



Full wwPDB EM Validation Report ⓘ

Apr 5, 2026 – 02:29 PM UTC

PDB ID : 6KSP / pdb_00006ksp
EMDB ID : EMD-0773
Title : Rat GluD1 receptor(splayed conformation) in complex with 7-CKA and Calcium ions
Authors : Burada, A.P.; Kumar, J.
Deposited on : 2019-08-24
Resolution : 8.10 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>
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:

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

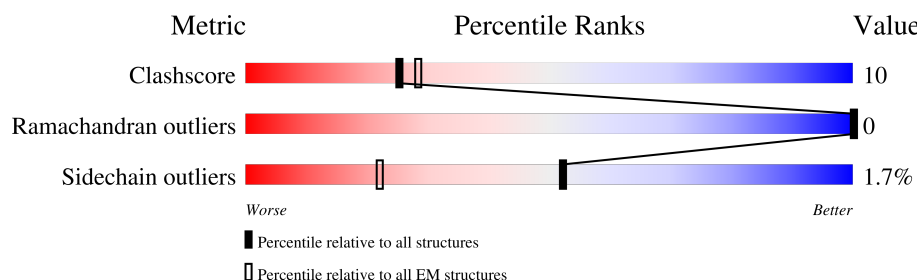
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 8.10 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)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	856	65% 22% • 12%
1	B	856	63% 23% • 12%
1	C	856	64% 23% • 12%
1	D	856	67% 21% 12%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 23744 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Glutamate receptor ionotropic, delta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	753	Total	C	N	O	S	0	0
			5936	3785	1010	1109	32		
1	B	753	Total	C	N	O	S	0	0
			5936	3785	1010	1109	32		
1	C	753	Total	C	N	O	S	0	0
			5936	3785	1010	1109	32		
1	D	753	Total	C	N	O	S	0	0
			5936	3785	1010	1109	32		

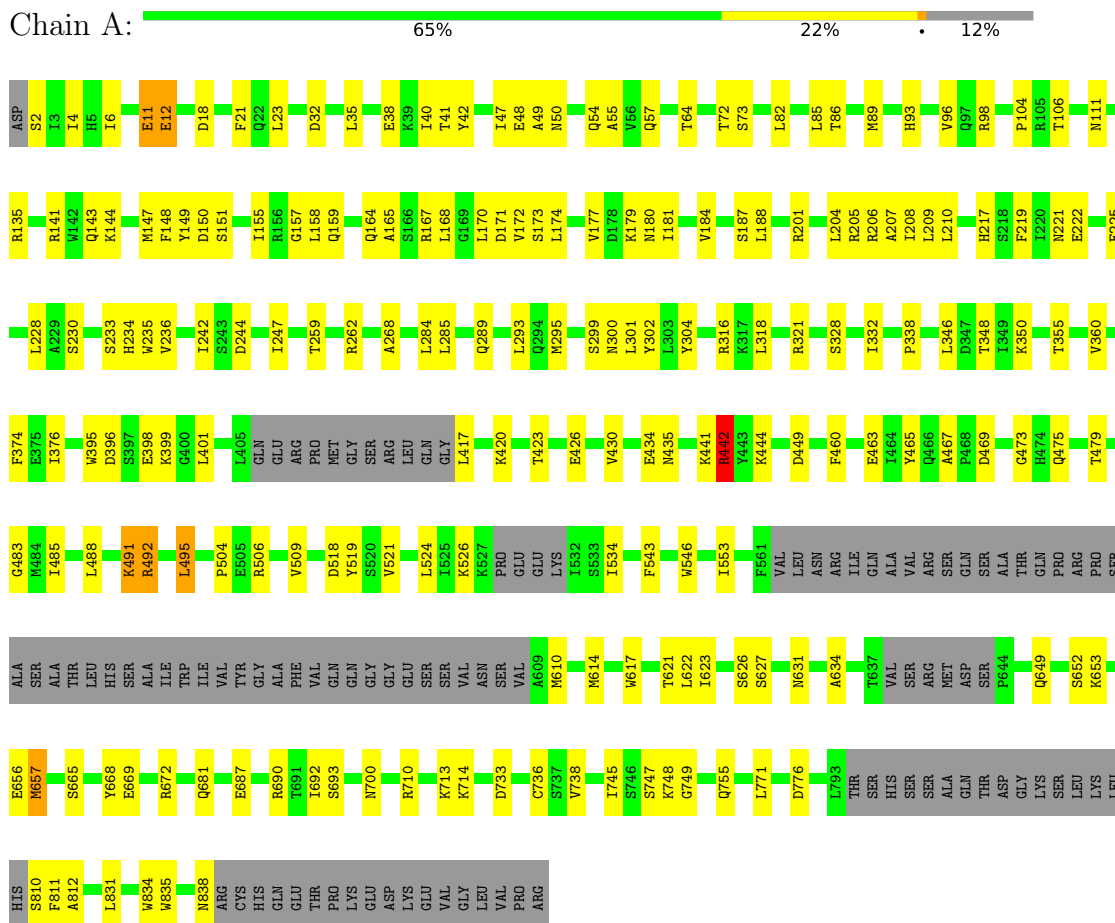
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	852	GLY	-	See sequence details	UNP Q62640
A	853	LEU	-	See sequence details	UNP Q62640
A	854	VAL	-	See sequence details	UNP Q62640
A	855	PRO	-	See sequence details	UNP Q62640
A	856	ARG	-	See sequence details	UNP Q62640
B	852	GLY	-	See sequence details	UNP Q62640
B	853	LEU	-	See sequence details	UNP Q62640
B	854	VAL	-	See sequence details	UNP Q62640
B	855	PRO	-	See sequence details	UNP Q62640
B	856	ARG	-	See sequence details	UNP Q62640
C	852	GLY	-	See sequence details	UNP Q62640
C	853	LEU	-	See sequence details	UNP Q62640
C	854	VAL	-	See sequence details	UNP Q62640
C	855	PRO	-	See sequence details	UNP Q62640
C	856	ARG	-	See sequence details	UNP Q62640
D	852	GLY	-	See sequence details	UNP Q62640
D	853	LEU	-	See sequence details	UNP Q62640
D	854	VAL	-	See sequence details	UNP Q62640
D	855	PRO	-	See sequence details	UNP Q62640
D	856	ARG	-	See sequence details	UNP Q62640

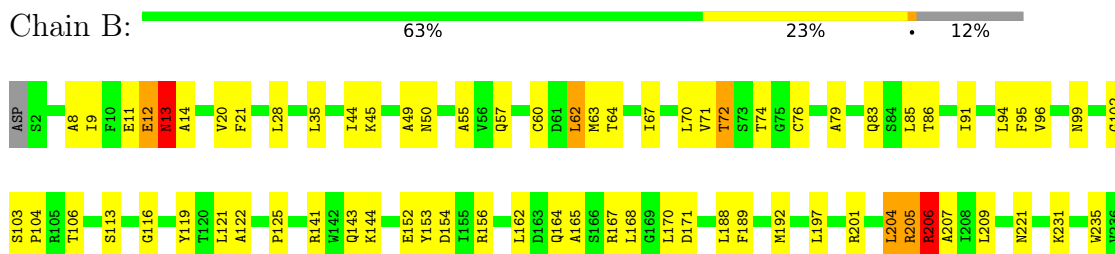
3 Residue-property plots

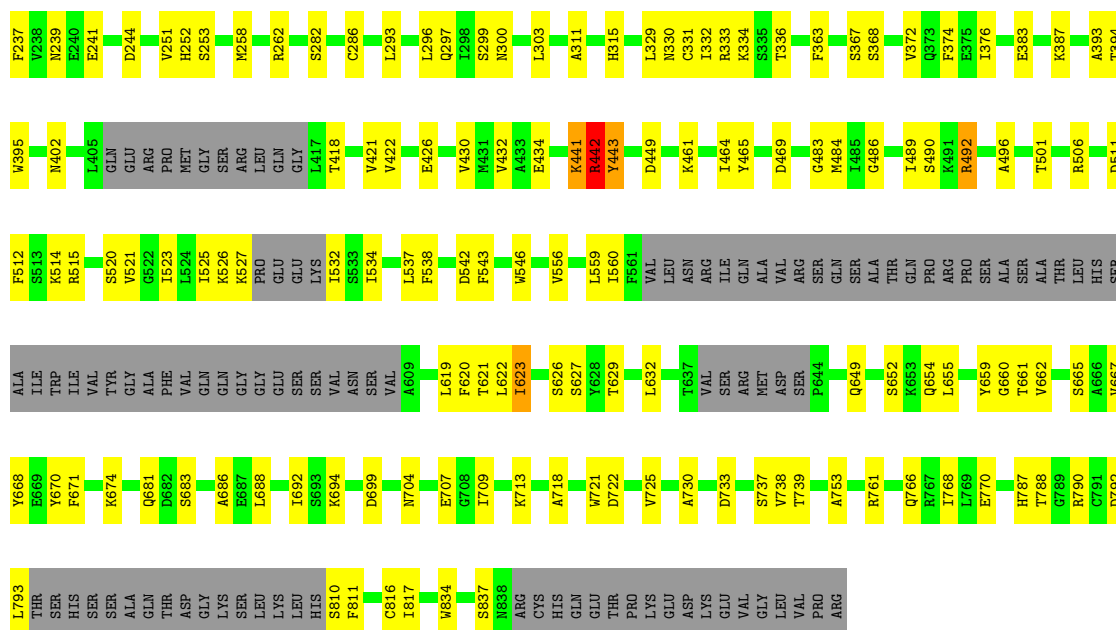
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

- Molecule 1: Glutamate receptor ionotropic, delta-1



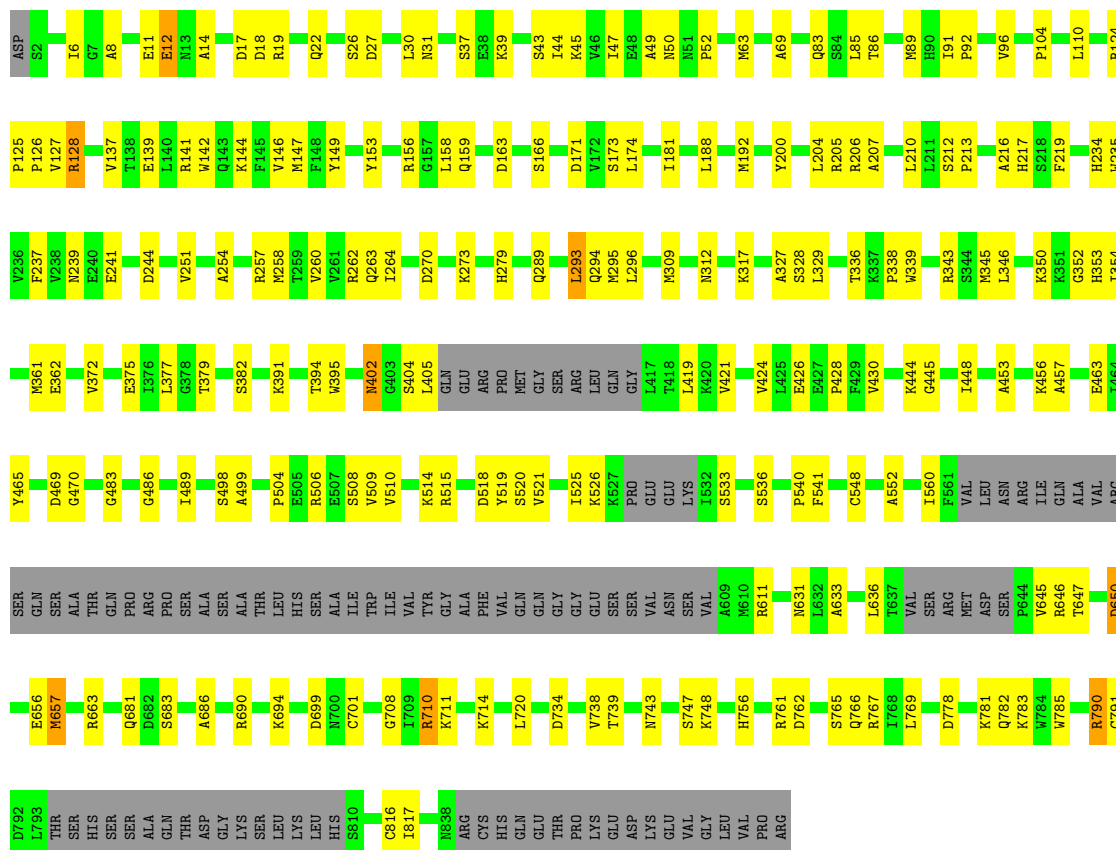
- Molecule 1: Glutamate receptor ionotropic, delta-1





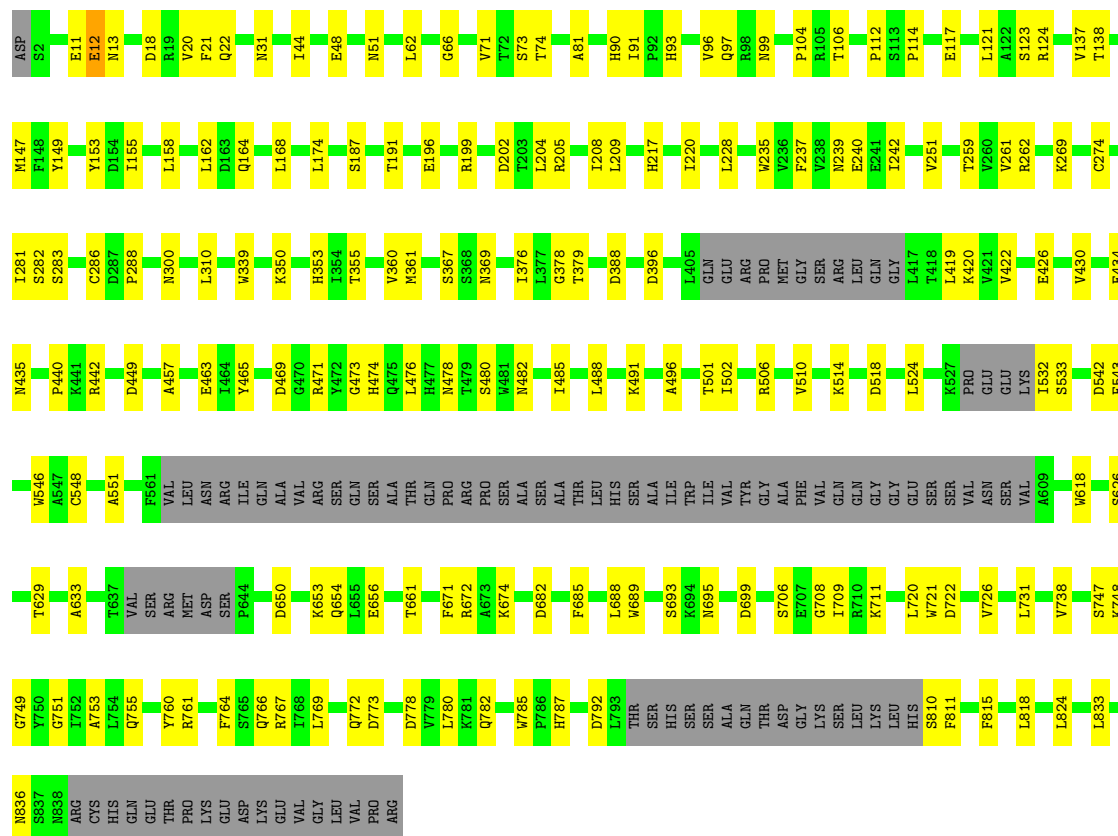
- Molecule 1: Glutamate receptor ionotropic, delta-1

Chain C: 64% 23% 12%



- Molecule 1: Glutamate receptor ionotropic, delta-1

Chain D:  67% 21% 12%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	14939	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40.38	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.26	0/6059	0.59	1/8208 (0.0%)
1	B	0.27	0/6059	0.63	4/8208 (0.0%)
1	C	0.26	0/6059	0.60	2/8208 (0.0%)
1	D	0.23	0/6059	0.58	0/8208
All	All	0.25	0/24236	0.60	7/32832 (0.0%)

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	2
1	B	0	4
1	C	0	1
All	All	0	7

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	13	ASN	N-CA-C	-8.40	92.91	110.80
1	C	650	ASP	CA-C-N	5.43	131.91	121.54
1	C	650	ASP	C-N-CA	5.43	131.91	121.54
1	A	657	MET	CB-CA-C	-5.30	110.48	116.63
1	B	441	LYS	CB-CG-CD	5.23	123.34	111.30
1	B	13	ASN	CB-CA-C	5.19	120.74	110.42
1	B	442	ARG	N-CA-C	-5.14	99.86	110.80

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	442	ARG	Sidechain
1	A	492	ARG	Sidechain
1	B	205	ARG	Sidechain
1	B	206	ARG	Sidechain
1	B	333	ARG	Sidechain
1	B	492	ARG	Sidechain
1	C	128	ARG	Sidechain

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	5936	0	5898	129	0
1	B	5936	0	5900	138	0
1	C	5936	0	5898	142	0
1	D	5936	0	5900	114	0
All	All	23744	0	23596	491	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 10.

All (491) 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:12:GLU:HG2	1:B:49:ALA:HB2	1.44	0.99
1:B:442:ARG:HA	1:B:442:ARG:NH2	1.90	0.86
1:A:434:GLU:HB2	1:A:442:ARG:HB3	1.61	0.82
1:C:445:GLY:H	1:C:783:LYS:HZ3	1.36	0.73
1:B:62:LEU:HD12	1:B:62:LEU:O	1.91	0.71
1:A:653:LYS:HD2	1:A:656:GLU:HG3	1.73	0.71
1:B:206:ARG:HB2	1:B:206:ARG:HH21	1.58	0.68
1:C:137:VAL:HG12	1:C:142:TRP:HB2	1.75	0.68
1:C:426:GLU:HG2	1:C:428:PRO:HD2	1.76	0.67
1:D:532:ILE:HG13	1:D:811:PHE:HB2	1.76	0.67
1:D:695:ASN:HD22	1:D:699:ASP:HB2	1.60	0.67
1:D:104:PRO:HB3	1:D:240:GLU:H	1.60	0.66
1:B:559:LEU:O	1:B:559:LEU:HG	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:ILE:HG12	1:B:45:LYS:HD3	1.76	0.65
1:C:681:GLN:HG3	1:C:683:SER:H	1.61	0.65
1:A:236:VAL:HA	1:A:259:THR:HB	1.79	0.64
1:B:12:GLU:CG	1:B:49:ALA:HB2	2.24	0.64
1:D:672:ARG:HG2	1:D:689:TRP:HE1	1.62	0.64
1:D:209:LEU:HB2	1:D:237:PHE:HA	1.78	0.64
1:C:19:ARG:NH2	1:C:289:GLN:OE1	2.31	0.63
1:A:11:GLU:HB2	1:A:47:ILE:HG13	1.79	0.63
1:B:206:ARG:HB2	1:B:206:ARG:NH2	2.14	0.63
1:B:209:LEU:HB2	1:B:237:PHE:HA	1.81	0.63
1:B:671:PHE:HA	1:B:674:LYS:HB2	1.81	0.63
1:C:293:LEU:O	1:C:293:LEU:HD22	1.99	0.63
1:A:518:ASP:HA	1:A:749:GLY:HA2	1.79	0.62
1:A:420:LYS:HA	1:A:463:GLU:HB2	1.81	0.62
1:A:614:MET:SD	1:A:614:MET:N	2.72	0.62
1:B:556:VAL:HA	1:B:559:LEU:HB3	1.82	0.62
1:C:317:LYS:NZ	1:C:345:MET:SD	2.74	0.61
1:A:149:TYR:HA	1:A:210:LEU:HB2	1.83	0.61
1:B:426:GLU:H	1:B:430:VAL:HB	1.66	0.61
1:B:374:PHE:HB2	1:B:395:TRP:HB3	1.83	0.61
1:A:181:ILE:HG21	1:A:219:PHE:HA	1.82	0.61
1:B:221:ASN:HD21	1:B:253:SER:HB2	1.66	0.60
1:A:150:ASP:HA	1:A:177:VAL:HB	1.82	0.60
1:A:111:ASN:HD22	1:B:50:ASN:HB3	1.67	0.60
1:A:155:ILE:O	1:A:159:GLN:NE2	2.32	0.60
1:B:442:ARG:HA	1:B:442:ARG:CZ	2.31	0.60
1:A:217:HIS:O	1:A:221:ASN:ND2	2.33	0.60
1:C:264:ILE:HD12	1:C:375:GLU:HB2	1.84	0.60
1:D:420:LYS:HA	1:D:463:GLU:HB2	1.83	0.60
1:B:62:LEU:HD11	1:B:67:ILE:HG23	1.84	0.59
1:A:188:LEU:HD11	1:B:167:ARG:HH21	1.67	0.59
1:C:8:ALA:HB3	1:C:44:ILE:HA	1.84	0.59
1:D:259:THR:HA	1:D:378:GLY:HA2	1.84	0.59
1:B:102:GLY:O	1:B:239:ASN:ND2	2.35	0.59
1:A:230:SER:HB3	1:A:233:SER:HB3	1.85	0.59
1:C:504:PRO:O	1:D:766:GLN:NE2	2.36	0.59
1:B:730:ALA:O	1:B:790:ARG:NH1	2.36	0.59
1:A:4:ILE:HB	1:A:40:ILE:HA	1.84	0.59
1:C:251:VAL:HA	1:C:254:ALA:HB3	1.83	0.59
1:A:295:MET:O	1:A:300:ASN:ND2	2.36	0.59
1:A:426:GLU:O	1:A:430:VAL:N	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:ILE:H	1:B:72:THR:HB	1.68	0.59
1:C:234:HIS:HA	1:C:257:ARG:HB2	1.85	0.59
1:C:128:ARG:HB3	1:C:395:TRP:HH2	1.67	0.58
1:D:518:ASP:HA	1:D:749:GLY:HA2	1.85	0.58
1:D:242:ILE:O	1:D:262:ARG:NH1	2.35	0.58
1:C:181:ILE:HG12	1:C:219:PHE:HB2	1.86	0.58
1:A:374:PHE:HB2	1:A:395:TRP:HB3	1.85	0.58
1:B:106:THR:HG22	1:B:156:ARG:HH21	1.68	0.58
1:C:457:ALA:O	1:C:767:ARG:NH1	2.36	0.58
1:D:833:LEU:HD23	1:D:836:ASN:HD22	1.69	0.58
1:B:201:ARG:HA	1:B:204:LEU:HB3	1.85	0.57
1:B:704:ASN:H	1:B:707:GLU:HB2	1.69	0.57
1:B:660:GLY:HA2	1:B:692:ILE:HG22	1.85	0.57
1:C:762:ASP:O	1:C:766:GLN:HB2	2.04	0.57
1:B:60:CYS:O	1:B:330:ASN:ND2	2.37	0.57
1:D:422:VAL:HA	1:D:465:TYR:HB2	1.85	0.57
1:B:442:ARG:HA	1:B:442:ARG:HH21	1.66	0.57
1:D:426:GLU:O	1:D:430:VAL:N	2.35	0.57
1:B:11:GLU:HB2	1:B:74:THR:HG22	1.87	0.57
1:B:788:THR:OG1	1:B:790:ARG:NH2	2.36	0.57
1:B:104:PRO:HG2	1:B:153:TYR:HB2	1.87	0.57
1:D:496:ALA:HB3	1:D:753:ALA:HB3	1.87	0.57
1:C:309:MET:HE1	1:C:352:GLY:HA3	1.86	0.56
1:C:645:VAL:HG22	1:C:647:THR:H	1.69	0.56
1:B:521:VAL:HG11	1:B:667:VAL:HG21	1.86	0.56
1:B:282:SER:HA	1:B:286:CYS:HB2	1.87	0.56
1:D:533:SER:O	1:D:810:SER:N	2.38	0.56
1:A:141:ARG:O	1:A:206:ARG:NH2	2.39	0.56
1:B:713:LYS:NZ	1:B:718:ALA:O	2.37	0.56
1:D:501:THR:O	1:D:506:ARG:NH2	2.39	0.56
1:C:37:SER:O	1:C:39:LYS:NZ	2.38	0.56
1:A:346:LEU:HG	1:A:350:LYS:HE2	1.88	0.56
1:D:31:ASN:ND2	1:D:282:SER:O	2.38	0.56
1:D:350:LYS:O	1:D:353:HIS:NE2	2.39	0.56
1:D:469:ASP:OD1	1:D:471:ARG:NH1	2.39	0.56
1:A:23:LEU:HD11	1:A:289:GLN:HE21	1.71	0.55
1:C:213:PRO:O	1:C:217:HIS:ND1	2.39	0.55
1:C:469:ASP:HB3	1:C:483:GLY:HA2	1.88	0.55
1:C:636:LEU:O	1:C:646:ARG:NH1	2.40	0.55
1:B:704:ASN:HB2	1:B:707:GLU:H	1.71	0.55
1:C:11:GLU:HG3	1:C:47:ILE:HD11	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:GLU:HB3	1:B:74:THR:HA	1.88	0.55
1:C:734:ASP:OD1	1:C:790:ARG:NH1	2.40	0.55
1:B:79:ALA:O	1:B:83:GLN:N	2.40	0.54
1:A:485:ILE:HG21	1:A:506:ARG:HE	1.72	0.54
1:D:71:VAL:HG11	1:D:310:LEU:HD11	1.89	0.54
1:A:6:ILE:HB	1:A:42:TYR:HB3	1.89	0.54
1:B:99:ASN:ND2	1:B:241:GLU:OE1	2.41	0.54
1:C:85:LEU:O	1:C:89:MET:N	2.40	0.54
1:D:473:GLY:HA3	1:D:485:ILE:HG12	1.89	0.54
1:A:151:SER:OG	1:A:179:LYS:NZ	2.39	0.54
1:D:761:ARG:O	1:D:764:PHE:HB2	2.08	0.54
1:B:79:ALA:HB3	1:B:95:PHE:HE1	1.72	0.54
1:B:144:LYS:HG2	1:B:171:ASP:HB2	1.90	0.54
1:C:353:HIS:ND1	1:C:362:GLU:OE2	2.41	0.54
1:C:790:ARG:NH2	1:C:791:CYS:SG	2.78	0.54
1:C:89:MET:HA	1:C:338:PRO:HG3	1.90	0.54
1:C:258:MET:H	1:C:379:THR:HB	1.73	0.54
1:C:525:ILE:O	1:C:739:THR:N	2.36	0.54
1:D:121:LEU:HA	1:D:367:SER:HB3	1.90	0.54
1:D:769:LEU:O	1:D:773:ASP:N	2.39	0.54
1:A:82:LEU:O	1:A:85:LEU:HB3	2.08	0.54
1:A:519:TYR:HB3	1:A:748:LYS:HB2	1.90	0.54
1:B:311:ALA:O	1:B:315:HIS:ND1	2.36	0.54
1:B:538:PHE:HB3	1:B:542:ASP:HB3	1.90	0.54
1:D:654:GLN:HE22	1:D:688:LEU:HD12	1.73	0.54
1:A:48:GLU:H	1:A:54:GLN:HE21	1.56	0.53
1:A:473:GLY:O	1:A:506:ARG:NH2	2.39	0.53
1:C:14:ALA:O	1:C:18:ASP:N	2.39	0.53
1:C:128:ARG:HH11	1:C:372:VAL:HG23	1.74	0.53
1:C:156:ARG:HA	1:D:155:ILE:HD11	1.90	0.53
1:D:73:SER:HA	1:D:96:VAL:HB	1.91	0.53
1:B:99:ASN:O	1:B:297:GLN:NE2	2.41	0.53
1:D:158:LEU:O	1:D:162:LEU:N	2.38	0.53
1:D:524:LEU:HD11	1:D:738:VAL:HB	1.90	0.53
1:B:164:GLN:O	1:B:168:LEU:N	2.36	0.53
1:C:11:GLU:HG2	1:C:49:ALA:HA	1.91	0.53
1:C:548:CYS:O	1:C:552:ALA:N	2.41	0.53
1:B:665:SER:H	1:B:668:TYR:HB3	1.73	0.53
1:B:733:ASP:O	1:B:790:ARG:NH1	2.41	0.53
1:D:488:LEU:O	1:D:755:GLN:NE2	2.42	0.53
1:D:671:PHE:HA	1:D:674:LYS:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:816:CYS:SG	1:C:817:ILE:N	2.82	0.53
1:D:543:PHE:HA	1:D:546:TRP:HB2	1.89	0.53
1:B:523:ILE:N	1:B:659:TYR:OH	2.39	0.53
1:A:164:GLN:OE1	1:A:167:ARG:NH1	2.40	0.52
1:C:514:LYS:HZ2	1:D:514:LYS:HD3	1.75	0.52
1:B:76:CYS:HA	1:B:95:PHE:HZ	1.74	0.52
1:A:12:GLU:HG2	1:A:49:ALA:HB2	1.91	0.52
1:B:376:ILE:HB	1:B:393:ALA:HB3	1.92	0.52
1:B:739:THR:HG23	1:B:793:LEU:HG	1.90	0.52
1:C:52:PRO:HB2	1:D:81:ALA:HB2	1.92	0.52
1:C:125:PRO:HG2	1:C:263:GLN:HG3	1.92	0.52
1:B:9:ILE:HA	1:B:45:LYS:HB2	1.91	0.52
1:B:363:PHE:HB3	1:B:367:SER:HA	1.92	0.52
1:C:778:ASP:OD2	1:C:782:GLN:NE2	2.42	0.52
1:A:106:THR:HA	1:A:157:GLY:HA3	1.92	0.52
1:A:209:LEU:HD11	1:A:235:TRP:HB3	1.90	0.52
1:C:163:ASP:OD2	1:D:149:TYR:OH	2.27	0.52
1:D:90:HIS:HB2	1:D:339:TRP:HB3	1.91	0.52
1:C:104:PRO:HG3	1:C:239:ASN:HA	1.91	0.52
1:A:623:ILE:O	1:A:626:SER:OG	2.23	0.52
1:C:92:PRO:HA	1:C:350:LYS:HE3	1.91	0.52
1:C:663:ARG:HB3	1:C:701:CYS:HB3	1.91	0.52
1:B:512:PHE:HA	1:B:753:ALA:HA	1.92	0.51
1:C:237:PHE:HD2	1:C:260:VAL:HG13	1.75	0.51
1:C:521:VAL:HG13	1:C:720:LEU:HD11	1.91	0.51
1:A:526:LYS:HA	1:A:738:VAL:HA	1.92	0.51
1:D:48:GLU:HB2	1:D:51:ASN:HD22	1.76	0.51
1:B:694:LYS:NZ	1:B:699:ASP:O	2.41	0.51
1:A:623:ILE:O	1:A:627:SER:OG	2.25	0.51
1:A:73:SER:HA	1:A:96:VAL:HB	1.92	0.51
1:B:231:LYS:HB3	1:B:383:GLU:HG3	1.93	0.51
1:D:251:VAL:HG11	1:D:379:THR:HG21	1.93	0.51
1:A:104:PRO:HG2	1:A:158:LEU:HD11	1.92	0.51
1:A:649:GLN:H	1:A:652:SER:HB2	1.75	0.51
1:B:20:VAL:HG11	1:B:300:ASN:HB3	1.91	0.51
1:B:206:ARG:HB2	1:B:206:ARG:CZ	2.39	0.51
1:B:434:GLU:HB3	1:B:442:ARG:CG	2.40	0.51
1:C:6:ILE:HG22	1:C:69:ALA:H	1.76	0.51
1:D:196:GLU:HA	1:D:199:ARG:HE	1.76	0.51
1:B:125:PRO:HG3	1:B:372:VAL:HB	1.93	0.50
1:B:501:THR:O	1:B:506:ARG:NH2	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:618:TRP:H	1:D:618:TRP:CD1	2.28	0.50
1:B:654:GLN:HE22	1:B:688:LEU:HD13	1.76	0.50
1:B:418:THR:HG22	1:B:461:LYS:HB2	1.94	0.50
1:D:261:VAL:HG13	1:D:376:ILE:HG13	1.93	0.50
1:A:50:ASN:O	1:B:113:SER:OG	2.28	0.50
1:A:681:GLN:NE2	1:C:690:ARG:O	2.41	0.50
1:B:252:HIS:HA	1:B:387:LYS:HE3	1.94	0.50
1:C:244:ASP:OD1	1:C:262:ARG:NH2	2.44	0.50
1:A:355:THR:HG22	1:A:360:VAL:HG22	1.92	0.50
1:A:810:SER:OG	1:A:811:PHE:N	2.45	0.50
1:C:139:GLU:O	1:C:141:ARG:NH1	2.44	0.50
1:A:148:PHE:N	1:A:208:ILE:O	2.44	0.50
1:A:174:LEU:HD21	1:B:162:LEU:HB3	1.94	0.50
1:D:476:LEU:HD13	1:D:478:ASN:HD22	1.76	0.50
1:A:12:GLU:N	1:A:12:GLU:OE2	2.44	0.49
1:A:710:ARG:O	1:A:714:LYS:N	2.45	0.49
1:B:164:GLN:HE21	1:B:168:LEU:HG	1.75	0.49
1:C:205:ARG:HD3	1:C:206:ARG:HG3	1.94	0.49
1:A:147:MET:N	1:A:173:SER:O	2.41	0.49
1:B:12:GLU:O	1:B:13:ASN:C	2.55	0.49
1:D:650:ASP:N	1:D:650:ASP:OD1	2.44	0.49
1:B:94:LEU:HA	1:B:121:LEU:HB2	1.94	0.49
1:C:293:LEU:C	1:C:293:LEU:HD13	2.38	0.49
1:C:657:MET:SD	1:C:657:MET:N	2.75	0.49
1:A:144:LYS:HA	1:A:171:ASP:HB3	1.93	0.49
1:A:534:ILE:HG23	1:A:812:ALA:HB3	1.94	0.49
1:C:159:GLN:HE21	1:D:155:ILE:HD13	1.77	0.49
1:C:631:ASN:ND2	1:D:629:THR:OG1	2.45	0.49
1:D:449:ASP:HB3	1:D:780:LEU:HD22	1.94	0.49
1:C:747:SER:O	1:D:772:GLN:NE2	2.45	0.49
1:D:22:GLN:HE21	1:D:44:ILE:HD12	1.78	0.49
1:B:165:ALA:O	1:B:170:LEU:N	2.46	0.49
1:B:421:VAL:O	1:B:465:TYR:N	2.46	0.49
1:C:156:ARG:HG2	1:D:155:ILE:HG13	1.95	0.49
1:C:43:SER:HB3	1:C:45:LYS:HE3	1.95	0.48
1:C:510:VAL:HA	1:C:756:HIS:HB2	1.95	0.48
1:C:12:GLU:HA	1:C:49:ALA:HB2	1.95	0.48
1:D:457:ALA:O	1:D:767:ARG:NH1	2.45	0.48
1:A:89:MET:HA	1:A:338:PRO:HA	1.94	0.48
1:B:534:ILE:O	1:B:537:LEU:HB3	2.13	0.48
1:C:11:GLU:HG3	1:C:47:ILE:HG13	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:149:TYR:HD2	1:C:153:TYR:HB3	1.78	0.48
1:C:560:ILE:HG13	1:C:611:ARG:HH11	1.77	0.48
1:A:519:TYR:O	1:A:747:SER:OG	2.32	0.48
1:B:681:GLN:HE22	1:B:686:ALA:H	1.62	0.48
1:C:11:GLU:HG3	1:C:47:ILE:CD1	2.44	0.48
1:C:63:MET:O	1:C:328:SER:OG	2.32	0.48
1:A:47:ILE:HD13	1:A:55:ALA:HB2	1.94	0.48
1:C:144:LYS:HG2	1:C:171:ASP:HB2	1.95	0.48
1:C:163:ASP:O	1:C:166:SER:OG	2.24	0.48
1:D:689:TRP:O	1:D:693:SER:N	2.47	0.48
1:A:72:THR:OG1	1:A:93:HIS:NE2	2.41	0.48
1:A:316:ARG:HH21	1:A:348:THR:HG22	1.78	0.48
1:B:116:GLY:O	1:B:119:TYR:OH	2.31	0.48
1:D:396:ASP:OD1	1:D:396:ASP:N	2.45	0.48
1:C:31:ASN:O	1:C:279:HIS:ND1	2.43	0.48
1:A:442:ARG:HG2	1:A:444:LYS:NZ	2.29	0.48
1:C:506:ARG:HA	1:C:509:VAL:HG22	1.96	0.48
1:D:164:GLN:O	1:D:168:LEU:HB2	2.14	0.48
1:B:74:THR:HG21	1:B:79:ALA:HB2	1.96	0.47
1:A:150:ASP:N	1:A:150:ASP:OD1	2.46	0.47
1:C:710:ARG:HH21	1:C:714:LYS:HE2	1.78	0.47
1:A:12:GLU:CG	1:A:49:ALA:HB2	2.44	0.47
1:B:623:ILE:O	1:B:627:SER:N	2.46	0.47
1:C:207:ALA:HB3	1:C:235:TRP:HA	1.97	0.47
1:C:769:LEU:HD11	1:D:748:LYS:HG2	1.95	0.47
1:A:135:ARG:HE	1:A:401:LEU:HD23	1.78	0.47
1:B:8:ALA:HB3	1:B:44:ILE:HA	1.96	0.47
1:D:283:SER:H	1:D:286:CYS:HB2	1.78	0.47
1:D:355:THR:HG22	1:D:360:VAL:HG22	1.96	0.47
1:A:244:ASP:HA	1:A:247:ILE:HD12	1.97	0.47
1:A:398:GLU:HG3	1:A:399:LYS:HG2	1.96	0.47
1:A:488:LEU:HA	1:A:755:GLN:HE21	1.80	0.47
1:C:251:VAL:HG12	1:C:258:MET:HE2	1.95	0.47
1:C:525:ILE:HG13	1:C:739:THR:HG23	1.97	0.47
1:A:86:THR:HG21	1:A:93:HIS:HB2	1.96	0.47
1:A:771:LEU:O	1:A:776:ASP:N	2.44	0.47
1:B:122:ALA:HB3	1:B:368:SER:HA	1.95	0.47
1:C:86:THR:HG23	1:C:91:ILE:H	1.79	0.47
1:C:212:SER:O	1:C:216:ALA:N	2.46	0.47
1:C:508:SER:OG	1:D:766:GLN:NE2	2.47	0.47
1:C:521:VAL:HA	1:C:720:LEU:HD21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:474:HIS:N	1:D:482:ASN:OD1	2.36	0.47
1:D:760:TYR:O	1:D:764:PHE:N	2.45	0.47
1:C:270:ASP:HA	1:C:273:LYS:HD3	1.95	0.47
1:D:114:PRO:HB2	1:D:117:GLU:HB2	1.96	0.47
1:D:502:ILE:HD11	1:D:751:GLY:H	1.79	0.47
1:B:511:ASP:OD1	1:B:761:ARG:NH1	2.47	0.47
1:C:200:TYR:O	1:C:204:LEU:HB2	2.14	0.46
1:C:631:ASN:HD21	1:D:626:SER:HA	1.81	0.46
1:D:12:GLU:HG2	1:D:99:ASN:HD21	1.79	0.46
1:A:207:ALA:O	1:A:236:VAL:N	2.49	0.46
1:B:141:ARG:NH2	1:B:143:GLN:OE1	2.48	0.46
1:B:526:LYS:H	1:B:713:LYS:HG3	1.79	0.46
1:C:128:ARG:HG3	1:C:372:VAL:CG2	2.45	0.46
1:C:694:LYS:HG3	1:C:699:ASP:HB3	1.97	0.46
1:D:269:LYS:HE2	1:D:288:PRO:HB3	1.97	0.46
1:B:63:MET:HE1	1:B:85:LEU:HD11	1.97	0.46
1:B:681:GLN:HG3	1:B:683:SER:H	1.79	0.46
1:B:421:VAL:HG13	1:B:496:ALA:HA	1.98	0.46
1:C:241:GLU:HA	1:C:296:LEU:HD23	1.96	0.46
1:A:614:MET:HG3	1:D:824:LEU:HG	1.97	0.46
1:B:9:ILE:HD13	1:B:55:ALA:HB1	1.98	0.46
1:B:192:MET:HE1	1:B:197:LEU:HD13	1.98	0.46
1:B:188:LEU:HB3	1:B:192:MET:HE3	1.98	0.46
1:B:422:VAL:HG13	1:B:484:MET:HG3	1.97	0.46
1:B:811:PHE:HB3	1:C:541:PHE:HZ	1.79	0.46
1:B:623:ILE:N	1:B:623:ILE:CD1	2.78	0.46
1:D:434:GLU:HB2	1:D:442:ARG:HB3	1.97	0.46
1:A:665:SER:H	1:A:668:TYR:HB3	1.80	0.46
1:C:128:ARG:HG3	1:C:372:VAL:HG21	1.97	0.46
1:B:619:LEU:HA	1:B:622:LEU:HB3	1.96	0.45
1:A:332:ILE:HG21	1:B:60:CYS:HB3	1.98	0.45
1:B:12:GLU:O	1:B:14:ALA:N	2.49	0.45
1:B:207:ALA:HB3	1:B:235:TRP:HD1	1.82	0.45
1:B:293:LEU:HD12	1:B:296:LEU:HD12	1.98	0.45
1:B:543:PHE:H	1:B:546:TRP:HD1	1.65	0.45
1:D:661:THR:HB	1:D:720:LEU:HB3	1.97	0.45
1:D:187:SER:O	1:D:191:THR:OG1	2.29	0.45
1:D:202:ASP:OD1	1:D:205:ARG:NH2	2.48	0.45
1:D:542:ASP:OD1	1:D:542:ASP:N	2.50	0.45
1:D:731:LEU:HB2	1:D:785:TRP:HD1	1.81	0.45
1:B:244:ASP:OD1	1:B:262:ARG:NH1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:649:GLN:HB2	1:B:652:SER:H	1.80	0.45
1:A:164:GLN:O	1:A:168:LEU:N	2.41	0.45
1:A:504:PRO:O	1:B:766:GLN:NE2	2.48	0.45
1:A:733:ASP:OD2	1:A:736:CYS:N	2.50	0.45
1:B:96:VAL:HG11	1:B:299:SER:HA	1.98	0.45
1:B:670:TYR:O	1:B:674:LYS:N	2.42	0.45
1:C:533:SER:O	1:C:536:SER:OG	2.32	0.45
1:D:62:LEU:O	1:D:66:GLY:N	2.45	0.45
1:D:653:LYS:HD2	1:D:656:GLU:HB2	1.98	0.45
1:C:394:THR:O	1:C:402:ASN:N	2.50	0.45
1:A:268:ALA:HA	1:A:285:LEU:HD22	1.97	0.45
1:A:417:LEU:HB2	1:A:460:PHE:HB3	1.99	0.45
1:A:423:THR:O	1:A:467:ALA:N	2.49	0.45
1:A:442:ARG:HG2	1:A:444:LYS:HZ2	1.82	0.45
1:A:521:VAL:H	1:A:745:ILE:HD11	1.81	0.45
1:B:556:VAL:HG13	1:B:559:LEU:HD23	1.99	0.45
1:C:27:ASP:HA	1:C:30:LEU:HB3	1.98	0.45
1:B:60:CYS:SG	1:B:330:ASN:ND2	2.90	0.45
1:C:11:GLU:HG3	1:C:47:ILE:CG1	2.46	0.45
1:C:166:SER:OG	1:D:174:LEU:O	2.34	0.45
1:C:486:GLY:HA2	1:C:489:ILE:HB	1.97	0.45
1:D:274:CYS:HB3	1:D:281:ILE:HB	1.99	0.45
1:A:653:LYS:HD2	1:A:656:GLU:CG	2.45	0.45
1:C:354:ILE:HD11	1:C:361:MET:HE3	1.99	0.45
1:A:396:ASP:OD1	1:A:396:ASP:N	2.50	0.44
1:B:251:VAL:HG12	1:B:258:MET:HG2	1.99	0.44
1:C:83:GLN:HE22	1:C:110:LEU:HB3	1.82	0.44
1:A:687:GLU:OE1	1:A:690:ARG:NH1	2.50	0.44
1:C:520:SER:OG	1:C:743:ASN:O	2.28	0.44
1:D:93:HIS:O	1:D:121:LEU:N	2.48	0.44
1:A:710:ARG:HD2	1:A:713:LYS:HD3	1.99	0.44
1:B:204:LEU:C	1:B:204:LEU:HD23	2.43	0.44
1:C:128:ARG:HB3	1:C:395:TRP:CH2	2.49	0.44
1:C:329:LEU:HB2	1:C:336:THR:HB	1.97	0.44
1:A:147:MET:HA	1:A:208:ILE:HB	1.98	0.44
1:A:621:THR:OG1	1:A:622:LEU:N	2.50	0.44
1:A:831:LEU:HA	1:A:834:TRP:HB3	1.99	0.44
1:A:18:ASP:HA	1:A:21:PHE:HB3	2.00	0.44
1:A:318:LEU:O	1:A:321:ARG:NH2	2.38	0.44
1:B:443:TYR:HE2	1:B:464:ILE:HG21	1.83	0.44
1:A:332:ILE:HG22	1:B:57:GLN:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:22:GLN:O	1:C:26:SER:HB3	2.17	0.44
1:B:189:PHE:HA	1:B:192:MET:HG2	1.99	0.44
1:D:792:ASP:OD1	1:D:792:ASP:N	2.51	0.44
1:C:312:ASN:ND2	1:C:353:HIS:O	2.51	0.44
1:C:453:ALA:HA	1:C:456:LYS:HE3	2.00	0.44
1:C:681:GLN:HE22	1:C:686:ALA:H	1.65	0.44
1:D:20:VAL:HG21	1:D:300:ASN:HB3	2.00	0.44
1:A:64:THR:HA	1:A:328:SER:HB3	1.98	0.44
1:C:188:LEU:O	1:C:192:MET:N	2.51	0.44
1:D:361:MET:HA	1:D:369:ASN:HD21	1.82	0.44
1:D:682:ASP:HB3	1:D:685:PHE:HB2	2.00	0.44
1:B:810:SER:OG	1:B:811:PHE:N	2.47	0.43
1:C:402:ASN:ND2	1:C:402:ASN:H	2.16	0.43
1:C:426:GLU:H	1:C:430:VAL:HB	1.83	0.43
1:C:761:ARG:NH1	1:C:765:SER:OG	2.51	0.43
1:D:204:LEU:HD23	1:D:228:LEU:HD11	2.00	0.43
1:D:239:ASN:HB3	1:D:242:ILE:HB	2.00	0.43
1:B:165:ALA:HB1	1:B:170:LEU:HB2	2.00	0.43
1:B:661:THR:OG1	1:B:662:VAL:N	2.43	0.43
1:C:377:LEU:HG	1:C:391:LYS:HA	1.99	0.43
1:C:419:LEU:O	1:C:463:GLU:N	2.41	0.43
1:C:424:VAL:HG22	1:C:470:GLY:HA2	2.00	0.43
1:D:510:VAL:HG11	1:D:753:ALA:HB1	2.00	0.43
1:A:435:ASN:HA	1:A:441:LYS:H	1.83	0.43
1:C:137:VAL:HA	1:C:142:TRP:H	1.82	0.43
1:C:237:PHE:HB3	1:C:260:VAL:HG22	1.99	0.43
1:C:514:LYS:HA	1:D:514:LYS:HZ2	1.83	0.43
1:A:32:ASP:HB2	1:A:35:LEU:HB2	2.00	0.43
1:A:143:GLN:H	1:A:206:ARG:HH12	1.67	0.43
1:A:284:LEU:HD22	1:A:293:LEU:HD11	2.01	0.43
1:D:518:ASP:OD2	1:D:747:SER:N	2.51	0.43
1:A:524:LEU:HD11	1:A:738:VAL:HB	1.99	0.43
1:B:432:VAL:HG22	1:B:443:TYR:HE1	1.83	0.43
1:D:97:GLN:HB3	1:D:124:ARG:HD3	2.01	0.43
1:B:525:ILE:HG12	1:B:709:ILE:HA	1.99	0.43
1:C:402:ASN:H	1:C:402:ASN:HD22	1.66	0.43
1:A:201:ARG:HH11	1:A:228:LEU:HD23	1.84	0.43
1:D:708:GLY:HA2	1:D:711:LYS:HB2	2.00	0.43
1:A:2:SER:N	1:A:38:GLU:OE1	2.52	0.43
1:A:553:ILE:H	1:A:553:ILE:HG13	1.66	0.43
1:B:62:LEU:HD12	1:B:62:LEU:C	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:394:THR:HG23	1:B:402:ASN:HB2	2.00	0.43
1:B:721:TRP:CD1	1:B:725:VAL:HB	2.54	0.43
1:C:526:LYS:HA	1:C:738:VAL:HG12	2.00	0.43
1:D:106:THR:HG1	1:D:153:TYR:HH	1.58	0.43
1:C:444:LYS:HA	1:C:448:ILE:HG13	2.00	0.43
1:D:476:LEU:N	1:D:480:SER:O	2.44	0.43
1:D:524:LEU:HD22	1:D:726:VAL:HA	2.01	0.43
1:B:28:LEU:HB3	1:B:35:LEU:HD13	2.01	0.42
1:B:331:CYS:HA	1:B:336:THR:HG21	1.99	0.42
1:D:12:GLU:HB3	1:D:13:ASN:H	1.70	0.42
1:A:692:ILE:O	1:A:700:ASN:ND2	2.52	0.42
1:B:792:ASP:OD1	1:B:792:ASP:N	2.52	0.42
1:C:540:PRO:HD3	1:C:633:ALA:HB2	2.01	0.42
1:C:708:GLY:HA2	1:C:711:LYS:HB2	2.02	0.42
1:A:374:PHE:HB3	1:A:376:ILE:HD11	2.01	0.42
1:B:486:GLY:HA2	1:B:489:ILE:HD12	2.01	0.42
1:C:128:ARG:HG3	1:C:372:VAL:HB	2.01	0.42
1:C:205:ARG:HH11	1:C:206:ARG:HG3	1.84	0.42
1:C:650:ASP:OD1	1:C:650:ASP:N	2.52	0.42
1:D:147:MET:HA	1:D:208:ILE:HB	2.02	0.42
1:D:476:LEU:HB3	1:D:478:ASN:H	1.84	0.42
1:D:18:ASP:HA	1:D:21:PHE:HB3	2.01	0.42
1:D:419:LEU:O	1:D:463:GLU:N	2.52	0.42
1:D:435:ASN:HA	1:D:440:PRO:HA	2.01	0.42
1:A:205:ARG:O	1:A:234:HIS:N	2.52	0.42
1:A:423:THR:N	1:A:465:TYR:O	2.49	0.42
1:A:518:ASP:OD2	1:B:515:ARG:NH2	2.51	0.42
1:B:816:CYS:SG	1:B:817:ILE:HG13	2.59	0.42
1:C:96:VAL:HG13	1:C:124:ARG:HD2	2.01	0.42
1:C:498:SER:OG	1:C:499:ALA:N	2.52	0.42
1:D:91:ILE:HD13	1:D:91:ILE:HA	1.93	0.42
1:B:86:THR:HB	1:B:91:ILE:HB	2.00	0.42
1:A:506:ARG:HA	1:A:509:VAL:HG22	2.00	0.42
1:B:103:SER:OG	1:B:152:GLU:OE1	2.30	0.42
1:C:515:ARG:HD3	1:C:518:ASP:HB3	2.02	0.42
1:D:11:GLU:OE1	1:D:74:THR:HA	2.19	0.42
1:A:201:ARG:HA	1:A:204:LEU:HB2	2.01	0.42
1:A:491:LYS:HA	1:A:755:GLN:HE22	1.85	0.42
1:B:434:GLU:HB3	1:B:442:ARG:HG2	2.02	0.42
1:D:721:TRP:CD1	1:D:722:ASP:H	2.36	0.42
1:A:180:ASN:HD22	1:A:181:ILE:H	1.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:VAL:HG23	1:D:138:THR:HG23	2.02	0.42
1:B:527:LYS:HG3	1:B:737:SER:HB2	2.02	0.41
1:A:634:ALA:HB1	1:B:632:LEU:HD13	2.02	0.41
1:C:404:SER:OG	1:C:405:LEU:N	2.53	0.41
1:A:4:ILE:O	1:A:41:THR:N	2.39	0.41
1:B:532:ILE:HG13	1:C:541:PHE:HB2	2.02	0.41
1:D:228:LEU:HG	1:D:235:TRP:HH2	1.86	0.41
1:A:143:GLN:HA	1:A:170:LEU:HD22	2.03	0.41
1:A:242:ILE:O	1:A:262:ARG:NE	2.48	0.41
1:A:835:TRP:HA	1:A:838:ASN:HB2	2.03	0.41
1:C:128:ARG:NH1	1:C:372:VAL:HG23	2.34	0.41
1:C:781:LYS:O	1:C:785:TRP:N	2.39	0.41
1:D:96:VAL:HG22	1:D:123:SER:HB3	2.01	0.41
1:A:222:GLU:HA	1:A:225:GLU:HB2	2.02	0.41
1:A:669:GLU:HB3	1:A:672:ARG:HH11	1.86	0.41
1:C:453:ALA:HA	1:C:456:LYS:HG2	2.02	0.41
1:D:379:THR:HA	1:D:388:ASP:O	2.21	0.41
1:C:125:PRO:HA	1:C:126:PRO:HD3	1.88	0.41
1:C:146:VAL:HB	1:C:173:SER:HB2	2.02	0.41
1:D:548:CYS:HA	1:D:551:ALA:HB3	2.02	0.41
1:B:8:ALA:O	1:B:45:LYS:N	2.52	0.41
1:B:469:ASP:HB2	1:B:483:GLY:HA3	2.03	0.41
1:C:327:ALA:HB3	1:C:339:TRP:CD1	2.56	0.41
1:A:631:ASN:HD21	1:D:633:ALA:HA	1.86	0.41
1:C:343:ARG:HA	1:C:346:LEU:HD12	2.03	0.41
1:C:421:VAL:N	1:C:463:GLU:O	2.40	0.41
1:D:815:PHE:HA	1:D:818:LEU:HB3	2.02	0.41
1:A:165:ALA:HB1	1:A:170:LEU:HB2	2.03	0.41
1:A:184:VAL:O	1:A:187:SER:OG	2.32	0.41
1:A:610:MET:SD	1:A:610:MET:N	2.89	0.41
1:B:64:THR:HA	1:B:329:LEU:H	1.85	0.41
1:C:50:ASN:HD22	1:D:112:PRO:HD2	1.85	0.41
1:C:519:TYR:N	1:C:748:LYS:O	2.53	0.41
1:A:164:GLN:HE21	1:A:168:LEU:HG	1.86	0.41
1:A:469:ASP:HB3	1:A:483:GLY:HA2	2.03	0.41
1:C:14:ALA:HA	1:C:17:ASP:HB2	2.02	0.41
1:D:706:SER:HA	1:D:709:ILE:HD12	2.01	0.41
1:D:782:GLN:HE22	1:D:787:HIS:HB3	1.86	0.41
1:A:57:GLN:HE22	1:B:332:ILE:HA	1.85	0.40
1:A:299:SER:HA	1:A:302:TYR:HD2	1.86	0.40
1:A:301:LEU:HA	1:A:304:TYR:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:514:LYS:HB2	1:B:768:ILE:HG21	2.03	0.40
1:C:104:PRO:HD2	1:C:210:LEU:HD23	2.01	0.40
1:C:421:VAL:O	1:C:465:TYR:N	2.35	0.40
1:C:521:VAL:HB	1:C:743:ASN:HB3	2.02	0.40
1:D:217:HIS:HA	1:D:220:ILE:HG22	2.03	0.40
1:D:471:ARG:O	1:D:474:HIS:NE2	2.54	0.40
1:D:778:ASP:O	1:D:782:GLN:HG2	2.21	0.40
1:B:11:GLU:CB	1:B:74:THR:HA	2.49	0.40
1:C:139:GLU:OE1	1:C:404:SER:OG	2.33	0.40
1:C:147:MET:HE2	1:C:174:LEU:HD22	2.02	0.40
1:A:475:GLN:HE22	1:A:479:THR:HA	1.87	0.40
1:A:492:ARG:HE	1:A:492:ARG:HB2	1.66	0.40
1:A:504:PRO:HB3	1:B:770:GLU:HG2	2.04	0.40
1:A:543:PHE:HA	1:A:546:TRP:HB2	2.04	0.40
1:B:520:SER:H	1:B:722:ASP:HB2	1.86	0.40
1:C:158:LEU:HA	1:C:158:LEU:HD23	1.89	0.40
1:C:708:GLY:HA2	1:C:711:LYS:HD3	2.02	0.40
1:A:181:ILE:HD12	1:A:219:PHE:HD1	1.85	0.40
1:A:495:LEU:HD22	1:A:495:LEU:HA	1.95	0.40
1:A:622:LEU:HD21	1:B:620:PHE:HE2	1.87	0.40
1:B:834:TRP:O	1:B:837:SER:OG	2.30	0.40
1:D:731:LEU:HB2	1:D:785:TRP:CD1	2.57	0.40
1:A:11:GLU:O	1:A:98:ARG:NH2	2.55	0.40
1:A:690:ARG:HA	1:A:693:SER:HB2	2.03	0.40
1:B:21:PHE:HB2	1:B:303:LEU:HD21	2.04	0.40
1:B:490:SER:OG	1:B:492:ARG:HG3	2.22	0.40
1:D:491:LYS:HG2	1:D:755:GLN:HE22	1.86	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 PDB entries followed by that with respect to all EM entries.

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	741/856 (87%)	682 (92%)	59 (8%)	0	100	100
1	B	741/856 (87%)	671 (91%)	70 (9%)	0	100	100
1	C	741/856 (87%)	682 (92%)	59 (8%)	0	100	100
1	D	741/856 (87%)	686 (93%)	55 (7%)	0	100	100
All	All	2964/3424 (87%)	2721 (92%)	243 (8%)	0	100	100

There are no Ramachandran outliers to report.

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 PDB entries followed by that with respect to all EM entries.

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	650/739 (88%)	641 (99%)	9 (1%)	59	72
1	B	650/739 (88%)	627 (96%)	23 (4%)	32	53
1	C	650/739 (88%)	639 (98%)	11 (2%)	53	69
1	D	650/739 (88%)	649 (100%)	1 (0%)	87	87
All	All	2600/2956 (88%)	2556 (98%)	44 (2%)	52	69

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	GLU
1	A	12	GLU
1	A	172	VAL
1	A	442	ARG
1	A	449	ASP
1	A	491	LYS
1	A	495	LEU
1	A	617	TRP
1	A	657	MET
1	B	12	GLU
1	B	13	ASN
1	B	62	LEU
1	B	70	LEU

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Mol	Chain	Res	Type
1	B	71	VAL
1	B	72	THR
1	B	154	ASP
1	B	204	LEU
1	B	205	ARG
1	B	206	ARG
1	B	334	LYS
1	B	441	LYS
1	B	442	ARG
1	B	443	TYR
1	B	449	ASP
1	B	560	ILE
1	B	621	THR
1	B	623	ILE
1	B	626	SER
1	B	629	THR
1	B	655	LEU
1	B	738	VAL
1	B	787	HIS
1	C	12	GLU
1	C	127	VAL
1	C	293	LEU
1	C	294	GLN
1	C	295	MET
1	C	382	SER
1	C	402	ASN
1	C	656	GLU
1	C	657	MET
1	C	710	ARG
1	C	790	ARG
1	D	12	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (65) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	180	ASN
1	A	214	GLN
1	A	221	ASN
1	A	271	ASN
1	A	277	ASN
1	A	289	GLN

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Mol	Chain	Res	Type
1	A	294	GLN
1	A	312	ASN
1	A	466	GLN
1	A	474	HIS
1	A	475	GLN
1	A	477	HIS
1	A	478	ASN
1	A	631	ASN
1	A	654	GLN
1	A	695	ASN
1	A	700	ASN
1	A	743	ASN
1	A	755	GLN
1	B	13	ASN
1	B	51	ASN
1	B	164	GLN
1	B	175	GLN
1	B	221	ASN
1	B	227	ASN
1	B	263	GLN
1	B	277	ASN
1	B	278	ASN
1	B	312	ASN
1	B	324	HIS
1	B	330	ASN
1	B	340	ASN
1	B	369	ASN
1	B	654	GLN
1	B	681	GLN
1	B	700	ASN
1	B	716	ASN
1	B	772	GLN
1	C	31	ASN
1	C	50	ASN
1	C	83	GLN
1	C	111	ASN
1	C	143	GLN
1	C	159	GLN
1	C	324	HIS
1	C	330	ASN
1	C	631	ASN
1	C	649	GLN

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Mol	Chain	Res	Type
1	C	681	GLN
1	D	22	GLN
1	D	36	GLN
1	D	50	ASN
1	D	51	ASN
1	D	217	HIS
1	D	272	GLN
1	D	324	HIS
1	D	369	ASN
1	D	373	GLN
1	D	478	ASN
1	D	654	GLN
1	D	677	ASN
1	D	695	ASN
1	D	766	GLN
1	D	836	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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 ⓘ

There are no chain breaks in this entry.