



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:43 PM UTC

PDB ID : 1KXV / pdb_00001kxv
Title : Camelid VHH Domains in Complex with Porcine Pancreatic alpha-Amylase
Authors : Desmyter, A.; Spinelli, S.; Payan, F.; Lauwereys, M.; Wyns, L.; Muyldermans, S.; Cambillau, C.
Deposited on : 2002-02-01
Resolution : 1.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

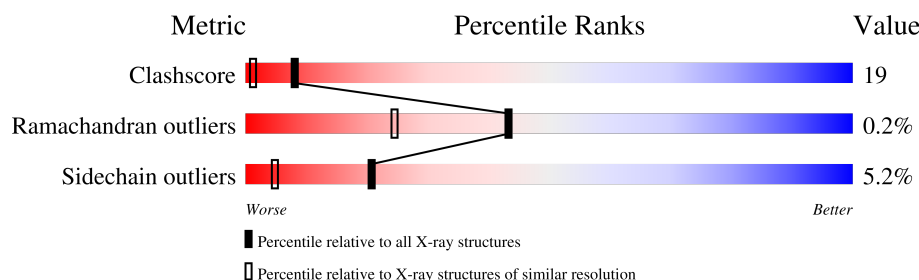
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	496	 74% 22% .
1	B	496	 74% 21% .
2	C	121	 80% 16% ...
2	D	121	 69% 21% 9% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10951 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 ALPHA-AMYLASE, PANCREATIC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	496	Total	C	N	O	S	0	0	0
			3900	2465	684	730	21			
1	B	496	Total	C	N	O	S	0	0	0
			3894	2459	685	729	21			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	243	LYS	GLN	SEE REMARK 999	UNP P00690
A	310	SER	ALA	SEE REMARK 999	UNP P00690
A	323	ILE	VAL	SEE REMARK 999	UNP P00690
A	404	GLN	GLU	SEE REMARK 999	UNP P00690
B	243	LYS	GLN	SEE REMARK 999	UNP P00690
B	310	SER	ALA	SEE REMARK 999	UNP P00690
B	323	ILE	VAL	SEE REMARK 999	UNP P00690
B	404	GLN	GLU	SEE REMARK 999	UNP P00690

- Molecule 2 is a protein called CAMELID VHH DOMAIN CAB10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	119	Total	C	N	O	S	0	0	0
			882	548	152	175	7			
2	D	121	Total	C	N	O	S	0	0	0
			898	556	155	180	7			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	590	Total	O	0	0
			590	590		
3	B	484	Total	O	0	0
			484	484		

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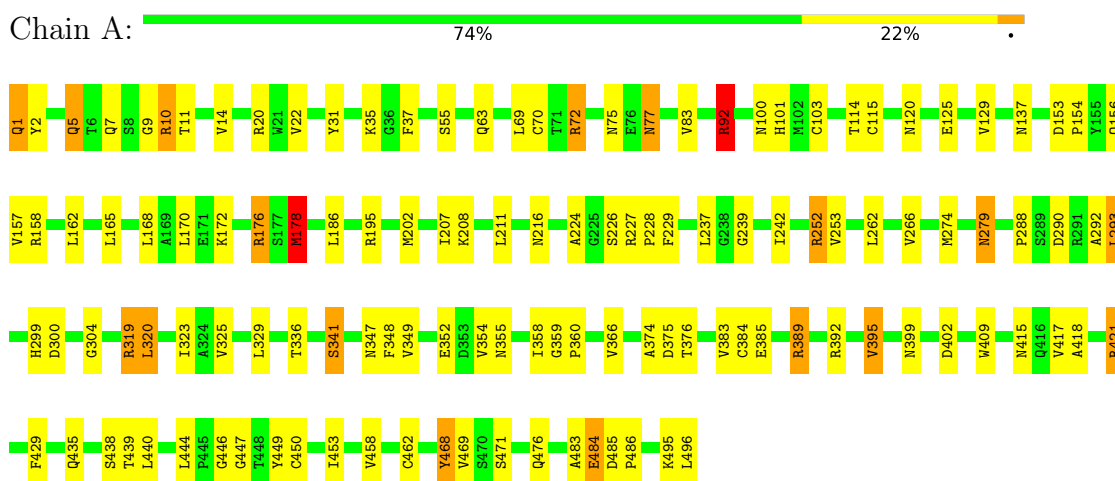
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	147	Total 147	O 147	0	0
3	D	156	Total 156	O 156	0	0

3 Residue-property plots

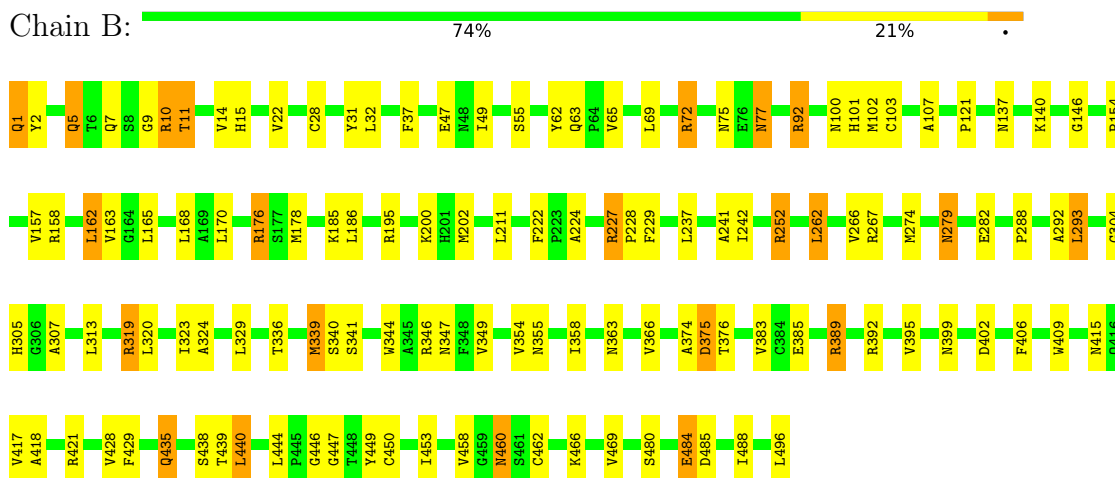
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

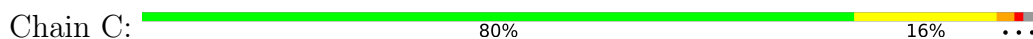
• Molecule 1: ALPHA-AMYLASE, PANCREATIC



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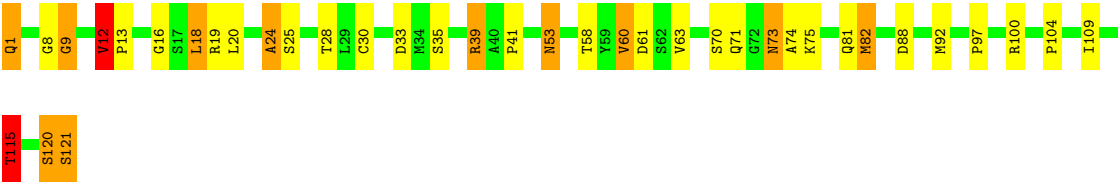


• Molecule 2: CAMELID VHH DOMAIN CAB10





● Molecule 2: CAMELID VHH DOMAIN CAB10



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.57Å 60.99Å 107.33Å 98.78° 100.88° 101.14°	Depositor
Resolution (Å)	29.34 – 1.60	Depositor
% Data completeness (in resolution range)	95.1 (29.34-1.60)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.205 , 0.229	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10951	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	1.06	3/4010 (0.1%)	1.34	46/5450 (0.8%)
1	B	1.00	5/4003 (0.1%)	1.23	29/5439 (0.5%)
2	C	1.00	2/901 (0.2%)	1.32	9/1225 (0.7%)
2	D	1.90	9/917 (1.0%)	1.76	21/1245 (1.7%)
All	All	1.14	19/9831 (0.2%)	1.34	105/13359 (0.8%)

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
2	C	0	2
All	All	0	3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	121	SER	C-OXT	31.62	1.86	1.23
2	D	8	GLY	C-N	-29.33	0.90	1.33
2	D	1	GLN	C-N	-11.73	1.18	1.33
2	D	120	SER	C-N	8.20	1.44	1.33
1	A	178	MET	SD-CE	-7.74	1.60	1.79
1	B	339	MET	SD-CE	-7.13	1.61	1.79
2	D	12	VAL	CA-CB	-6.88	1.45	1.54
1	B	102	MET	SD-CE	6.79	1.96	1.79
1	B	421	ARG	CD-NE	-6.74	1.36	1.46
2	D	8	GLY	C-O	-6.26	1.15	1.23
2	D	121	SER	N-CA	6.04	1.57	1.46
1	B	178	MET	SD-CE	-6.00	1.64	1.79
1	A	395	VAL	CA-CB	5.97	1.62	1.54
1	A	253	VAL	CA-CB	5.83	1.63	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	33	ASP	CA-C	5.53	1.60	1.52
2	D	58	THR	C-O	-5.49	1.17	1.23
2	C	46	ASP	CA-C	-5.17	1.46	1.52
2	D	82	MET	SD-CE	-5.17	1.66	1.79
1	B	421	ARG	NE-CZ	-5.00	1.27	1.33

All (105) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	121	SER	N-CA-CB	17.69	140.58	110.50
2	D	8	GLY	CA-C-N	16.00	152.78	121.41
2	D	8	GLY	C-N-CA	16.00	152.78	121.41
2	D	60	VAL	CG1-CB-CG2	15.44	144.77	110.80
2	D	8	GLY	O-C-N	-14.21	104.22	122.70
1	A	195	ARG	NE-CZ-NH2	12.68	130.61	119.20
1	A	252	ARG	NE-CZ-NH2	12.27	130.24	119.20
1	B	252	ARG	NE-CZ-NH2	11.95	129.95	119.20
1	A	195	ARG	NE-CZ-NH1	-11.71	109.79	121.50
1	B	176	ARG	NE-CZ-NH2	11.70	129.73	119.20
2	D	12	VAL	N-CA-CB	-11.07	95.71	111.21
1	B	252	ARG	NE-CZ-NH1	-10.90	110.60	121.50
1	A	252	ARG	NE-CZ-NH1	-10.79	110.70	121.50
1	A	72	ARG	NE-CZ-NH2	10.22	128.40	119.20
1	A	176	ARG	NE-CZ-NH2	10.22	128.40	119.20
1	A	10	ARG	NE-CZ-NH2	9.93	128.14	119.20
1	A	176	ARG	NE-CZ-NH1	-9.73	111.77	121.50
1	A	10	ARG	NE-CZ-NH1	-9.59	111.91	121.50
1	A	72	ARG	NE-CZ-NH1	-9.55	111.95	121.50
1	A	92	ARG	NE-CZ-NH2	9.35	127.61	119.20
1	A	176	ARG	CB-CG-CD	9.28	132.63	111.30
1	A	252	ARG	CD-NE-CZ	9.25	137.35	124.40
1	A	252	ARG	CB-CG-CD	9.02	132.04	111.30
2	D	120	SER	CA-C-N	-9.01	105.48	121.70
2	D	120	SER	C-N-CA	-9.01	105.48	121.70
1	A	224	ALA	CA-C-N	-8.84	104.09	121.41
1	A	224	ALA	C-N-CA	-8.84	104.09	121.41
1	A	92	ARG	NE-CZ-NH1	-8.80	112.70	121.50
1	B	176	ARG	NE-CZ-NH1	-8.77	112.73	121.50
1	B	72	ARG	NE-CZ-NH2	8.52	126.87	119.20
1	A	421	ARG	NE-CZ-NH2	8.30	126.67	119.20
1	B	252	ARG	CB-CG-CD	8.29	130.36	111.30
1	A	92	ARG	CB-CG-CD	8.24	130.26	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	72	ARG	NE-CZ-NH1	-8.04	113.46	121.50
1	B	252	ARG	CD-NE-CZ	7.95	135.52	124.40
1	B	304	GLY	N-CA-C	7.92	125.51	115.21
2	D	53	ASN	CA-CB-CG	-7.90	104.70	112.60
1	A	226	SER	N-CA-C	7.86	121.46	108.96
1	B	195	ARG	NE-CZ-NH2	7.85	126.27	119.20
1	A	389	ARG	NE-CZ-NH2	7.74	126.17	119.20
1	B	176	ARG	CB-CG-CD	7.67	128.94	111.30
1	A	389	ARG	NE-CZ-NH1	-7.58	113.92	121.50
1	B	10	ARG	NE-CZ-NH2	7.27	125.75	119.20
1	A	389	ARG	CG-CD-NE	-7.25	96.05	112.00
1	A	421	ARG	NE-CZ-NH1	-7.09	114.41	121.50
1	A	195	ARG	CD-NE-CZ	6.86	134.00	124.40
2	D	71	GLN	CB-CG-CD	6.82	124.19	112.60
2	D	115	THR	CB-CA-C	-6.75	94.57	109.56
1	B	389	ARG	NE-CZ-NH1	-6.75	114.75	121.50
1	B	389	ARG	NE-CZ-NH2	6.72	125.25	119.20
1	B	222	PHE	CA-C-N	-6.66	112.80	119.92
1	B	222	PHE	C-N-CA	-6.66	112.80	119.92
1	B	195	ARG	NE-CZ-NH1	-6.63	114.87	121.50
1	A	304	GLY	N-CA-C	6.62	123.81	115.21
1	A	224	ALA	O-C-N	-6.56	114.55	122.22
1	A	292	ALA	N-CA-C	6.52	120.46	109.76
1	A	92	ARG	CD-NE-CZ	6.46	133.45	124.40
2	C	108	ILE	N-CA-C	-6.42	100.42	108.89
1	A	55	SER	N-CA-C	6.34	119.14	108.23
1	A	293	LEU	CA-CB-CG	-6.30	94.24	116.30
1	A	72	ARG	CD-NE-CZ	6.16	133.03	124.40
2	D	121	SER	N-CA-C	-6.12	93.86	111.00
2	C	33	ASP	N-CA-C	-6.10	100.63	110.32
2	D	25	SER	N-CA-C	6.08	119.56	109.46
1	A	70	CYS	N-CA-CB	-6.07	101.43	110.90
1	B	293	LEU	CA-CB-CG	-6.02	95.24	116.30
1	A	336	THR	N-CA-C	5.91	119.27	109.46
2	D	35	SER	CB-CA-C	-5.90	101.06	110.14
1	B	176	ARG	CD-NE-CZ	5.89	132.65	124.40
1	A	293	LEU	N-CA-C	-5.85	98.98	108.52
1	A	341	SER	N-CA-C	5.75	117.97	109.69
1	B	55	SER	N-CA-C	5.74	118.11	108.23
1	B	292	ALA	N-CA-C	5.74	119.18	109.76
1	B	146	GLY	N-CA-C	-5.74	107.75	115.21
1	A	360	PRO	N-CA-C	5.71	115.95	110.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	293	LEU	N-CA-C	-5.65	99.31	108.52
2	D	24	ALA	N-CA-C	-5.64	96.02	107.41
2	D	8	GLY	N-CA-C	5.64	126.54	113.18
1	A	359	GLY	CA-C-N	5.60	123.75	119.66
1	A	359	GLY	C-N-CA	5.60	123.75	119.66
1	B	336	THR	N-CA-C	5.59	119.36	109.96
2	D	71	GLN	CA-C-N	-5.58	116.75	122.27
2	D	71	GLN	C-N-CA	-5.58	116.75	122.27
2	D	71	GLN	CA-CB-CG	-5.53	103.04	114.10
2	D	39	ARG	NE-CZ-NH2	5.53	124.17	119.20
1	A	137	ASN	N-CA-C	5.50	119.62	113.02
2	C	43	LYS	CA-C-N	-5.48	115.36	121.83
2	C	43	LYS	C-N-CA	-5.48	115.36	121.83
1	A	114	THR	N-CA-C	5.45	119.19	112.54
1	A	239	GLY	N-CA-C	5.43	122.30	114.10
2	C	39	ARG	NE-CZ-NH2	5.34	124.00	119.20
2	D	115	THR	CA-CB-CG2	5.29	119.48	110.50
1	B	406	PHE	N-CA-C	-5.28	102.23	110.10
1	B	137	ASN	N-CA-C	5.28	119.36	113.02
1	A	10	ARG	CD-NE-CZ	5.23	131.72	124.40
1	A	471	SER	N-CA-C	5.23	117.72	111.71
2	C	25	SER	N-CA-C	5.20	118.47	110.42
1	A	176	ARG	CD-NE-CZ	5.17	131.64	124.40
1	B	11	THR	N-CA-C	5.10	120.62	113.02
2	C	119	VAL	CA-C-N	5.10	130.88	121.70
2	C	119	VAL	C-N-CA	5.10	130.88	121.70
1	B	10	ARG	NE-CZ-NH1	-5.09	116.41	121.50
1	A	103	CYS	N-CA-C	5.09	117.37	110.35
1	B	241	ALA	N-CA-C	5.08	116.82	111.28
2	C	73	ASN	N-CA-C	5.03	117.41	111.33

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	468	TYR	Sidechain
2	C	39	ARG	Sidechain
2	C	45	ARG	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	3900	0	3671	145	0
1	B	3894	0	3666	178	0
2	C	882	0	848	16	0
2	D	898	0	863	24	0
3	A	590	0	0	28	0
3	B	484	0	0	25	0
3	C	147	0	0	4	0
3	D	156	0	0	9	0
All	All	10951	0	9048	363	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 19.

All (363) 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:1:GLN:NE2	1:B:1:GLN:H1	1.18	1.41
1:B:1:GLN:NE2	1:B:1:GLN:N	1.90	1.19
1:B:282:GLU:OE1	3:B:626:HOH:O	1.66	1.14
1:A:1:GLN:OE1	1:A:1:GLN:CA	2.00	1.09
1:A:409:TRP:CH2	1:A:417:VAL:HG11	1.90	1.06
1:B:341:SER:O	1:B:383:VAL:HG22	1.55	1.05
1:A:229:PHE:HE2	1:A:252:ARG:HD3	1.17	1.04
1:B:170:LEU:O	1:B:176:ARG:HD3	1.58	1.04
1:B:229:PHE:HE2	1:B:252:ARG:HD3	1.19	1.01
1:B:157:VAL:HG21	1:B:242:ILE:HD11	1.42	1.01
1:B:484:GLU:HG2	1:B:485:ASP:N	1.77	1.00
1:B:409:TRP:CH2	1:B:417:VAL:HG11	1.96	0.99
1:A:1:GLN:OE1	1:A:1:GLN:HA	1.13	0.99
1:A:207:ILE:O	1:A:211:LEU:HD13	1.63	0.98
1:B:266:VAL:HG13	1:B:323:ILE:HD12	1.45	0.98
1:B:2:TYR:O	1:B:252:ARG:HD2	1.64	0.97
1:B:1:GLN:N	1:B:1:GLN:CD	2.23	0.96
1:B:69:LEU:HD21	1:B:185:LYS:HE3	1.48	0.96
1:B:320:LEU:O	1:B:323:ILE:HG13	1.63	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:SER:O	1:A:383:VAL:HG22	1.66	0.95
1:A:2:TYR:O	1:A:252:ARG:HD2	1.68	0.94
1:B:358:ILE:HB	3:B:826:HOH:O	1.66	0.94
1:B:170:LEU:O	1:B:176:ARG:CD	2.15	0.94
2:D:61:ASP:OD2	3:D:159:HOH:O	1.87	0.92
2:D:16:GLY:HA3	3:D:226:HOH:O	1.68	0.92
1:B:1:GLN:H1	1:B:1:GLN:HE21	1.04	0.92
1:A:170:LEU:O	1:A:176:ARG:HD3	1.68	0.91
1:A:1:GLN:N	1:A:228:PRO:O	2.03	0.91
1:A:170:LEU:O	1:A:176:ARG:CD	2.19	0.91
1:A:229:PHE:CE2	1:A:252:ARG:HD3	2.05	0.90
1:A:325:VAL:O	1:A:329:LEU:HD23	1.71	0.89
1:B:154:PRO:HA	1:B:157:VAL:HG22	1.55	0.88
1:A:484:GLU:CD	1:A:485:ASP:H	1.83	0.86
1:A:100:ASN:HD22	1:A:101:HIS:HD2	1.23	0.85
1:B:1:GLN:CD	1:B:1:GLN:H3	1.84	0.83
1:B:49:ILE:HG21	1:B:103:CYS:CB	2.08	0.83
1:B:100:ASN:HD22	1:B:101:HIS:HD2	1.23	0.82
1:A:484:GLU:OE2	1:A:484:GLU:N	2.13	0.82
1:B:392:ARG:O	1:B:395:VAL:HG22	1.80	0.82
2:C:12:VAL:HG23	2:C:13:PRO:HD2	1.62	0.81
1:B:229:PHE:CE2	1:B:252:ARG:HD3	2.10	0.81
1:B:157:VAL:HG21	1:B:242:ILE:CD1	2.12	0.80
1:B:266:VAL:CG1	1:B:323:ILE:HD12	2.11	0.80
1:A:300:ASP:HB2	3:A:737:HOH:O	1.80	0.80
1:A:72:ARG:CD	3:A:545:HOH:O	2.28	0.79
1:A:349:VAL:HG13	3:A:838:HOH:O	1.82	0.79
1:B:339:MET:HE3	1:B:339:MET:HA	1.62	0.79
1:B:439:THR:C	1:B:440:LEU:HD12	2.08	0.78
1:B:355:ASN:ND2	3:B:805:HOH:O	2.15	0.78
1:B:5:GLN:HA	1:B:5:GLN:HE21	1.48	0.78
1:A:157:VAL:HG11	1:A:242:ILE:HD11	1.64	0.77
2:D:13:PRO:HD2	3:D:226:HOH:O	1.86	0.75
1:B:157:VAL:HG23	1:B:158:ARG:HD2	1.68	0.75
1:B:341:SER:C	1:B:383:VAL:HG22	2.11	0.75
1:A:5:GLN:HA	1:A:5:GLN:HE21	1.50	0.75
1:A:323:ILE:HD11	1:A:486:PRO:HG2	1.68	0.75
1:B:49:ILE:HG22	1:B:65:VAL:CG1	2.16	0.74
1:A:72:ARG:HD3	3:A:545:HOH:O	1.88	0.74
1:B:154:PRO:HA	1:B:157:VAL:CG2	2.18	0.74
1:B:69:LEU:HD22	1:B:69:LEU:N	2.02	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:ARG:HH21	1:A:319:ARG:HG2	1.53	0.74
2:D:12:VAL:HG13	3:D:226:HOH:O	1.86	0.74
2:C:12:VAL:CG2	2:C:13:PRO:HD2	2.18	0.73
1:A:383:VAL:HG23	1:A:385:GLU:OE2	1.88	0.73
1:A:72:ARG:HD2	3:A:545:HOH:O	1.88	0.73
1:A:207:ILE:O	1:A:211:LEU:CD1	2.36	0.73
1:B:313:LEU:HD11	3:B:831:HOH:O	1.87	0.73
1:B:72:ARG:HD2	3:B:566:HOH:O	1.87	0.73
1:B:440:LEU:HD12	1:B:440:LEU:N	2.04	0.72
1:B:49:ILE:HG21	1:B:103:CYS:HB2	1.71	0.72
1:B:383:VAL:HG23	1:B:385:GLU:OE2	1.89	0.72
1:B:266:VAL:CG1	1:B:323:ILE:CD1	2.67	0.72
1:A:288:PRO:HG3	3:C:208:HOH:O	1.88	0.72
1:B:69:LEU:HD21	1:B:185:LYS:CE	2.19	0.72
1:B:347:ASN:HB2	3:B:805:HOH:O	1.89	0.72
1:B:409:TRP:CH2	1:B:417:VAL:CG1	2.71	0.72
1:A:5:GLN:O	1:A:92:ARG:CD	2.38	0.71
1:A:5:GLN:O	1:A:92:ARG:HD2	1.91	0.71
1:B:267:ARG:HG3	3:B:831:HOH:O	1.89	0.70
1:B:446:GLY:HA2	1:B:469:VAL:HG13	1.73	0.70
1:A:447:GLY:HA3	1:A:495:LYS:HZ3	1.54	0.70
1:B:266:VAL:HG13	1:B:323:ILE:CD1	2.21	0.70
1:B:15:HIS:HB3	1:B:339:MET:HE3	1.74	0.70
1:A:417:VAL:HG12	1:A:418:ALA:N	2.06	0.70
1:A:208:LYS:HE3	3:C:205:HOH:O	1.92	0.70
1:B:484:GLU:CG	1:B:485:ASP:N	2.53	0.69
1:B:319:ARG:HG2	1:B:319:ARG:HH21	1.57	0.69
1:B:15:HIS:HB3	1:B:339:MET:CE	2.23	0.69
1:B:15:HIS:CD2	1:B:339:MET:HE1	2.28	0.69
1:A:447:GLY:HA3	1:A:495:LYS:NZ	2.07	0.69
1:B:227:ARG:CD	3:B:799:HOH:O	2.40	0.69
1:A:154:PRO:HA	1:A:157:VAL:HG12	1.74	0.68
1:A:383:VAL:CG2	1:A:385:GLU:OE2	2.41	0.68
1:A:154:PRO:HA	1:A:157:VAL:CG1	2.24	0.68
1:A:439:THR:C	1:A:440:LEU:HD12	2.19	0.68
1:B:279:ASN:H	1:B:279:ASN:HD22	1.41	0.68
2:C:60:VAL:HB	3:C:208:HOH:O	1.93	0.68
1:A:389:ARG:HD2	1:A:453:ILE:O	1.94	0.68
1:B:392:ARG:O	1:B:395:VAL:CG2	2.41	0.68
1:B:438:SER:HB2	1:B:440:LEU:HD11	1.76	0.68
1:B:363:ASN:O	1:B:366:VAL:HG22	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:ILE:CG2	1:B:103:CYS:CB	2.72	0.67
1:A:157:VAL:HG11	1:A:242:ILE:CD1	2.25	0.67
1:B:346:ARG:C	3:B:805:HOH:O	2.37	0.67
1:B:72:ARG:CD	3:B:566:HOH:O	2.43	0.67
1:A:170:LEU:O	1:A:176:ARG:HD2	1.94	0.67
1:A:409:TRP:CH2	1:A:417:VAL:CG1	2.75	0.66
1:B:320:LEU:O	1:B:323:ILE:CG1	2.39	0.66
2:C:13:PRO:HA	2:C:120:SER:HA	1.77	0.66
1:A:447:GLY:O	1:A:469:VAL:HG22	1.96	0.66
2:C:47:PHE:HD2	3:C:208:HOH:O	1.79	0.65
1:B:320:LEU:HA	1:B:323:ILE:HG12	1.78	0.65
1:A:274:MET:HE2	1:A:415:ASN:O	1.96	0.65
2:D:33:ASP:OD2	2:D:100:ARG:HD2	1.96	0.65
1:A:279:ASN:HD22	1:A:279:ASN:H	1.45	0.65
1:B:2:TYR:O	1:B:252:ARG:CD	2.43	0.65
1:B:1:GLN:NE2	3:B:799:HOH:O	2.29	0.65
1:B:339:MET:CE	1:B:340:SER:H	2.10	0.64
1:B:320:LEU:HA	1:B:323:ILE:HD11	1.80	0.64
1:B:170:LEU:O	1:B:176:ARG:HD2	1.96	0.64
1:B:15:HIS:CG	1:B:339:MET:HE1	2.33	0.64
1:A:449:TYR:HE2	1:A:469:VAL:HG21	1.62	0.64
1:B:320:LEU:HA	1:B:323:ILE:CG1	2.27	0.64
1:B:227:ARG:HD3	3:B:799:HOH:O	1.95	0.63
1:B:320:LEU:HA	1:B:323:ILE:CD1	2.29	0.63
1:A:349:VAL:HG23	1:A:349:VAL:O	1.99	0.62
1:B:100:ASN:HD22	1:B:101:HIS:CD2	2.12	0.62
2:D:28:THR:HG22	3:D:138:HOH:O	1.97	0.62
1:B:49:ILE:HD11	1:B:63:GLN:OE1	2.00	0.62
1:A:20:ARG:CZ	3:A:673:HOH:O	2.47	0.62
1:B:462:CYS:SG	1:B:496:LEU:HD11	2.40	0.61
1:A:341:SER:C	1:A:383:VAL:HG22	2.24	0.61
1:A:484:GLU:CG	3:A:745:HOH:O	2.48	0.61
1:B:355:ASN:O	3:B:826:HOH:O	2.15	0.61
1:B:339:MET:HE2	1:B:340:SER:H	1.66	0.60
1:A:496:LEU:HD12	1:A:496:LEU:N	2.16	0.60
1:B:449:TYR:HE2	1:B:469:VAL:HG11	1.64	0.60
1:A:266:VAL:HG12	1:A:320:LEU:HG	1.84	0.60
1:A:83:VAL:HG12	3:A:971:HOH:O	2.02	0.60
1:A:100:ASN:HD22	1:A:101:HIS:CD2	2.14	0.60
2:D:19:ARG:HE	2:D:81:GLN:HE21	1.49	0.59
1:A:446:GLY:HA2	1:A:469:VAL:HG23	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:ILE:CD1	3:B:817:HOH:O	2.50	0.59
1:A:5:GLN:O	1:A:92:ARG:HD3	2.02	0.59
1:A:484:GLU:HG2	3:A:745:HOH:O	2.02	0.59
1:B:154:PRO:CA	1:B:157:VAL:HG22	2.30	0.59
1:B:458:VAL:HG23	1:B:458:VAL:O	2.01	0.59
1:B:389:ARG:HD2	1:B:453:ILE:O	2.02	0.59
1:B:460:ASN:CG	1:B:460:ASN:O	2.44	0.59
1:B:49:ILE:CG2	1:B:103:CYS:HB3	2.32	0.59
1:A:347:ASN:OD1	3:A:838:HOH:O	2.16	0.58
1:B:383:VAL:CG2	1:B:385:GLU:OE2	2.51	0.58
1:A:323:ILE:HD11	1:A:486:PRO:CG	2.32	0.58
1:B:69:LEU:HD22	1:B:69:LEU:H	1.67	0.58
1:B:237:LEU:HD21	1:B:307:ALA:HB1	1.84	0.58
2:D:120:SER:O	2:D:121:SER:C	2.40	0.58
1:A:290:ASP:HB3	3:A:772:HOH:O	2.02	0.58
1:A:449:TYR:CE2	1:A:469:VAL:HG21	2.38	0.58
1:B:227:ARG:HG3	1:B:228:PRO:HD2	1.84	0.58
2:D:24:ALA:HB1	2:D:109:ILE:HD11	1.84	0.58
1:A:129:VAL:CG1	1:A:178:MET:HG2	2.33	0.58
2:D:30:CYS:O	2:D:75:LYS:HE2	2.02	0.58
1:B:274:MET:HE1	1:B:429:PHE:CB	2.34	0.57
2:D:19:ARG:HE	2:D:81:GLN:NE2	2.03	0.57
1:B:31:TYR:OH	1:B:392:ARG:HG3	2.04	0.57
1:B:320:LEU:N	1:B:320:LEU:HD12	2.19	0.57
1:B:417:VAL:HG12	1:B:418:ALA:N	2.19	0.57
1:A:323:ILE:HG23	1:A:429:PHE:CD2	2.39	0.57
1:A:476:GLN:CD	3:A:640:HOH:O	2.47	0.57
2:D:33:ASP:OD1	3:D:128:HOH:O	2.18	0.57
1:A:374:ALA:C	1:A:376:THR:H	2.11	0.57
1:B:447:GLY:O	1:B:469:VAL:HG12	2.04	0.57
1:B:450:CYS:SG	1:B:496:LEU:HD11	2.45	0.57
1:B:282:GLU:OE2	1:B:288:PRO:HA	2.05	0.57
1:B:440:LEU:N	1:B:440:LEU:CD1	2.68	0.56
1:B:157:VAL:CG2	1:B:242:ILE:HD11	2.27	0.56
1:A:211:LEU:HD12	1:A:211:LEU:N	2.21	0.56
2:C:12:VAL:CG2	2:C:13:PRO:CD	2.84	0.56
1:A:409:TRP:CZ3	1:A:417:VAL:HG11	2.39	0.56
1:A:483:ALA:HB1	1:A:484:GLU:OE2	2.05	0.56
1:A:63:GLN:HG2	1:A:165:LEU:HD21	1.88	0.56
1:B:447:GLY:H	1:B:469:VAL:HG13	1.71	0.56
1:A:14:VAL:HG11	1:A:37:PHE:CE2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:MET:CE	1:A:415:ASN:O	2.53	0.56
1:A:349:VAL:CG1	3:A:838:HOH:O	2.47	0.56
1:A:438:SER:CB	1:A:440:LEU:HD11	2.36	0.56
1:B:274:MET:CE	1:B:429:PHE:HB2	2.36	0.56
2:D:97:PRO:HB2	2:D:109:ILE:HD12	1.88	0.55
1:B:447:GLY:N	1:B:469:VAL:HG13	2.20	0.55
1:B:75:ASN:HB2	1:B:77:ASN:OD1	2.06	0.55
2:D:60:VAL:HG12	2:D:63:VAL:HG22	1.87	0.55
1:B:341:SER:HB2	1:B:383:VAL:HG21	1.88	0.55
1:B:344:TRP:CD1	3:B:826:HOH:O	2.60	0.55
1:A:22:VAL:HG12	3:A:769:HOH:O	2.08	0.54
1:B:274:MET:HE1	1:B:429:PHE:HB2	1.90	0.54
2:C:119:VAL:O	2:C:120:SER:HB2	2.07	0.54
1:A:447:GLY:H	1:A:469:VAL:HG23	1.73	0.54
1:B:49:ILE:CD1	1:B:63:GLN:OE1	2.56	0.53
1:A:438:SER:HB2	1:A:440:LEU:HD11	1.89	0.53
1:B:154:PRO:O	1:B:157:VAL:HG22	2.09	0.53
1:B:349:VAL:HG22	1:B:354:VAL:HG13	1.90	0.53
1:A:172:LYS:HD2	3:A:775:HOH:O	2.07	0.53
2:C:12:VAL:HG21	2:C:16:GLY:HA3	1.91	0.53
1:A:458:VAL:HG23	1:A:458:VAL:O	2.08	0.53
1:A:476:GLN:HG3	3:A:722:HOH:O	2.09	0.53
1:B:14:VAL:HG11	1:B:37:PHE:CE2	2.44	0.53
1:B:157:VAL:HG23	1:B:158:ARG:CD	2.39	0.53
1:B:366:VAL:HG23	1:B:366:VAL:O	2.09	0.53
1:A:323:ILE:HG23	1:A:429:PHE:HD2	1.74	0.52
1:A:484:GLU:CD	1:A:485:ASP:N	2.63	0.52
1:B:200:LYS:HE3	3:B:703:HOH:O	2.08	0.52
2:C:12:VAL:HG22	2:C:13:PRO:N	2.24	0.52
1:B:274:MET:HE2	1:B:415:ASN:O	2.09	0.52
1:B:392:ARG:HD2	3:B:637:HOH:O	2.09	0.52
1:A:157:VAL:CG1	1:A:242:ILE:HD11	2.38	0.52
1:A:469:VAL:HG23	1:A:469:VAL:O	2.09	0.52
1:A:129:VAL:HG13	1:A:178:MET:HG2	1.92	0.52
1:B:154:PRO:C	1:B:157:VAL:HG22	2.34	0.52
1:A:348:PHE:C	3:A:838:HOH:O	2.52	0.52
1:B:449:TYR:CE2	1:B:469:VAL:HG11	2.44	0.51
1:A:349:VAL:O	1:A:349:VAL:CG2	2.57	0.51
1:B:1:GLN:HB3	3:B:731:HOH:O	2.10	0.51
1:A:2:TYR:O	1:A:252:ARG:CD	2.52	0.51
2:D:20:LEU:HG	2:D:82:MET:HE2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:39:ARG:NH2	2:C:94:TYR:OH	2.43	0.51
1:B:374:ALA:C	1:B:376:THR:H	2.16	0.51
1:A:227:ARG:NH2	3:A:713:HOH:O	2.44	0.50
1:A:417:VAL:CG1	1:A:418:ALA:N	2.72	0.50
1:A:462:CYS:SG	1:A:496:LEU:HD11	2.52	0.50
1:B:319:ARG:HG2	1:B:319:ARG:NH2	2.23	0.50
1:A:170:LEU:CD2	1:A:202:MET:SD	2.99	0.50
1:A:299:HIS:HE1	3:A:694:HOH:O	1.95	0.50
1:B:11:THR:H	1:B:399:ASN:HD21	1.60	0.49
2:C:13:PRO:HA	2:C:120:SER:CA	2.42	0.49
1:A:409:TRP:CZ2	1:A:417:VAL:HG11	2.45	0.49
2:C:74:ALA:O	2:C:77:THR:OG1	2.23	0.49
1:A:389:ARG:CD	1:A:453:ILE:O	2.60	0.49
1:A:178:MET:HA	1:A:178:MET:CE	2.41	0.49
1:A:319:ARG:HG2	1:A:319:ARG:NH2	2.26	0.49
2:D:88:ASP:C	2:D:88:ASP:OD1	2.55	0.49
1:B:349:VAL:HG23	1:B:349:VAL:O	2.13	0.48
1:A:383:VAL:HG23	1:A:384:CYS:N	2.28	0.48
1:A:484:GLU:CG	1:A:485:ASP:N	2.76	0.48
1:B:435:GLN:HG2	1:B:480:SER:HA	1.95	0.48
1:A:374:ALA:C	1:A:376:THR:N	2.70	0.48
1:B:428:VAL:HG13	1:B:428:VAL:O	2.12	0.48
1:A:157:VAL:HG13	1:A:158:ARG:HD2	1.95	0.48
1:A:349:VAL:HG22	1:A:354:VAL:HG13	1.95	0.48
1:A:375:ASP:O	1:A:376:THR:OG1	2.22	0.48
1:A:347:ASN:C	3:A:838:HOH:O	2.55	0.48
1:B:47:GLU:HG2	1:B:65:VAL:CG2	2.43	0.48
2:D:73:ASN:HD22	2:D:74:ALA:N	2.11	0.48
1:A:409:TRP:CZ2	1:A:417:VAL:CG1	2.97	0.48
1:B:5:GLN:O	1:B:92:ARG:HD2	2.14	0.48
1:B:15:HIS:HB3	1:B:339:MET:HE1	1.96	0.48
2:D:9:GLY:HA3	2:D:115:THR:CG2	2.44	0.48
1:A:468:TYR:CD1	1:A:468:TYR:C	2.90	0.47
1:A:7:GLN:OE1	1:A:10:ARG:HD2	2.14	0.47
1:A:440:LEU:HD12	1:A:440:LEU:N	2.29	0.47
1:B:170:LEU:CD2	1:B:202:MET:SD	3.02	0.47
1:B:274:MET:HE1	1:B:429:PHE:HB3	1.96	0.47
1:A:153:ASP:HB3	1:A:156:GLN:HG2	1.96	0.47
1:B:62:TYR:O	1:B:101:HIS:HE1	1.98	0.47
1:B:227:ARG:NE	3:B:691:HOH:O	2.45	0.47
1:B:339:MET:HE3	1:B:339:MET:CA	2.38	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:GLN:HA	1:B:5:GLN:NE2	2.24	0.46
1:B:170:LEU:HD21	1:B:202:MET:SD	2.55	0.46
1:A:229:PHE:HE2	1:A:252:ARG:CD	2.06	0.46
1:A:237:LEU:N	1:A:237:LEU:HD22	2.30	0.46
1:A:348:PHE:HA	1:A:352:GLU:O	2.15	0.46
1:B:69:LEU:N	1:B:69:LEU:CD2	2.74	0.46
1:A:476:GLN:CG	3:A:722:HOH:O	2.63	0.46
2:D:1:GLN:HA	3:D:192:HOH:O	2.14	0.46
2:C:73:ASN:OD1	2:C:74:ALA:N	2.48	0.46
1:B:49:ILE:HD13	1:B:103:CYS:HB2	1.97	0.46
1:B:49:ILE:HG23	1:B:103:CYS:HB3	1.96	0.46
1:B:22:VAL:HG13	3:B:591:HOH:O	2.16	0.46
1:B:266:VAL:CG1	1:B:323:ILE:HD11	2.45	0.45
1:A:447:GLY:N	1:A:469:VAL:HG23	2.31	0.45
1:B:279:ASN:HD22	1:B:279:ASN:N	2.09	0.45
1:A:496:LEU:N	1:A:496:LEU:CD1	2.80	0.45
1:A:349:VAL:HG22	1:A:354:VAL:CG1	2.46	0.45
1:A:31:TYR:OH	1:A:392:ARG:HG3	2.16	0.45
1:A:207:ILE:C	1:A:211:LEU:HD13	2.35	0.45
1:B:262:LEU:HD13	1:B:324:ALA:HB1	1.99	0.45
1:B:28:CYS:HA	1:B:32:LEU:HB2	1.98	0.45
2:D:104:PRO:HG2	3:D:243:HOH:O	2.17	0.45
1:B:447:GLY:N	1:B:469:VAL:CG1	2.80	0.45
1:A:7:GLN:HB3	1:A:10:ARG:HD3	1.98	0.45
1:A:319:ARG:HD2	3:A:809:HOH:O	2.17	0.45
1:B:9:GLY:HA2	3:B:729:HOH:O	2.16	0.45
1:A:154:PRO:CA	1:A:157:VAL:HG12	2.45	0.45
1:B:354:VAL:HA	3:B:954:HOH:O	2.17	0.44
1:A:211:LEU:CD1	1:A:211:LEU:H	2.30	0.44
1:B:320:LEU:CA	1:B:323:ILE:HG12	2.46	0.44
1:A:347:ASN:CG	3:A:838:HOH:O	2.60	0.44
1:A:216:ASN:OD1	1:A:227:ARG:HD2	2.18	0.44
1:A:440:LEU:N	1:A:440:LEU:CD1	2.80	0.44
1:B:15:HIS:HD2	3:B:671:HOH:O	2.00	0.44
1:B:375:ASP:O	1:B:376:THR:OG1	2.29	0.44
1:B:402:ASP:C	1:B:402:ASP:OD2	2.61	0.44
1:A:1:GLN:NE2	3:A:633:HOH:O	2.51	0.44
1:B:107:ALA:HB3	1:B:121:PRO:HG2	2.00	0.44
1:B:14:VAL:CG1	1:B:37:PHE:CE2	3.01	0.44
1:B:72:ARG:HD3	3:B:566:HOH:O	2.16	0.44
1:B:392:ARG:C	1:B:395:VAL:HG22	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:GLN:HG2	1:B:165:LEU:CD2	2.47	0.43
1:B:409:TRP:CZ2	1:B:417:VAL:CG1	3.01	0.43
1:B:462:CYS:HB2	1:B:466:LYS:NZ	2.33	0.43
1:A:211:LEU:CD1	1:A:211:LEU:N	2.81	0.43
1:A:22:VAL:HG11	3:A:944:HOH:O	2.18	0.43
1:A:323:ILE:CD1	1:A:486:PRO:HG2	2.45	0.43
1:B:428:VAL:CG1	1:B:488:ILE:HB	2.48	0.43
2:D:73:ASN:HD22	2:D:73:ASN:C	2.26	0.43
1:A:9:GLY:HA2	3:A:818:HOH:O	2.18	0.43
1:A:435:GLN:HG2	3:A:940:HOH:O	2.18	0.43
1:B:469:VAL:HG13	1:B:469:VAL:O	2.17	0.43
2:D:121:SER:C	3:D:153:HOH:O	2.60	0.43
2:C:6:GLU:OE2	2:C:95:CYS:HB3	2.18	0.43
1:B:355:ASN:HB3	1:B:358:ILE:HD12	2.00	0.43
1:A:170:LEU:HD21	1:A:202:MET:SD	2.59	0.43
1:B:417:VAL:CG1	1:B:418:ALA:N	2.82	0.43
1:B:339:MET:HE3	1:B:340:SER:H	1.80	0.42
2:D:12:VAL:HG21	2:D:18:LEU:HG	2.01	0.42
1:B:69:LEU:HD21	1:B:185:LYS:NZ	2.34	0.42
2:C:73:ASN:OD1	2:C:77:THR:HG21	2.19	0.42
1:A:383:VAL:HG21	1:A:385:GLU:OE2	2.16	0.42
1:B:409:TRP:CZ3	1:B:417:VAL:HG11	2.52	0.42
1:B:435:GLN:HE21	1:B:435:GLN:HB3	1.52	0.42
1:B:449:TYR:HE2	1:B:469:VAL:CG1	2.29	0.42
1:B:162:LEU:HD22	1:B:163:VAL:HG13	2.02	0.42
1:B:374:ALA:C	1:B:376:THR:N	2.76	0.42
1:A:120:ASN:HB3	1:A:125:GLU:HB2	2.00	0.42
2:D:53:ASN:HD22	2:D:53:ASN:HA	1.55	0.42
1:A:11:THR:H	1:A:399:ASN:HD21	1.67	0.42
1:A:211:LEU:HD12	1:A:211:LEU:H	1.85	0.42
1:A:329:LEU:N	1:A:329:LEU:HD22	2.35	0.42
1:B:49:ILE:HD11	3:B:817:HOH:O	2.18	0.42
1:A:14:VAL:CG1	1:A:37:PHE:CE2	3.02	0.41
1:A:450:CYS:SG	1:A:496:LEU:HD11	2.60	0.41
1:B:49:ILE:CG2	1:B:65:VAL:CG1	2.95	0.41
1:A:438:SER:HB3	1:A:440:LEU:HD11	2.01	0.41
1:B:2:TYR:C	1:B:252:ARG:HD2	2.41	0.41
1:B:409:TRP:CZ2	1:B:417:VAL:HG13	2.56	0.41
2:C:12:VAL:CG2	2:C:16:GLY:HA3	2.50	0.41
1:A:101:HIS:CE1	1:A:165:LEU:HD13	2.55	0.41
1:B:320:LEU:CA	1:B:323:ILE:CG1	2.99	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:ARG:HH21	1:A:319:ARG:CG	2.21	0.41
1:B:49:ILE:HG21	1:B:103:CYS:SG	2.61	0.41
1:B:458:VAL:O	1:B:458:VAL:CG2	2.65	0.41
1:A:5:GLN:HA	1:A:5:GLN:NE2	2.28	0.41
1:A:483:ALA:O	1:A:484:GLU:C	2.62	0.41
1:B:7:GLN:OE1	1:B:10:ARG:HD2	2.20	0.40
1:B:49:ILE:HG22	1:B:65:VAL:HG11	2.02	0.40
1:B:63:GLN:HG2	1:B:165:LEU:HD21	2.03	0.40
1:B:339:MET:HE2	1:B:340:SER:O	2.20	0.40
1:A:402:ASP:HA	3:A:652:HOH:O	2.20	0.40
1:B:69:LEU:H	1:B:69:LEU:CD2	2.33	0.40
1:B:266:VAL:HG11	1:B:323:ILE:CD1	2.50	0.40
1:A:75:ASN:OD1	1:A:77:ASN:OD1	2.39	0.40
1:B:320:LEU:N	1:B:320:LEU:CD1	2.83	0.40
1:A:355:ASN:HB3	1:A:358:ILE:HD12	2.03	0.40
1:B:462:CYS:HB2	1:B:466:LYS:HZ1	1.87	0.40

There are no symmetry-related clashes.

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	494/496 (100%)	482 (98%)	12 (2%)	0	100	100
1	B	494/496 (100%)	480 (97%)	13 (3%)	1 (0%)	43	24
2	C	117/121 (97%)	112 (96%)	5 (4%)	0	100	100
2	D	119/121 (98%)	114 (96%)	4 (3%)	1 (1%)	16	5
All	All	1224/1234 (99%)	1188 (97%)	34 (3%)	2 (0%)	43	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	224	ALA
2	D	9	GLY

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	411/413 (100%)	390 (95%)	21 (5%)	21	5
1	B	410/413 (99%)	388 (95%)	22 (5%)	20	4
2	C	95/97 (98%)	93 (98%)	2 (2%)	47	23
2	D	97/97 (100%)	89 (92%)	8 (8%)	10	2
All	All	1013/1020 (99%)	960 (95%)	53 (5%)	21	5

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
1	A	5	GLN
1	A	35	LYS
1	A	69	LEU
1	A	77	ASN
1	A	92	ARG
1	A	115	CYS
1	A	162	LEU
1	A	168	LEU
1	A	178	MET
1	A	186	LEU
1	A	262	LEU
1	A	279	ASN
1	A	293	LEU
1	A	319	ARG
1	A	320	LEU
1	A	366	VAL
1	A	395	VAL
1	A	421	ARG
1	A	444	LEU

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Mol	Chain	Res	Type
1	A	484	GLU
1	B	1	GLN
1	B	5	GLN
1	B	77	ASN
1	B	92	ARG
1	B	140	LYS
1	B	162	LEU
1	B	168	LEU
1	B	186	LEU
1	B	211	LEU
1	B	227	ARG
1	B	262	LEU
1	B	279	ASN
1	B	293	LEU
1	B	305	HIS
1	B	319	ARG
1	B	329	LEU
1	B	375	ASP
1	B	435	GLN
1	B	440	LEU
1	B	444	LEU
1	B	460	ASN
1	B	484	GLU
2	C	35	SER
2	C	88	ASP
2	D	12	VAL
2	D	18	LEU
2	D	39	ARG
2	D	41	PRO
2	D	70	SER
2	D	73	ASN
2	D	92	MET
2	D	115	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	15	HIS
1	A	100	ASN
1	A	101	HIS
1	A	201	HIS

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Mol	Chain	Res	Type
1	A	279	ASN
1	A	299	HIS
1	A	350	ASN
1	A	364	ASN
1	A	399	ASN
1	B	1	GLN
1	B	5	GLN
1	B	15	HIS
1	B	75	ASN
1	B	101	HIS
1	B	201	HIS
1	B	279	ASN
1	B	299	HIS
1	B	350	ASN
1	B	373	ASN
1	B	399	ASN
1	B	435	GLN
1	B	476	GLN
2	C	27	ASN
2	C	81	GLN
2	D	53	ASN
2	D	73	ASN
2	D	81	GLN
2	D	116	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	D	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	1:GLN	C	2:VAL	N	1.18
1	D	8:GLY	C	9:GLY	N	0.90

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.