GLN
1	B	117	GLN
1	B	118	ASN
1	B	170	HIS
1	B	184	GLN
1	B	258	ASN
1	B	282	HIS
1	B	285	GLN
1	B	303	HIS
1	B	335	ASN
1	B	368	ASN
1	B	503	GLN
1	C	28	ASN
1	C	76	HIS
1	C	89	GLN
1	C	117	GLN
1	C	118	ASN
1	C	170	HIS
1	C	184	GLN
1	C	227	HIS
1	C	282	HIS
1	C	285	GLN
1	C	303	HIS
1	C	335	ASN
1	C	368	ASN
1	C	428	GLN
1	C	503	GLN
1	D	17	GLN
1	D	28	ASN
1	D	75	GLN
1	D	76	HIS

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Mol	Chain	Res	Type
1	D	89	GLN
1	D	117	GLN
1	D	118	ASN
1	D	170	HIS
1	D	184	GLN
1	D	227	HIS
1	D	282	HIS
1	D	285	GLN
1	D	303	HIS
1	D	335	ASN
1	D	368	ASN

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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	ANP	A	1528	2	29,33,33	1.65	9 (31%)	28,52,52	1.41	2 (7%)
3	ANP	B	2528	2	29,33,33	1.56	8 (27%)	28,52,52	1.49	2 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ANP	C	3528	2	29,33,33	1.60	8 (27%)	28,52,52	1.44	3 (10%)
3	ANP	D	4528	2	29,33,33	1.60	9 (31%)	28,52,52	1.46	2 (7%)

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	ANP	A	1528	2	-	0/13/38/38	0/3/3/3
3	ANP	B	2528	2	-	1/13/38/38	0/3/3/3
3	ANP	C	3528	2	-	0/13/38/38	0/3/3/3
3	ANP	D	4528	2	-	0/13/38/38	0/3/3/3

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1528	ANP	PB-O2B	-3.29	1.47	1.56
3	D	4528	ANP	PB-O2B	-3.26	1.47	1.56
3	B	2528	ANP	PB-O2B	-3.24	1.47	1.56
3	C	3528	ANP	PB-O2B	-3.23	1.47	1.56
3	C	3528	ANP	PG-O2G	-2.89	1.48	1.56
3	D	4528	ANP	PG-O2G	-2.82	1.49	1.56
3	B	2528	ANP	PG-O2G	-2.77	1.49	1.56
3	A	1528	ANP	PG-O2G	-2.72	1.49	1.56
3	A	1528	ANP	PG-O3G	-2.66	1.49	1.56
3	C	3528	ANP	PG-O3G	-2.61	1.49	1.56
3	B	2528	ANP	PG-O3G	-2.40	1.50	1.56
3	D	4528	ANP	PG-O3G	-2.36	1.50	1.56
3	D	4528	ANP	PB-O1B	2.03	1.48	1.46
3	A	1528	ANP	PB-O1B	2.09	1.48	1.46
3	B	2528	ANP	C4-N3	2.10	1.38	1.35
3	C	3528	ANP	C4-N3	2.12	1.38	1.35
3	A	1528	ANP	C4-N3	2.15	1.38	1.35
3	D	4528	ANP	C4-N3	2.19	1.38	1.35
3	A	1528	ANP	C2-N3	2.31	1.36	1.32
3	B	2528	ANP	O4'-C1'	2.39	1.44	1.41
3	D	4528	ANP	C2-N3	2.40	1.36	1.32
3	D	4528	ANP	O4'-C1'	2.42	1.44	1.41
3	B	2528	ANP	C2-N3	2.49	1.36	1.32
3	A	1528	ANP	O4'-C1'	2.51	1.44	1.41
3	C	3528	ANP	C2-N3	2.56	1.36	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	3528	ANP	O4'-C1'	2.60	1.44	1.41
3	D	4528	ANP	PG-O1G	2.88	1.49	1.46
3	C	3528	ANP	PB-O3A	2.98	1.62	1.59
3	B	2528	ANP	PG-O1G	3.01	1.49	1.46
3	B	2528	ANP	PB-O3A	3.02	1.62	1.59
3	A	1528	ANP	PG-O1G	3.04	1.49	1.46
3	C	3528	ANP	PG-O1G	3.29	1.49	1.46
3	D	4528	ANP	PB-O3A	3.47	1.63	1.59
3	A	1528	ANP	PB-O3A	3.71	1.63	1.59

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4528	ANP	O1G-PG-N3B	-2.82	107.58	111.79
3	C	3528	ANP	O1G-PG-N3B	-2.62	107.88	111.79
3	A	1528	ANP	O1B-PB-N3B	-2.19	108.51	111.79
3	C	3528	ANP	O1B-PB-N3B	-2.04	108.75	111.79
3	B	2528	ANP	O1G-PG-N3B	2.95	116.20	111.79
3	C	3528	ANP	O2B-PB-O1B	4.34	118.90	109.87
3	D	4528	ANP	O2B-PB-O1B	4.57	119.37	109.87
3	B	2528	ANP	O2B-PB-O1B	4.63	119.50	109.87
3	A	1528	ANP	O2B-PB-O1B	4.73	119.71	109.87

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	2528	ANP	O1G-PG-N3B-PB

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1528	ANP	2	0
3	B	2528	ANP	2	0
3	C	3528	ANP	2	0
3	D	4528	ANP	2	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

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	518/548 (94%)	0.37	13 (2%) 58 65	9, 23, 41, 83	0
1	B	518/548 (94%)	0.45	20 (3%) 40 47	11, 24, 40, 82	0
1	C	518/548 (94%)	0.61	30 (5%) 24 30	8, 22, 41, 77	0
1	D	518/548 (94%)	0.36	18 (3%) 44 51	11, 24, 41, 77	0
All	All	2072/2192 (94%)	0.45	81 (3%) 40 47	8, 23, 41, 83	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	526	ALA	5.4
1	A	9	VAL	5.3
1	D	9	VAL	5.1
1	A	189	LYS	4.9
1	C	526	ALA	4.7
1	B	189	LYS	4.7
1	C	55	SER	4.6
1	D	56	LEU	4.5
1	C	190	ASP	4.4
1	B	190	ASP	4.4
1	C	56	LEU	4.3
1	C	58	ASP	4.3
1	A	190	ASP	4.2
1	C	189	LYS	4.1
1	D	191	GLY	4.0
1	C	191	GLY	4.0
1	C	9	VAL	4.0
1	D	55	SER	3.8
1	D	525	LYS	3.7
1	D	192	LYS	3.7
1	B	9	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	192	LYS	3.6
1	D	190	ASP	3.5
1	A	56	LEU	3.5
1	C	308	TYR	3.5
1	D	235	GLU	3.4
1	A	525	LYS	3.3
1	B	308	TYR	3.3
1	C	14	GLU	3.2
1	C	525	LYS	3.1
1	B	14	GLU	3.0
1	C	349	VAL	3.0
1	A	55	SER	3.0
1	B	526	ALA	2.9
1	C	57	GLY	2.9
1	B	72	ILE	2.9
1	D	189	LYS	2.9
1	C	346	TYR	2.9
1	C	357	GLY	2.9
1	B	55	SER	2.8
1	A	15	GLY	2.7
1	B	483	GLU	2.7
1	A	192	LYS	2.7
1	D	367	LYS	2.7
1	B	304	TYR	2.6
1	B	56	LEU	2.6
1	C	235	GLU	2.6
1	D	58	ASP	2.6
1	A	191	GLY	2.6
1	B	424	GLU	2.6
1	B	525	LYS	2.6
1	B	58	ASP	2.5
1	D	13	PRO	2.5
1	C	231	PRO	2.5
1	A	235	GLU	2.5
1	D	14	GLU	2.4
1	C	229	ARG	2.4
1	C	355	LEU	2.4
1	C	356	ALA	2.4
1	B	235	GLU	2.4
1	A	166	ASN	2.4
1	C	367	LYS	2.4
1	B	367	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	225	VAL	2.3
1	A	17	GLN	2.3
1	D	483	GLU	2.3
1	D	526	ALA	2.3
1	D	491	GLU	2.3
1	C	72	ILE	2.3
1	C	329	GLY	2.2
1	B	472	ASN	2.2
1	C	380	THR	2.2
1	C	335	ASN	2.1
1	B	372	VAL	2.1
1	D	427	LYS	2.1
1	B	346	TYR	2.1
1	D	268	PHE	2.1
1	B	475	LEU	2.0
1	C	218	GLY	2.0
1	C	304	TYR	2.0
1	C	483	GLU	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(Å ²)	Q<0.9
3	ANP	C	3528	31/31	0.81	0.28	3.63	46,58,60,62	0
3	ANP	B	2528	31/31	0.83	0.22	2.39	50,59,65,66	0
3	ANP	D	4528	31/31	0.84	0.20	2.00	46,54,57,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ANP	A	1528	31/31	0.85	0.20	1.61	43,47,56,58	0
2	MG	D	4527	1/1	0.88	0.18	-	57,57,57,57	0
2	MG	A	1527	1/1	0.87	0.30	-	71,71,71,71	0
2	MG	C	3527	1/1	0.37	0.41	-	85,85,85,85	0
2	MG	B	2527	1/1	0.85	0.25	-	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.