



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:19 PM GMT

PDB ID : 3TKK
Title : Crystal Structure Analysis of a recombinant predicted acetamidase/ formamida-
dase from the thermophile thermoanaerobacter tengcongensis
Authors : Qian, M.; Huang, Q.; Wu, G.; Lai, L.; Tang, Y.; Pei, J.; Kusunoki, M.
Deposited on : 2011-08-26
Resolution : 1.99 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

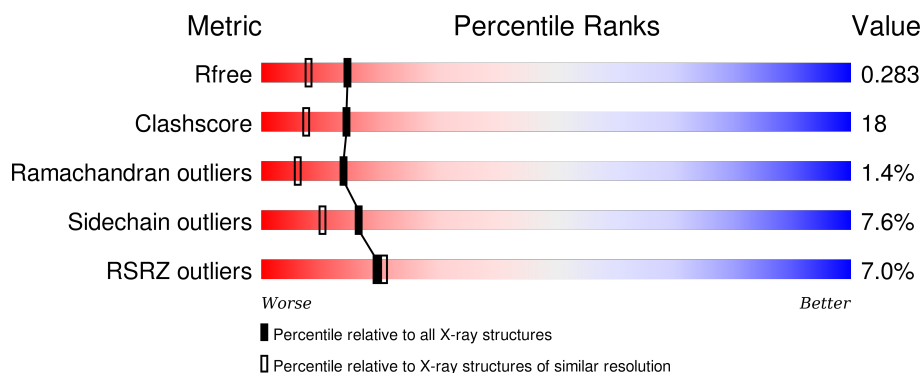
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	301	<div> <div>6%</div> <div>67%</div> <div>28%</div> <div>• •</div> </div>
1	B	301	<div> <div>6%</div> <div>73%</div> <div>23%</div> <div>•</div> </div>
1	C	301	<div> <div>6%</div> <div>72%</div> <div>24%</div> <div>5%</div> </div>
1	D	301	<div> <div>9%</div> <div>67%</div> <div>27%</div> <div>5%</div> </div>

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

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CA	A	300	-	-	-	X
2	CA	B	304	-	-	-	X
3	ZN	B	303	-	-	-	X
3	ZN	D	307	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9872 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 Predicted acetamidase/formamidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2257	1436	369	442	10			
1	B	301	Total	C	N	O	S	0	0	0
			2257	1436	369	442	10			
1	C	301	Total	C	N	O	S	0	0	0
			2257	1436	369	442	10			
1	D	301	Total	C	N	O	S	0	0	0
			2257	1436	369	442	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q8R8S5
A	-1	HIS	-	EXPRESSION TAG	UNP Q8R8S5
B	-2	SER	-	EXPRESSION TAG	UNP Q8R8S5
B	-1	HIS	-	EXPRESSION TAG	UNP Q8R8S5
C	-2	SER	-	EXPRESSION TAG	UNP Q8R8S5
C	-1	HIS	-	EXPRESSION TAG	UNP Q8R8S5
D	-2	SER	-	EXPRESSION TAG	UNP Q8R8S5
D	-1	HIS	-	EXPRESSION TAG	UNP Q8R8S5

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

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

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Zn 1	0	0
3	A	1	Total 1	Zn 1	0	0
3	D	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0

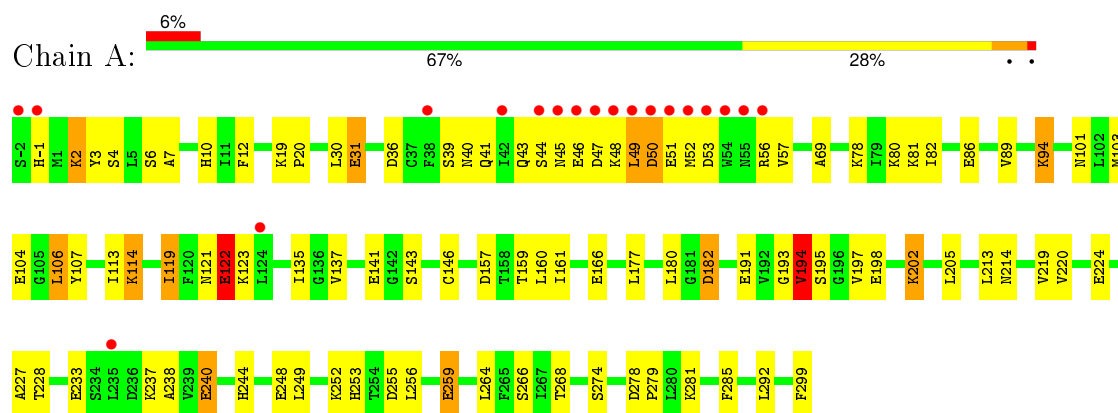
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	292	Total 292	O 292	0	0
4	B	153	Total 153	O 153	0	0
4	C	254	Total 254	O 254	0	0
4	D	136	Total 136	O 136	0	0

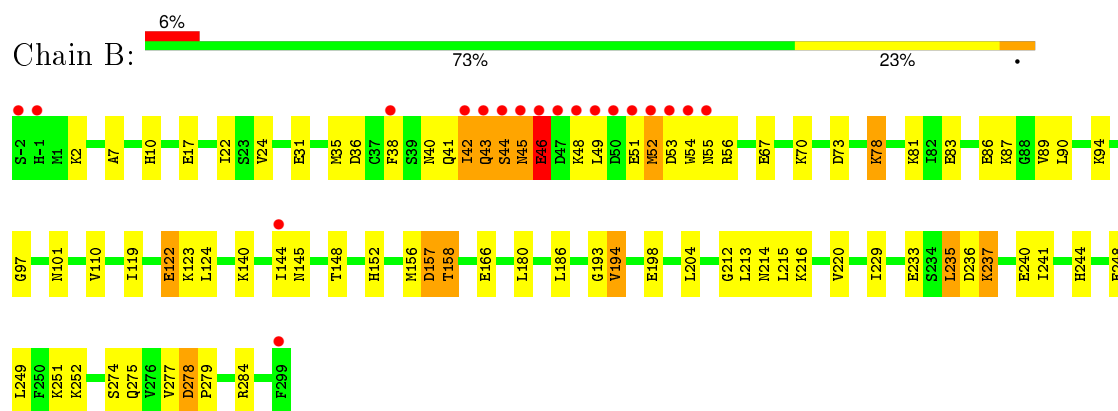
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

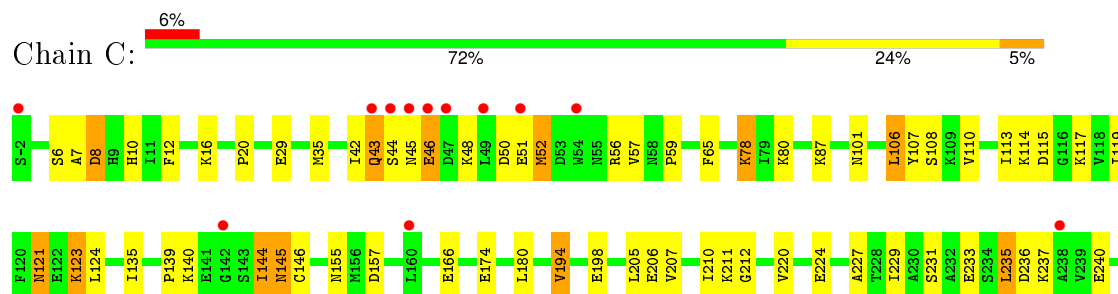
• Molecule 1: Predicted acetamidase/formamidase



• Molecule 1: Predicted acetamidase/formamidase

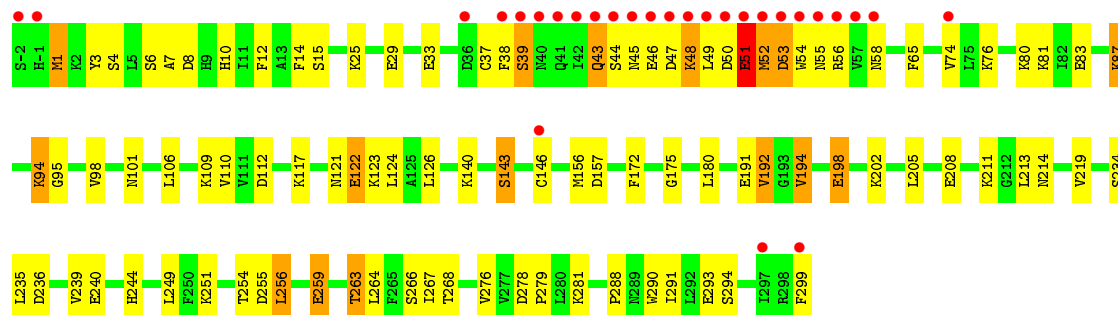


• Molecule 1: Predicted acetamidase/formamidase





- Molecule 1: Predicted acetamidase/formamidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	41.23Å 152.88Å 100.25Å 90.00° 99.49° 90.00°	Depositor
Resolution (Å)	50.00 – 1.99 40.67 – 1.99	Depositor EDS
% Data completeness (in resolution range)	94.7 (50.00-1.99) 94.7 (40.67-1.99)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.4.0077	Depositor
R, R_{free}	0.174 , 0.237 0.240 , 0.283	Depositor DCC
R_{free} test set	3985 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.692	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.1	EDS
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 79240 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9872	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CSD, ZN, CA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.52	25/2285 (1.1%)	1.22	6/3088 (0.2%)
1	B	1.03	0/2285	0.98	5/3088 (0.2%)
1	C	1.29	5/2285 (0.2%)	1.11	7/3088 (0.2%)
1	D	1.04	3/2285 (0.1%)	0.95	1/3088 (0.0%)
All	All	1.24	33/9140 (0.4%)	1.07	19/12352 (0.2%)

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

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

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	192	VAL	CB-CG1	-11.70	1.28	1.52
1	A	137	VAL	CB-CG2	6.98	1.67	1.52
1	A	69	ALA	CA-CB	6.56	1.66	1.52
1	C	198	GLU	CG-CD	6.54	1.61	1.51
1	C	146	CYS	CB-SG	-6.31	1.71	1.82
1	D	198	GLU	CG-CD	6.31	1.61	1.51
1	A	274	SER	CB-OG	6.11	1.50	1.42
1	A	233	GLU	CG-CD	6.11	1.61	1.51
1	A	202	LYS	CE-NZ	6.07	1.64	1.49
1	A	86	GLU	CB-CG	-5.98	1.40	1.52
1	A	195	SER	CB-OG	5.84	1.49	1.42
1	A	122	GLU	CD-OE2	5.77	1.31	1.25
1	A	202	LYS	CD-CE	5.73	1.65	1.51
1	A	220	VAL	CB-CG2	5.61	1.64	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	219	VAL	CB-CG1	5.55	1.64	1.52
1	C	259	GLU	CD-OE1	5.54	1.31	1.25
1	A	166	GLU	CD-OE2	-5.49	1.19	1.25
1	A	31	GLU	CG-CD	5.48	1.60	1.51
1	C	224	GLU	CG-CD	5.38	1.60	1.51
1	A	240	GLU	CD-OE2	5.34	1.31	1.25
1	A	281	LYS	CB-CG	5.33	1.67	1.52
1	A	224	GLU	CB-CG	5.29	1.62	1.52
1	A	281	LYS	N-CA	5.23	1.56	1.46
1	C	231	SER	CB-OG	5.23	1.49	1.42
1	D	198	GLU	CB-CG	5.19	1.62	1.52
1	A	159	THR	CB-CG2	5.18	1.69	1.52
1	A	240	GLU	CG-CD	5.17	1.59	1.51
1	A	198	GLU	CD-OE2	5.16	1.31	1.25
1	A	2	LYS	CE-NZ	5.16	1.61	1.49
1	A	191	GLU	CB-CG	5.14	1.61	1.52
1	A	197	VAL	CB-CG1	5.05	1.63	1.52
1	A	259	GLU	CB-CG	-5.05	1.42	1.52
1	A	89	VAL	CB-CG2	5.03	1.63	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	8	ASP	CB-CG-OD1	7.59	125.14	118.30
1	A	202	LYS	CD-CE-NZ	6.92	127.61	111.70
1	B	284	ARG	NE-CZ-NH1	-6.66	116.97	120.30
1	C	235	LEU	CB-CG-CD2	-6.41	100.11	111.00
1	A	157	ASP	CB-CG-OD1	6.16	123.85	118.30
1	C	284	ARG	NE-CZ-NH2	6.15	123.37	120.30
1	B	278	ASP	CB-CG-OD2	5.96	123.66	118.30
1	C	157	ASP	CB-CG-OD2	5.89	123.60	118.30
1	A	146	CYS	CA-CB-SG	-5.82	103.53	114.00
1	B	157	ASP	CB-CG-OD2	5.74	123.47	118.30
1	B	213	LEU	CA-CB-CG	5.64	128.28	115.30
1	D	124	LEU	CA-CB-CG	5.52	127.99	115.30
1	C	210	ILE	CB-CA-C	-5.37	100.86	111.60
1	C	292	LEU	CB-CG-CD1	5.35	120.09	111.00
1	A	182	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	177	LEU	CB-CG-CD2	5.27	119.96	111.00
1	C	8	ASP	CB-CG-OD2	-5.17	113.64	118.30
1	A	94	LYS	CD-CE-NZ	-5.16	99.84	111.70
1	B	198	GLU	OE1-CD-OE2	5.06	129.38	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	160	LEU	Mainchain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2257	0	2294	74	1
1	B	2257	0	2292	90	0
1	C	2257	0	2293	75	0
1	D	2257	0	2293	107	0
2	A	2	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	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	292	0	0	14	1
4	B	153	0	0	9	0
4	C	254	0	0	14	0
4	D	136	0	0	11	0
All	All	9872	0	9172	327	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 18.

All (327) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:ASN:HB3	1:B:110:VAL:CG2	1.22	1.57
1:A:135:ILE:CG1	1:A:135:ILE:CD1	1.78	1.56
1:A:161:ILE:CD1	1:A:161:ILE:CG1	1.76	1.56
1:B:45:ASN:CB	1:B:110:VAL:HG23	1.54	1.33

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:ASN:CB	1:B:110:VAL:CG2	2.05	1.33
1:B:216:LYS:HD2	4:B:506:HOH:O	1.45	1.12
1:A:31:GLU:OE1	1:A:202:LYS:HE2	1.49	1.11
1:B:45:ASN:HB2	1:B:110:VAL:HG23	1.31	1.09
1:B:45:ASN:HB3	1:B:110:VAL:HG21	1.10	1.09
1:B:43:GLN:HG3	1:B:87:LYS:NZ	1.68	1.06
1:B:43:GLN:O	1:B:110:VAL:HG21	1.53	1.06
1:C:51:GLU:HB2	4:C:786:HOH:O	1.57	1.05
1:A:143:SER:O	4:A:455:HOH:O	1.76	1.02
1:D:58:ASN:HB3	1:D:146:CYS:SG	2.02	1.00
1:B:44:SER:O	1:B:45:ASN:ND2	1.96	0.98
1:D:254:THR:HG22	1:D:256:LEU:H	1.27	0.98
1:D:15:SER:HB2	1:D:143:SER:HB2	1.46	0.97
1:D:44:SER:HB2	1:D:48:LYS:HG2	1.46	0.97
1:D:45:ASN:HA	4:D:651:HOH:O	1.63	0.96
1:A:49:LEU:HD12	1:A:50:ASP:H	1.28	0.95
1:D:45:ASN:HB2	1:D:110:VAL:HG23	1.48	0.95
1:C:29:GLU:OE2	1:C:80:LYS:NZ	2.00	0.94
1:A:31:GLU:OE1	1:A:202:LYS:CE	2.17	0.93
1:B:45:ASN:HB3	1:B:110:VAL:CB	2.01	0.91
1:A:114:LYS:HE3	4:A:398:HOH:O	1.71	0.91
1:B:43:GLN:HG3	1:B:87:LYS:HZ3	1.28	0.90
1:C:144:ILE:H	1:C:144:ILE:HD12	1.37	0.90
1:A:7:ALA:O	1:A:10:HIS:HD2	1.54	0.90
1:A:41:GLN:HE22	1:A:53:ASP:H	1.14	0.90
1:C:212:GLY:O	1:D:46:GLU:HG3	1.72	0.89
1:A:31:GLU:HB3	1:A:202:LYS:HE3	1.55	0.88
1:C:220:VAL:HG23	1:C:229:ILE:HD11	1.56	0.87
1:B:53:ASP:O	1:B:55:ASN:N	2.08	0.86
1:A:44:SER:O	1:A:48:LYS:HB2	1.74	0.86
1:D:299:PHE:OXT	4:D:658:HOH:O	1.92	0.86
1:C:121:ASN:HD22	1:C:121:ASN:C	1.79	0.86
1:D:44:SER:O	1:D:45:ASN:HB3	1.76	0.85
1:D:12:PHE:O	1:D:146:CYS:HB2	1.75	0.85
1:A:43:GLN:H	1:A:48:LYS:HD3	1.41	0.84
1:D:50:ASP:HB3	1:D:106:LEU:HD13	1.59	0.84
1:D:58:ASN:CB	1:D:146:CYS:SG	2.65	0.83
1:A:256:LEU:HD21	4:C:466:HOH:O	1.79	0.82
1:B:53:ASP:O	1:B:53:ASP:OD1	1.98	0.81
1:D:276:VAL:HB	1:D:281:LYS:HE3	1.60	0.81
1:B:278:ASP:HB3	1:B:279:PRO:HD2	1.60	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:45:ASN:HB2	1:D:110:VAL:CG2	2.10	0.81
1:D:249:LEU:HD21	1:D:299:PHE:HZ	1.46	0.80
1:C:121:ASN:ND2	1:C:123:LYS:H	1.80	0.80
1:B:248:GLU:OE2	1:B:252:LYS:HE3	1.81	0.80
1:B:43:GLN:HG3	1:B:87:LYS:HZ1	1.47	0.79
1:C:121:ASN:ND2	1:C:124:LEU:H	1.79	0.79
1:C:45:ASN:OD1	4:C:603:HOH:O	2.00	0.79
1:A:255:ASP:OD2	4:A:414:HOH:O	2.01	0.79
1:D:299:PHE:HB2	4:D:799:HOH:O	1.83	0.79
1:C:211:LYS:O	4:C:706:HOH:O	2.02	0.78
1:B:43:GLN:CG	1:B:87:LYS:NZ	2.46	0.77
1:D:249:LEU:HD21	1:D:299:PHE:CZ	2.19	0.77
1:D:50:ASP:O	1:D:51:GLU:HB3	1.84	0.76
1:B:235:LEU:HB2	1:D:236:ASP:OD1	1.86	0.76
1:A:104:GLU:OE2	4:A:392:HOH:O	2.03	0.75
1:B:237:LYS:HE2	1:B:241:ILE:HG13	1.70	0.74
1:B:45:ASN:CB	1:B:110:VAL:HG21	1.96	0.74
1:A:135:ILE:CB	1:A:135:ILE:CD1	2.66	0.74
1:C:139:PRO:HG3	1:C:144:ILE:HD11	1.69	0.74
1:C:220:VAL:CG2	1:C:229:ILE:HD11	2.19	0.73
1:C:117:LYS:NZ	4:C:627:HOH:O	2.14	0.73
1:B:278:ASP:HB3	1:B:279:PRO:CD	2.19	0.73
1:D:44:SER:HB2	1:D:48:LYS:CG	2.18	0.73
1:C:139:PRO:CG	1:C:144:ILE:HD11	2.20	0.72
1:C:44:SER:HB2	1:C:48:LYS:HE2	1.72	0.71
1:D:254:THR:CG2	1:D:256:LEU:HB2	2.19	0.71
1:A:279:PRO:O	4:A:334:HOH:O	2.07	0.71
1:A:41:GLN:NE2	1:A:53:ASP:H	1.86	0.71
1:C:220:VAL:HG23	1:C:229:ILE:CD1	2.20	0.71
1:D:254:THR:HG22	1:D:256:LEU:N	2.04	0.70
1:C:7:ALA:O	1:C:10:HIS:HD2	1.74	0.70
1:D:264:LEU:O	1:D:268:THR:HG22	1.90	0.70
1:D:46:GLU:OE1	1:D:46:GLU:HA	1.90	0.70
1:B:2:LYS:HE2	1:B:31:GLU:HG3	1.73	0.69
1:C:144:ILE:N	1:C:144:ILE:HD12	2.02	0.69
1:B:42:ILE:HD13	1:B:48:LYS:O	1.91	0.69
1:B:43:GLN:CG	1:B:87:LYS:HZ1	2.05	0.68
1:C:121:ASN:OD1	4:C:466:HOH:O	2.12	0.68
1:C:248:GLU:OE2	1:C:252:LYS:HD2	1.92	0.68
1:B:24:VAL:O	1:B:67:GLU:HG3	1.94	0.68
1:D:239:VAL:CG2	4:D:597:HOH:O	2.41	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:LYS:HE3	1:B:237:LYS:O	1.95	0.67
1:B:78:LYS:HG3	1:B:166:GLU:HG2	1.76	0.67
1:B:237:LYS:HE3	1:B:237:LYS:HA	1.77	0.66
1:B:123:LYS:NZ	1:D:294:SER:O	2.27	0.66
1:D:214:ASN:HB3	4:D:481:HOH:O	1.95	0.66
1:D:44:SER:O	1:D:45:ASN:CB	2.43	0.65
1:A:49:LEU:HD12	1:A:50:ASP:N	2.07	0.65
1:D:254:THR:HG21	1:D:256:LEU:HB2	1.76	0.65
1:C:121:ASN:HD21	1:C:124:LEU:H	1.41	0.65
1:C:44:SER:O	1:C:46:GLU:N	2.26	0.64
1:B:45:ASN:CG	1:B:46:GLU:N	2.51	0.64
1:D:276:VAL:CB	1:D:281:LYS:HE3	2.28	0.64
1:C:43:GLN:OE1	4:C:819:HOH:O	2.14	0.64
1:D:112:ASP:OD2	4:D:329:HOH:O	2.14	0.64
1:A:248:GLU:OE2	1:A:252:LYS:HD2	1.98	0.64
1:B:237:LYS:HE3	1:B:237:LYS:CA	2.28	0.63
1:A:12:PHE:CE1	1:A:56:ARG:HA	2.33	0.63
1:D:240:GLU:O	1:D:244:HIS:HD2	1.81	0.63
1:C:278:ASP:HB3	1:C:279:PRO:HD2	1.80	0.63
1:D:1:MET:HG2	1:D:3:TYR:CZ	2.34	0.62
1:D:140:LYS:HG2	1:D:172:PHE:HB3	1.81	0.62
1:D:94:LYS:HE2	1:D:95:GLY:N	2.15	0.62
1:A:41:GLN:HE22	1:A:53:ASP:N	1.93	0.61
1:D:48:LYS:CE	4:D:666:HOH:O	2.49	0.61
1:C:240:GLU:O	1:C:244:HIS:HD2	1.83	0.61
1:C:121:ASN:ND2	1:C:121:ASN:C	2.51	0.60
1:A:122:GLU:H	1:A:122:GLU:CD	2.04	0.60
1:C:78:LYS:HD3	1:C:166:GLU:HG2	1.82	0.60
1:D:7:ALA:H	1:D:33:GLU:CG	2.14	0.60
1:D:259:GLU:O	1:D:263:THR:HB	2.00	0.60
1:B:56:ARG:HG2	1:B:56:ARG:HH11	1.67	0.60
1:D:15:SER:HB2	1:D:143:SER:CB	2.29	0.59
1:A:278:ASP:HB3	1:A:279:PRO:CD	2.32	0.59
1:D:278:ASP:HB3	1:D:279:PRO:CD	2.33	0.59
1:B:144:ILE:HG12	1:B:152:HIS:CE1	2.37	0.59
1:C:10:HIS:HE1	4:C:417:HOH:O	1.87	0.58
1:D:47:ASP:O	1:D:49:LEU:N	2.29	0.58
1:C:236:ASP:O	1:C:240:GLU:HG3	2.04	0.57
1:A:141:GLU:HB2	4:A:337:HOH:O	2.04	0.57
1:A:43:GLN:HB2	1:A:48:LYS:HZ2	1.69	0.57
1:B:43:GLN:H	1:B:48:LYS:HE3	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:276:VAL:CA	1:D:281:LYS:HE3	2.35	0.57
1:B:240:GLU:O	1:B:244:HIS:HD2	1.86	0.57
1:A:-1:HIS:N	4:A:543:HOH:O	2.37	0.57
1:A:205:LEU:C	1:A:205:LEU:HD12	2.25	0.57
1:B:97:GLY:HA2	1:B:277:VAL:HG12	1.87	0.57
1:A:7:ALA:O	1:A:10:HIS:CD2	2.46	0.56
1:D:278:ASP:HB3	1:D:279:PRO:HD2	1.85	0.56
1:B:216:LYS:CD	4:B:506:HOH:O	2.24	0.56
1:C:121:ASN:ND2	1:C:123:LYS:N	2.51	0.56
1:A:113:ILE:C	1:A:114:LYS:HG2	2.26	0.56
1:C:42:ILE:HG22	1:C:42:ILE:O	2.06	0.56
1:C:212:GLY:O	1:D:46:GLU:C	2.44	0.56
1:A:240:GLU:O	1:A:244:HIS:HD2	1.89	0.55
1:A:36:ASP:H	1:A:40:ASN:ND2	2.04	0.55
1:B:2:LYS:HE2	1:B:31:GLU:CG	2.36	0.55
1:C:44:SER:C	1:C:46:GLU:H	2.10	0.55
1:D:7:ALA:N	1:D:33:GLU:HG2	2.22	0.55
1:D:290:TRP:HA	1:D:293:GLU:HG3	1.87	0.55
1:B:157:ASP:HB2	1:B:275:GLN:HG2	1.89	0.55
1:D:7:ALA:HB2	1:D:33:GLU:HG3	1.89	0.55
1:D:7:ALA:H	1:D:33:GLU:HG2	1.72	0.55
1:B:67:GLU:OE1	4:B:363:HOH:O	2.18	0.55
1:B:122:GLU:H	1:B:122:GLU:CD	2.09	0.55
1:B:36:ASP:H	1:B:40:ASN:HD21	1.55	0.54
1:B:119:ILE:H	1:B:119:ILE:HD12	1.71	0.54
1:A:78:LYS:HB2	1:A:78:LYS:NZ	2.23	0.54
1:A:6:SER:HB3	4:A:360:HOH:O	2.07	0.54
1:D:37:CSD:HB3	1:D:198:GLU:OE1	2.07	0.54
1:C:121:ASN:HD21	1:C:124:LEU:N	2.04	0.54
1:A:51:GLU:O	1:A:51:GLU:HG3	2.06	0.54
1:D:239:VAL:HG23	1:D:240:GLU:N	2.21	0.54
1:D:251:LYS:O	4:D:556:HOH:O	2.18	0.54
1:C:248:GLU:OE2	1:C:252:LYS:CD	2.55	0.54
1:A:119:ILE:N	1:A:119:ILE:HD13	2.23	0.53
1:C:212:GLY:C	1:D:46:GLU:HG3	2.28	0.53
1:C:272:GLN:HE21	1:C:284:ARG:HE	1.55	0.53
1:D:7:ALA:O	1:D:10:HIS:HD2	1.91	0.53
1:A:213:LEU:HD22	1:A:299:PHE:HB3	1.91	0.53
1:B:36:ASP:H	1:B:40:ASN:ND2	2.07	0.53
1:D:94:LYS:HE2	1:D:95:GLY:H	1.73	0.52
1:A:264:LEU:O	1:A:268:THR:HG22	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:ILE:HG23	1:C:135:ILE:O	2.07	0.52
1:C:237:LYS:HD2	1:C:237:LYS:O	2.09	0.52
1:C:205:LEU:C	1:C:205:LEU:HD12	2.30	0.52
1:D:157:ASP:OD2	1:D:191:GLU:OE1	2.28	0.52
1:B:144:ILE:HD13	1:B:148:THR:O	2.10	0.52
1:B:41:GLN:CD	1:B:52:MET:HB3	2.30	0.51
1:D:117:LYS:HA	1:D:126:LEU:O	2.11	0.51
1:D:38:PHE:O	1:D:56:ARG:O	2.29	0.51
1:A:278:ASP:HB3	1:A:279:PRO:HD2	1.92	0.51
1:A:244:HIS:HE1	1:C:101:ASN:OD1	1.93	0.51
1:C:206:GLU:HG2	1:C:207:VAL:N	2.25	0.51
1:D:50:ASP:O	1:D:51:GLU:CB	2.58	0.51
1:D:38:PHE:O	1:D:39:SER:HB2	2.10	0.50
1:D:205:LEU:HD12	1:D:205:LEU:C	2.31	0.50
1:B:70:LYS:HB2	1:B:73:ASP:OD2	2.11	0.50
1:B:7:ALA:O	1:B:10:HIS:HD2	1.95	0.50
1:D:29:GLU:OE2	1:D:80:LYS:NZ	2.33	0.50
1:B:42:ILE:HG23	1:B:89:VAL:HG21	1.94	0.50
1:C:44:SER:HB2	1:C:48:LYS:HG2	1.94	0.50
1:B:49:LEU:HB2	1:B:51:GLU:HG2	1.94	0.50
1:A:193:GLY:O	1:A:194:VAL:HB	2.11	0.50
1:D:44:SER:C	1:D:46:GLU:H	2.14	0.50
1:D:46:GLU:HB2	4:D:649:HOH:O	2.10	0.50
1:B:101:ASN:OD1	1:D:244:HIS:HE1	1.95	0.49
1:A:78:LYS:HD2	1:A:80:LYS:HE2	1.93	0.49
1:D:276:VAL:HB	1:D:281:LYS:CE	2.37	0.49
1:C:272:GLN:NE2	1:C:284:ARG:HE	2.09	0.49
1:B:45:ASN:HB3	1:B:110:VAL:HB	1.90	0.49
1:D:43:GLN:OE1	1:D:43:GLN:N	2.45	0.49
1:C:278:ASP:HB3	1:C:279:PRO:CD	2.43	0.49
1:D:122:GLU:CD	1:D:122:GLU:H	2.15	0.49
1:D:52:MET:O	1:D:53:ASP:HB2	2.12	0.49
1:B:42:ILE:HG23	4:B:356:HOH:O	2.12	0.49
1:D:239:VAL:HG21	4:D:597:HOH:O	2.07	0.49
1:B:144:ILE:HD12	1:B:144:ILE:O	2.11	0.49
1:D:38:PHE:O	1:D:39:SER:CB	2.61	0.49
1:A:122:GLU:N	1:A:122:GLU:CD	2.66	0.49
1:B:220:VAL:CG1	1:B:229:ILE:HD11	2.42	0.49
1:C:46:GLU:C	1:C:48:LYS:H	2.15	0.49
1:D:87:LYS:O	1:D:87:LYS:HG2	2.13	0.49
1:D:45:ASN:OD1	1:D:109:LYS:HA	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:299:PHE:HD1	4:C:335:HOH:O	1.95	0.48
1:C:20:PRO:HB3	1:C:65:PHE:HB2	1.95	0.48
1:D:254:THR:HG22	1:D:256:LEU:HB2	1.94	0.48
1:D:219:VAL:HG11	1:D:299:PHE:CE1	2.49	0.48
1:C:50:ASP:HB2	1:C:106:LEU:HD12	1.94	0.48
1:D:98:VAL:HB	1:D:192:VAL:HG13	1.96	0.48
1:A:-1:HIS:HE1	4:A:634:HOH:O	1.95	0.48
1:A:237:LYS:O	1:A:238:ALA:C	2.50	0.47
1:A:43:GLN:CB	1:A:48:LYS:HZ2	2.27	0.47
1:D:58:ASN:HB2	1:D:146:CYS:SG	2.51	0.47
1:C:113:ILE:O	1:C:114:LYS:HD3	2.14	0.47
1:B:158:THR:HG22	1:B:274:SER:CB	2.44	0.47
1:A:161:ILE:CB	1:A:161:ILE:CD1	2.79	0.47
1:D:254:THR:HG22	1:D:255:ASP:N	2.29	0.47
1:B:90:LEU:HD13	1:D:263:THR:HG23	1.97	0.47
1:B:251:LYS:O	1:B:251:LYS:HG2	2.13	0.47
1:D:121:ASN:ND2	4:D:529:HOH:O	2.45	0.47
1:A:46:GLU:CD	1:B:212:GLY:HA2	2.35	0.47
1:B:193:GLY:H	1:D:266:SER:HB2	1.78	0.47
1:C:87:LYS:CE	4:C:373:HOH:O	2.60	0.47
1:A:114:LYS:CE	4:A:398:HOH:O	2.46	0.46
1:B:237:LYS:O	1:B:237:LYS:CE	2.63	0.46
1:C:6:SER:HB2	1:C:8:ASP:OD1	2.16	0.46
1:A:113:ILE:O	1:A:114:LYS:HG2	2.14	0.46
1:C:240:GLU:O	1:C:244:HIS:CD2	2.67	0.46
1:D:156:MET:O	1:D:157:ASP:C	2.54	0.46
1:C:12:PHE:CE1	1:C:56:ARG:HA	2.50	0.46
1:A:253:HIS:HE1	4:A:408:HOH:O	1.99	0.46
1:B:220:VAL:HG12	1:B:229:ILE:HD11	1.98	0.46
1:B:46:GLU:C	1:B:48:LYS:H	2.19	0.46
1:A:94:LYS:HG3	1:A:103:MET:O	2.16	0.46
1:C:52:MET:HE1	1:C:57:VAL:HG12	1.96	0.46
1:C:139:PRO:HG2	1:C:144:ILE:HD11	1.96	0.46
1:C:174:GLU:HG2	4:C:741:HOH:O	2.15	0.46
1:C:43:GLN:HB3	1:C:43:GLN:HE21	1.52	0.46
1:B:94:LYS:NZ	4:B:646:HOH:O	2.48	0.46
1:B:244:HIS:HE1	1:D:101:ASN:OD1	1.98	0.45
1:A:36:ASP:H	1:A:40:ASN:HD21	1.62	0.45
1:B:193:GLY:H	1:D:266:SER:CB	2.29	0.45
1:B:119:ILE:HD12	1:B:119:ILE:N	2.31	0.45
1:C:240:GLU:OE2	4:C:362:HOH:O	2.21	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:GLU:HG3	4:B:430:HOH:O	2.16	0.45
1:A:46:GLU:HG2	1:B:214:ASN:ND2	2.31	0.45
1:D:65:PHE:CE1	1:D:175:GLY:HA3	2.52	0.45
1:D:44:SER:CB	1:D:48:LYS:HG2	2.33	0.45
1:D:7:ALA:O	1:D:10:HIS:CD2	2.69	0.44
1:A:51:GLU:O	1:A:51:GLU:CG	2.65	0.44
1:C:212:GLY:O	1:D:46:GLU:O	2.36	0.44
1:A:94:LYS:HE2	1:A:104:GLU:OE1	2.16	0.44
1:A:214:ASN:ND2	4:A:769:HOH:O	2.41	0.44
1:B:144:ILE:HD11	1:B:152:HIS:NE2	2.32	0.44
1:A:82:ILE:HG21	1:A:82:ILE:HD13	1.75	0.44
1:A:180:LEU:HD23	1:A:180:LEU:N	2.32	0.44
1:B:7:ALA:HB1	1:B:35:MET:HG3	1.99	0.44
1:A:228:THR:HB	4:A:304:HOH:O	2.18	0.44
1:B:237:LYS:C	1:B:237:LYS:HE3	2.38	0.44
1:A:101:ASN:OD1	1:C:244:HIS:HE1	2.00	0.44
1:A:39:SER:O	1:A:40:ASN:HB2	2.18	0.43
1:C:248:GLU:HG3	1:C:252:LYS:HD3	1.99	0.43
1:C:114:LYS:HD3	1:C:114:LYS:HA	1.73	0.43
1:D:1:MET:HG2	1:D:3:TYR:CE2	2.53	0.43
1:B:41:GLN:OE1	1:B:52:MET:HB3	2.18	0.43
1:B:158:THR:HG22	1:B:274:SER:HB3	2.00	0.43
1:A:31:GLU:CB	1:A:202:LYS:HE3	2.38	0.43
1:B:186:LEU:HG	1:D:267:ILE:O	2.19	0.43
1:B:216:LYS:HB3	4:B:432:HOH:O	2.18	0.43
1:D:45:ASN:CG	1:D:110:VAL:H	2.21	0.43
1:D:6:SER:HA	1:D:33:GLU:HG2	2.01	0.43
1:C:35:MET:HB2	1:C:59:PRO:HG2	2.00	0.43
1:A:2:LYS:HE2	1:A:31:GLU:OE2	2.19	0.43
1:C:206:GLU:OE1	4:C:336:HOH:O	2.21	0.43
1:D:14:PHE:O	1:D:143:SER:HA	2.18	0.43
1:B:156:MET:O	1:B:158:THR:HG22	2.18	0.43
1:B:45:ASN:CG	1:B:110:VAL:HB	2.38	0.42
1:B:81:LYS:HB2	1:B:204:LEU:HB3	2.00	0.42
1:D:12:PHE:CE1	1:D:56:ARG:HA	2.54	0.42
1:C:7:ALA:O	1:C:10:HIS:CD2	2.63	0.42
1:D:54:TRP:C	1:D:56:ARG:H	2.22	0.42
1:D:1:MET:HG2	1:D:3:TYR:OH	2.20	0.42
1:D:81:LYS:HE3	1:D:81:LYS:HB2	1.56	0.42
1:D:83:GLU:OE2	1:D:202:LYS:HE3	2.20	0.42
1:C:45:ASN:HA	1:C:110:VAL:HG23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:288:PRO:HD2	1:D:291:ILE:HG13	2.02	0.42
1:D:76:LYS:HE2	1:D:208:GLU:OE2	2.20	0.42
1:A:-1:HIS:CE1	4:A:634:HOH:O	2.72	0.42
1:A:227:ALA:HA	1:A:285:PHE:O	2.20	0.42
1:B:45:ASN:CB	1:B:110:VAL:HB	2.49	0.42
1:B:144:ILE:HD12	1:B:145:ASN:O	2.19	0.42
1:B:193:GLY:HA2	4:B:319:HOH:O	2.20	0.42
1:B:215:LEU:HD23	1:B:215:LEU:HA	1.91	0.42
1:B:236:ASP:OD2	1:D:234:SER:HB2	2.19	0.41
1:C:123:LYS:HB2	1:C:123:LYS:NZ	2.35	0.41
1:B:236:ASP:OD1	1:D:281:LYS:NZ	2.38	0.41
1:A:43:GLN:H	1:A:48:LYS:CD	2.20	0.41
1:C:227:ALA:HA	1:C:285:PHE:O	2.20	0.41
1:B:237:LYS:NZ	1:B:240:GLU:OE1	2.52	0.41
1:D:74:VAL:HG21	1:D:213:LEU:HD23	2.01	0.41
1:D:37:CSD:O	1:D:58:ASN:N	2.47	0.41
1:B:236:ASP:OD1	1:D:235:LEU:HB2	2.20	0.41
1:A:121:ASN:HB2	1:A:122:GLU:OE1	2.20	0.41
1:B:122:GLU:N	1:B:122:GLU:CD	2.74	0.41
1:C:264:LEU:HD21	1:C:291:ILE:CG2	2.50	0.41
1:C:43:GLN:C	1:C:45:ASN:N	2.74	0.41
1:D:240:GLU:O	1:D:244:HIS:CD2	2.68	0.41
1:A:119:ILE:N	1:A:119:ILE:CD1	2.84	0.41
1:B:38:PHE:O	1:B:41:GLN:HG3	2.21	0.41
1:C:274:SER:HA	4:C:304:HOH:O	2.21	0.41
1:C:123:LYS:CB	1:C:123:LYS:NZ	2.84	0.40
1:A:182:ASP:O	1:A:182:ASP:CG	2.58	0.40
1:D:254:THR:CG2	1:D:255:ASP:N	2.83	0.40
1:C:114:LYS:HB2	1:C:119:ILE:HD11	2.03	0.40
1:B:67:GLU:HB3	4:B:363:HOH:O	2.20	0.40
1:C:235:LEU:HA	1:C:235:LEU:HD23	1.81	0.40
1:A:19:LYS:HA	1:A:20:PRO:HD3	1.91	0.40
1:D:1:MET:HE3	1:D:1:MET:HB2	1.89	0.40
1:A:193:GLY:O	1:A:194:VAL:CB	2.69	0.40
1:A:106:LEU:HD13	1:A:107:TYR:H	1.86	0.40
1:A:3:TYR:O	1:A:30:LEU:HA	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:LYS:NZ	4:A:455:HOH:O[1_455]	1.55	0.65

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	298/301 (99%)	276 (93%)	19 (6%)	3 (1%)	19	11
1	B	298/301 (99%)	273 (92%)	21 (7%)	4 (1%)	15	7
1	C	298/301 (99%)	277 (93%)	17 (6%)	4 (1%)	15	7
1	D	298/301 (99%)	276 (93%)	16 (5%)	6 (2%)	9	3
All	All	1192/1204 (99%)	1102 (92%)	73 (6%)	17 (1%)	14	6

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	VAL
1	B	44	SER
1	B	194	VAL
1	C	52	MET
1	C	194	VAL
1	D	39	SER
1	D	48	LYS
1	D	51	GLU
1	D	53	ASP
1	D	194	VAL
1	A	50	ASP
1	B	46	GLU
1	B	54	TRP
1	C	46	GLU
1	C	145	ASN
1	A	45	ASN
1	D	55	ASN

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	245/245 (100%)	229 (94%)	16 (6%)	21	15
1	B	245/245 (100%)	226 (92%)	19 (8%)	16	10
1	C	245/245 (100%)	224 (91%)	21 (9%)	13	7
1	D	245/245 (100%)	227 (93%)	18 (7%)	17	11
All	All	980/980 (100%)	906 (92%)	74 (8%)	16	10

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	47	ASP
1	A	49	LEU
1	A	52	MET
1	A	57	VAL
1	A	81	LYS
1	A	106	LEU
1	A	114	LYS
1	A	119	ILE
1	A	122	GLU
1	A	123	LYS
1	A	194	VAL
1	A	249	LEU
1	A	259	GLU
1	A	266	SER
1	A	292	LEU
1	B	17	GLU
1	B	22	ILE
1	B	42	ILE
1	B	43	GLN
1	B	45	ASN
1	B	46	GLU
1	B	52	MET
1	B	78	LYS

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Mol	Chain	Res	Type
1	B	83	GLU
1	B	86	GLU
1	B	122	GLU
1	B	124	LEU
1	B	140	LYS
1	B	158	THR
1	B	180	LEU
1	B	194	VAL
1	B	235	LEU
1	B	237	LYS
1	B	249	LEU
1	C	16	LYS
1	C	43	GLN
1	C	78	LYS
1	C	106	LEU
1	C	107	TYR
1	C	108	SER
1	C	115	ASP
1	C	121	ASN
1	C	123	LYS
1	C	140	LYS
1	C	144	ILE
1	C	145	ASN
1	C	155	ASN
1	C	180	LEU
1	C	194	VAL
1	C	233	GLU
1	C	249	LEU
1	C	256	LEU
1	C	266	SER
1	C	281	LYS
1	C	292	LEU
1	D	1	MET
1	D	4	SER
1	D	8	ASP
1	D	25	LYS
1	D	43	GLN
1	D	51	GLU
1	D	52	MET
1	D	87	LYS
1	D	94	LYS
1	D	122	GLU

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Mol	Chain	Res	Type
1	D	123	LYS
1	D	143	SER
1	D	180	LEU
1	D	194	VAL
1	D	211	LYS
1	D	256	LEU
1	D	259	GLU
1	D	263	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	HIS
1	A	40	ASN
1	A	41	GLN
1	A	244	HIS
1	B	10	HIS
1	B	40	ASN
1	B	43	GLN
1	B	214	ASN
1	B	244	HIS
1	B	253	HIS
1	C	10	HIS
1	C	43	GLN
1	C	45	ASN
1	C	101	ASN
1	C	121	ASN
1	C	145	ASN
1	C	244	HIS
1	C	272	GLN
1	D	10	HIS
1	D	244	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	CSD	A	37	1,3	3,7,8	2.17	1 (33%)	3,8,10	5.25	1 (33%)
1	CSD	B	37	1,3	3,7,8	0.73	0	3,8,10	4.07	2 (66%)
1	CSD	C	37	1,3	3,7,8	0.60	0	3,8,10	6.61	2 (66%)
1	CSD	D	37	1,3	3,7,8	2.27	1 (33%)	3,8,10	4.37	1 (33%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSD	A	37	1,3	-	0/2/6/8	0/0/0/0
1	CSD	B	37	1,3	-	0/2/6/8	0/0/0/0
1	CSD	C	37	1,3	-	0/2/6/8	0/0/0/0
1	CSD	D	37	1,3	-	0/2/6/8	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	37	CSD	CB-SG	3.19	1.97	1.79
1	D	37	CSD	CB-SG	3.71	1.99	1.79

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	37	CSD	O-C-CA	-2.29	119.52	125.49
1	B	37	CSD	O-C-CA	-2.24	119.65	125.49
1	B	37	CSD	OD1-SG-CB	6.68	116.54	105.40
1	D	37	CSD	OD1-SG-CB	7.47	117.85	105.40
1	A	37	CSD	OD1-SG-CB	8.90	120.23	105.40
1	C	37	CSD	OD1-SG-CB	11.20	124.07	105.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	37	CSD	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	300/301 (99%)	0.65	19 (6%) 23 24	15, 27, 61, 95	0
1	B	300/301 (99%)	0.52	19 (6%) 23 24	21, 42, 68, 103	0
1	C	300/301 (99%)	0.63	18 (6%) 25 27	17, 31, 69, 118	0
1	D	300/301 (99%)	0.67	28 (9%) 11 11	24, 46, 80, 110	0
All	All	1200/1204 (99%)	0.62	84 (7%) 19 21	15, 38, 70, 118	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	49	LEU	12.1
1	D	52	MET	11.5
1	D	49	LEU	10.0
1	D	44	SER	9.4
1	D	55	ASN	9.2
1	B	44	SER	8.8
1	A	47	ASP	8.5
1	D	53	ASP	7.7
1	D	47	ASP	7.0
1	B	45	ASN	7.0
1	D	54	TRP	6.9
1	C	47	ASP	6.8
1	A	49	LEU	6.7
1	A	52	MET	6.7
1	D	38	PHE	6.6
1	B	47	ASP	6.4
1	B	54	TRP	6.3
1	B	43	GLN	6.2
1	B	52	MET	5.9
1	B	42	ILE	5.7
1	A	54	TRP	5.7

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Mol	Chain	Res	Type	RSRZ
1	A	53	ASP	5.7
1	D	45	ASN	5.6
1	D	299	PHE	5.6
1	C	44	SER	5.6
1	A	50	ASP	5.5
1	B	46	GLU	5.5
1	A	46	GLU	5.0
1	B	50	ASP	4.8
1	D	43	GLN	4.4
1	B	-1	HIS	4.2
1	D	46	GLU	4.1
1	B	51	GLU	4.0
1	C	45	ASN	4.0
1	D	57	VAL	4.0
1	D	39	SER	4.0
1	D	41	GLN	3.9
1	C	54	TRP	3.9
1	A	-2	SER	3.8
1	A	42	ILE	3.7
1	B	-2	SER	3.7
1	D	58	ASN	3.3
1	C	51	GLU	3.3
1	D	-2	SER	3.3
1	D	51	GLU	3.2
1	B	48	LYS	3.2
1	A	55	ASN	3.1
1	D	146	CYS	3.1
1	D	74	VAL	3.1
1	A	45	ASN	3.1
1	D	40	ASN	3.1
1	A	51	GLU	3.1
1	D	36	ASP	3.1
1	D	42	ILE	3.1
1	B	53	ASP	3.0
1	A	56	ARG	3.0
1	D	48	LYS	2.9
1	A	48	LYS	2.8
1	C	271	ALA	2.8
1	A	38	PHE	2.8
1	C	265	PHE	2.7
1	C	46	GLU	2.7
1	B	38	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	50	ASP	2.6
1	D	-1	HIS	2.6
1	C	287	LEU	2.5
1	D	56	ARG	2.5
1	B	55	ASN	2.5
1	D	297	ILE	2.5
1	A	-1	HIS	2.4
1	C	-2	SER	2.4
1	A	124	LEU	2.3
1	A	235	LEU	2.3
1	B	144	ILE	2.2
1	C	264	LEU	2.2
1	B	299	PHE	2.1
1	C	49	LEU	2.1
1	A	44	SER	2.1
1	C	238	ALA	2.0
1	C	291	ILE	2.0
1	C	160	LEU	2.0
1	C	262	ALA	2.0
1	C	43	GLN	2.0
1	C	142	GLY	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	CSD	A	37	8/9	0.69	0.18	-	36,39,46,49	3
1	CSD	B	37	8/9	0.88	0.10	-	51,54,63,64	0
1	CSD	C	37	8/9	0.83	0.18	-	42,44,55,55	3
1	CSD	D	37	8/9	0.69	0.19	-	56,60,67,67	0

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	ZN	B	303	1/1	0.99	0.18	7.58	8,8,8,8	0
3	ZN	D	307	1/1	0.99	0.23	7.39	12,12,12,12	0
2	CA	B	304	1/1	0.98	0.18	6.23	5,5,5,5	0
2	CA	A	300	1/1	0.97	0.17	2.27	5,5,5,5	0
2	CA	D	308	1/1	0.99	0.14	0.11	7,7,7,7	0
2	CA	C	306	1/1	0.99	0.09	-2.47	4,4,4,4	0
3	ZN	C	305	1/1	0.99	0.06	-3.16	2,2,2,2	1
2	CA	A	302	1/1	0.99	0.03	-7.33	6,6,6,6	0
3	ZN	A	301	1/1	0.99	0.02	-8.49	7,7,7,7	1

6.5 Other polymers [i](#)

There are no such residues in this entry.