



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:36 AM UTC

PDB ID : 1KXQ / pdb_00001kxq
Title : Camelid VHH Domain 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

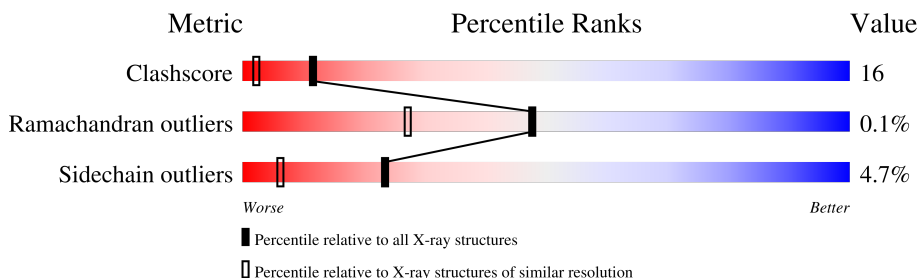
1 Overall quality at a glance

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	
1	B	496	
1	C	496	
1	D	496	
2	E	120	
2	F	120	
2	G	120	
2	H	120	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 22119 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
			3910	2471	688	730	21			
1	B	496	Total	C	N	O	S	0	0	0
			3910	2471	688	730	21			
1	C	496	Total	C	N	O	S	0	0	0
			3910	2471	688	730	21			
1	D	496	Total	C	N	O	S	0	0	0
			3910	2471	688	730	21			

There are 16 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
C	243	LYS	GLN	SEE REMARK 999	UNP P00690
C	310	SER	ALA	SEE REMARK 999	UNP P00690
C	323	ILE	VAL	SEE REMARK 999	UNP P00690
C	404	GLN	GLU	SEE REMARK 999	UNP P00690
D	243	LYS	GLN	SEE REMARK 999	UNP P00690
D	310	SER	ALA	SEE REMARK 999	UNP P00690
D	323	ILE	VAL	SEE REMARK 999	UNP P00690
D	404	GLN	GLU	SEE REMARK 999	UNP P00690

- Molecule 2 is a protein called antibody VHH fragment CABAMD9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	120	Total	C	N	O	S	0	0	0
			916	570	158	185	3			
2	F	120	Total	C	N	O	S	0	0	0
			916	570	158	185	3			
2	G	120	Total	C	N	O	S	0	0	0
			916	570	158	185	3			
2	H	120	Total	C	N	O	S	0	0	0
			916	570	158	185	3			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	B	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		

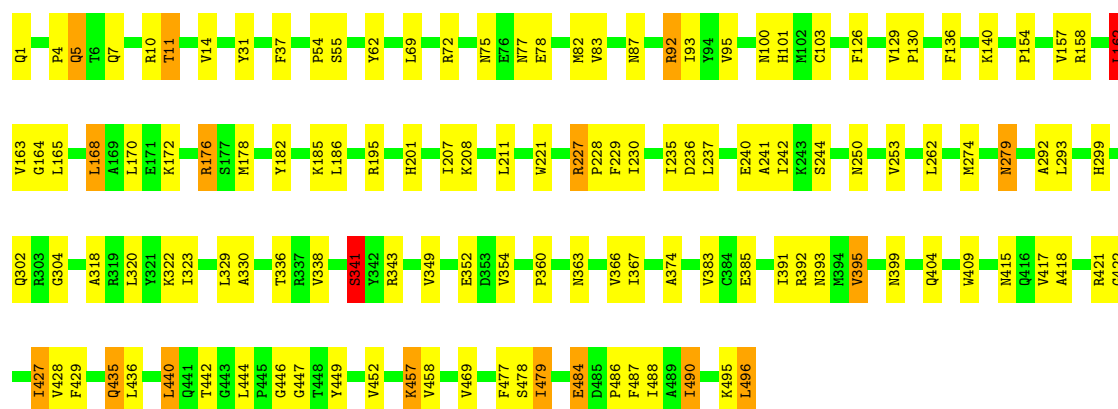
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	586	Total	O	0	0
			586	586		
5	B	540	Total	O	0	0
			540	540		
5	C	508	Total	O	0	0
			508	508		

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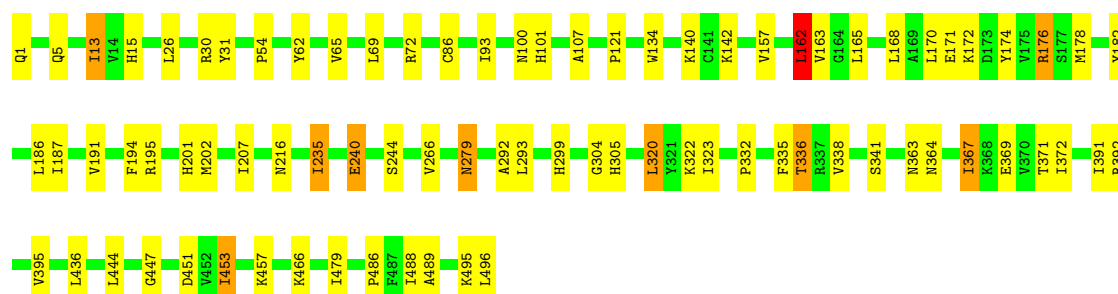
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	635	Total 635	O 635	0	0
5	E	143	Total 143	O 143	0	0
5	F	137	Total 137	O 137	0	0
5	G	144	Total 144	O 144	0	0
5	H	114	Total 114	O 114	0	0



- Molecule 1: alpha-amylase, pancreatic

Chain D: 83% 15% .



- Molecule 2: antibody VHH fragment CABAMD9

Chain E: 81% 16% .



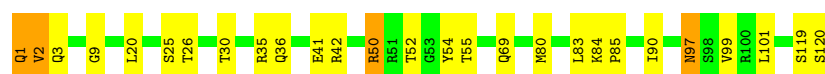
- Molecule 2: antibody VHH fragment CABAMD9

Chain F: 78% 18% .



- Molecule 2: antibody VHH fragment CABAMD9

Chain G: 78% 19% .



- Molecule 2: antibody VHH fragment CABAMD9

Chain H: 71% 26% .

Q1	V2	Q3	G8	G9	Q13	L18	S19	L20	S25	D29	T30	Q36	A37	P38	G39	K40	E41	I48	Y49	R50	S57	V61	R64	F65	Q69	M80	L83	K84	P85	E86	D87	N97	S98	V99	R100	L101	F108	Q112	G113	T114	S119	S120
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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	65.17Å 100.92Å 103.74Å 79.31° 72.63° 86.11°	Depositor
Resolution (Å)	29.79 – 1.60	Depositor
% Data completeness (in resolution range)	96.8 (29.79-1.60)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.197 , 0.219	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	22119	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

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: CL, CA

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.52	0/4020	0.99	13/5461 (0.2%)
1	B	0.49	0/4020	0.98	15/5461 (0.3%)
1	C	0.47	0/4020	0.97	15/5461 (0.3%)
1	D	0.50	0/4020	0.97	13/5461 (0.2%)
2	E	0.43	0/935	0.88	2/1267 (0.2%)
2	F	0.46	0/935	0.87	1/1267 (0.1%)
2	G	0.46	0/935	0.88	2/1267 (0.2%)
2	H	0.42	0/935	0.85	3/1267 (0.2%)
All	All	0.49	0/19820	0.96	64/26912 (0.2%)

There are no bond length outliers.

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	304	GLY	N-CA-C	8.68	125.02	114.69
1	C	304	GLY	N-CA-C	7.86	124.05	114.69
1	B	304	GLY	N-CA-C	7.61	123.75	114.69
1	C	11	THR	N-CA-C	7.39	122.42	112.88
1	A	318	ALA	N-CA-C	7.35	119.93	111.11
1	C	336	THR	N-CA-C	7.11	121.27	109.46
1	D	292	ALA	N-CA-C	7.09	121.38	109.76
1	D	336	THR	N-CA-C	6.91	120.94	109.46
1	B	11	THR	N-CA-C	6.82	122.55	112.94
1	C	55	SER	N-CA-C	6.77	119.46	108.63
1	B	336	THR	N-CA-C	6.68	121.19	109.96
1	A	6	THR	N-CA-C	-6.66	101.93	110.53
1	A	55	SER	N-CA-C	6.58	119.54	108.23
1	B	292	ALA	N-CA-C	6.55	120.50	109.76
1	D	304	GLY	N-CA-C	6.54	123.72	115.21
1	C	292	ALA	N-CA-C	6.53	120.48	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	395	VAL	N-CA-C	-6.50	104.15	110.72
1	B	195	ARG	N-CA-C	-6.44	98.24	108.73
1	A	292	ALA	N-CA-C	6.33	120.13	109.76
1	B	306	GLY	N-CA-C	-6.33	105.20	112.79
1	C	195	ARG	N-CA-C	-6.25	98.55	108.73
1	A	336	THR	N-CA-C	6.18	120.34	109.96
1	B	194	PHE	N-CA-C	5.99	118.52	109.23
1	A	195	ARG	N-CA-C	-5.98	99.45	108.96
1	A	11	THR	N-CA-C	5.96	121.35	112.94
1	D	293	LEU	N-CA-C	-5.94	98.84	108.52
1	A	216	ASN	N-CA-C	-5.92	102.72	110.53
2	H	50	ARG	N-CA-C	5.91	120.10	113.01
1	D	195	ARG	N-CA-C	-5.89	99.14	108.73
1	D	216	ASN	N-CA-C	-5.87	102.78	110.53
1	B	318	ALA	N-CA-C	5.78	118.05	111.11
1	C	341	SER	N-CA-C	5.72	117.92	109.69
1	B	293	LEU	N-CA-C	-5.70	99.24	108.52
2	H	99	VAL	N-CA-C	-5.67	107.21	112.83
1	C	162	LEU	N-CA-C	-5.57	101.14	109.15
2	G	50	ARG	N-CA-C	5.56	119.69	113.01
1	B	341	SER	N-CA-C	5.54	117.67	109.69
1	B	143	THR	N-CA-C	5.54	118.69	110.48
1	D	395	VAL	N-CA-C	-5.53	105.13	110.72
1	C	457	LYS	N-CA-C	-5.51	100.20	108.96
1	C	302	GLN	N-CA-C	-5.46	106.46	113.01
1	C	293	LEU	N-CA-C	-5.45	99.53	108.41
2	E	50	ARG	N-CA-C	5.44	119.40	112.87
1	B	185	LYS	N-CA-C	-5.44	105.25	111.07
2	F	50	ARG	N-CA-C	5.33	119.04	112.54
1	C	103	CYS	N-CA-C	5.31	116.78	110.19
1	A	293	LEU	N-CA-C	-5.30	99.87	108.52
1	D	162	LEU	N-CA-C	-5.29	101.53	109.15
1	D	65	VAL	N-CA-C	-5.29	108.69	113.71
1	A	194	PHE	N-CA-C	5.28	117.41	109.23
1	D	194	PHE	N-CA-C	5.26	117.38	109.23
2	H	29	ASP	N-CA-C	5.25	117.08	111.36
1	A	341	SER	N-CA-C	5.24	117.24	109.69
2	G	99	VAL	N-CA-C	-5.23	107.65	112.83
1	D	341	SER	N-CA-C	5.22	117.20	109.69
2	E	99	VAL	N-CA-C	-5.20	107.68	112.83
1	C	95	VAL	N-CA-C	5.18	115.77	108.36
1	D	332	PRO	N-CA-C	5.15	120.30	113.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(^o)	Ideal(^o)
1	B	59	TRP	N-CA-C	5.11	119.15	113.01
1	B	41	GLN	N-CA-C	-5.11	100.84	108.96
1	B	103	CYS	N-CA-C	5.11	116.53	110.19
1	A	102	MET	N-CA-C	-5.06	102.16	109.29
1	D	457	LYS	N-CA-C	-5.03	100.96	108.96
1	C	318	ALA	N-CA-C	5.00	119.62	111.37

There are no chirality outliers.

There are no planarity outliers.

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	3910	0	3693	131	0
1	B	3910	0	3693	138	0
1	C	3910	0	3693	154	0
1	D	3910	0	3693	87	0
2	E	916	0	864	16	0
2	F	916	0	864	38	0
2	G	916	0	864	18	0
2	H	916	0	864	30	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	586	0	0	16	0
5	B	540	0	0	17	0
5	C	508	0	0	12	0
5	D	635	0	0	16	0
5	E	143	0	0	2	0
5	F	137	0	0	15	0
5	G	144	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	H	114	0	0	8	0
All	All	22119	0	18228	599	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 16.

All (599) 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)
2:F:3:GLN:HG2	2:F:25:SER:HB3	1.29	1.12
1:A:26:LEU:HD21	5:A:4489:HOH:O	1.52	1.09
1:A:352:GLU:OE2	2:H:97:ASN:OD1	1.75	1.04
1:C:1:GLN:HG3	1:C:230:ILE:HD12	1.36	1.03
1:D:453:ILE:HD11	1:D:489:ALA:HB3	1.39	1.02
1:A:341:SER:O	1:A:383:VAL:HG12	1.60	1.01
1:D:100:ASN:HD22	1:D:101:HIS:HD2	1.12	0.98
1:A:409:TRP:CH2	1:A:417:VAL:HG11	2.03	0.93
1:B:409:TRP:CH2	1:B:417:VAL:HG11	2.04	0.92
1:A:5:GLN:HA	1:A:5:GLN:HE21	1.34	0.92
1:B:427:ILE:HD13	1:B:428:VAL:N	1.83	0.92
1:C:83:VAL:HG22	1:C:93:ILE:HD12	1.49	0.92
1:C:341:SER:O	1:C:383:VAL:HG12	1.70	0.91
1:A:100:ASN:HD22	1:A:101:HIS:HD2	1.17	0.91
1:C:427:ILE:HD13	1:C:428:VAL:N	1.87	0.90
1:C:170:LEU:O	1:C:176:ARG:HD2	1.72	0.89
1:C:352:GLU:HG3	2:F:97:ASN:OD1	1.72	0.89
1:B:100:ASN:HD22	1:B:101:HIS:HD2	1.16	0.88
5:A:4492:HOH:O	1:B:460:ASN:HB3	1.71	0.88
1:A:392:ARG:O	1:A:395:VAL:HG22	1.73	0.88
1:D:322:LYS:HG2	1:D:391:ILE:HD11	1.56	0.88
1:A:363:ASN:O	1:A:366:VAL:HG22	1.75	0.87
2:H:57:SER:HB3	5:H:1202:HOH:O	1.73	0.86
1:C:322:LYS:HG2	1:C:391:ILE:HD11	1.56	0.86
1:C:363:ASN:O	1:C:366:VAL:HG22	1.76	0.86
1:B:5:GLN:HA	1:B:5:GLN:HE21	1.39	0.86
1:B:274:MET:HE1	1:B:429:PHE:HB2	1.58	0.86
1:B:446:GLY:HA2	1:B:469:VAL:HG23	1.56	0.86
1:B:325:VAL:HG11	1:B:391:ILE:HD12	1.58	0.85
1:B:392:ARG:O	1:B:395:VAL:HG22	1.78	0.84
1:B:182:TYR:O	1:B:186:LEU:HD13	1.79	0.82
1:B:343:ARG:HG3	1:B:383:VAL:HG12	1.62	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:LEU:O	1:B:176:ARG:HD2	1.80	0.81
1:A:325:VAL:HG11	1:A:391:ILE:HD12	1.63	0.81
2:E:3:GLN:HG2	2:E:25:SER:HB3	1.63	0.81
1:A:341:SER:C	1:A:383:VAL:HG12	2.06	0.80
2:F:62:LYS:HB3	5:F:159:HOH:O	1.82	0.79
1:A:182:TYR:O	1:A:186:LEU:HD13	1.82	0.79
1:D:453:ILE:HD11	1:D:489:ALA:CB	2.12	0.79
1:A:279:ASN:HD22	1:A:279:ASN:H	1.31	0.79
1:B:323:ILE:HD11	1:B:486:PRO:HG2	1.63	0.79
1:C:100:ASN:HD22	1:C:101:HIS:HD2	1.31	0.79
1:C:383:VAL:HG13	1:C:385:GLU:OE2	1.82	0.79
2:F:3:GLN:CG	2:F:25:SER:HB3	2.12	0.79
1:B:355:ASN:O	1:B:358:ILE:HD13	1.83	0.78
1:B:363:ASN:O	1:B:366:VAL:HG22	1.84	0.78
1:B:279:ASN:H	1:B:279:ASN:HD22	1.29	0.77
1:D:86:CYS:HB2	1:D:93:ILE:HD11	1.66	0.77
1:C:392:ARG:O	1:C:395:VAL:HG22	1.84	0.77
1:B:86:CYS:HB2	1:B:93:ILE:HD11	1.66	0.77
1:A:392:ARG:O	1:A:395:VAL:CG2	2.32	0.77
1:C:182:TYR:O	1:C:186:LEU:HD13	1.85	0.76
1:D:13:ILE:HD12	1:D:335:PHE:CE2	2.20	0.76
1:A:446:GLY:HA2	1:A:469:VAL:HG23	1.67	0.76
1:D:279:ASN:HD22	1:D:279:ASN:H	1.33	0.76
1:B:457:LYS:HZ3	1:B:496:LEU:HD12	1.49	0.76
2:H:3:GLN:HG2	2:H:25:SER:HB3	1.66	0.76
2:H:80:MET:HE2	2:H:83:LEU:HD21	1.67	0.76
1:C:279:ASN:H	1:C:279:ASN:HD22	1.32	0.76
1:A:92:ARG:C	1:A:93:ILE:HD12	2.12	0.75
2:F:3:GLN:HG2	2:F:25:SER:CB	2.15	0.74
2:F:13:GLN:HG3	5:F:195:HOH:O	1.87	0.74
2:H:38:PRO:HA	5:H:1310:HOH:O	1.88	0.74
1:A:383:VAL:HG13	1:A:385:GLU:OE2	1.86	0.74
2:H:61:VAL:HG23	5:H:1202:HOH:O	1.87	0.74
1:C:490:ILE:HD13	1:C:490:ILE:N	2.03	0.73
1:B:69:LEU:HD21	1:B:185:LYS:HE3	1.70	0.73
1:C:487:PHE:C	1:C:488:ILE:HD12	2.14	0.73
1:B:427:ILE:HD13	1:B:428:VAL:H	1.51	0.73
1:A:100:ASN:HD22	1:A:101:HIS:CD2	2.05	0.73
1:C:427:ILE:HD12	1:C:429:PHE:CE1	2.23	0.73
1:B:142:LYS:HG3	5:B:4320:HOH:O	1.89	0.73
1:C:5:GLN:HE21	1:C:5:GLN:HA	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ASN:HD22	1:B:101:HIS:CD2	2.03	0.72
1:B:427:ILE:HD12	1:B:429:PHE:CE1	2.23	0.72
1:C:490:ILE:HD13	1:C:490:ILE:H	1.53	0.72
1:C:323:ILE:HD11	1:C:486:PRO:HG2	1.72	0.72
1:C:392:ARG:O	1:C:395:VAL:CG2	2.37	0.72
1:D:170:LEU:O	1:D:176:ARG:HD2	1.88	0.72
1:A:227:ARG:HG3	1:A:228:PRO:HD2	1.72	0.72
1:A:484:GLU:HG3	5:A:4514:HOH:O	1.89	0.72
1:A:93:ILE:HD12	1:A:93:ILE:N	2.05	0.71
1:D:100:ASN:HD22	1:D:101:HIS:CD2	2.02	0.71
1:B:13:ILE:HD12	1:B:335:PHE:HE2	1.55	0.71
1:C:1:GLN:HE22	1:C:227:ARG:HG3	1.56	0.71
2:F:4:LEU:HD21	2:F:95:THR:CG2	2.21	0.71
1:B:417:VAL:HG12	1:B:418:ALA:N	2.04	0.71
1:C:72:ARG:HD3	5:C:4277:HOH:O	1.91	0.70
1:C:484:GLU:CD	1:C:484:GLU:H	2.00	0.70
1:B:123:SER:HB2	1:B:125:GLU:HG3	1.72	0.70
1:B:325:VAL:CG1	1:B:391:ILE:HD12	2.22	0.70
2:F:112:GLN:CD	5:F:126:HOH:O	2.33	0.70
2:F:4:LEU:HD21	2:F:95:THR:HG23	1.74	0.70
2:G:3:GLN:HG2	2:G:25:SER:HB3	1.73	0.70
1:C:341:SER:C	1:C:383:VAL:HG12	2.15	0.69
2:E:9:GLY:HA2	2:E:18:LEU:HD13	1.74	0.69
1:C:1:GLN:HG2	5:C:4414:HOH:O	1.91	0.69
1:C:436:LEU:HD22	1:C:479:ILE:HD13	1.74	0.69
1:A:366:VAL:HG23	5:A:4147:HOH:O	1.91	0.69
1:C:72:ARG:CD	5:C:4277:HOH:O	2.40	0.69
1:D:453:ILE:HD12	1:D:453:ILE:N	2.07	0.69
1:A:325:VAL:CG1	1:A:391:ILE:HD12	2.23	0.69
1:B:323:ILE:HD11	1:B:486:PRO:CG	2.22	0.69
1:C:1:GLN:CG	1:C:230:ILE:HD12	2.19	0.69
1:A:391:ILE:HD13	5:A:4280:HOH:O	1.94	0.68
1:D:367:ILE:HD12	1:D:367:ILE:N	2.06	0.68
1:C:427:ILE:HD13	1:C:428:VAL:H	1.55	0.68
1:A:170:LEU:O	1:A:176:ARG:HD2	1.93	0.68
2:G:42:ARG:HD2	5:G:261:HOH:O	1.93	0.68
1:D:240:GLU:HG2	5:D:4253:HOH:O	1.93	0.68
1:C:449:TYR:HE2	1:C:469:VAL:HG11	1.59	0.67
1:D:171:GLU:HG2	5:D:4251:HOH:O	1.94	0.67
1:A:323:ILE:HD11	1:A:486:PRO:HG2	1.76	0.67
1:D:13:ILE:HD12	1:D:335:PHE:HE2	1.57	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:ARG:CZ	1:A:189:ILE:HD11	2.24	0.67
1:B:274:MET:HE1	1:B:429:PHE:CB	2.24	0.67
1:A:157:VAL:HG23	1:A:158:ARG:CD	2.25	0.67
2:G:50:ARG:HE	2:G:69:GLN:NE2	1.92	0.67
1:A:86:CYS:HB2	1:A:93:ILE:HD11	1.77	0.67
1:A:80:ARG:NH2	1:A:189:ILE:CD1	2.58	0.66
1:C:83:VAL:CG2	1:C:93:ILE:HD12	2.24	0.66
1:A:5:GLN:HA	1:A:5:GLN:NE2	2.10	0.66
1:B:457:LYS:NZ	1:B:496:LEU:HD12	2.10	0.66
1:B:31:TYR:OH	1:B:392:ARG:HG3	1.95	0.66
1:D:13:ILE:HD13	1:D:336:THR:O	1.95	0.66
1:A:157:VAL:HG23	1:A:158:ARG:HD2	1.79	0.65
1:D:13:ILE:HD13	1:D:13:ILE:H	1.60	0.65
1:A:26:LEU:HD11	5:A:4362:HOH:O	1.96	0.65
1:A:244:SER:HB2	1:A:253:VAL:HG11	1.78	0.65
1:C:201:HIS:HE1	5:C:4413:HOH:O	1.79	0.65
1:A:409:TRP:CH2	1:A:417:VAL:CG1	2.80	0.65
1:C:154:PRO:HA	1:C:157:VAL:HG22	1.79	0.65
1:C:447:GLY:HA3	1:C:495:LYS:HE3	1.77	0.65
1:D:391:ILE:N	1:D:391:ILE:HD12	2.12	0.65
2:F:50:ARG:HE	2:F:69:GLN:NE2	1.95	0.64
1:C:360:PRO:HG2	1:C:367:ILE:HD12	1.78	0.64
2:F:80:MET:HE2	2:F:83:LEU:HD21	1.79	0.64
1:C:352:GLU:CG	2:F:97:ASN:OD1	2.44	0.64
1:D:31:TYR:OH	1:D:392:ARG:HG3	1.98	0.64
1:B:366:VAL:HG23	5:B:4399:HOH:O	1.95	0.64
1:B:367:ILE:HD12	1:B:367:ILE:N	2.13	0.64
1:D:453:ILE:CD1	1:D:489:ALA:HB3	2.21	0.64
1:B:240:GLU:HG3	5:B:4145:HOH:O	1.97	0.64
1:C:479:ILE:HD12	1:C:479:ILE:N	2.13	0.63
1:C:343:ARG:NH2	1:C:343:ARG:HB3	2.13	0.63
2:E:48:ILE:HD13	2:E:49:TYR:C	2.22	0.63
1:B:1:GLN:HE22	1:B:227:ARG:HG3	1.63	0.63
1:B:438:SER:HB2	1:B:440:LEU:HD21	1.80	0.63
1:C:360:PRO:HG2	1:C:367:ILE:CD1	2.27	0.63
1:D:72:ARG:HB3	1:D:367:ILE:HD13	1.80	0.63
1:A:417:VAL:HG12	1:A:418:ALA:N	2.14	0.63
1:D:72:ARG:HD2	1:D:367:ILE:CD1	2.28	0.63
1:A:26:LEU:HD23	1:A:85:ARG:HH12	1.63	0.63
1:D:201:HIS:HE1	5:D:4380:HOH:O	1.82	0.63
2:E:50:ARG:HE	2:E:69:GLN:NE2	1.96	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:446:GLY:HA2	1:C:469:VAL:HG13	1.81	0.63
2:H:30:THR:HG23	2:H:101:LEU:HD21	1.81	0.62
1:B:86:CYS:CB	1:B:93:ILE:HD11	2.30	0.62
1:B:142:LYS:HE3	5:B:4258:HOH:O	1.98	0.62
1:B:383:VAL:CG2	1:B:385:GLU:OE2	2.47	0.62
1:C:11:THR:H	1:C:399:ASN:HD21	1.47	0.62
1:D:453:ILE:N	1:D:453:ILE:CD1	2.63	0.62
1:D:235:ILE:HD12	1:D:235:ILE:N	2.15	0.62
1:D:367:ILE:HD12	1:D:367:ILE:H	1.64	0.62
1:B:343:ARG:CG	1:B:383:VAL:HG12	2.30	0.62
1:B:75:ASN:HB2	1:B:77:ASN:OD1	2.00	0.62
2:F:65:PHE:CE2	2:F:80:MET:HE3	2.35	0.62
1:B:409:TRP:CH2	1:B:417:VAL:CG1	2.82	0.62
1:C:391:ILE:HD12	1:C:391:ILE:N	2.15	0.62
1:A:458:VAL:HG23	1:A:458:VAL:O	1.99	0.61
1:C:349:VAL:HG23	1:C:349:VAL:O	1.99	0.61
1:B:22:VAL:HG12	5:B:4103:HOH:O	1.99	0.61
1:B:69:LEU:HD21	1:B:185:LYS:CE	2.30	0.61
1:A:308:GLY:HA3	1:A:312:ILE:HD11	1.83	0.61
1:C:14:VAL:HG21	1:C:37:PHE:CE2	2.35	0.61
1:D:367:ILE:HD13	5:D:4076:HOH:O	2.00	0.61
1:C:83:VAL:HA	1:C:93:ILE:HD12	1.83	0.61
2:E:42:ARG:CZ	5:E:195:HOH:O	2.49	0.61
2:F:3:GLN:HE21	2:F:25:SER:HB3	1.65	0.60
2:F:36:GLN:HG3	5:F:172:HOH:O	2.01	0.60
1:D:72:ARG:HD2	1:D:367:ILE:HD11	1.83	0.60
1:B:72:ARG:NE	1:B:367:ILE:HD11	2.16	0.60
2:H:50:ARG:HE	2:H:69:GLN:NE2	1.99	0.60
1:B:274:MET:CE	1:B:429:PHE:HB2	2.30	0.60
1:B:69:LEU:HD22	1:B:69:LEU:N	2.16	0.60
1:C:484:GLU:CD	1:C:484:GLU:N	2.59	0.60
1:C:409:TRP:CH2	1:C:417:VAL:HG11	2.36	0.60
1:D:72:ARG:HB3	1:D:367:ILE:CD1	2.32	0.60
1:B:107:ALA:HB3	1:B:121:PRO:HG2	1.82	0.59
1:B:1:GLN:HG3	1:B:230:ILE:CG1	2.31	0.59
1:B:13:ILE:HD12	1:B:335:PHE:CE2	2.36	0.59
1:A:14:VAL:HG11	1:A:37:PHE:CE2	2.38	0.59
1:B:5:GLN:HA	1:B:5:GLN:NE2	2.15	0.59
1:B:231:PHE:CD1	1:B:293:LEU:HD22	2.38	0.59
1:A:26:LEU:HD23	1:A:85:ARG:NH1	2.18	0.59
1:C:208:LYS:HD3	1:C:250:ASN:OD1	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:372:ILE:N	1:D:372:ILE:HD12	2.17	0.59
2:G:50:ARG:HE	2:G:69:GLN:HE21	1.50	0.59
1:B:440:LEU:N	1:B:440:LEU:HD22	2.18	0.59
1:C:279:ASN:HD22	1:C:279:ASN:N	2.01	0.59
1:C:447:GLY:O	1:C:469:VAL:HG12	2.03	0.59
1:C:100:ASN:HD22	1:C:101:HIS:CD2	2.16	0.58
1:C:320:LEU:HD13	5:C:414:HOH:O	2.03	0.58
1:B:274:MET:HE3	1:B:417:VAL:HB	1.86	0.58
1:A:201:HIS:HE1	5:A:455:HOH:O	1.85	0.58
1:A:449:TYR:HE2	1:A:469:VAL:HG21	1.66	0.58
1:C:458:VAL:HG23	1:C:458:VAL:O	2.03	0.58
1:C:237:LEU:N	1:C:237:LEU:HD22	2.18	0.58
1:D:451:ASP:CG	1:D:488:ILE:HD12	2.28	0.58
1:D:170:LEU:HD11	1:D:202:MET:SD	2.43	0.58
2:E:59:ASP:HA	2:E:62:LYS:HE2	1.85	0.58
1:D:162:LEU:HD22	1:D:163:VAL:HG22	1.85	0.57
1:D:367:ILE:CD1	1:D:367:ILE:H	2.17	0.57
1:C:207:ILE:O	1:C:211:LEU:HD13	2.03	0.57
1:B:92:ARG:C	1:B:93:ILE:HD12	2.30	0.57
1:C:83:VAL:HA	1:C:93:ILE:CD1	2.34	0.57
2:F:50:ARG:HE	2:F:69:GLN:HE21	1.52	0.57
1:A:383:VAL:CG1	1:A:385:GLU:OE2	2.53	0.57
1:D:171:GLU:HG3	5:D:412:HOH:O	2.03	0.57
1:A:308:GLY:CA	1:A:312:ILE:HD11	2.35	0.56
1:D:202:MET:CE	1:D:207:ILE:HD11	2.35	0.56
1:C:1:GLN:HG3	1:C:230:ILE:CD1	2.25	0.56
1:C:75:ASN:HB2	1:C:77:ASN:OD1	2.05	0.56
1:C:383:VAL:CG1	1:C:385:GLU:OE2	2.51	0.56
1:A:447:GLY:O	1:A:469:VAL:HG22	2.05	0.56
2:H:30:THR:CG2	2:H:101:LEU:HD21	2.35	0.56
1:A:484:GLU:H	1:A:484:GLU:CD	2.14	0.56
1:A:93:ILE:N	1:A:93:ILE:CD1	2.69	0.56
1:B:11:THR:H	1:B:399:ASN:HD21	1.53	0.56
1:C:31:TYR:OH	1:C:392:ARG:HG3	2.06	0.56
1:A:153:ASP:HB3	1:A:156:GLN:HG2	1.88	0.56
1:C:449:TYR:CE2	1:C:469:VAL:HG11	2.41	0.56
1:C:274:MET:HE2	1:C:415:ASN:O	2.05	0.55
1:B:93:ILE:HD12	1:B:93:ILE:N	2.21	0.55
1:B:417:VAL:CG1	1:B:418:ALA:N	2.69	0.55
1:C:447:GLY:N	1:C:469:VAL:HG13	2.21	0.55
1:C:478:SER:C	1:C:479:ILE:HD12	2.32	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:LEU:N	1:B:237:LEU:HD22	2.21	0.55
1:B:367:ILE:N	1:B:367:ILE:CD1	2.70	0.55
1:B:458:VAL:HG23	1:B:458:VAL:O	2.07	0.55
1:B:123:SER:CB	1:B:125:GLU:HG3	2.36	0.55
2:H:20:LEU:HG	5:H:3124:HOH:O	2.06	0.55
1:A:460:ASN:CG	1:A:460:ASN:O	2.49	0.55
2:H:13:GLN:HG3	5:H:3162:HOH:O	2.05	0.55
1:A:5:GLN:O	1:A:92:ARG:HD2	2.07	0.55
1:B:391:ILE:HD13	5:B:4221:HOH:O	2.06	0.55
1:C:457:LYS:HZ2	1:C:496:LEU:HD12	1.71	0.55
1:B:14:VAL:HG21	1:B:37:PHE:CE2	2.42	0.55
1:B:162:LEU:O	1:B:165:LEU:HG	2.07	0.55
1:D:363:ASN:O	1:D:364:ASN:OD1	2.25	0.55
1:A:235:ILE:N	1:A:235:ILE:HD12	2.21	0.55
1:A:312:ILE:HD13	5:A:4011:HOH:O	2.07	0.55
1:B:162:LEU:HD22	1:B:163:VAL:HG22	1.88	0.55
1:C:479:ILE:HG12	1:C:488:ILE:HD13	1.89	0.55
1:D:323:ILE:HD11	1:D:486:PRO:HG2	1.88	0.54
2:G:30:THR:HG23	2:G:101:LEU:HD21	1.88	0.54
1:A:352:GLU:OE2	2:H:97:ASN:CG	2.50	0.54
1:A:316:TRP:CH2	1:A:343:ARG:HG3	2.42	0.54
1:B:69:LEU:HD22	1:B:69:LEU:H	1.69	0.54
2:G:26:THR:HB	5:G:252:HOH:O	2.07	0.54
1:A:341:SER:HB2	1:A:383:VAL:HG11	1.87	0.54
1:A:442:THR:O	1:A:444:LEU:HD13	2.07	0.54
1:D:1:GLN:HG2	5:D:4626:HOH:O	2.08	0.54
1:C:374:ALA:HB2	1:D:369:GLU:OE2	2.07	0.54
1:C:7:GLN:HB3	1:C:10:ARG:HD2	1.89	0.54
1:A:207:ILE:O	1:A:211:LEU:HD13	2.08	0.54
1:C:164:GLY:HA3	2:F:54:TYR:HB2	1.89	0.54
1:C:274:MET:HE3	1:C:417:VAL:HB	1.90	0.54
1:D:322:LYS:HG2	1:D:391:ILE:CD1	2.36	0.54
1:C:14:VAL:HG22	1:C:338:VAL:HB	1.90	0.54
2:H:3:GLN:CG	2:H:25:SER:HB3	2.38	0.54
2:H:61:VAL:CG2	5:H:1202:HOH:O	2.52	0.54
2:H:64:ARG:NH2	2:H:87:ASP:OD2	2.32	0.54
1:A:268:LYS:HE2	5:A:4464:HOH:O	2.08	0.53
1:C:274:MET:HE1	1:C:429:PHE:HB2	1.90	0.53
1:C:479:ILE:CG1	1:C:488:ILE:HD13	2.39	0.53
1:D:5:GLN:HB3	5:D:4100:HOH:O	2.07	0.53
1:B:72:ARG:HB3	1:B:367:ILE:HD13	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:343:ARG:HB3	1:C:343:ARG:HH21	1.72	0.53
2:F:3:GLN:NE2	5:F:178:HOH:O	2.41	0.53
1:A:479:ILE:HD11	1:A:488:ILE:HG13	1.90	0.53
1:D:170:LEU:CD1	1:D:202:MET:SD	2.97	0.53
1:D:187:ILE:HD13	1:D:191:VAL:HG23	1.89	0.53
1:A:440:LEU:HD23	1:A:477:PHE:HD2	1.73	0.53
1:A:80:ARG:NH2	1:A:189:ILE:HD11	2.23	0.53
1:A:392:ARG:C	1:A:395:VAL:HG22	2.34	0.53
1:A:428:VAL:CG1	1:A:488:ILE:HB	2.39	0.53
1:B:77:ASN:ND2	1:B:78:GLU:HG3	2.23	0.53
1:B:440:LEU:HD23	1:B:477:PHE:HD2	1.73	0.53
1:B:367:ILE:CD1	1:B:367:ILE:H	2.21	0.52
1:C:11:THR:H	1:C:399:ASN:ND2	2.08	0.52
1:A:438:SER:HB2	1:A:440:LEU:HD21	1.91	0.52
1:C:154:PRO:HA	1:C:157:VAL:CG2	2.40	0.52
2:H:8:GLY:O	2:H:114:THR:HG21	2.09	0.52
1:B:367:ILE:HD13	5:B:4126:HOH:O	2.08	0.52
2:F:37:ALA:HB2	5:F:183:HOH:O	2.10	0.52
1:A:447:GLY:N	1:A:469:VAL:HG23	2.24	0.52
1:C:409:TRP:CH2	1:C:417:VAL:CG1	2.92	0.52
1:A:478:SER:O	1:A:479:ILE:HD13	2.10	0.52
1:B:75:ASN:CB	1:B:77:ASN:OD1	2.57	0.52
1:D:367:ILE:N	1:D:367:ILE:CD1	2.72	0.52
1:D:453:ILE:HD13	1:D:489:ALA:H	1.74	0.52
1:C:235:ILE:N	1:C:235:ILE:HD12	2.25	0.51
2:G:55:THR:HB	5:G:203:HOH:O	2.09	0.51
1:B:392:ARG:O	1:B:395:VAL:CG2	2.56	0.51
1:C:447:GLY:H	1:C:469:VAL:HG13	1.75	0.51
1:A:80:ARG:HG2	1:A:189:ILE:HD12	1.92	0.51
1:D:72:ARG:CD	1:D:367:ILE:HD11	2.39	0.51
2:H:3:GLN:HG2	2:H:25:SER:CB	2.38	0.51
1:A:484:GLU:CD	1:A:484:GLU:N	2.69	0.51
1:D:13:ILE:HD13	1:D:13:ILE:N	2.26	0.51
2:E:97:ASN:C	2:E:97:ASN:HD22	2.19	0.51
1:C:82:MET:HE3	1:C:93:ILE:HD13	1.92	0.50
1:D:436:LEU:C	1:D:436:LEU:HD23	2.36	0.50
2:E:50:ARG:HE	2:E:69:GLN:HE21	1.57	0.50
1:B:13:ILE:HD13	1:B:13:ILE:H	1.76	0.50
1:C:62:TYR:O	1:C:101:HIS:HE1	1.94	0.50
1:C:366:VAL:HG23	1:C:366:VAL:O	2.11	0.50
2:F:4:LEU:HD21	2:F:95:THR:HG21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:GLY:C	1:A:312:ILE:CD1	2.85	0.50
1:C:72:ARG:HD2	5:C:4277:HOH:O	2.07	0.50
1:D:235:ILE:N	1:D:235:ILE:CD1	2.73	0.50
1:D:305:HIS:CG	1:D:305:HIS:O	2.65	0.50
2:G:84:LYS:HB2	2:G:85:PRO:HD2	1.94	0.50
1:A:77:ASN:HD22	1:B:458:VAL:HG23	1.76	0.50
1:A:349:VAL:O	1:A:349:VAL:HG13	2.12	0.50
1:B:77:ASN:CG	1:B:78:GLU:N	2.69	0.50
1:B:357:TRP:CE2	1:B:358:ILE:HD12	2.46	0.50
1:C:440:LEU:N	1:C:440:LEU:HD22	2.27	0.50
2:F:3:GLN:HB3	5:F:249:HOH:O	2.11	0.50
2:H:97:ASN:C	2:H:97:ASN:HD22	2.20	0.50
1:A:157:VAL:HG23	1:A:158:ARG:HD3	1.93	0.50
1:B:1:GLN:HG3	1:B:230:ILE:HG12	1.94	0.50
1:A:349:VAL:HG11	2:H:108:PHE:CE1	2.47	0.50
1:C:77:ASN:CG	1:C:78:GLU:N	2.70	0.50
1:C:101:HIS:CE1	1:C:165:LEU:HD13	2.47	0.50
1:D:182:TYR:O	1:D:186:LEU:HD13	2.11	0.50
1:A:120:ASN:HB3	1:A:125:GLU:HB2	1.94	0.50
1:C:4:PRO:HA	1:C:229:PHE:CG	2.47	0.50
2:F:64:ARG:HD3	5:F:234:HOH:O	2.11	0.49
1:A:279:ASN:HD22	1:A:279:ASN:N	1.98	0.49
1:C:322:LYS:HG2	1:C:391:ILE:CD1	2.36	0.49
1:A:237:LEU:N	1:A:237:LEU:HD22	2.27	0.49
1:B:440:LEU:HD23	1:B:477:PHE:CD2	2.48	0.49
1:D:93:ILE:HD12	1:D:93:ILE:N	2.27	0.49
2:G:1:GLN:HG3	2:G:1:GLN:O	2.10	0.49
1:A:274:MET:CE	1:A:429:PHE:HB2	2.42	0.49
1:B:236:ASP:O	1:B:257:LYS:NZ	2.42	0.49
2:H:13:GLN:HA	2:H:119:SER:HG	1.77	0.49
1:A:274:MET:HE2	1:A:415:ASN:O	2.13	0.49
1:A:417:VAL:CG1	1:A:418:ALA:N	2.76	0.49
1:C:75:ASN:ND2	5:C:4145:HOH:O	2.44	0.49
1:B:343:ARG:HB3	1:B:343:ARG:NH2	2.28	0.49
1:A:449:TYR:CE2	1:A:469:VAL:HG21	2.46	0.49
1:C:469:VAL:HG13	1:C:469:VAL:O	2.13	0.49
1:D:391:ILE:N	1:D:391:ILE:CD1	2.76	0.49
2:F:2:VAL:HG13	5:F:221:HOH:O	2.11	0.49
2:F:64:ARG:NH1	5:F:182:HOH:O	2.44	0.49
2:G:97:ASN:C	2:G:97:ASN:HD22	2.19	0.49
1:A:31:TYR:OH	1:A:392:ARG:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:PRO:O	1:C:157:VAL:HG22	2.13	0.48
1:B:358:ILE:HD12	1:B:358:ILE:N	2.27	0.48
1:B:383:VAL:HG23	1:B:385:GLU:OE2	2.13	0.48
2:F:3:GLN:HE21	2:F:25:SER:CB	2.25	0.48
1:A:440:LEU:HD23	1:A:477:PHE:CD2	2.47	0.48
1:B:75:ASN:ND2	5:B:4187:HOH:O	2.45	0.48
1:B:305:HIS:HD2	5:B:4313:HOH:O	1.96	0.48
1:C:186:LEU:N	1:C:186:LEU:CD1	2.76	0.48
1:C:330:ALA:O	1:C:421:ARG:NE	2.46	0.48
1:C:490:ILE:N	1:C:490:ILE:CD1	2.74	0.48
1:A:11:THR:H	1:A:399:ASN:HD21	1.62	0.48
1:A:86:CYS:CB	1:A:93:ILE:HD11	2.43	0.48
1:A:421:ARG:NH1	5:A:4015:HOH:O	2.46	0.48
1:B:77:ASN:CG	1:B:78:GLU:H	2.21	0.48
1:B:201:HIS:HE1	5:B:4431:HOH:O	1.95	0.48
1:C:1:GLN:OE1	1:C:227:ARG:HG2	2.14	0.48
1:A:274:MET:HE1	1:A:429:PHE:CB	2.44	0.48
2:E:14:ALA:HB2	2:E:119:SER:O	2.13	0.48
2:F:3:GLN:CD	5:F:179:HOH:O	2.56	0.48
1:A:140:LYS:HZ2	1:A:140:LYS:HB2	1.78	0.48
1:C:449:TYR:HE2	1:C:469:VAL:CG1	2.27	0.48
2:E:1:GLN:HG3	2:E:1:GLN:O	2.14	0.48
1:B:366:VAL:HG23	1:B:366:VAL:O	2.12	0.48
1:C:242:ILE:HD12	1:C:242:ILE:N	2.29	0.48
1:C:392:ARG:C	1:C:395:VAL:HG22	2.39	0.48
1:D:142:LYS:HE2	5:D:4491:HOH:O	2.12	0.48
2:H:1:GLN:O	2:H:1:GLN:HG3	2.13	0.48
2:H:13:GLN:HA	2:H:119:SER:OG	2.14	0.48
2:H:84:LYS:HB2	2:H:85:PRO:HD2	1.95	0.48
1:A:440:LEU:HD22	1:A:440:LEU:N	2.29	0.48
1:D:371:THR:C	1:D:372:ILE:HD12	2.39	0.48
2:E:12:VAL:HG23	5:E:184:HOH:O	2.13	0.48
1:A:154:PRO:HA	1:A:157:VAL:HG22	1.96	0.48
1:B:279:ASN:H	1:B:279:ASN:ND2	2.06	0.48
1:A:186:LEU:N	1:A:186:LEU:CD1	2.76	0.47
1:B:484:GLU:CD	1:B:485:ASP:H	2.22	0.47
1:C:5:GLN:HA	1:C:5:GLN:NE2	2.26	0.47
1:C:236:ASP:HA	1:C:240:GLU:OE2	2.13	0.47
2:F:2:VAL:HG13	2:F:2:VAL:O	2.14	0.47
1:A:140:LYS:HB2	1:A:140:LYS:NZ	2.30	0.47
1:B:13:ILE:HD13	1:B:336:THR:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:360:PRO:CG	1:C:367:ILE:HD12	2.44	0.47
1:D:479:ILE:HD11	1:D:488:ILE:CG1	2.43	0.47
1:B:357:TRP:CD2	1:B:358:ILE:HD12	2.50	0.47
1:C:1:GLN:HE22	1:C:227:ARG:CG	2.25	0.47
1:D:338:VAL:HG13	5:D:4390:HOH:O	2.14	0.47
1:A:349:VAL:HG12	1:A:354:VAL:HG13	1.95	0.47
1:B:305:HIS:CE1	5:B:4436:HOH:O	2.68	0.47
1:D:299:HIS:HD2	5:D:4021:HOH:O	1.97	0.47
1:B:1:GLN:HG3	1:B:230:ILE:HG13	1.95	0.47
1:B:279:ASN:HD22	1:B:279:ASN:N	1.99	0.47
1:B:447:GLY:HA3	1:B:495:LYS:HE3	1.97	0.47
1:C:366:VAL:HG23	5:C:4473:HOH:O	2.14	0.47
1:D:279:ASN:HD22	1:D:279:ASN:N	2.00	0.47
1:D:488:ILE:HD13	5:D:4055:HOH:O	2.14	0.47
2:H:48:ILE:HD13	2:H:49:TYR:C	2.39	0.47
1:A:312:ILE:HD12	1:A:312:ILE:H	1.79	0.47
1:D:202:MET:HE3	1:D:207:ILE:HD11	1.96	0.47
2:H:50:ARG:HE	2:H:69:GLN:HE21	1.62	0.47
1:D:447:GLY:HA3	1:D:495:LYS:HE3	1.96	0.47
1:A:449:TYR:HE2	1:A:469:VAL:CG2	2.28	0.47
1:A:366:VAL:HG23	1:A:366:VAL:O	2.15	0.46
2:F:84:LYS:HB2	2:F:85:PRO:HD2	1.97	0.46
2:F:2:VAL:CG1	5:F:221:HOH:O	2.62	0.46
2:E:48:ILE:HD13	2:E:48:ILE:C	2.41	0.46
1:B:1:GLN:HG2	5:B:4370:HOH:O	2.15	0.46
1:B:436:LEU:C	1:B:436:LEU:HD23	2.41	0.46
2:G:119:SER:O	2:G:120:SER:CB	2.64	0.46
1:C:488:ILE:HD12	1:C:488:ILE:N	2.30	0.46
1:C:338:VAL:HG13	5:C:4241:HOH:O	2.16	0.46
1:A:279:ASN:H	1:A:279:ASN:ND2	2.07	0.46
2:G:2:VAL:HG13	2:G:2:VAL:O	2.16	0.46
1:A:274:MET:HE1	1:A:429:PHE:HB2	1.97	0.45
1:A:447:GLY:H	1:A:469:VAL:HG23	1.80	0.45
1:C:435:GLN:HE21	1:C:435:GLN:HB3	1.60	0.45
1:C:440:LEU:HD23	1:C:477:PHE:HD2	1.80	0.45
2:F:3:GLN:NE2	2:F:25:SER:HB3	2.30	0.45
2:G:2:VAL:HG13	5:G:256:HOH:O	2.15	0.45
2:H:36:GLN:OE1	2:H:37:ALA:O	2.32	0.45
1:C:54:PRO:HD3	2:F:27:TYR:CE1	2.51	0.45
1:B:404:GLN:OE1	1:B:422:GLY:HA3	2.15	0.45
1:A:80:ARG:CZ	1:A:189:ILE:CD1	2.93	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:ALA:O	1:A:312:ILE:HD11	2.16	0.45
1:B:142:LYS:HB2	5:B:4258:HOH:O	2.17	0.45
1:B:208:LYS:HD3	1:B:250:ASN:OD1	2.15	0.45
1:D:86:CYS:CB	1:D:93:ILE:HD11	2.42	0.45
2:E:48:ILE:HD13	2:E:49:TYR:N	2.32	0.45
2:H:9:GLY:HA2	2:H:18:LEU:HD13	1.99	0.45
1:A:185:LYS:O	1:A:189:ILE:HG12	2.16	0.45
1:A:242:ILE:HD12	1:A:242:ILE:N	2.31	0.45
1:A:404:GLN:OE1	1:A:422:GLY:HA3	2.17	0.45
1:B:4:PRO:HA	1:B:229:PHE:CG	2.52	0.45
1:C:5:GLN:O	1:C:92:ARG:HD2	2.17	0.45
1:C:157:VAL:HG23	1:C:158:ARG:HG3	1.99	0.45
1:C:299:HIS:HD2	5:C:4041:HOH:O	1.99	0.45
2:G:9:GLY:N	2:G:20:LEU:HD23	2.31	0.45
1:A:469:VAL:HG23	1:A:469:VAL:O	2.16	0.45
1:B:72:ARG:CZ	1:B:367:ILE:HD11	2.47	0.45
1:B:158:ARG:HD3	1:B:203:TRP:CD2	2.52	0.45
1:C:442:THR:O	1:C:444:LEU:HD13	2.16	0.45
2:G:80:MET:HE2	2:G:83:LEU:HD21	1.99	0.45
1:B:241:ALA:HB3	5:B:4496:HOH:O	2.16	0.44
1:C:136:PHE:CE1	1:C:168:LEU:HD13	2.52	0.44
1:C:162:LEU:HD22	1:C:163:VAL:HG22	1.99	0.44
5:A:4486:HOH:O	1:B:77:ASN:HB3	2.16	0.44
1:B:457:LYS:NZ	1:B:496:LEU:HA	2.32	0.44
1:C:391:ILE:N	1:C:391:ILE:CD1	2.81	0.44
2:F:13:GLN:CG	5:F:195:HOH:O	2.58	0.44
2:F:112:GLN:CG	5:F:126:HOH:O	2.63	0.44
1:C:7:GLN:HB3	1:C:10:ARG:CD	2.46	0.44
1:C:77:ASN:CG	1:C:78:GLU:H	2.25	0.44
1:D:466:LYS:HE2	5:D:4399:HOH:O	2.16	0.44
1:B:11:THR:H	1:B:399:ASN:ND2	2.16	0.44
1:B:15:HIS:HD2	5:B:4008:HOH:O	2.01	0.44
1:B:170:LEU:CD1	1:B:202:MET:SD	3.06	0.44
1:A:1:GLN:OE1	1:A:227:ARG:HG2	2.17	0.44
1:A:14:VAL:HG12	1:A:338:VAL:HB	2.00	0.44
1:B:229:PHE:CE2	1:B:252:ARG:HD3	2.53	0.44
1:C:392:ARG:O	1:C:395:VAL:HG23	2.14	0.44
2:E:65:PHE:HA	2:E:79:GLN:O	2.18	0.44
1:D:299:HIS:HE1	5:D:4216:HOH:O	1.99	0.44
1:C:393:ASN:HB3	1:C:452:VAL:HB	2.00	0.44
1:B:126:PHE:O	1:B:130:PRO:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:439:THR:C	1:B:440:LEU:HD22	2.43	0.44
1:C:436:LEU:HD23	1:C:436:LEU:C	2.42	0.44
1:D:30:ARG:NH1	1:D:372:ILE:HD11	2.32	0.44
1:D:479:ILE:HD11	1:D:488:ILE:HG13	2.00	0.44
1:A:101:HIS:CE1	1:A:165:LEU:HD13	2.52	0.43
1:C:10:ARG:HD3	5:C:4280:HOH:O	2.18	0.43
1:C:77:ASN:HA	5:D:4047:HOH:O	2.16	0.43
1:C:323:ILE:HD11	1:C:486:PRO:CG	2.45	0.43
1:A:458:VAL:O	1:A:458:VAL:CG2	2.66	0.43
1:B:409:TRP:CZ3	1:B:417:VAL:HG11	2.50	0.43
1:C:185:LYS:HB3	1:C:185:LYS:HE2	1.86	0.43
1:A:129:VAL:CG1	1:A:178:MET:HG2	2.48	0.43
1:C:154:PRO:C	1:C:157:VAL:HG22	2.43	0.43
1:C:417:VAL:HG12	1:C:418:ALA:N	2.33	0.43
1:D:54:PRO:HD3	2:E:27:TYR:CE1	2.53	0.43
1:A:26:LEU:CD2	1:A:85:ARG:HH12	2.30	0.43
1:A:227:ARG:HG3	1:A:228:PRO:CD	2.46	0.43
2:F:64:ARG:NH2	2:F:87:ASP:OD2	2.32	0.43
1:A:126:PHE:O	1:A:130:PRO:HA	2.18	0.43
1:A:154:PRO:HA	1:A:157:VAL:CG2	2.49	0.43
1:B:427:ILE:HD13	1:B:427:ILE:C	2.42	0.43
1:C:427:ILE:HD12	1:C:429:PHE:CD1	2.54	0.43
1:A:14:VAL:HG11	1:A:37:PHE:CD2	2.53	0.43
1:A:235:ILE:N	1:A:235:ILE:CD1	2.81	0.43
1:A:447:GLY:N	1:A:469:VAL:CG2	2.81	0.43
1:B:427:ILE:HG12	1:B:489:ALA:HB2	2.01	0.43
1:D:170:LEU:O	1:D:176:ARG:CD	2.63	0.43
1:B:13:ILE:HD13	1:B:13:ILE:N	2.34	0.43
1:B:484:GLU:CD	1:B:484:GLU:N	2.77	0.43
1:C:129:VAL:CG1	1:C:178:MET:HG2	2.49	0.43
1:C:211:LEU:CD2	1:C:230:ILE:HD13	2.49	0.43
1:D:172:LYS:HE2	1:D:172:LYS:HB3	1.68	0.43
1:B:358:ILE:HD12	1:B:358:ILE:H	1.83	0.43
1:C:274:MET:CE	1:C:429:PHE:HB2	2.48	0.43
1:D:451:ASP:OD1	1:D:488:ILE:CD1	2.66	0.43
1:A:14:VAL:CG1	1:A:37:PHE:CD2	3.02	0.42
1:A:446:GLY:CA	1:A:469:VAL:HG23	2.42	0.42
1:B:22:VAL:HG11	5:B:4327:HOH:O	2.19	0.42
1:B:72:ARG:HB3	1:B:367:ILE:CD1	2.48	0.42
2:F:12:VAL:HG22	2:F:13:GLN:N	2.34	0.42
1:A:142:LYS:HE3	5:A:4434:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:449:TYR:CE1	1:C:495:LYS:HB2	2.54	0.42
1:A:15:HIS:HD2	5:A:4023:HOH:O	2.02	0.42
1:A:435:GLN:HG2	1:A:436:LEU:N	2.35	0.42
1:C:417:VAL:CG1	1:C:418:ALA:N	2.82	0.42
2:H:3:GLN:CD	5:H:1746:HOH:O	2.63	0.42
1:C:241:ALA:C	1:C:242:ILE:HD12	2.44	0.42
1:C:440:LEU:HD23	1:C:477:PHE:CD2	2.54	0.42
1:B:164:GLY:HA3	2:G:54:TYR:HB2	2.02	0.42
1:B:323:ILE:HG23	1:B:429:PHE:CD2	2.55	0.42
1:C:126:PHE:O	1:C:130:PRO:HA	2.20	0.42
1:C:349:VAL:HG11	2:F:2:VAL:HG11	2.02	0.42
1:B:427:ILE:CD1	1:B:428:VAL:N	2.69	0.42
1:C:352:GLU:OE2	1:C:354:VAL:HG12	2.20	0.42
1:D:279:ASN:H	1:D:279:ASN:ND2	2.09	0.42
1:D:451:ASP:CG	1:D:488:ILE:CD1	2.91	0.42
2:H:40:LYS:HB3	5:H:3153:HOH:O	2.18	0.42
2:H:65:PHE:CE2	2:H:80:MET:HE3	2.55	0.42
1:C:274:MET:HE1	1:C:429:PHE:CB	2.49	0.42
1:D:62:TYR:O	1:D:101:HIS:HE1	2.03	0.42
1:B:161:GLN:HE21	1:B:164:GLY:HA2	1.85	0.42
1:C:14:VAL:CG2	1:C:37:PHE:CE2	3.03	0.42
1:D:15:HIS:HD2	5:D:4011:HOH:O	2.03	0.42
1:C:172:LYS:O	1:C:176:ARG:HD3	2.20	0.41
1:C:244:SER:HB2	1:C:253:VAL:HG11	2.01	0.41
1:C:299:HIS:HE1	5:C:4070:HOH:O	2.01	0.41
1:C:349:VAL:O	1:C:349:VAL:CG2	2.68	0.41
1:A:236:ASP:O	1:A:257:LYS:NZ	2.52	0.41
1:B:157:VAL:O	1:B:201:HIS:HD2	2.02	0.41
1:D:451:ASP:OD2	1:D:453:ILE:HD13	2.20	0.41
1:D:30:ARG:CZ	1:D:372:ILE:HD11	2.51	0.41
1:D:174:TYR:CZ	1:D:178:MET:HE3	2.55	0.41
1:C:427:ILE:HD13	1:C:427:ILE:C	2.45	0.41
1:D:101:HIS:CE1	1:D:165:LEU:HD13	2.55	0.41
1:D:157:VAL:O	1:D:201:HIS:HD2	2.03	0.41
2:E:112:GLN:HE21	2:E:112:GLN:HB3	1.62	0.41
1:B:14:VAL:CG2	1:B:37:PHE:CD2	3.03	0.41
1:B:31:TYR:O	1:B:35:LYS:HB2	2.21	0.41
1:D:187:ILE:HD13	1:D:191:VAL:CG2	2.50	0.41
1:A:261:LYS:HA	1:A:261:LYS:HD3	1.94	0.41
1:B:87:ASN:HB3	1:B:221:TRP:CD1	2.56	0.41
1:C:154:PRO:CA	1:C:157:VAL:HG22	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:72:ARG:NE	1:D:367:ILE:HD11	2.36	0.41
1:D:107:ALA:HB3	1:D:121:PRO:HG2	2.03	0.41
1:D:266:VAL:HG12	1:D:320:LEU:HG	2.02	0.41
1:A:479:ILE:HD11	1:A:488:ILE:CG1	2.51	0.41
1:C:87:ASN:HB3	1:C:221:TRP:CD1	2.56	0.41
1:C:352:GLU:OE2	1:C:354:VAL:CG1	2.69	0.41
1:C:427:ILE:CD1	1:C:428:VAL:N	2.72	0.41
1:D:26:LEU:HD21	5:D:4485:HOH:O	2.20	0.41
1:D:186:LEU:C	1:D:191:VAL:HG22	2.46	0.41
2:F:3:GLN:CB	5:F:249:HOH:O	2.68	0.41
2:F:59:ASP:O	2:F:62:LYS:HG2	2.20	0.41
1:A:323:ILE:HD11	1:A:486:PRO:CG	2.45	0.41
5:A:4486:HOH:O	1:B:77:ASN:CB	2.69	0.41
1:C:227:ARG:HG3	1:C:228:PRO:HD2	2.03	0.41
1:C:428:VAL:O	1:C:428:VAL:HG13	2.21	0.41
1:D:162:LEU:O	1:D:165:LEU:HG	2.21	0.41
1:A:1:GLN:OE1	1:A:227:ARG:CG	2.70	0.40
1:A:142:LYS:HE3	1:A:142:LYS:HB2	1.90	0.40
1:A:299:HIS:HE1	5:A:4114:HOH:O	2.04	0.40
1:B:229:PHE:HE2	1:B:252:ARG:CD	2.34	0.40
1:B:349:VAL:HG13	1:B:349:VAL:O	2.21	0.40
1:B:14:VAL:HG21	1:B:37:PHE:CD2	2.56	0.40
1:B:244:SER:HB2	1:B:253:VAL:HG11	2.02	0.40
1:B:449:TYR:CE1	1:B:495:LYS:HB2	2.56	0.40
1:C:186:LEU:N	1:C:186:LEU:HD12	2.35	0.40
1:A:189:ILE:HD13	1:A:189:ILE:HA	1.85	0.40
1:A:488:ILE:HD13	5:A:4063:HOH:O	2.21	0.40
1:B:165:LEU:HD23	2:G:52:THR:HG21	2.03	0.40
1:C:75:ASN:CB	1:C:77:ASN:OD1	2.69	0.40
1:B:299:HIS:HD2	5:B:4005:HOH:O	2.04	0.40
1:C:360:PRO:HG2	1:C:367:ILE:HD11	2.01	0.40
1:C:404:GLN:OE1	1:C:422:GLY:HA3	2.21	0.40
2:G:35:ARG:HA	2:G:90:ILE:O	2.22	0.40
1:A:120:ASN:ND2	1:A:123:SER:HB2	2.37	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%)	485 (98%)	9 (2%)	0	100	100
1	B	494/496 (100%)	483 (98%)	11 (2%)	0	100	100
1	C	494/496 (100%)	483 (98%)	11 (2%)	0	100	100
1	D	494/496 (100%)	484 (98%)	10 (2%)	0	100	100
2	E	118/120 (98%)	114 (97%)	4 (3%)	0	100	100
2	F	118/120 (98%)	114 (97%)	2 (2%)	2 (2%)	7	1
2	G	118/120 (98%)	115 (98%)	2 (2%)	1 (1%)	16	5
2	H	118/120 (98%)	111 (94%)	7 (6%)	0	100	100
All	All	2448/2464 (99%)	2389 (98%)	56 (2%)	3 (0%)	48	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	2	VAL
2	G	2	VAL
2	F	38	PRO

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	413/413 (100%)	393 (95%)	20 (5%)	23	6
1	B	413/413 (100%)	391 (95%)	22 (5%)	20	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	413/413 (100%)	394 (95%)	19 (5%)	24	6
1	D	413/413 (100%)	397 (96%)	16 (4%)	28	9
2	E	95/96 (99%)	90 (95%)	5 (5%)	20	4
2	F	95/96 (99%)	91 (96%)	4 (4%)	26	7
2	G	95/96 (99%)	91 (96%)	4 (4%)	26	7
2	H	95/96 (99%)	90 (95%)	5 (5%)	20	4
All	All	2032/2036 (100%)	1937 (95%)	95 (5%)	23	6

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	92	ARG
1	A	93	ILE
1	A	140	LYS
1	A	162	LEU
1	A	168	LEU
1	A	170	LEU
1	A	186	LEU
1	A	227	ARG
1	A	235	ILE
1	A	262	LEU
1	A	279	ASN
1	A	293	LEU
1	A	317	ASP
1	A	320	LEU
1	A	329	LEU
1	A	352	GLU
1	A	440	LEU
1	A	460	ASN
1	A	496	LEU
1	B	5	GLN
1	B	13	ILE
1	B	49	ILE
1	B	92	ARG
1	B	134	TRP
1	B	140	LYS
1	B	157	VAL
1	B	162	LEU
1	B	168	LEU

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Mol	Chain	Res	Type
1	B	211	LEU
1	B	227	ARG
1	B	240	GLU
1	B	262	LEU
1	B	279	ASN
1	B	320	LEU
1	B	329	LEU
1	B	349	VAL
1	B	367	ILE
1	B	383	VAL
1	B	427	ILE
1	B	440	LEU
1	B	496	LEU
1	C	5	GLN
1	C	69	LEU
1	C	92	ARG
1	C	140	LYS
1	C	162	LEU
1	C	168	LEU
1	C	176	ARG
1	C	227	ARG
1	C	262	LEU
1	C	279	ASN
1	C	329	LEU
1	C	341	SER
1	C	427	ILE
1	C	435	GLN
1	C	440	LEU
1	C	479	ILE
1	C	484	GLU
1	C	490	ILE
1	C	496	LEU
1	D	13	ILE
1	D	69	LEU
1	D	134	TRP
1	D	140	LYS
1	D	162	LEU
1	D	168	LEU
1	D	176	ARG
1	D	235	ILE
1	D	240	GLU
1	D	244	SER

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Mol	Chain	Res	Type
1	D	279	ASN
1	D	320	LEU
1	D	367	ILE
1	D	444	LEU
1	D	453	ILE
1	D	496	LEU
2	E	36	GLN
2	E	41	GLU
2	E	48	ILE
2	E	59	ASP
2	E	97	ASN
2	F	36	GLN
2	F	41	GLU
2	F	97	ASN
2	F	112	GLN
2	G	1	GLN
2	G	36	GLN
2	G	41	GLU
2	G	97	ASN
2	H	36	GLN
2	H	41	GLU
2	H	48	ILE
2	H	97	ASN
2	H	112	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (68) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	7	GLN
1	A	15	HIS
1	A	75	ASN
1	A	101	HIS
1	A	161	GLN
1	A	201	HIS
1	A	279	ASN
1	A	299	HIS
1	A	350	ASN
1	A	373	ASN
1	A	399	ASN
1	A	435	GLN
1	A	476	GLN

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Mol	Chain	Res	Type
1	B	5	GLN
1	B	7	GLN
1	B	15	HIS
1	B	75	ASN
1	B	101	HIS
1	B	161	GLN
1	B	201	HIS
1	B	279	ASN
1	B	299	HIS
1	B	350	ASN
1	B	363	ASN
1	B	373	ASN
1	B	399	ASN
1	B	435	GLN
1	B	476	GLN
1	C	5	GLN
1	C	15	HIS
1	C	75	ASN
1	C	101	HIS
1	C	161	GLN
1	C	201	HIS
1	C	279	ASN
1	C	299	HIS
1	C	350	ASN
1	C	363	ASN
1	C	373	ASN
1	C	399	ASN
1	C	435	GLN
1	C	476	GLN
1	D	1	GLN
1	D	15	HIS
1	D	101	HIS
1	D	161	GLN
1	D	201	HIS
1	D	279	ASN
1	D	299	HIS
1	D	363	ASN
1	D	373	ASN
1	D	435	GLN
1	D	476	GLN
2	E	3	GLN
2	E	69	GLN

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Mol	Chain	Res	Type
2	E	97	ASN
2	E	112	GLN
2	F	3	GLN
2	F	69	GLN
2	F	79	GLN
2	F	112	GLN
2	G	3	GLN
2	G	69	GLN
2	G	97	ASN
2	G	112	GLN
2	H	69	GLN
2	H	112	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 ⓘ

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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.