



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:44 AM GMT

PDB ID : 3MZ3
Title : Crystal structure of Co2+ HDAC8 complexed with M344
Authors : Dowling, D.P.; Gattis, S.G.; Fierke, C.A.; Christianson, D.W.
Deposited on : 2010-05-11
Resolution : 3.20 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

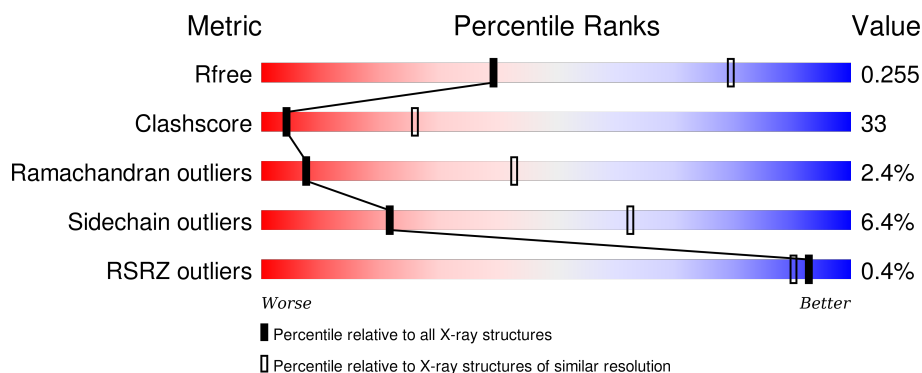
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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	389	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 44%, yellow 43%, orange 5%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 44% 43% 5% 8% </div> </div>
1	B	389	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 43%, yellow 43%, orange 5%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 43% 43% 5% 8% </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5647 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 Histone deacetylase 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	17	0	0
			2785	1789	463	514	19			
1	B	356	Total	C	N	O	S	17	0	0
			2777	1785	462	511	19			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	378	ILE	-	EXPRESSION TAG	UNP Q9BY41
A	379	GLU	-	EXPRESSION TAG	UNP Q9BY41
A	380	GLY	-	EXPRESSION TAG	UNP Q9BY41
A	381	ARG	-	EXPRESSION TAG	UNP Q9BY41
A	382	GLY	-	EXPRESSION TAG	UNP Q9BY41
A	383	SER	-	EXPRESSION TAG	UNP Q9BY41
A	384	HIS	-	EXPRESSION TAG	UNP Q9BY41
A	385	HIS	-	EXPRESSION TAG	UNP Q9BY41
A	386	HIS	-	EXPRESSION TAG	UNP Q9BY41
A	387	HIS	-	EXPRESSION TAG	UNP Q9BY41
A	388	HIS	-	EXPRESSION TAG	UNP Q9BY41
A	389	HIS	-	EXPRESSION TAG	UNP Q9BY41
B	378	ILE	-	EXPRESSION TAG	UNP Q9BY41
B	379	GLU	-	EXPRESSION TAG	UNP Q9BY41
B	380	GLY	-	EXPRESSION TAG	UNP Q9BY41
B	381	ARG	-	EXPRESSION TAG	UNP Q9BY41
B	382	GLY	-	EXPRESSION TAG	UNP Q9BY41
B	383	SER	-	EXPRESSION TAG	UNP Q9BY41
B	384	HIS	-	EXPRESSION TAG	UNP Q9BY41
B	385	HIS	-	EXPRESSION TAG	UNP Q9BY41
B	386	HIS	-	EXPRESSION TAG	UNP Q9BY41
B	387	HIS	-	EXPRESSION TAG	UNP Q9BY41
B	388	HIS	-	EXPRESSION TAG	UNP Q9BY41
B	389	HIS	-	EXPRESSION TAG	UNP Q9BY41

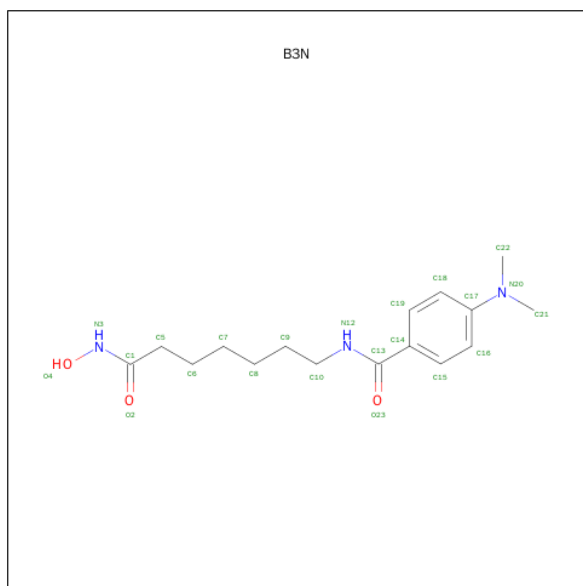
- Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Co	0	0
			1	1		
2	A	1	Total	Co	0	0
			1	1		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	K	0	0
			2	2		
3	A	2	Total	K	0	0
			2	2		

- Molecule 4 is 4-(DIMETHYLAMINO)-N-[7-(HYDROXYAMINO)-7-OXOHEPTYL]BENZAMIDE (three-letter code: B3N) (formula: C₁₆H₂₅N₃O₃).

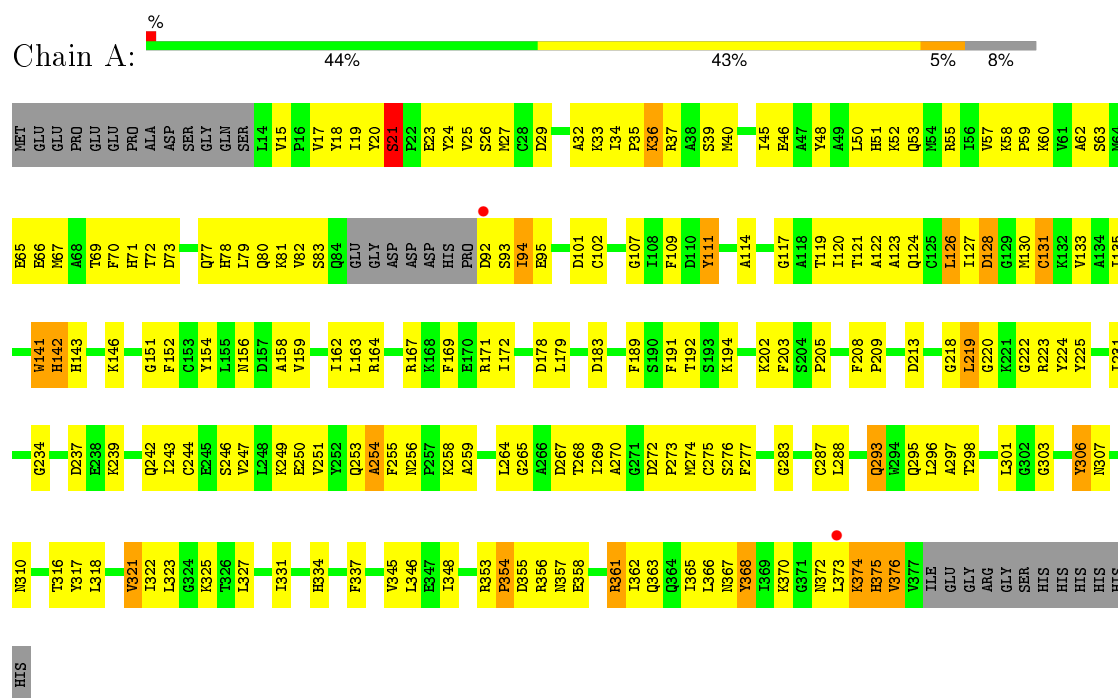


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	15	Total 15	O 15	0	0
5	B	20	Total 20	O 20	0	0

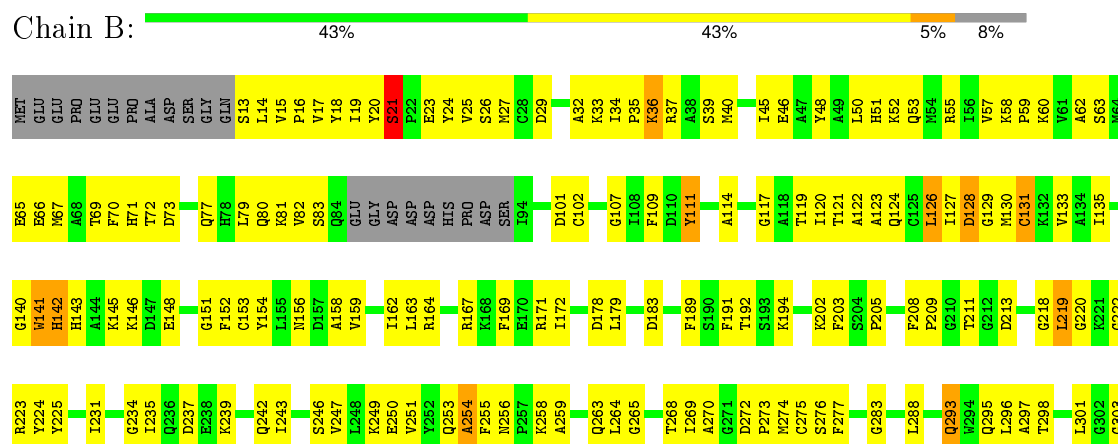
3 Residue-property plots

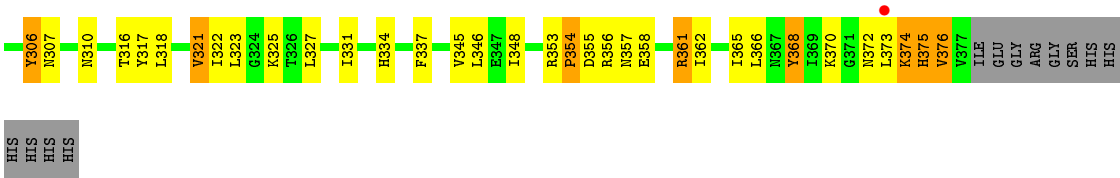
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Histone deacetylase 8



• Molecule 1: Histone deacetylase 8





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.63Å 86.14Å 94.51Å 90.00° 94.07° 90.00°	Depositor
Resolution (Å)	50.00 – 3.20 49.37 – 3.19	Depositor EDS
% Data completeness (in resolution range)	92.2 (50.00-3.20) 91.7 (49.37-3.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 3.19Å)	Xtriage
Refinement program	CNS 2.1	Depositor
R, R_{free}	0.204 , 0.256 0.204 , 0.255	Depositor DCC
R_{free} test set	684 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	63.2	Xtriage
Anisotropy	0.953	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 61.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 14720 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5647	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K, CO, B3N

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	0.41	0/2853	0.64	0/3869
1	B	0.42	0/2845	0.65	0/3858
All	All	0.42	0/5698	0.65	0/7727

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
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	306	TYR	Sidechain
1	B	306	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	2785	0	2752	186	0
1	B	2777	0	2748	185	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	22	0	24	4	0
4	B	22	0	24	4	0
5	A	15	0	0	1	0
5	B	20	0	0	1	0
All	All	5647	0	5548	367	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 (367) 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:272:ASP:OD2	1:B:274:MET:HB2	1.69	0.93
1:A:272:ASP:OD2	1:A:274:MET:HB2	1.70	0.90
1:B:127:ILE:HG22	1:B:169:PHE:HE1	1.37	0.90
1:B:13:SER:HB2	1:B:15:VAL:HG12	1.53	0.89
1:A:127:ILE:HG22	1:A:169:PHE:HE1	1.37	0.88
1:A:83:SER:HB2	1:A:107:GLY:H	1.42	0.84
1:B:83:SER:HB2	1:B:107:GLY:H	1.43	0.83
1:A:218:GLY:C	1:A:219:LEU:HD13	2.03	0.78
1:B:55:ARG:HH11	1:B:55:ARG:HG3	1.48	0.78
1:B:218:GLY:C	1:B:219:LEU:HD13	2.04	0.77
1:A:169:PHE:HD2	1:A:258:LYS:HB3	1.51	0.76
1:B:254:ALA:HB1	1:B:376:VAL:CG2	2.16	0.75
1:A:254:ALA:HB1	1:A:376:VAL:CG2	2.17	0.75
1:A:55:ARG:HG3	1:A:55:ARG:HH11	1.49	0.75
1:B:169:PHE:HD2	1:B:258:LYS:HB3	1.51	0.74
1:B:254:ALA:HB1	1:B:376:VAL:HG21	1.71	0.73
1:B:18:TYR:CE1	1:B:20:TYR:HB2	2.23	0.73
1:A:18:TYR:CE1	1:A:20:TYR:HB2	2.24	0.73
1:A:254:ALA:HB1	1:A:376:VAL:HG21	1.71	0.73
1:A:63:SER:OG	1:A:66:GLU:HG3	1.89	0.72
1:B:69:THR:HG22	1:B:163:LEU:HD13	1.72	0.72
1:A:73:ASP:O	1:A:77:GLN:HB2	1.90	0.71
1:A:83:SER:HB2	1:A:107:GLY:N	2.04	0.71
1:A:219:LEU:HD22	1:A:219:LEU:N	2.05	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:THR:HG22	1:A:163:LEU:HD13	1.71	0.71
1:B:63:SER:OG	1:B:66:GLU:HG3	1.90	0.71
1:B:83:SER:HB2	1:B:107:GLY:N	2.04	0.71
1:A:78:HIS:CE1	1:A:93:SER:HA	2.25	0.71
1:B:73:ASP:O	1:B:77:GLN:HB2	1.91	0.70
1:A:375:HIS:H	1:A:375:HIS:CD2	2.10	0.70
1:B:219:LEU:N	1:B:219:LEU:HD22	2.07	0.69
1:B:223:ARG:O	1:B:224:TYR:HB2	1.93	0.69
1:A:223:ARG:O	1:A:224:TYR:HB2	1.92	0.69
1:B:375:HIS:H	1:B:375:HIS:CD2	2.09	0.68
1:B:179:LEU:HD23	1:B:274:MET:HG2	1.76	0.68
1:B:374:LYS:C	1:B:376:VAL:H	1.97	0.67
1:A:127:ILE:HG22	1:A:169:PHE:CE1	2.27	0.67
1:A:179:LEU:HD23	1:A:274:MET:HG2	1.77	0.66
1:A:218:GLY:O	1:A:219:LEU:HD13	1.95	0.66
1:A:307:ASN:HB3	1:A:310:ASN:HB2	1.76	0.66
1:B:124:GLN:NE2	1:B:127:ILE:HD11	2.11	0.66
1:B:55:ARG:NH1	1:B:55:ARG:HG3	2.10	0.65
1:A:32:ALA:O	1:A:35:PRO:HD3	1.96	0.65
1:B:32:ALA:O	1:B:35:PRO:HD3	1.96	0.65
1:B:218:GLY:O	1:B:219:LEU:HD13	1.95	0.65
1:A:124:GLN:NE2	1:A:127:ILE:HD11	2.11	0.65
1:A:152:PHE:HE2	4:B:390:B3N:H211	1.61	0.65
1:A:239:LYS:HA	1:A:242:GLN:HE21	1.62	0.65
1:A:158:ALA:O	1:A:162:ILE:HG13	1.97	0.65
1:B:13:SER:CB	1:B:15:VAL:HG12	2.26	0.65
1:B:307:ASN:HB3	1:B:310:ASN:HB2	1.77	0.65
1:B:239:LYS:HA	1:B:242:GLN:HE21	1.62	0.64
1:B:50:LEU:HD22	1:B:323:LEU:HD12	1.80	0.64
1:B:158:ALA:O	1:B:162:ILE:HG13	1.97	0.64
1:A:374:LYS:C	1:A:376:VAL:H	1.97	0.64
1:A:247:VAL:O	1:A:251:VAL:HG23	1.98	0.63
1:A:50:LEU:HD22	1:A:323:LEU:HD12	1.81	0.63
1:A:55:ARG:NH1	1:A:55:ARG:HG3	2.10	0.63
1:B:247:VAL:O	1:B:251:VAL:HG23	1.98	0.63
1:B:20:TYR:CG	1:B:21:SER:N	2.67	0.63
1:B:18:TYR:HE1	1:B:20:TYR:HB2	1.62	0.62
1:A:20:TYR:CG	1:A:21:SER:N	2.67	0.62
1:A:18:TYR:HE1	1:A:20:TYR:HB2	1.62	0.62
1:B:127:ILE:HG22	1:B:169:PHE:CE1	2.26	0.62
4:A:390:B3N:H211	1:B:152:PHE:HE2	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:ASP:OD1	1:B:357:ASN:HB2	2.00	0.62
1:A:191:PHE:CE2	1:A:219:LEU:HD23	2.35	0.61
1:B:19:ILE:CG2	1:B:59:PRO:HB3	2.30	0.61
1:A:19:ILE:CG2	1:A:59:PRO:HB3	2.29	0.61
1:B:191:PHE:CE2	1:B:219:LEU:HD23	2.35	0.61
1:A:355:ASP:OD1	1:A:357:ASN:HB2	1.99	0.61
1:B:355:ASP:O	1:B:357:ASN:N	2.34	0.61
1:A:223:ARG:HG2	1:A:224:TYR:CD2	2.36	0.60
1:A:355:ASP:O	1:A:357:ASN:N	2.34	0.60
1:A:208:PHE:CD2	1:A:209:PRO:HA	2.36	0.60
1:B:223:ARG:HG2	1:B:224:TYR:CD2	2.36	0.60
1:B:375:HIS:CD2	1:B:375:HIS:N	2.70	0.59
1:B:14:LEU:O	1:B:16:PRO:HD3	2.02	0.59
1:A:69:THR:HG22	1:A:163:LEU:CD1	2.33	0.59
1:B:208:PHE:CD2	1:B:209:PRO:HA	2.37	0.59
1:B:374:LYS:HZ2	1:B:374:LYS:HB2	1.69	0.58
1:B:331:ILE:HD11	1:B:346:LEU:HA	1.86	0.58
1:A:19:ILE:HG23	1:A:59:PRO:HB3	1.86	0.58
1:B:83:SER:CB	1:B:107:GLY:H	2.16	0.57
1:A:375:HIS:N	1:A:375:HIS:CD2	2.70	0.57
1:B:143:HIS:NE2	4:B:390:B3N:N3	2.50	0.57
1:B:239:LYS:O	1:B:243:ILE:HG13	2.05	0.57
1:A:178:ASP:OD1	1:A:265:GLY:HA3	2.04	0.57
1:B:17:VAL:HG13	1:B:57:VAL:CG2	2.35	0.57
1:B:69:THR:HG22	1:B:163:LEU:CD1	2.34	0.57
1:A:331:ILE:HD11	1:A:346:LEU:HA	1.86	0.57
1:A:239:LYS:O	1:A:243:ILE:HG13	2.04	0.57
1:B:178:ASP:OD1	1:B:265:GLY:HA3	2.04	0.57
1:B:225:TYR:CZ	1:B:376:VAL:HG12	2.40	0.56
1:B:254:ALA:HB1	1:B:376:VAL:HG23	1.87	0.56
1:B:19:ILE:HG23	1:B:59:PRO:HB3	1.87	0.56
1:A:225:TYR:CZ	1:A:376:VAL:HG12	2.40	0.56
1:B:203:PHE:CE2	1:B:213:ASP:HB2	2.41	0.56
1:A:63:SER:HG	1:A:66:GLU:HG3	1.68	0.56
1:B:36:LYS:O	1:B:40:MET:HG3	2.05	0.56
1:B:143:HIS:CD2	4:B:390:B3N:H24	2.24	0.56
1:A:202:LYS:HD2	1:A:276:SER:OG	2.06	0.56
1:A:156:ASN:OD1	1:A:159:VAL:HG23	2.06	0.56
1:A:128:ASP:C	1:A:130:MET:H	2.09	0.56
1:A:29:ASP:OD1	1:A:39:SER:OG	2.25	0.55
1:A:83:SER:CB	1:A:107:GLY:H	2.15	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:374:LYS:O	1:B:376:VAL:N	2.40	0.55
1:A:172:ILE:N	1:A:172:ILE:HD12	2.22	0.55
1:B:288:LEU:HD23	1:B:321:VAL:CG2	2.37	0.55
1:B:29:ASP:OD1	1:B:39:SER:OG	2.24	0.55
1:A:203:PHE:CE2	1:A:213:ASP:HB2	2.41	0.55
1:A:264:LEU:HD13	1:A:318:LEU:HD13	1.88	0.55
1:B:202:LYS:HD2	1:B:276:SER:OG	2.06	0.55
1:B:264:LEU:HD13	1:B:318:LEU:HD13	1.88	0.55
1:B:128:ASP:C	1:B:130:MET:H	2.09	0.55
1:A:124:GLN:HE22	1:A:127:ILE:HD11	1.71	0.55
1:B:33:LYS:HE3	1:B:152:PHE:CE1	2.42	0.55
1:B:250:GLU:OE1	1:B:370:LYS:HG3	2.07	0.54
1:B:172:ILE:N	1:B:172:ILE:HD12	2.22	0.54
1:B:171:ARG:HH21	1:B:171:ARG:HG3	1.73	0.54
1:A:79:LEU:O	1:A:82:VAL:HG12	2.07	0.54
1:A:254:ALA:HB1	1:A:376:VAL:HG23	1.88	0.54
1:A:171:ARG:HG3	1:A:171:ARG:HH21	1.72	0.54
1:B:124:GLN:HE22	1:B:127:ILE:HD11	1.72	0.54
1:A:83:SER:HB2	1:A:107:GLY:CA	2.37	0.54
1:A:33:LYS:HE3	1:A:152:PHE:CE1	2.42	0.54
1:A:250:GLU:OE1	1:A:370:LYS:HG3	2.07	0.54
1:A:374:LYS:CB	1:A:374:LYS:NZ	2.70	0.54
1:B:250:GLU:HB2	1:B:373:LEU:HD23	1.88	0.54
1:A:62:ALA:HB2	1:A:117:GLY:N	2.22	0.54
1:B:34:ILE:HG21	1:B:37:ARG:NH1	2.23	0.54
1:A:111:TYR:CD1	1:A:111:TYR:C	2.81	0.54
1:B:111:TYR:C	1:B:111:TYR:CD1	2.81	0.54
1:A:36:LYS:O	1:A:40:MET:HG3	2.07	0.54
1:A:202:LYS:HB2	1:A:277:PHE:CE2	2.43	0.54
1:A:288:LEU:HD23	1:A:321:VAL:CG2	2.38	0.54
1:B:374:LYS:CB	1:B:374:LYS:NZ	2.70	0.54
1:A:111:TYR:O	1:A:114:ALA:HB3	2.08	0.54
1:A:374:LYS:O	1:A:376:VAL:N	2.40	0.53
1:B:224:TYR:HD1	1:B:372:ASN:HB3	1.73	0.53
1:B:79:LEU:O	1:B:82:VAL:HG12	2.06	0.53
1:B:83:SER:HB2	1:B:107:GLY:CA	2.38	0.53
1:A:250:GLU:HB2	1:A:373:LEU:HD23	1.89	0.53
1:B:202:LYS:HB2	1:B:277:PHE:CE2	2.44	0.53
1:B:62:ALA:HB2	1:B:117:GLY:N	2.24	0.53
1:A:17:VAL:HG13	1:A:57:VAL:CG2	2.37	0.53
1:B:24:TYR:O	1:B:27:MET:HB3	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:TYR:HD1	1:A:372:ASN:HB3	1.73	0.53
1:B:71:HIS:HA	1:B:146:LYS:O	2.08	0.53
1:B:374:LYS:C	1:B:376:VAL:N	2.62	0.53
1:A:269:ILE:O	1:A:270:ALA:C	2.48	0.52
1:B:251:VAL:CG2	1:B:373:LEU:HD21	2.39	0.52
1:A:24:TYR:O	1:A:27:MET:HB3	2.08	0.52
1:A:374:LYS:C	1:A:376:VAL:N	2.63	0.52
1:B:19:ILE:HG13	1:B:122:ALA:HB2	1.92	0.52
1:B:334:HIS:H	1:B:337:PHE:HB2	1.75	0.52
1:A:102:CYS:HB3	1:A:154:TYR:CE2	2.45	0.52
1:A:189:PHE:O	1:A:192:THR:HG22	2.10	0.52
1:A:20:TYR:HE1	1:A:25:VAL:HG21	1.75	0.52
1:A:18:TYR:CD1	1:A:20:TYR:HB2	2.45	0.51
1:A:70:PHE:CD1	1:A:159:VAL:HG11	2.44	0.51
1:A:71:HIS:HA	1:A:146:LYS:O	2.10	0.51
1:B:156:ASN:OD1	1:B:159:VAL:HG23	2.09	0.51
1:A:34:ILE:HG21	1:A:37:ARG:NH1	2.24	0.51
1:B:269:ILE:O	1:B:270:ALA:C	2.48	0.51
1:B:251:VAL:HG22	1:B:373:LEU:HD21	1.93	0.51
1:B:18:TYR:CD1	1:B:20:TYR:HB2	2.45	0.51
1:B:20:TYR:HE1	1:B:25:VAL:HG21	1.74	0.51
1:B:375:HIS:H	1:B:375:HIS:HD2	1.55	0.51
1:B:46:GLU:C	1:B:48:TYR:H	2.13	0.51
1:B:111:TYR:O	1:B:114:ALA:HB3	2.10	0.51
1:B:70:PHE:CD1	1:B:159:VAL:HG11	2.45	0.51
1:B:17:VAL:HG13	1:B:57:VAL:HG23	1.93	0.51
1:A:375:HIS:H	1:A:375:HIS:HD2	1.55	0.51
1:B:135:ILE:HD11	1:B:322:ILE:HD13	1.92	0.51
1:A:53:GLN:NE2	1:A:325:LYS:HE3	2.25	0.51
1:B:374:LYS:NZ	1:B:374:LYS:HB2	2.25	0.51
1:B:189:PHE:O	1:B:192:THR:HG22	2.10	0.51
1:A:251:VAL:CG2	1:A:373:LEU:HD21	2.41	0.51
1:B:126:LEU:CD1	1:B:297:ALA:HB1	2.42	0.50
1:A:334:HIS:H	1:A:337:PHE:HB2	1.75	0.50
1:A:169:PHE:CD2	1:A:259:ALA:HB2	2.45	0.50
1:A:135:ILE:HD11	1:A:322:ILE:HD13	1.92	0.50
1:A:135:ILE:HD11	1:A:322:ILE:CD1	2.42	0.50
1:B:169:PHE:CD2	1:B:259:ALA:HB2	2.46	0.50
1:A:374:LYS:HB2	1:A:374:LYS:NZ	2.26	0.50
1:A:19:ILE:HG13	1:A:122:ALA:HB2	1.92	0.50
1:B:102:CYS:HB3	1:B:154:TYR:CE2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:GLU:C	1:A:48:TYR:H	2.14	0.49
1:A:21:SER:HB2	1:A:58:LYS:HG3	1.94	0.49
1:A:223:ARG:HG3	1:A:223:ARG:HH11	1.78	0.49
1:B:101:ASP:OD2	1:B:151:GLY:N	2.44	0.49
1:B:53:GLN:NE2	1:B:325:LYS:HE3	2.27	0.49
1:B:368:TYR:O	1:B:368:TYR:HD1	1.95	0.49
1:A:368:TYR:HD1	1:A:368:TYR:O	1.96	0.49
1:A:59:PRO:HA	1:A:121:THR:HG21	1.95	0.48
1:A:189:PHE:HB3	1:A:192:THR:CG2	2.42	0.48
1:B:223:ARG:HG3	1:B:223:ARG:HH11	1.78	0.48
1:B:224:TYR:CD1	1:B:372:ASN:HB3	2.48	0.48
1:A:251:VAL:HG22	1:A:373:LEU:HD21	1.94	0.48
1:A:93:SER:C	1:A:95:GLU:H	2.17	0.48
1:B:63:SER:HG	1:B:66:GLU:HG3	1.74	0.48
1:A:53:GLN:HE22	1:A:325:LYS:CE	2.27	0.48
1:A:205:PRO:HA	5:A:403:HOH:O	2.13	0.48
1:B:355:ASP:C	1:B:357:ASN:H	2.17	0.48
1:A:53:GLN:HE22	1:A:325:LYS:HE3	1.79	0.48
1:B:189:PHE:HB3	1:B:192:THR:CG2	2.43	0.48
1:B:59:PRO:HA	1:B:121:THR:HG21	1.95	0.48
1:B:288:LEU:HD23	1:B:321:VAL:HG21	1.95	0.48
1:A:288:LEU:HD23	1:A:321:VAL:HG21	1.96	0.48
1:B:135:ILE:HD11	1:B:322:ILE:CD1	2.42	0.48
1:B:21:SER:HB2	1:B:58:LYS:HG3	1.94	0.48
1:A:224:TYR:CD1	1:A:372:ASN:HB3	2.48	0.48
1:A:288:LEU:HD22	1:A:318:LEU:HD22	1.94	0.48
1:B:237:ASP:OD1	1:B:283:GLY:HA3	2.14	0.47
1:B:67:MET:C	1:B:69:THR:H	2.18	0.47
1:B:288:LEU:HD22	1:B:318:LEU:HD22	1.96	0.47
1:B:55:ARG:O	1:B:57:VAL:HG23	2.14	0.47
1:A:361:ARG:O	1:A:365:ILE:HG13	2.14	0.47
1:A:355:ASP:C	1:A:357:ASN:H	2.16	0.47
1:A:130:MET:O	1:A:131:CYS:HB3	2.14	0.47
1:A:101:ASP:OD2	1:A:151:GLY:N	2.44	0.47
1:B:130:MET:O	1:B:131:CYS:HB3	2.14	0.47
1:A:141:TRP:HA	1:A:141:TRP:CE3	2.50	0.47
1:A:126:LEU:CD1	1:A:297:ALA:HB1	2.45	0.47
1:A:374:LYS:HB2	1:A:374:LYS:HZ2	1.78	0.47
1:A:17:VAL:HG13	1:A:57:VAL:HG23	1.95	0.47
1:B:141:TRP:HA	1:B:141:TRP:CE3	2.50	0.47
1:B:13:SER:HB2	1:B:15:VAL:H	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:LYS:HA	1:B:242:GLN:NE2	2.28	0.47
1:B:179:LEU:CD2	1:B:274:MET:HG2	2.45	0.46
1:A:219:LEU:HD22	1:A:219:LEU:H	1.79	0.46
1:B:361:ARG:O	1:B:365:ILE:HG13	2.15	0.46
1:B:80:GLN:HE21	1:B:109:PHE:HE2	1.64	0.46
1:B:101:ASP:O	1:B:152:PHE:CD1	2.68	0.46
1:A:142:HIS:CD2	1:A:142:HIS:H	2.33	0.46
1:B:53:GLN:HE22	1:B:325:LYS:HE3	1.80	0.46
1:B:293:GLN:O	1:B:293:GLN:HG2	2.15	0.46
1:A:239:LYS:HA	1:A:242:GLN:NE2	2.27	0.46
1:A:133:VAL:HG22	1:A:298:THR:HB	1.98	0.46
1:A:237:ASP:OD1	1:A:283:GLY:HA3	2.15	0.46
1:A:143:HIS:CD2	4:A:390:B3N:H24	2.33	0.46
1:A:101:ASP:O	1:A:152:PHE:CD1	2.68	0.46
1:B:220:GLY:C	1:B:222:GLY:N	2.69	0.46
1:A:124:GLN:C	1:A:126:LEU:H	2.19	0.46
1:B:220:GLY:C	1:B:222:GLY:H	2.20	0.46
1:B:142:HIS:CD2	1:B:142:HIS:H	2.33	0.45
1:B:124:GLN:C	1:B:126:LEU:H	2.20	0.45
1:B:53:GLN:HE22	1:B:325:LYS:CE	2.28	0.45
1:B:133:VAL:HG22	1:B:298:THR:HB	1.99	0.45
1:A:220:GLY:C	1:A:222:GLY:N	2.69	0.45
1:A:293:GLN:HG2	1:A:293:GLN:O	2.16	0.45
1:B:246:SER:O	1:B:249:LYS:HB3	2.15	0.45
1:A:55:ARG:O	1:A:57:VAL:HG23	2.15	0.45
1:A:203:PHE:C	1:A:203:PHE:CD1	2.90	0.45
1:A:273:PRO:C	1:A:275:CYS:N	2.70	0.45
1:A:244:CYS:HG	1:A:287:CYS:HB3	1.81	0.45
1:A:78:HIS:HE1	1:A:93:SER:HA	1.77	0.45
1:A:23:GLU:H	1:A:23:GLU:CD	2.20	0.45
1:B:23:GLU:H	1:B:23:GLU:CD	2.20	0.45
1:B:269:ILE:HG21	1:B:348:ILE:CD1	2.47	0.45
1:A:67:MET:C	1:A:69:THR:H	2.20	0.45
1:A:143:HIS:NE2	4:A:390:B3N:N3	2.65	0.45
1:B:269:ILE:HG21	1:B:348:ILE:HD12	1.99	0.45
1:A:205:PRO:HD2	1:B:353:ARG:HB3	1.98	0.45
1:A:358:GLU:HB3	1:A:361:ARG:HB2	1.98	0.45
1:B:50:LEU:HD21	1:B:327:LEU:HD21	2.00	0.44
1:B:203:PHE:CD1	1:B:203:PHE:C	2.90	0.44
1:B:128:ASP:C	1:B:130:MET:N	2.70	0.44
1:A:246:SER:O	1:A:249:LYS:HB3	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:ILE:HG21	1:A:348:ILE:HD12	1.98	0.44
1:A:358:GLU:O	1:A:362:ILE:HG12	2.18	0.44
1:A:269:ILE:HG21	1:A:348:ILE:CD1	2.47	0.44
1:A:189:PHE:HB3	1:A:192:THR:HG22	1.99	0.44
1:B:317:TYR:O	1:B:321:VAL:HG13	2.18	0.44
1:B:361:ARG:HH21	1:B:365:ILE:CG1	2.31	0.44
1:A:128:ASP:C	1:A:130:MET:N	2.70	0.44
1:A:183:ASP:N	1:A:183:ASP:OD2	2.51	0.44
1:A:223:ARG:O	1:A:224:TYR:CB	2.64	0.44
1:A:119:THR:HG23	1:A:301:LEU:HD21	2.00	0.44
1:A:94:ILE:O	1:A:94:ILE:HG22	2.17	0.44
1:A:80:GLN:HE21	1:A:109:PHE:HE2	1.65	0.44
1:B:258:LYS:HA	1:B:258:LYS:HD2	1.82	0.43
1:B:358:GLU:HB3	1:B:361:ARG:HB2	1.99	0.43
1:A:15:VAL:HG13	1:A:15:VAL:O	2.18	0.43
1:B:362:ILE:O	1:B:366:LEU:HG	2.18	0.43
1:B:169:PHE:CD2	1:B:258:LYS:HB3	2.43	0.43
1:B:234:GLY:HA3	1:B:355:ASP:HA	2.00	0.43
1:A:317:TYR:O	1:A:321:VAL:HG13	2.18	0.43
1:A:220:GLY:C	1:A:222:GLY:H	2.20	0.43
1:B:273:PRO:C	1:B:275:CYS:N	2.71	0.43
1:B:358:GLU:O	1:B:362:ILE:HG12	2.17	0.43
1:A:72:THR:HG23	1:A:146:LYS:HG3	2.01	0.43
1:A:50:LEU:HD21	1:A:327:LEU:HD21	1.99	0.43
4:A:390:B3N:H221	1:B:152:PHE:CE2	2.53	0.43
1:A:219:LEU:CD2	1:A:219:LEU:N	2.77	0.43
1:A:362:ILE:O	1:A:366:LEU:HG	2.19	0.43
1:A:37:ARG:NH2	1:A:303:GLY:O	2.52	0.43
1:B:189:PHE:HB3	1:B:192:THR:HG22	2.00	0.43
1:B:45:ILE:HA	1:B:316:THR:HG23	2.01	0.43
1:B:183:ASP:N	1:B:183:ASP:OD2	2.50	0.43
1:A:353:ARG:HD3	1:B:205:PRO:O	2.18	0.42
1:B:171:ARG:HA	1:B:194:LYS:O	2.18	0.42
1:B:72:THR:HG23	1:B:146:LYS:HG3	2.00	0.42
1:A:353:ARG:HB2	1:A:354:PRO:HD2	2.01	0.42
1:A:267:ASP:OD1	1:A:267:ASP:N	2.47	0.42
1:A:124:GLN:C	1:A:126:LEU:N	2.73	0.42
1:B:243:ILE:HG23	1:B:366:LEU:HD21	2.01	0.42
1:B:60:LYS:HG2	1:B:121:THR:OG1	2.19	0.42
1:A:179:LEU:HB2	1:A:268:THR:HG22	2.02	0.42
1:A:288:LEU:CD2	1:A:318:LEU:HD22	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ARG:NH1	1:A:55:ARG:CG	2.78	0.42
1:A:361:ARG:HH21	1:A:365:ILE:CG1	2.32	0.42
1:B:211:THR:HG21	5:B:398:HOH:O	2.19	0.42
1:A:231:ILE:CG2	1:A:357:ASN:HD21	2.33	0.42
1:A:205:PRO:O	1:B:353:ARG:HD3	2.20	0.42
1:B:15:VAL:HG13	1:B:15:VAL:O	2.19	0.42
1:A:171:ARG:HA	1:A:194:LYS:O	2.19	0.42
1:B:219:LEU:H	1:B:219:LEU:HD22	1.80	0.41
1:A:243:ILE:HG23	1:A:366:LEU:HD21	2.02	0.41
1:A:60:LYS:HG2	1:A:121:THR:OG1	2.20	0.41
1:A:273:PRO:HD2	1:A:306:TYR:HB2	2.02	0.41
1:B:164:ARG:HD3	1:B:167:ARG:HD2	2.02	0.41
1:B:19:ILE:HG22	1:B:19:ILE:O	2.20	0.41
1:B:124:GLN:C	1:B:126:LEU:N	2.73	0.41
1:B:288:LEU:CD2	1:B:318:LEU:HD22	2.50	0.41
1:B:119:THR:HG23	1:B:301:LEU:HD21	2.01	0.41
1:A:164:ARG:HD3	1:A:167:ARG:HD2	2.02	0.41
1:A:19:ILE:HG22	1:A:19:ILE:O	2.21	0.41
1:B:37:ARG:NH2	1:B:303:GLY:O	2.54	0.41
1:B:191:PHE:CZ	1:B:219:LEU:HD23	2.56	0.41
1:A:234:GLY:HA3	1:A:355:ASP:HA	2.01	0.41
1:B:171:ARG:HB3	1:B:255:PHE:CZ	2.55	0.41
1:B:34:ILE:HG21	1:B:37:ARG:HH11	1.85	0.41
1:A:169:PHE:CD2	1:A:258:LYS:HB3	2.42	0.41
1:A:152:PHE:CE2	4:B:390:B3N:H221	2.56	0.41
1:B:293:GLN:C	1:B:295:GLN:OE1	2.58	0.41
1:A:92:ASP:O	1:A:92:ASP:CG	2.59	0.41
1:A:51:HIS:CE1	1:A:52:LYS:HE2	2.55	0.41
1:B:129:GLY:C	1:B:131:CYS:H	2.24	0.41
1:B:353:ARG:HB2	1:B:354:PRO:HD2	2.01	0.41
1:B:243:ILE:O	1:B:247:VAL:HG23	2.20	0.41
1:B:178:ASP:HB2	1:B:263:GLN:OE1	2.20	0.41
1:A:171:ARG:HB3	1:A:255:PHE:CE2	2.56	0.41
1:B:140:GLY:O	1:B:142:HIS:N	2.54	0.41
1:B:51:HIS:CE1	1:B:52:LYS:HE2	2.56	0.41
1:A:45:ILE:HA	1:A:316:THR:HG23	2.02	0.41
1:B:145:LYS:HB2	1:B:148:GLU:HB3	2.03	0.41
1:A:120:ILE:O	1:A:123:ALA:HB3	2.21	0.41
1:A:244:CYS:HG	1:A:287:CYS:CB	2.33	0.41
1:A:363:GLN:HG3	1:A:367:ASN:HD21	1.86	0.41
1:B:179:LEU:HB2	1:B:268:THR:HG22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:TYR:CE1	1:A:25:VAL:HG21	2.56	0.40
1:B:120:ILE:O	1:B:123:ALA:HB3	2.20	0.40
1:A:67:MET:C	1:A:69:THR:N	2.75	0.40
1:A:243:ILE:O	1:A:247:VAL:HG23	2.21	0.40
1:B:171:ARG:HB3	1:B:255:PHE:CE2	2.56	0.40
1:A:353:ARG:HB3	1:B:205:PRO:HD2	2.03	0.40
1:A:179:LEU:CD2	1:A:274:MET:HG2	2.48	0.40
1:B:67:MET:C	1:B:69:THR:N	2.74	0.40
1:A:171:ARG:HB3	1:A:255:PHE:CZ	2.56	0.40
1:B:141:TRP:HE3	1:B:153:CYS:HG	1.68	0.40
1:B:273:PRO:HD2	1:B:306:TYR:HB2	2.02	0.40
1:B:231:ILE:CG2	1:B:357:ASN:HD21	2.35	0.40
1:B:234:GLY:O	1:B:235:ILE:C	2.60	0.40
1:A:19:ILE:HG12	1:A:135:ILE:O	2.21	0.40
1:A:293:GLN:C	1:A:295:GLN:OE1	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	353/389 (91%)	287 (81%)	57 (16%)	9 (2%)	7	41
1	B	352/389 (90%)	288 (82%)	56 (16%)	8 (2%)	8	44
All	All	705/778 (91%)	575 (82%)	113 (16%)	17 (2%)	7	43

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	SER
1	A	354	PRO
1	A	356	ARG

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Mol	Chain	Res	Type
1	B	21	SER
1	B	354	PRO
1	B	356	ARG
1	A	141	TRP
1	A	254	ALA
1	A	375	HIS
1	B	141	TRP
1	B	254	ALA
1	B	375	HIS
1	A	374	LYS
1	B	374	LYS
1	A	131	CYS
1	B	131	CYS
1	A	94	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	299/326 (92%)	280 (94%)	19 (6%)	22	62
1	B	298/326 (91%)	279 (94%)	19 (6%)	22	62
All	All	597/652 (92%)	559 (94%)	38 (6%)	22	62

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	SER
1	A	26	SER
1	A	36	LYS
1	A	65	GLU
1	A	81	LYS
1	A	111	TYR
1	A	126	LEU
1	A	128	ASP
1	A	142	HIS

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Mol	Chain	Res	Type
1	A	219	LEU
1	A	253	GLN
1	A	256	ASN
1	A	293	GLN
1	A	296	LEU
1	A	321	VAL
1	A	345	VAL
1	A	361	ARG
1	A	368	TYR
1	A	376	VAL
1	B	21	SER
1	B	26	SER
1	B	36	LYS
1	B	65	GLU
1	B	81	LYS
1	B	111	TYR
1	B	126	LEU
1	B	128	ASP
1	B	142	HIS
1	B	219	LEU
1	B	253	GLN
1	B	256	ASN
1	B	293	GLN
1	B	296	LEU
1	B	321	VAL
1	B	345	VAL
1	B	361	ARG
1	B	368	TYR
1	B	376	VAL

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	53	GLN
1	A	80	GLN
1	A	84	GLN
1	A	242	GLN
1	A	256	ASN
1	A	363	GLN
1	A	367	ASN
1	A	372	ASN
1	A	375	HIS

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Mol	Chain	Res	Type
1	B	53	GLN
1	B	80	GLN
1	B	84	GLN
1	B	242	GLN
1	B	256	ASN
1	B	363	GLN
1	B	367	ASN
1	B	372	ASN
1	B	375	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 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	$\# Z > 2$
4	B3N	A	390	2	22,22,22	2.90	9 (40%)	26,27,27	1.13	2 (7%)
4	B3N	B	390	2	22,22,22	2.97	9 (40%)	26,27,27	1.14	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
4	B3N	A	390	2	-	0/20/20/20	0/1/1/1
4	B3N	B	390	2	-	0/20/20/20	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	390	B3N	O4-N3	-8.93	1.23	1.39
4	A	390	B3N	O4-N3	-8.27	1.24	1.39
4	B	390	B3N	O23-C13	2.07	1.27	1.23
4	B	390	B3N	C19-C18	2.08	1.42	1.38
4	A	390	B3N	C19-C18	2.23	1.42	1.38
4	A	390	B3N	O23-C13	2.31	1.28	1.23
4	A	390	B3N	C17-N20	2.91	1.44	1.37
4	B	390	B3N	C17-N20	3.00	1.44	1.37
4	B	390	B3N	C18-C17	3.16	1.45	1.39
4	A	390	B3N	C18-C17	3.47	1.46	1.39
4	A	390	B3N	C16-C15	3.72	1.45	1.38
4	A	390	B3N	C15-C14	3.97	1.46	1.39
4	B	390	B3N	C15-C14	4.02	1.46	1.39
4	B	390	B3N	C19-C14	4.02	1.46	1.39
4	B	390	B3N	C16-C15	4.05	1.46	1.38
4	A	390	B3N	C19-C14	4.75	1.47	1.39
4	A	390	B3N	C16-C17	5.24	1.49	1.39
4	B	390	B3N	C16-C17	5.67	1.50	1.39

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	390	B3N	O2-C1-N3	-2.57	120.49	123.53
4	B	390	B3N	C18-C17-N20	-2.55	118.00	121.64
4	A	390	B3N	C18-C17-N20	-2.16	118.55	121.64
4	B	390	B3N	O2-C1-N3	-2.11	121.04	123.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	390	B3N	4	0
4	B	390	B3N	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	357/389 (91%)	-0.08	2 (0%) 90 84	45, 81, 109, 141	15 (4%)
1	B	356/389 (91%)	-0.06	1 (0%) 94 93	44, 81, 109, 127	15 (4%)
All	All	713/778 (91%)	-0.07	3 (0%) 93 90	44, 81, 109, 141	30 (4%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	92	ASP	2.7
1	A	373	LEU	2.5
1	B	373	LEU	2.2

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	B3N	B	390	22/22	0.95	0.24	0.38	73,86,97,98	0
4	B3N	A	390	22/22	0.96	0.21	-0.13	60,72,76,76	0
3	K	A	402	1/1	0.91	0.13	-1.31	53,53,53,53	0
3	K	A	404	1/1	0.90	0.11	-1.90	67,67,67,67	0
2	CO	A	400	1/1	0.99	0.12	-2.27	74,74,74,74	0
3	K	B	403	1/1	0.93	0.08	-2.99	53,53,53,53	0
2	CO	B	401	1/1	0.97	0.10	-3.14	70,70,70,70	0
3	K	B	405	1/1	0.97	0.06	-3.34	72,72,72,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.