



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:12 AM GMT

PDB ID : 2G47
Title : Crystal structure of human insulin-degrading enzyme in complex with amyloid-beta (1-40)
Authors : Shen, Y.; Tang, W.-J.
Deposited on : 2006-02-21
Resolution : 2.10 Å(reported)

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

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

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

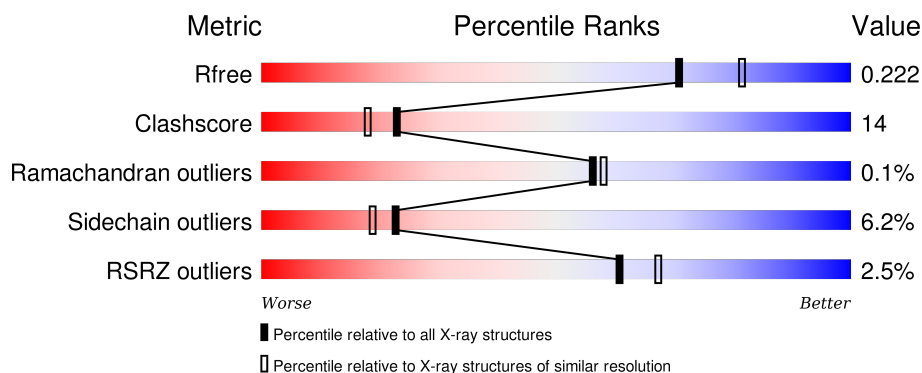
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	990	<div> <div>2%</div> <div>70%</div> <div>25%</div> <div>• •</div> </div>
1	B	990	<div> <div>2%</div> <div>71%</div> <div>24%</div> <div>• •</div> </div>
2	C	40	<div> <div>10%</div> <div>15%</div> <div>10%</div> <div>•</div> <div>73%</div> </div>
2	D	40	<div> <div>8%</div> <div>13%</div> <div>13%</div> <div>•</div> <div>73%</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
3	DIO	A	2000	-	-	-	X
3	DIO	B	2001	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16962 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 Insulin-degrading enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	964	Total	C	N	O	S	0	0	0
			7853	5053	1319	1447	34			
1	B	965	Total	C	N	O	S	0	0	0
			7855	5055	1319	1447	34			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MET	-	INITIATING METHIONINE	UNP P14735
A	31	HIS	-	EXPRESSION TAG	UNP P14735
A	32	HIS	-	EXPRESSION TAG	UNP P14735
A	33	HIS	-	EXPRESSION TAG	UNP P14735
A	34	HIS	-	EXPRESSION TAG	UNP P14735
A	35	HIS	-	EXPRESSION TAG	UNP P14735
A	36	HIS	-	EXPRESSION TAG	UNP P14735
A	37	ALA	-	CLONING ARTIFACT	UNP P14735
A	38	ALA	-	CLONING ARTIFACT	UNP P14735
A	39	GLY	-	CLONING ARTIFACT	UNP P14735
A	40	ILE	-	CLONING ARTIFACT	UNP P14735
A	41	PRO	-	CLONING ARTIFACT	UNP P14735
A	111	GLN	GLU	ENGINEERED	UNP P14735
B	30	MET	-	INITIATING METHIONINE	UNP P14735
B	31	HIS	-	EXPRESSION TAG	UNP P14735
B	32	HIS	-	EXPRESSION TAG	UNP P14735
B	33	HIS	-	EXPRESSION TAG	UNP P14735
B	34	HIS	-	EXPRESSION TAG	UNP P14735
B	35	HIS	-	EXPRESSION TAG	UNP P14735
B	36	HIS	-	EXPRESSION TAG	UNP P14735
B	37	ALA	-	CLONING ARTIFACT	UNP P14735
B	38	ALA	-	CLONING ARTIFACT	UNP P14735
B	39	GLY	-	CLONING ARTIFACT	UNP P14735
B	40	ILE	-	CLONING ARTIFACT	UNP P14735
B	41	PRO	-	CLONING ARTIFACT	UNP P14735

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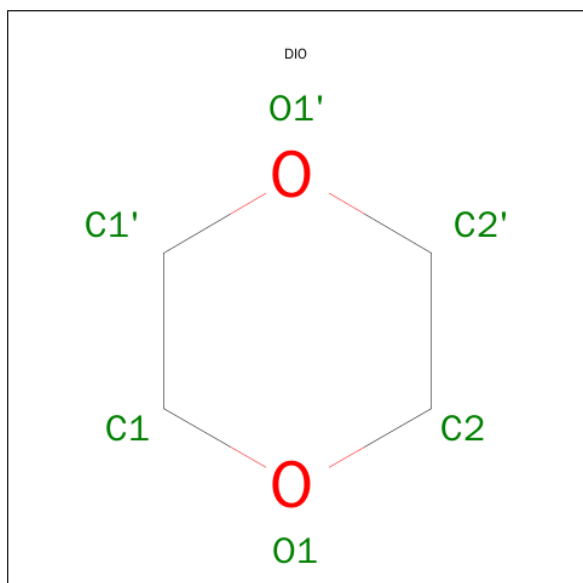
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Chain	Residue	Modelled	Actual	Comment	Reference
B	111	GLN	GLU	ENGINEERED	UNP P14735

- Molecule 2 is a protein called amyloid protein beta A4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	11	Total	C	N	O	0	0	0
			86	57	12	17			
2	D	11	Total	C	N	O	0	0	0
			90	59	12	19			

- Molecule 3 is 1,4-DIETHYLENE DIOXIDE (three-letter code: DIO) (formula: C₄H₈O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			6	4	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	564	Total	O	0	0
			564	564		
4	B	497	Total	O	0	0
			497	497		

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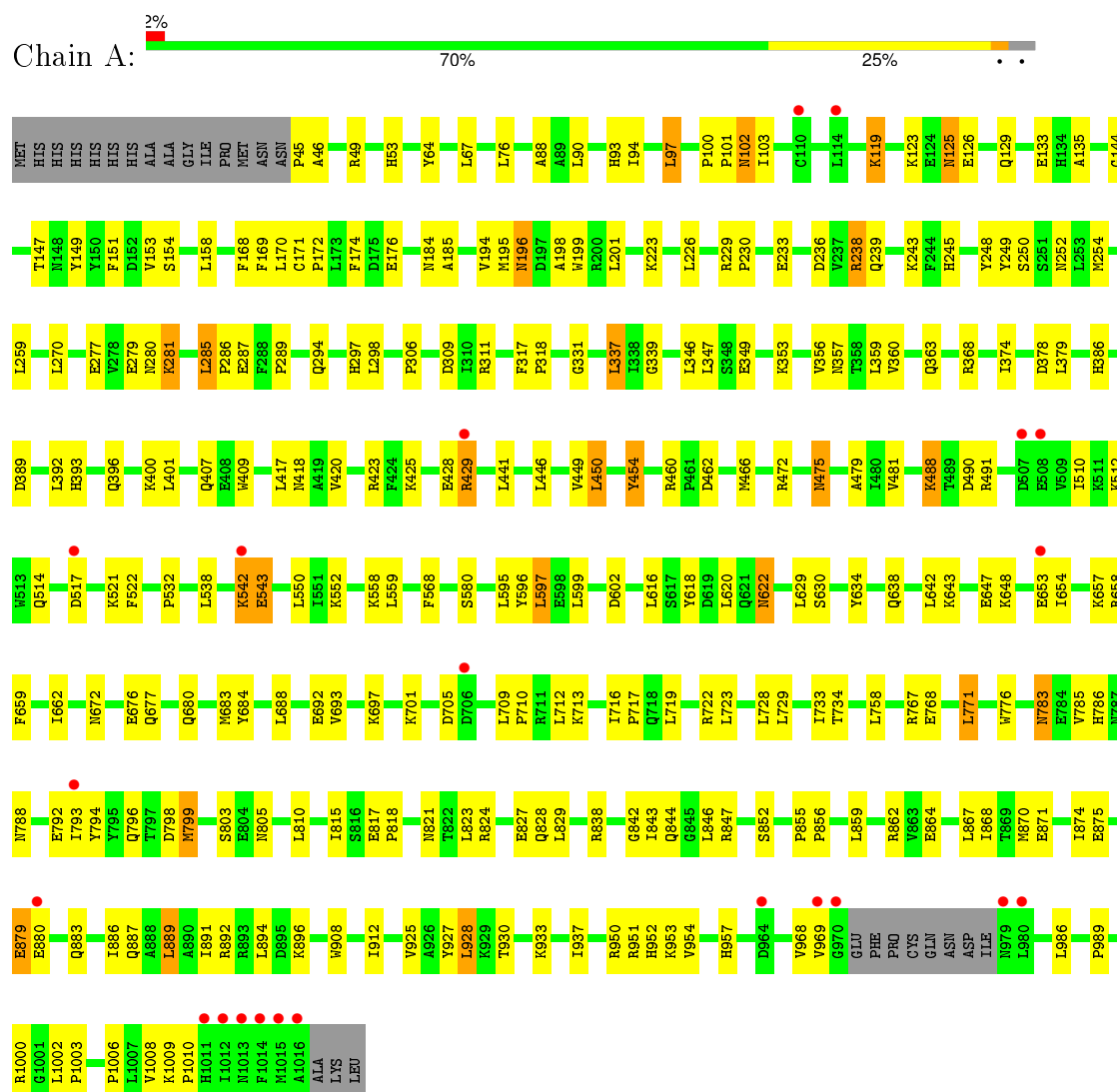
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	2	Total	O	0	0
			2	2		
4	D	3	Total	O	0	0
			3	3		

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: Insulin-degrading enzyme



• Molecule 1: Insulin-degrading enzyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	262.43 Å 262.43 Å 90.71 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.77 – 2.10 29.77 – 2.10	Depositor EDS
% Data completeness (in resolution range)	90.5 (29.77-2.10) 90.6 (29.77-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.10 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.203 , 0.223 0.202 , 0.222	Depositor DCC
R_{free} test set	18633 reflections (9.95%)	DCC
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.3	EDS
Estimated twinning fraction	0.018 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 194459 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16962	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.78% 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: DIO

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	0.38	0/8049	0.60	0/10887
1	B	0.38	0/8051	0.64	2/10891 (0.0%)
2	C	1.32	1/86 (1.2%)	1.19	1/113 (0.9%)
2	D	1.65	1/90 (1.1%)	1.31	1/118 (0.8%)
All	All	0.41	2/16276 (0.0%)	0.63	4/22009 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	18	VAL	CB-CG2	-5.76	1.40	1.52
2	D	18	VAL	CB-CG2	-5.03	1.42	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	23	ASP	CB-CG-OD2	8.15	125.64	118.30
1	B	67	LEU	CA-CB-CG	7.30	132.08	115.30
2	C	23	ASP	N-CA-C	5.25	125.18	111.00
1	B	259	LEU	CA-CB-CG	5.04	126.89	115.30

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	7853	0	7773	225	1
1	B	7855	0	7767	197	1
2	C	86	0	78	3	0
2	D	90	0	82	8	0
3	A	6	0	8	1	0
3	B	6	0	8	2	0
4	A	564	0	0	28	0
4	B	497	0	0	24	0
4	C	2	0	0	0	0
4	D	3	0	0	0	0
All	All	16962	0	15716	427	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 14.

All (427) 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:782:ARG:NH2	1:B:961:ARG:O	1.65	1.28
1:B:892:ARG:HD2	4:B:2433:HOH:O	1.50	1.12
2:D:17:LEU:O	2:D:17:LEU:HD23	1.62	0.99
1:B:208:THR:HG23	1:B:477:ARG:HH22	1.28	0.96
1:B:817:GLU:HG3	1:B:818:PRO:HD3	1.49	0.94
2:D:17:LEU:H	2:D:17:LEU:HD22	1.32	0.94
1:A:102:ASN:HD22	1:A:102:ASN:H	1.14	0.94
1:B:864:GLU:HG2	1:B:986:LEU:HD21	1.50	0.93
1:B:674:ARG:HD3	4:B:2199:HOH:O	1.69	0.92
2:D:17:LEU:C	2:D:17:LEU:HD23	1.86	0.92
1:A:705:ASP:HB2	4:A:2459:HOH:O	1.68	0.92
1:A:879:GLU:HG2	4:A:2307:HOH:O	1.69	0.91
1:A:425:LYS:HE2	1:A:428:GLU:OE2	1.71	0.89
1:A:309:ASP:H	1:A:672:ASN:HD21	1.17	0.88
1:A:512:LYS:NZ	4:A:2554:HOH:O	2.07	0.88
1:A:287:GLU:HG2	1:A:289:PRO:HD3	1.55	0.88
2:D:17:LEU:N	2:D:17:LEU:HD22	1.85	0.87
1:A:864:GLU:HG2	1:A:986:LEU:HD21	1.58	0.86
1:A:441:LEU:HD23	1:A:449:VAL:HG11	1.55	0.86
1:B:309:ASP:H	1:B:672:ASN:HD21	1.19	0.85
1:A:874:ILE:O	1:A:933:LYS:HE3	1.78	0.84
1:B:521:LYS:NZ	4:B:2136:HOH:O	2.11	0.84
1:B:102:ASN:H	1:B:102:ASN:HD22	1.24	0.83
1:B:782:ARG:NH2	1:B:961:ARG:C	2.33	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:GLU:OE2	4:B:2440:HOH:O	1.97	0.81
1:A:1000:ARG:HD2	4:B:2259:HOH:O	1.79	0.81
1:A:123:LYS:HB3	1:A:126:GLU:HB2	1.62	0.80
1:B:950:ARG:HD2	4:B:2213:HOH:O	1.82	0.80
1:A:771:LEU:HD21	1:A:954:VAL:HG23	1.65	0.79
1:A:599:LEU:HD23	1:A:662:ILE:HD12	1.62	0.79
1:A:827:GLU:OE1	1:A:862:ARG:HD3	1.82	0.79
1:A:174:PHE:O	1:A:238:ARG:HD3	1.84	0.78
1:A:309:ASP:H	1:A:672:ASN:ND2	1.82	0.78
1:A:815:ILE:HA	1:A:870:MET:HE2	1.66	0.78
1:B:309:ASP:H	1:B:672:ASN:ND2	1.83	0.77
1:B:791:ILE:HD11	1:B:793:ILE:HD11	1.64	0.76
1:B:783:ASN:ND2	1:B:786:HIS:H	1.83	0.76
1:A:558:LYS:NZ	4:A:2216:HOH:O	2.17	0.76
1:A:294:GLN:H	1:A:297:HIS:HD2	1.34	0.76
2:D:17:LEU:N	2:D:17:LEU:CD2	2.49	0.75
1:B:658:ARG:NH2	4:B:2456:HOH:O	2.18	0.75
1:B:224:TYR:CZ	1:B:229:ARG:HD2	2.22	0.75
1:B:208:THR:HG23	1:B:477:ARG:NH2	2.02	0.74
1:A:558:LYS:NZ	1:A:558:LYS:HB2	2.02	0.74
1:B:294:GLN:H	1:B:297:HIS:HD2	1.33	0.74
1:A:135:ALA:HA	1:A:892:ARG:NH2	2.02	0.74
1:B:429:ARG:HD2	4:B:2318:HOH:O	1.87	0.74
1:A:102:ASN:HD22	1:A:102:ASN:N	1.85	0.74
1:B:425:LYS:HE2	1:B:428:GLU:OE2	1.87	0.74
1:A:821:ASN:HB2	4:A:2008:HOH:O	1.89	0.73
1:A:446:LEU:HD22	1:A:446:LEU:H	1.52	0.72
1:A:339:GLY:O	2:C:1:ASP:HA	1.90	0.72
1:B:886:ILE:HG23	1:B:928:LEU:HD13	1.70	0.72
1:B:591:ASN:O	1:B:595:LEU:HD23	1.90	0.71
1:B:299:LYS:HD2	1:B:510:ILE:HD13	1.71	0.71
1:A:491:ARG:HD2	4:A:2179:HOH:O	1.89	0.71
1:A:886:ILE:HG23	1:A:928:LEU:HD22	1.72	0.71
1:B:886:ILE:HG23	1:B:928:LEU:CD1	2.20	0.71
1:B:883:GLN:NE2	4:B:2432:HOH:O	2.23	0.71
1:B:94:ILE:HD11	1:B:244:PHE:CZ	2.26	0.70
1:B:49:ARG:NH2	1:B:446:LEU:HD23	2.07	0.70
1:A:53:HIS:H	1:A:53:HIS:CD2	2.07	0.70
1:B:760:PRO:HA	1:B:763:LEU:HD23	1.74	0.69
1:B:423:ARG:NE	4:B:2440:HOH:O	2.17	0.69
1:A:643:LYS:NZ	4:A:2560:HOH:O	2.21	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:968:VAL:O	1:A:969:VAL:HG22	1.93	0.69
1:B:817:GLU:HG3	1:B:818:PRO:CD	2.21	0.69
1:B:236:ASP:OD1	4:B:2074:HOH:O	2.12	0.68
1:A:93:HIS:HE1	1:A:368:ARG:HH21	1.41	0.68
1:A:102:ASN:ND2	1:A:102:ASN:H	1.90	0.67
1:A:446:LEU:CD2	4:A:2165:HOH:O	2.42	0.67
1:A:622:ASN:H	1:A:622:ASN:HD22	1.42	0.67
1:B:125:ASN:H	1:B:125:ASN:HD22	1.42	0.67
1:A:805:ASN:HD22	1:A:844:GLN:HE22	1.43	0.67
1:B:93:HIS:HE1	1:B:368:ARG:HH21	1.43	0.66
1:B:538:LEU:H	1:B:732:ASN:HD21	1.43	0.66
1:A:368:ARG:HD2	4:A:2150:HOH:O	1.94	0.66
1:A:538:LEU:HD12	1:A:538:LEU:N	2.10	0.66
1:A:337:LEU:HD23	1:A:401:LEU:HD22	1.78	0.65
1:B:102:ASN:N	1:B:102:ASN:HD22	1.90	0.65
1:B:94:ILE:HG12	1:B:95:GLY:N	2.12	0.65
1:A:783:ASN:ND2	1:A:785:VAL:H	1.95	0.65
1:B:229:ARG:HB3	1:B:230:PRO:HD3	1.78	0.64
1:A:88:ALA:HB3	1:A:151:PHE:CE2	2.32	0.64
1:A:783:ASN:ND2	1:A:786:HIS:H	1.96	0.64
1:B:84:ASP:OD2	1:B:896:LYS:HG2	1.97	0.64
1:A:184:ASN:HD21	1:A:223:LYS:NZ	1.96	0.64
1:B:196:ASN:HD22	1:B:199:TRP:H	1.43	0.64
1:B:815:ILE:HG22	1:B:870:MET:HE2	1.80	0.63
1:A:233:GLU:OE1	4:A:2534:HOH:O	2.15	0.62
1:A:950:ARG:HD2	4:A:2278:HOH:O	1.98	0.62
1:B:192:LYS:HD3	1:B:193:ASN:OD1	1.99	0.62
1:A:709:LEU:HB3	1:A:710:PRO:HD3	1.80	0.62
1:B:874:ILE:O	1:B:933:LYS:HD2	2.00	0.62
1:B:357:ASN:HB2	1:B:378:ASP:OD1	2.00	0.61
1:B:102:ASN:H	1:B:102:ASN:ND2	1.98	0.61
1:A:407:GLN:HG3	1:A:409:TRP:CD1	2.35	0.61
1:B:622:ASN:H	1:B:622:ASN:HD22	1.45	0.61
1:B:423:ARG:NH2	4:B:2440:HOH:O	2.27	0.61
1:B:126:GLU:HG3	4:B:2336:HOH:O	2.00	0.61
1:B:449:VAL:HG13	1:B:450:LEU:HD13	1.81	0.61
1:A:298:LEU:HD21	1:A:318:PRO:HG2	1.83	0.61
1:B:294:GLN:H	1:B:297:HIS:CD2	2.18	0.60
1:B:716:ILE:HB	1:B:717:PRO:HD3	1.83	0.60
1:A:783:ASN:HD22	1:A:785:VAL:H	1.50	0.60
1:A:1002:LEU:C	1:B:1006:PRO:HB3	2.22	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:ARG:NH1	1:A:462:ASP:OD2	2.35	0.60
1:A:542:LYS:NZ	1:A:542:LYS:HB2	2.17	0.60
1:A:97:LEU:HB2	1:A:144:GLY:O	2.01	0.59
1:A:517:ASP:HB2	4:A:2189:HOH:O	2.02	0.59
1:B:205:GLU:O	1:B:208:THR:HG22	2.01	0.59
1:A:490:ASP:OD1	1:A:491:ARG:HG3	2.02	0.59
1:A:968:VAL:HG23	1:A:969:VAL:HG23	1.84	0.59
1:B:679:HIS:HD2	1:B:851:GLN:OE1	1.85	0.59
2:D:17:LEU:H	2:D:17:LEU:CD2	2.09	0.59
1:B:285:LEU:HD13	1:B:286:PRO:O	2.02	0.59
1:A:684:TYR:OH	1:A:697:LYS:HG2	2.03	0.58
1:B:94:ILE:HD11	1:B:244:PHE:HZ	1.68	0.58
1:B:100:PRO:HG2	1:B:103:ILE:HB	1.84	0.58
1:A:716:ILE:HB	1:A:717:PRO:HD3	1.85	0.58
1:A:236:ASP:CG	1:A:239:GLN:HG2	2.23	0.58
1:B:805:ASN:HD22	1:B:844:GLN:HE22	1.51	0.58
1:A:889:LEU:HB3	1:A:928:LEU:HD11	1.84	0.58
1:A:168:PHE:O	1:A:172:PRO:HG3	2.04	0.58
1:B:999:LYS:NZ	4:B:2415:HOH:O	2.36	0.58
1:A:176:GLU:OE2	1:A:238:ARG:HG3	2.03	0.57
1:A:331:GLY:HA3	1:A:363:GLN:OE1	2.04	0.57
1:A:968:VAL:C	1:A:969:VAL:CG2	2.72	0.57
1:A:100:PRO:HG2	1:A:103:ILE:HB	1.85	0.57
1:B:479:ALA:HB2	3:B:2001:DIO:H12	1.85	0.57
1:B:49:ARG:HH22	1:B:446:LEU:HD23	1.67	0.57
1:B:709:LEU:N	1:B:710:PRO:HD2	2.20	0.57
1:A:475:ASN:ND2	1:A:475:ASN:C	2.59	0.57
1:B:196:ASN:HD21	1:B:198:ALA:HB3	1.70	0.56
1:A:883:GLN:NE2	4:A:2042:HOH:O	2.38	0.56
1:B:299:LYS:HD2	1:B:510:ILE:CD1	2.34	0.56
1:A:125:ASN:HD22	1:A:125:ASN:H	1.52	0.56
1:A:658:ARG:NH2	4:A:2147:HOH:O	2.39	0.56
1:A:558:LYS:HZ2	1:A:558:LYS:HB2	1.69	0.56
1:B:898:LYS:HE3	4:B:2040:HOH:O	2.06	0.56
1:B:402:ARG:HG2	1:B:468:LEU:HD13	1.88	0.56
1:A:223:LYS:HE2	4:A:2387:HOH:O	2.05	0.56
1:A:429:ARG:HB3	1:A:429:ARG:HH11	1.71	0.56
1:B:602:ASP:OD1	1:B:658:ARG:HD3	2.06	0.56
1:A:597:LEU:HG	1:A:620:LEU:HG	1.88	0.55
1:B:868:ILE:HD12	1:B:984:PRO:HD3	1.87	0.55
1:A:429:ARG:NH1	4:A:2033:HOH:O	2.38	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:LEU:CD2	1:A:449:VAL:HG11	2.31	0.55
1:B:49:ARG:NH2	4:B:2269:HOH:O	2.18	0.55
1:A:908:TRP:CZ2	1:A:912:ILE:HD11	2.42	0.55
1:A:722:ARG:HB3	1:A:758:LEU:HD13	1.87	0.55
1:A:767:ARG:NH1	1:A:1006:PRO:HA	2.22	0.55
1:B:339:GLY:O	2:D:1:ASP:HA	2.07	0.55
1:A:874:ILE:HG22	1:A:937:ILE:HD11	1.88	0.54
1:A:927:TYR:O	1:A:930:THR:HB	2.06	0.54
1:B:673:PHE:CD1	1:B:697:LYS:HE3	2.42	0.54
1:A:446:LEU:HD22	4:A:2165:HOH:O	2.03	0.54
1:B:654:ILE:HD13	1:B:712:LEU:HD13	1.89	0.54
1:A:950:ARG:NH1	4:A:2289:HOH:O	2.41	0.54
1:A:580:SER:HB2	1:A:723:LEU:HD23	1.90	0.54
1:B:961:ARG:HG2	1:B:961:ARG:O	2.06	0.54
1:A:538:LEU:CD1	1:A:538:LEU:N	2.71	0.54
1:A:803:SER:HA	1:A:927:TYR:CE2	2.43	0.54
1:A:243:LYS:NZ	1:A:243:LYS:HB2	2.22	0.54
1:A:692:GLU:HG2	1:A:693:VAL:HG23	1.90	0.54
1:A:407:GLN:HG3	1:A:409:TRP:NE1	2.23	0.54
1:B:191:GLU:OE2	4:B:2348:HOH:O	2.19	0.54
1:A:184:ASN:HD21	1:A:223:LYS:HZ1	1.54	0.53
1:A:771:LEU:HD21	1:A:954:VAL:CG2	2.35	0.53
1:B:461:PRO:O	1:B:465:GLU:HG3	2.08	0.53
1:A:868:ILE:O	1:A:871:GLU:HB3	2.08	0.53
1:A:558:LYS:HZ3	1:A:558:LYS:HB2	1.73	0.53
4:B:2406:HOH:O	2:D:1:ASP:HB2	2.07	0.53
1:B:622:ASN:H	1:B:622:ASN:ND2	2.07	0.53
1:A:185:ALA:HB2	1:A:828:GLN:HE22	1.74	0.53
1:B:93:HIS:CE1	1:B:368:ARG:HH21	2.25	0.53
1:A:887:GLN:OE1	4:A:2041:HOH:O	2.19	0.53
1:B:259:LEU:C	1:B:259:LEU:HD12	2.29	0.53
1:A:1006:PRO:HB3	1:B:1002:LEU:C	2.28	0.52
1:B:368:ARG:HD3	4:B:2313:HOH:O	2.09	0.52
1:A:475:ASN:ND2	1:A:475:ASN:O	2.43	0.52
1:A:119:LYS:HB2	1:A:171:CYS:HB2	1.91	0.52
1:B:147:THR:HG22	1:B:149:TYR:CE1	2.45	0.52
1:A:93:HIS:HE1	1:A:368:ARG:NH2	2.08	0.52
1:B:815:ILE:HG22	1:B:870:MET:CE	2.40	0.52
1:B:392:LEU:O	1:B:396:GLN:HG3	2.10	0.52
1:B:879:GLU:HB2	4:B:2488:HOH:O	2.09	0.52
1:A:285:LEU:HD23	1:A:286:PRO:HD2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:LEU:HB2	1:B:144:GLY:O	2.09	0.52
1:B:49:ARG:HG2	1:B:68:GLU:HB3	1.91	0.51
1:A:843:ILE:HG22	1:A:844:GLN:H	1.73	0.51
1:B:94:ILE:HG21	1:B:248:TYR:HB3	1.93	0.51
1:A:317:PHE:HD2	1:A:475:ASN:ND2	2.09	0.51
1:B:474:GLU:C	1:B:475:ASN:HD22	2.14	0.51
1:A:349:GLU:CD	1:A:521:LYS:HD3	2.31	0.51
1:B:309:ASP:N	1:B:672:ASN:HD21	1.99	0.51
1:A:653:GLU:N	4:A:2250:HOH:O	2.27	0.51
1:A:677:GLN:HB2	1:A:680:GLN:HG3	1.92	0.51
1:B:978:ILE:O	1:B:979:ASN:CB	2.59	0.51
1:B:782:ARG:CZ	1:B:961:ARG:O	2.51	0.51
1:A:53:HIS:HB3	4:A:2366:HOH:O	2.10	0.51
1:B:760:PRO:HA	1:B:763:LEU:CD2	2.41	0.51
1:B:616:LEU:HD21	1:B:638:GLN:HG2	1.93	0.50
1:A:676:GLU:OE2	1:A:676:GLU:HA	2.10	0.50
1:B:490:ASP:OD1	1:B:491:ARG:HG3	2.11	0.50
1:B:877:MET:HG3	1:B:881:ALA:HB3	1.93	0.50
1:B:238:ARG:O	1:B:242:LEU:HD23	2.12	0.50
1:A:194:VAL:HG12	1:A:195:MET:HE1	1.94	0.50
1:A:196:ASN:HD21	1:A:198:ALA:HB3	1.76	0.50
1:A:510:ILE:O	1:A:514:GLN:HG3	2.11	0.50
1:A:294:GLN:H	1:A:297:HIS:CD2	2.20	0.50
1:A:194:VAL:HG12	1:A:195:MET:CE	2.42	0.50
1:B:208:THR:HG21	1:B:477:ARG:HH12	1.77	0.50
1:A:968:VAL:O	1:A:969:VAL:CG2	2.59	0.49
1:B:968:VAL:HG23	1:B:969:VAL:HG22	1.92	0.49
1:B:733:ILE:HD13	1:B:733:ILE:N	2.28	0.49
1:A:879:GLU:OE1	1:A:879:GLU:HA	2.13	0.49
1:B:783:ASN:ND2	1:B:785:VAL:H	2.10	0.49
1:A:374:ILE:HD11	2:C:3:GLU:HG3	1.94	0.49
1:A:90:LEU:HD13	1:A:169:PHE:CE2	2.48	0.49
1:A:538:LEU:H	1:A:538:LEU:CD1	2.25	0.49
1:A:722:ARG:HH11	1:A:722:ARG:HG3	1.77	0.49
1:A:552:LYS:HB3	1:A:559:LEU:HB3	1.94	0.49
1:A:309:ASP:N	1:A:672:ASN:HD21	1.96	0.49
1:A:659:PHE:CE1	1:A:712:LEU:HD12	2.48	0.49
1:B:510:ILE:O	1:B:514:GLN:HG3	2.13	0.49
1:A:429:ARG:CB	1:A:429:ARG:HH11	2.24	0.49
1:A:135:ALA:HA	1:A:892:ARG:CZ	2.42	0.49
1:A:49:ARG:HH12	1:A:446:LEU:HD23	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:864:GLU:HG3	4:A:2301:HOH:O	2.13	0.48
1:A:862:ARG:NE	1:A:862:ARG:HA	2.28	0.48
1:B:291:HIS:ND1	1:B:292:PRO:HD2	2.28	0.48
1:A:602:ASP:OD1	1:A:658:ARG:HD3	2.14	0.48
1:A:359:LEU:HD23	1:A:360:VAL:N	2.28	0.48
1:A:94:ILE:HG13	1:A:248:TYR:HB3	1.94	0.48
1:A:543:GLU:H	1:A:543:GLU:CD	2.16	0.48
1:A:475:ASN:C	1:A:475:ASN:HD22	2.16	0.48
1:A:306:PRO:HB3	1:A:481:VAL:CG1	2.44	0.48
1:B:874:ILE:HG22	1:B:937:ILE:HD11	1.96	0.48
1:A:538:LEU:HD23	1:A:734:THR:HG23	1.94	0.47
1:B:799:MET:HE1	1:B:1008:VAL:HA	1.95	0.47
1:B:100:PRO:HB2	1:B:102:ASN:ND2	2.29	0.47
1:A:446:LEU:HD23	4:A:2165:HOH:O	2.09	0.47
1:A:311:ARG:HB3	1:A:379:LEU:HB2	1.96	0.47
1:B:224:TYR:CE2	1:B:229:ARG:HD2	2.48	0.47
1:A:968:VAL:C	1:A:969:VAL:HG23	2.34	0.47
1:A:532:PRO:HG3	1:A:634:TYR:CD2	2.49	0.47
1:A:259:LEU:HD23	1:A:259:LEU:C	2.35	0.47
1:A:67:LEU:N	1:A:67:LEU:HD23	2.29	0.47
1:A:776:TRP:HA	1:A:953:LYS:O	2.14	0.47
1:A:45:PRO:HG2	1:A:46:ALA:H	1.80	0.47
1:A:568:PHE:HA	4:A:2222:HOH:O	2.13	0.47
1:B:607:TYR:CE2	1:B:644:LYS:HG2	2.49	0.47
1:B:559:LEU:HD11	1:B:729:LEU:HD22	1.95	0.47
1:B:552:LYS:HB3	1:B:559:LEU:HB3	1.95	0.47
1:A:622:ASN:H	1:A:622:ASN:ND2	2.12	0.47
1:A:196:ASN:HD22	1:A:199:TRP:H	1.60	0.47
1:B:184:ASN:HD21	1:B:223:LYS:NZ	2.12	0.47
1:A:488:LYS:HD3	1:A:488:LYS:N	2.29	0.47
1:A:102:ASN:ND2	1:A:102:ASN:N	2.57	0.47
1:A:252:ASN:HB3	1:A:280:ASN:OD1	2.15	0.47
1:B:359:LEU:HD23	1:B:359:LEU:C	2.35	0.47
1:B:552:LYS:NZ	1:B:743:GLN:NE2	2.62	0.47
1:A:93:HIS:CE1	1:A:368:ARG:HH21	2.28	0.47
1:B:933:LYS:O	1:B:937:ILE:HG12	2.15	0.47
1:B:298:LEU:HD13	1:B:475:ASN:HA	1.96	0.47
1:A:64:TYR:CE2	1:A:450:LEU:HD21	2.49	0.47
1:B:552:LYS:HZ1	1:B:743:GLN:HE22	1.61	0.46
1:A:843:ILE:HG22	1:A:844:GLN:N	2.30	0.46
1:B:556:MET:O	1:B:556:MET:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:506:PRO:HG2	1:B:509:VAL:CG2	2.46	0.46
1:A:768:GLU:HB3	1:A:843:ILE:HG13	1.97	0.46
1:B:501:LYS:HE3	4:B:2166:HOH:O	2.15	0.46
1:B:638:GLN:HB2	1:B:639:PRO:HD3	1.98	0.46
1:B:597:LEU:HG	1:B:620:LEU:HG	1.97	0.46
1:B:93:HIS:HD2	1:B:145:GLU:O	1.98	0.46
1:B:404:GLU:HG2	1:B:524:LEU:CD2	2.46	0.46
1:B:371:MET:HE2	1:B:371:MET:HB3	1.73	0.46
1:B:102:ASN:N	1:B:102:ASN:ND2	2.61	0.46
1:B:285:LEU:HD22	1:B:286:PRO:HD2	1.98	0.46
1:B:843:ILE:HG22	1:B:844:GLN:N	2.31	0.46
1:B:311:ARG:HD2	1:B:379:LEU:O	2.16	0.46
1:B:67:LEU:HD13	1:B:75:VAL:HB	1.98	0.45
1:B:204:LEU:HD13	3:B:2001:DIO:H2'2	1.98	0.45
1:A:359:LEU:HD23	1:A:359:LEU:C	2.37	0.45
1:A:357:ASN:HB2	1:A:378:ASP:OD1	2.16	0.45
1:A:599:LEU:HD21	1:A:659:PHE:HA	1.99	0.45
1:B:722:ARG:HD2	1:B:756:LYS:HB2	1.98	0.45
1:A:151:PHE:C	1:A:151:PHE:CD1	2.89	0.45
1:A:49:ARG:HH22	1:A:446:LEU:HD21	1.81	0.45
1:A:349:GLU:O	1:A:353:LYS:HG3	2.16	0.45
1:B:529:GLU:O	1:B:637:LYS:HE3	2.17	0.45
1:A:129:GLN:O	1:A:133:GLU:HG3	2.17	0.45
1:A:618:TYR:HA	1:A:630:SER:O	2.16	0.45
1:A:776:TRP:CE3	1:A:989:PRO:HB3	2.52	0.45
1:B:787:ASN:OD1	1:B:962:GLU:HB2	2.17	0.45
1:B:224:TYR:OH	1:B:229:ARG:NH1	2.50	0.45
1:B:886:ILE:HG23	1:B:928:LEU:HD11	1.98	0.45
1:A:643:LYS:HE3	1:A:647:GLU:OE2	2.17	0.45
1:A:446:LEU:N	1:A:446:LEU:HD22	2.28	0.44
1:A:957:HIS:HD2	4:A:2281:HOH:O	2.00	0.44
1:B:824:ARG:O	1:B:828:GLN:HA	2.16	0.44
1:A:153:VAL:HG22	1:A:154:SER:N	2.32	0.44
1:A:815:ILE:HG22	1:A:870:MET:CE	2.47	0.44
1:B:94:ILE:HD12	1:B:248:TYR:CB	2.47	0.44
1:B:567:PHE:HE1	1:B:903:GLU:OE2	1.99	0.44
1:B:599:LEU:HD23	1:B:662:ILE:HD12	1.99	0.44
1:A:796:GLN:HB3	1:A:952:HIS:HB2	2.00	0.44
1:A:684:TYR:CZ	1:A:688:LEU:HD11	2.53	0.44
1:B:474:GLU:HG2	1:B:475:ASN:ND2	2.32	0.44
1:B:530:PHE:HA	1:B:637:LYS:HD2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:446:LEU:CD2	1:A:446:LEU:H	2.25	0.44
1:A:798:ASP:CG	1:A:799:MET:H	2.21	0.44
1:A:887:GLN:HE21	1:A:891:ILE:HD11	1.82	0.44
1:A:245:HIS:O	1:A:249:TYR:HB2	2.18	0.44
1:B:591:ASN:O	1:B:595:LEU:CD2	2.60	0.44
1:A:654:ILE:HD13	1:A:712:LEU:HD13	1.98	0.44
1:A:418:ASN:HB3	1:A:454:TYR:O	2.18	0.44
1:A:393:HIS:HE1	4:A:2144:HOH:O	2.00	0.44
1:A:64:TYR:HD2	1:A:76:LEU:HD11	1.83	0.43
1:A:374:ILE:HD12	1:A:374:ILE:C	2.38	0.43
1:B:810:LEU:HG	1:B:928:LEU:HD21	1.99	0.43
1:B:709:LEU:HG	1:B:713:LYS:HE3	2.00	0.43
1:A:799:MET:HE1	1:A:1006:PRO:HG2	2.00	0.43
1:A:346:LEU:HA	1:A:522:PHE:CE2	2.53	0.43
1:B:799:MET:SD	1:B:1006:PRO:HG2	2.58	0.43
1:A:908:TRP:CE2	1:A:912:ILE:HD11	2.54	0.43
1:A:871:GLU:O	1:A:875:GLU:HG3	2.18	0.43
1:B:90:LEU:C	1:B:90:LEU:HD23	2.39	0.43
1:A:229:ARG:HB3	1:A:230:PRO:HD3	2.01	0.43
1:B:934:GLU:HG2	1:B:938:LYS:HE2	2.01	0.43
1:B:950:ARG:CD	4:B:2213:HOH:O	2.54	0.43
1:A:196:ASN:ND2	1:A:199:TRP:H	2.17	0.43
1:B:776:TRP:CE3	1:B:989:PRO:HB3	2.53	0.43
1:B:542:LYS:HG2	1:B:543:GLU:OE2	2.19	0.43
1:B:400:LYS:HD3	4:B:2146:HOH:O	2.18	0.43
1:B:229:ARG:HB3	1:B:230:PRO:CD	2.48	0.43
1:B:123:LYS:HB3	1:B:126:GLU:HB2	2.01	0.43
1:B:54:ILE:HG22	1:B:450:LEU:HD22	2.01	0.43
1:B:767:ARG:NH1	1:B:1006:PRO:HA	2.34	0.43
1:A:196:ASN:ND2	1:A:198:ALA:H	2.16	0.43
1:A:719:LEU:HD13	1:A:719:LEU:C	2.39	0.43
1:A:654:ILE:HD11	1:A:712:LEU:HD22	2.00	0.43
1:B:88:ALA:HA	1:B:257:CYS:O	2.19	0.43
1:B:353:LYS:HD3	1:B:355:TRP:CZ3	2.54	0.43
1:B:309:ASP:O	1:B:668:ARG:HG2	2.19	0.42
1:A:596:TYR:CD2	1:A:597:LEU:HD13	2.54	0.42
1:A:158:LEU:O	1:A:158:LEU:HD23	2.19	0.42
1:B:460:ARG:HD2	1:B:462:ASP:OD2	2.19	0.42
2:C:16:LYS:HG2	2:C:16:LYS:HZ2	1.57	0.42
1:B:860:GLU:OE1	1:B:953:LYS:HE2	2.20	0.42
1:A:794:TYR:CE1	1:A:838:ARG:HD3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:542:LYS:HG3	1:A:543:GLU:HG3	2.01	0.42
1:A:472:ARG:HG2	1:A:472:ARG:HH11	1.84	0.42
1:A:683:MET:HA	1:A:792:GLU:OE2	2.19	0.42
1:A:712:LEU:HD23	1:A:712:LEU:O	2.20	0.42
1:A:542:LYS:HB2	1:A:542:LYS:HZ2	1.85	0.42
1:A:896:LYS:HD3	1:A:896:LYS:HA	1.84	0.42
1:B:208:THR:CG2	1:B:477:ARG:HH12	2.31	0.42
1:B:783:ASN:HD21	1:B:786:HIS:H	1.60	0.42
1:B:398:ILE:HG23	1:B:468:LEU:CD2	2.50	0.42
1:A:793:ILE:O	1:A:847:ARG:HA	2.20	0.42
1:B:306:PRO:HB3	1:B:481:VAL:CG1	2.49	0.42
1:B:446:LEU:HD23	4:B:2269:HOH:O	2.18	0.42
1:B:404:GLU:HG2	1:B:524:LEU:HD21	2.02	0.42
1:A:894:LEU:HG	1:A:925:VAL:HG21	2.02	0.42
1:A:392:LEU:O	1:A:396:GLN:HG3	2.19	0.42
1:A:170:LEU:HD21	1:A:277:GLU:HG3	2.01	0.42
1:A:1009:LYS:HA	1:A:1010:PRO:HD3	1.86	0.42
1:B:189:GLU:O	1:B:192:LYS:HD2	2.20	0.41
1:A:842:GLY:HA3	1:A:1008:VAL:HG23	2.02	0.41
1:B:618:TYR:HA	1:B:630:SER:O	2.20	0.41
1:B:437:ILE:HA	1:B:440:ILE:HG12	2.02	0.41
1:B:488:LYS:HD3	1:B:488:LYS:N	2.35	0.41
1:A:550:LEU:HD11	1:A:558:LYS:HG3	2.01	0.41
1:B:689:LEU:CD1	1:B:995:MET:HG2	2.50	0.41
1:A:479:ALA:HB2	3:A:2000:DIO:H12	2.02	0.41
1:A:386:HIS:HD2	1:A:389:ASP:OD2	2.03	0.41
1:A:817:GLU:HG3	1:A:818:PRO:CD	2.50	0.41
1:A:337:LEU:HD23	1:A:401:LEU:CD2	2.48	0.41
1:A:709:LEU:HG	1:A:713:LYS:HE3	2.03	0.41
1:A:634:TYR:O	1:A:638:GLN:HG3	2.21	0.41
1:B:179:LYS:HD2	1:B:237:VAL:HB	2.01	0.41
1:B:855:PRO:HB2	1:B:857:HIS:CE1	2.55	0.41
1:B:765:ARG:HD3	1:B:914:GLN:OE1	2.20	0.41
1:A:90:LEU:HD23	1:A:90:LEU:C	2.41	0.41
1:B:600:LEU:HD23	1:B:620:LEU:HD21	2.02	0.41
1:B:577:GLU:CD	1:B:912:ILE:HD13	2.41	0.41
1:A:400:LYS:HD3	4:A:2155:HOH:O	2.19	0.41
1:A:829:LEU:O	1:A:852:SER:HB2	2.21	0.41
1:B:793:ILE:O	1:B:847:ARG:HA	2.20	0.41
1:A:1003:PRO:HB3	1:B:1006:PRO:HD3	2.03	0.41
1:A:196:ASN:ND2	1:A:198:ALA:N	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:PHE:C	1:B:151:PHE:CD1	2.94	0.41
1:B:796:GLN:HB3	1:B:952:HIS:HB2	2.02	0.41
1:B:714:ALA:O	1:B:717:PRO:HD2	2.20	0.41
1:A:697:LYS:O	1:A:701:LYS:HG2	2.21	0.41
1:A:817:GLU:HG3	1:A:818:PRO:HD3	2.03	0.41
1:B:965:SER:O	1:B:966:CYS:HB2	2.21	0.41
1:A:100:PRO:HA	1:A:101:PRO:HD3	1.95	0.41
1:B:908:TRP:CZ2	1:B:912:ILE:HD11	2.55	0.41
1:B:374:ILE:C	1:B:374:ILE:HD12	2.41	0.41
1:A:90:LEU:HD13	1:A:169:PHE:CD2	2.55	0.41
1:A:559:LEU:HD11	1:A:729:LEU:HG	2.03	0.41
1:A:279:GLU:HB3	1:A:281:LYS:HE3	2.02	0.41
1:B:208:THR:CG2	1:B:477:ARG:HH22	2.14	0.40
1:B:196:ASN:ND2	1:B:198:ALA:HB3	2.36	0.40
1:B:843:ILE:HG22	1:B:844:GLN:H	1.86	0.40
1:B:259:LEU:HD12	1:B:260:GLY:N	2.36	0.40
1:A:776:TRP:CD2	1:A:989:PRO:HB3	2.56	0.40
1:B:552:LYS:HZ1	1:B:743:GLN:NE2	2.20	0.40
1:B:600:LEU:HD21	1:B:649:MET:CG	2.51	0.40
1:B:580:SER:HA	1:B:581:PRO:HD3	1.90	0.40
1:A:147:THR:HG22	1:A:149:TYR:CE1	2.56	0.40
1:B:600:LEU:CD2	1:B:649:MET:HG2	2.52	0.40
1:A:855:PRO:HA	1:A:856:PRO:HD3	1.90	0.40
1:B:512:LYS:HE3	1:B:512:LYS:HB2	1.85	0.40
1:A:423:ARG:NE	4:A:2544:HOH:O	2.53	0.40
1:B:986:LEU:HA	1:B:987:PRO:HD3	1.94	0.40
1:A:517:ASP:C	1:A:517:ASP:OD2	2.58	0.40
1:A:250:SER:O	1:A:254:MET:HG3	2.21	0.40
1:A:824:ARG:O	1:A:828:GLN:HA	2.22	0.40
1:A:420:VAL:HG12	1:A:634:TYR:OH	2.21	0.40
1:B:153:VAL:HG22	1:B:154:SER:N	2.36	0.40
1:B:86:SER:HB3	1:B:158:LEU:HG	2.03	0.40
1:B:94:ILE:HD12	1:B:248:TYR:HB3	2.02	0.40
1:B:580:SER:HB2	1:B:723:LEU:HD23	2.02	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:239:GLN:OE1	1:B:234:GLY:O[5_554]	2.19	0.01

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	960/990 (97%)	927 (97%)	33 (3%)	0	100	100
1	B	961/990 (97%)	934 (97%)	26 (3%)	1 (0%)	56	58
2	C	7/40 (18%)	7 (100%)	0	0	100	100
2	D	7/40 (18%)	7 (100%)	0	0	100	100
All	All	1935/2060 (94%)	1875 (97%)	59 (3%)	1 (0%)	56	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1012	I LE

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	855/883 (97%)	808 (94%)	47 (6%)	27	23
1	B	854/883 (97%)	798 (93%)	56 (7%)	21	17
2	C	8/31 (26%)	7 (88%)	1 (12%)	6	3
2	D	9/31 (29%)	6 (67%)	3 (33%)	0	0
All	All	1726/1828 (94%)	1619 (94%)	107 (6%)	23	19

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	LEU
1	A	102	ASN
1	A	119	LYS
1	A	125	ASN
1	A	196	ASN
1	A	201	LEU
1	A	226	LEU
1	A	238	ARG
1	A	270	LEU
1	A	281	LYS
1	A	285	LEU
1	A	337	LEU
1	A	347	LEU
1	A	356	VAL
1	A	417	LEU
1	A	429	ARG
1	A	450	LEU
1	A	454	TYR
1	A	466	MET
1	A	475	ASN
1	A	488	LYS
1	A	542	LYS
1	A	543	GLU
1	A	595	LEU
1	A	597	LEU
1	A	616	LEU
1	A	622	ASN
1	A	629	LEU
1	A	642	LEU
1	A	648	LYS
1	A	657	LYS
1	A	728	LEU
1	A	733	ILE
1	A	771	LEU
1	A	783	ASN
1	A	788	ASN
1	A	799	MET
1	A	810	LEU
1	A	823	LEU
1	A	846	LEU
1	A	859	LEU
1	A	867	LEU
1	A	879	GLU

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Mol	Chain	Res	Type
1	A	880	GLU
1	A	889	LEU
1	A	928	LEU
1	A	951	ARG
1	B	67	LEU
1	B	94	ILE
1	B	97	LEU
1	B	102	ASN
1	B	116	LEU
1	B	125	ASN
1	B	158	LEU
1	B	162	LEU
1	B	173	LEU
1	B	192	LYS
1	B	201	LEU
1	B	204	LEU
1	B	208	THR
1	B	226	LEU
1	B	259	LEU
1	B	324	LYS
1	B	337	LEU
1	B	417	LEU
1	B	450	LEU
1	B	468	LEU
1	B	470	LYS
1	B	488	LYS
1	B	556	MET
1	B	597	LEU
1	B	613	LEU
1	B	616	LEU
1	B	621	GLN
1	B	622	ASN
1	B	629	LEU
1	B	642	LEU
1	B	657	LYS
1	B	668	ARG
1	B	689	LEU
1	B	700	LEU
1	B	701	LYS
1	B	712	LEU
1	B	719	LEU
1	B	720	LEU

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Mol	Chain	Res	Type
1	B	728	LEU
1	B	733	ILE
1	B	771	LEU
1	B	783	ASN
1	B	788	ASN
1	B	810	LEU
1	B	817	GLU
1	B	823	LEU
1	B	846	LEU
1	B	859	LEU
1	B	861	SER
1	B	889	LEU
1	B	896	LYS
1	B	962	GLU
1	B	969	VAL
1	B	981	SER
1	B	993	GLN
1	B	1002	LEU
2	C	23	ASP
2	D	3	GLU
2	D	17	LEU
2	D	22	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (60) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	53	HIS
1	A	93	HIS
1	A	102	ASN
1	A	108	HIS
1	A	125	ASN
1	A	184	ASN
1	A	196	ASN
1	A	231	ASN
1	A	294	GLN
1	A	297	HIS
1	A	386	HIS
1	A	393	HIS
1	A	412	GLN
1	A	475	ASN
1	A	575	ASN

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Mol	Chain	Res	Type
1	A	622	ASN
1	A	672	ASN
1	A	736	GLN
1	A	783	ASN
1	A	805	ASN
1	A	828	GLN
1	A	841	ASN
1	A	887	GLN
1	A	922	ASN
1	A	957	HIS
1	B	93	HIS
1	B	102	ASN
1	B	125	ASN
1	B	129	GLN
1	B	157	HIS
1	B	184	ASN
1	B	196	ASN
1	B	232	GLN
1	B	294	GLN
1	B	297	HIS
1	B	300	GLN
1	B	332	HIS
1	B	475	ASN
1	B	499	GLN
1	B	502	GLN
1	B	589	HIS
1	B	605	ASN
1	B	621	GLN
1	B	622	ASN
1	B	672	ASN
1	B	679	HIS
1	B	718	GLN
1	B	730	HIS
1	B	732	ASN
1	B	743	GLN
1	B	770	GLN
1	B	781	GLN
1	B	783	ASN
1	B	805	ASN
1	B	828	GLN
1	B	883	GLN
1	B	922	ASN

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Mol	Chain	Res	Type
1	B	957	HIS
1	B	993	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 ligands 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	$\# Z > 2$
3	DIO	A	2000	-	6,6,6	0.75	0	6,6,6	0.21	0
3	DIO	B	2001	-	6,6,6	0.80	0	6,6,6	0.20	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DIO	A	2000	-	-	0/0/6/6	0/1/1/1
3	DIO	B	2001	-	-	0/0/6/6	0/1/1/1

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.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2000	DIO	1	0
3	B	2001	DIO	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	964/990 (97%)	-0.18	22 (2%) 64 70	22, 34, 50, 71	0
1	B	965/990 (97%)	-0.24	19 (1%) 68 73	18, 30, 48, 74	0
2	C	11/40 (27%)	1.61	4 (36%) 0 0	38, 45, 51, 56	0
2	D	11/40 (27%)	1.41	3 (27%) 1 1	33, 41, 50, 51	0
All	All	1951/2060 (94%)	-0.19	48 (2%) 61 67	18, 32, 50, 74	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1014	PHE	9.3
1	A	1014	PHE	8.4
1	A	1016	ALA	7.2
1	A	1015	MET	6.1
1	B	978	ILE	6.1
1	A	1012	ILE	5.8
2	D	23	ASP	5.6
1	A	964	ASP	5.3
1	B	1013	ASN	5.0
2	C	23	ASP	4.5
1	A	1013	ASN	4.3
1	B	1012	ILE	4.2
1	A	970	GLY	4.0
2	D	22	GLU	4.0
1	A	1011	HIS	4.0
2	C	17	LEU	3.8
1	B	1011	HIS	3.7
1	B	964	ASP	3.7
1	A	979	ASN	3.4
1	A	969	VAL	3.0
2	C	19	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	517	ASP	2.9
2	C	16	LYS	2.7
1	A	980	LEU	2.7
1	B	542	LYS	2.6
1	A	114	LEU	2.5
2	D	17	LEU	2.5
1	A	508	GLU	2.5
1	B	543	GLU	2.5
1	A	542	LYS	2.5
1	B	315	VAL	2.5
1	B	979	ASN	2.5
1	B	970	GLY	2.5
1	A	517	ASP	2.3
1	A	880	GLU	2.3
1	A	429	ARG	2.3
1	B	971	GLU	2.3
1	B	980	LEU	2.2
1	A	706	ASP	2.1
1	A	793	ILE	2.1
1	B	52	ASN	2.1
1	B	962	GLU	2.1
1	A	507	ASP	2.1
1	A	110	CYS	2.1
1	B	44	ASN	2.1
1	B	982	GLN	2.1
1	B	478	VAL	2.0
1	A	653	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(\AA^2)	Q<0.9
3	DIO	A	2000	6/6	0.92	0.26	4.44	55,56,57,57	0
3	DIO	B	2001	6/6	0.92	0.24	3.63	52,52,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.