



Full wwPDB EM Validation Report ⓘ

May 3, 2026 – 12:50 AM JST

PDB ID : 9UNR / pdb_00009unr
EMDB ID : EMD-64361
Title : native NMDA receptor-GluN1/N2A/N2B-S1 in the closed state
Authors : Yu, J.; Xu, R.S.; Ge, J.P.
Deposited on : 2025-04-24
Resolution : 2.98 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**
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

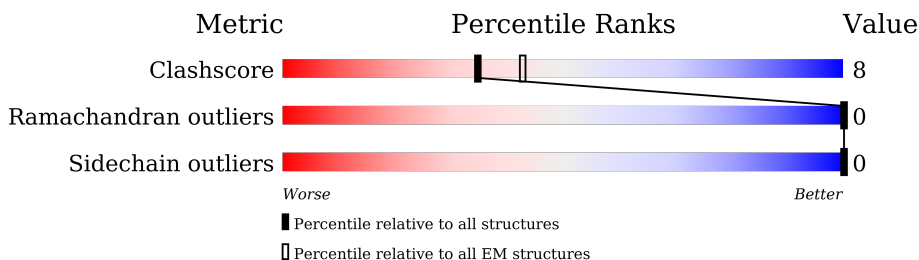
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.98 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	819	
1	C	819	
2	B	809	
3	D	808	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 23939 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, NMDA 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	794	Total	C	N	O	S	0	0
			5939	3804	1022	1078	35		
1	C	805	Total	C	N	O	S	0	0
			6085	3888	1054	1108	35		

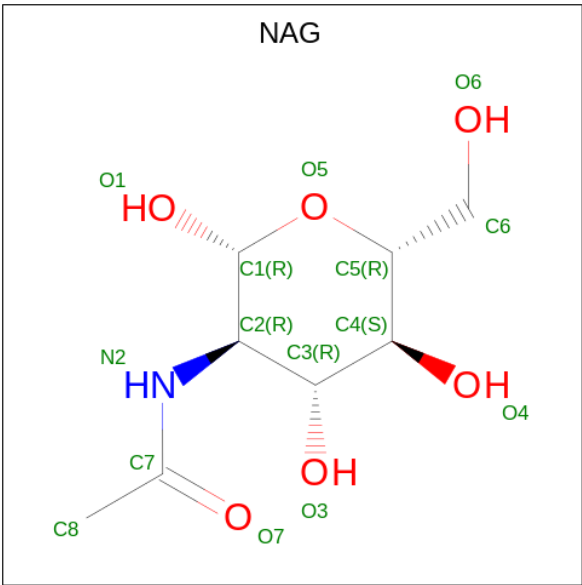
- Molecule 2 is a protein called Glutamate receptor ionotropic, NMDA 2A.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	792	Total	C	N	O	S	0	0
			5725	3697	949	1046	33		

- Molecule 3 is a protein called Glutamate receptor ionotropic, NMDA 2B.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	791	Total	C	N	O	S	0	0
			5882	3791	960	1090	41		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



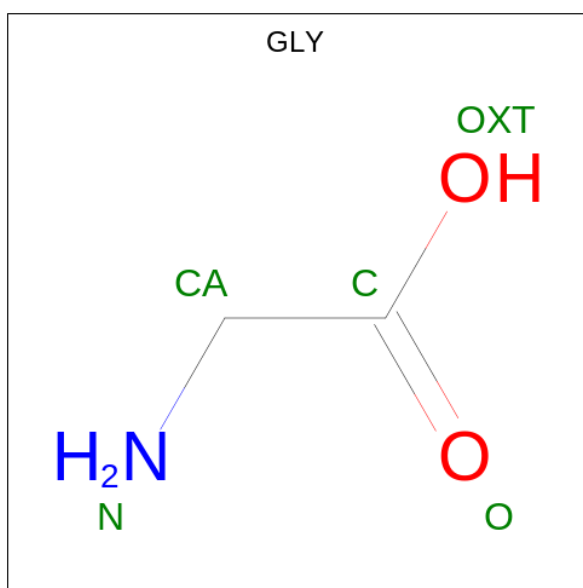
Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	

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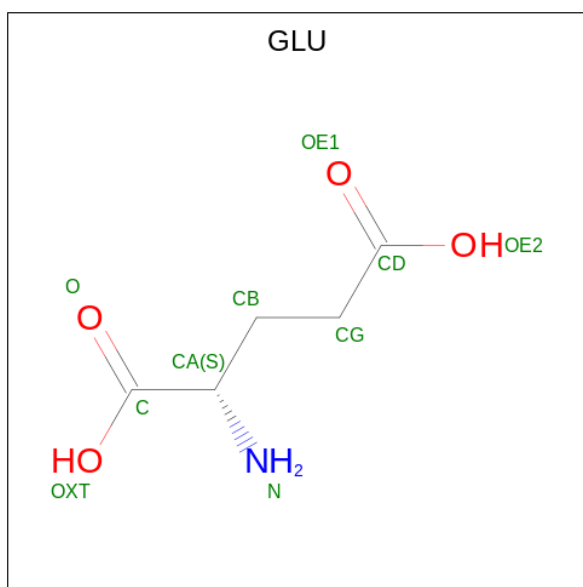
Mol	Chain	Residues	Atoms				AltConf
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	D	1	Total	C	N	O	0
			14	8	1	5	
4	D	1	Total	C	N	O	0
			14	8	1	5	
4	D	1	Total	C	N	O	0
			14	8	1	5	

- Molecule 5 is GLYCINE (CCD ID: GLY) (formula: $C_2H_5NO_2$).



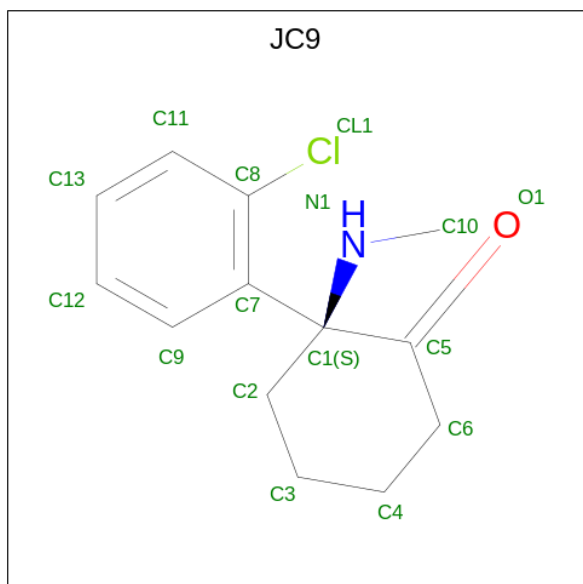
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			4	2	1	1	
5	C	1	Total	C	N	O	0
			4	2	1	1	

- Molecule 6 is GLUTAMIC ACID (CCD ID: GLU) (formula: $C_5H_9NO_4$).



Mol	Chain	Residues	Atoms				AltConf
6	B	1	Total	C	N	O	0
			9	5	1	3	
6	D	1	Total	C	N	O	0
			9	5	1	3	

- Molecule 7 is (2 {S})-2-(2-chlorophenyl)-2-(methylamino)cyclohexan-1-one (CCD ID: JC9) (formula: C₁₃H₁₆ClNO) (labeled as "Ligand of Interest" by depositor).

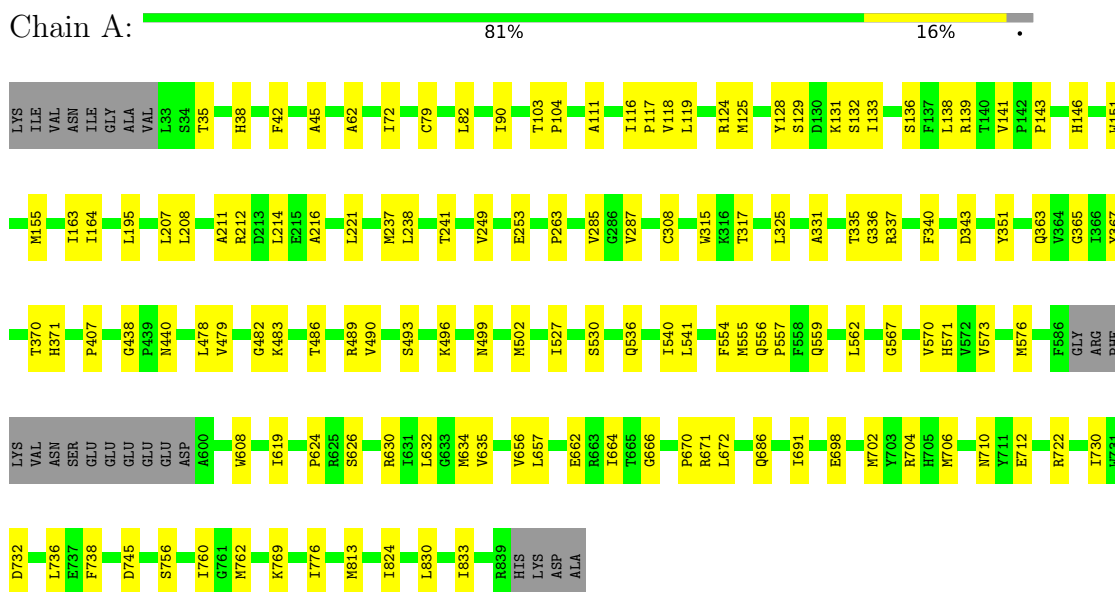


Mol	Chain	Residues	Atoms					AltConf
7	D	1	Total	C	Cl	N	O	0
			16	13	1	1	1	

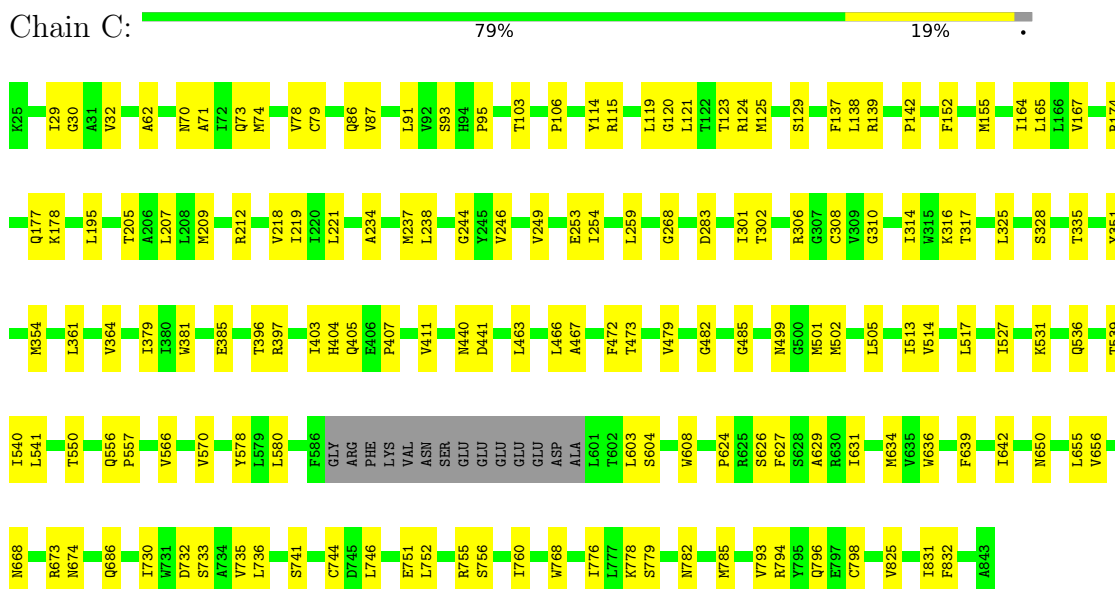
3 Residue-property plots

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, NMDA 1

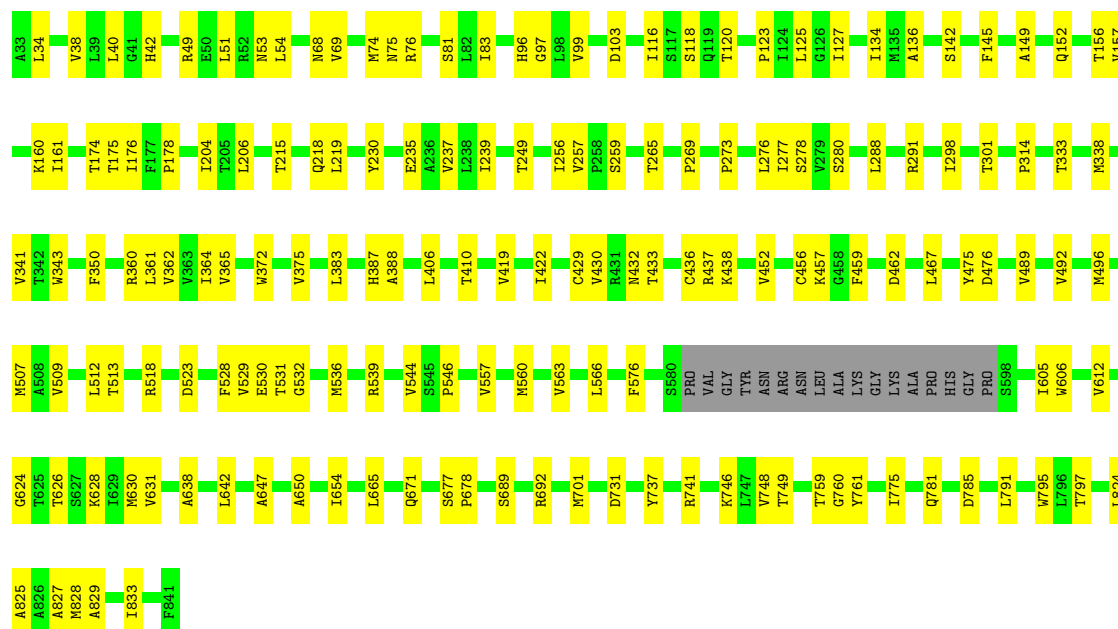


- Molecule 1: Glutamate receptor ionotropic, NMDA 1



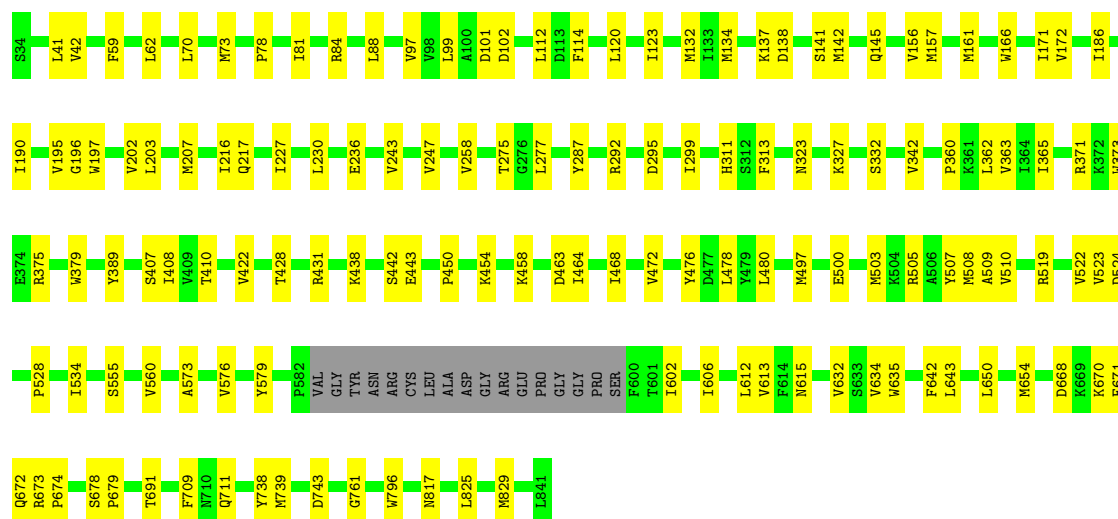
- Molecule 2: Glutamate receptor ionotropic, NMDA 2A

Chain B:  78% 20%



- Molecule 3: Glutamate receptor ionotropic, NMDA 2B

Chain D:  81% 17%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	343745	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	50	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor

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: JC9, NAG

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	0.09	0/6074	0.23	0/8272
1	C	0.09	0/6224	0.25	0/8469
2	B	0.09	0/5851	0.25	0/8009
3	D	0.09	0/6006	0.26	0/8176
All	All	0.09	0/24155	0.24	0/32926

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5939	0	5708	92	0
1	C	6085	0	5876	109	0
2	B	5725	0	5330	109	0
3	D	5882	0	5604	88	0
4	A	98	0	91	0	0
4	B	28	0	26	0	0
4	C	98	0	91	1	0
4	D	42	0	39	0	0
5	A	4	0	2	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	4	0	2	1	0
6	B	9	0	5	3	0
6	D	9	0	5	0	0
7	D	16	0	0	4	0
All	All	23939	0	22779	372	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 8.

All (372) 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)
7:D:905:JC9:C5	7:D:905:JC9:CL1	2.64	0.82
3:D:528:PRO:HB3	3:D:761:GLY:HA3	1.71	0.72
1:C:668:ASN:HA	1:C:673:ARG:HE	1.55	0.70
1:A:556:GLN:HG3	1:A:557:PRO:HD3	1.72	0.70
2:B:507:MET:HE2	2:B:775:ILE:HD11	1.75	0.69
2:B:160:LYS:HG3	2:B:383:LEU:HD22	1.75	0.68
1:C:86:GLN:HE21	1:C:306:ARG:HB2	1.57	0.68
2:B:626:THR:HG21	1:C:832:PHE:HB2	1.76	0.67
2:B:161:ILE:HD13	2:B:383:LEU:HD21	1.75	0.66
1:A:567:GLY:O	1:A:571:HIS:ND1	2.24	0.66
3:D:195:VAL:HG22	3:D:428:THR:HG21	1.77	0.65
2:B:419:VAL:HG12	2:B:456:CYS:HB3	1.77	0.65
1:C:124:ARG:NH1	1:C:142:PRO:O	2.30	0.65
2:B:175:THR:HG22	2:B:176:ILE:H	1.61	0.65
2:B:647:ALA:HB1	1:C:655:LEU:HD21	1.79	0.65
1:C:513:ILE:HG21	1:C:517:LEU:HD22	1.79	0.64
2:B:288:LEU:HA	2:B:291:ARG:HD3	1.78	0.64
2:B:219:LEU:HD11	2:B:239:ILE:HG23	1.79	0.64
1:C:93:SER:HB3	1:C:121:LEU:HD12	1.80	0.64
2:B:257:VAL:HB	2:B:276:LEU:HD11	1.80	0.64
2:B:116:ILE:O	2:B:120:THR:OG1	2.16	0.63
1:C:550:THR:O	1:C:556:GLN:NE2	2.29	0.63
3:D:534:ILE:HG22	3:D:691:THR:HG22	1.81	0.63
1:C:479:VAL:HG12	1:C:482:GLY:H	1.63	0.62
2:B:824:LEU:HG	2:B:828:MET:HE1	1.81	0.62
1:A:536:GLN:NE2	1:A:756:SER:O	2.32	0.62
2:B:422:ILE:HG13	2:B:429:CYS:HB3	1.79	0.62
1:C:174:ARG:O	1:C:178:LYS:HG3	2.00	0.61
3:D:408:ILE:HD13	3:D:476:TYR:HB2	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:483:LYS:NZ	1:A:712:GLU:O	2.34	0.61
2:B:430:VAL:HG12	2:B:432:ASN:H	1.66	0.61
2:B:576:PHE:HE1	2:B:624:GLY:H	1.49	0.61
2:B:364:ILE:HB	2:B:372:TRP:HB3	1.82	0.60
3:D:500:GLU:OE2	3:D:505:ARG:NH1	2.34	0.60
1:A:241:THR:HG23	1:A:263:PRO:HB3	1.82	0.60
3:D:138:ASP:HB3	3:D:141:SER:HB3	1.83	0.59
1:A:702:MET:O	1:A:706:MET:HG3	2.02	0.59
1:C:259:LEU:HD11	1:C:361:LEU:HD21	1.84	0.59
1:C:218:VAL:HG22	1:C:246:VAL:HB	1.84	0.59
3:D:166:TRP:HB3	3:D:227:ILE:HG13	1.82	0.59
2:B:406:LEU:H	2:B:475:TYR:HA	1.68	0.59
1:A:559:GLN:HB3	1:A:562:LEU:HB2	1.85	0.59
1:C:95:PRO:HG3	1:C:103:THR:HG21	1.85	0.59
2:B:665:LEU:HD22	2:B:701:MET:HE1	1.83	0.59
2:B:781:GLN:NE2	2:B:785:ASP:OD2	2.36	0.59
1:A:124:ARG:NH2	1:A:351:TYR:OH	2.35	0.59
1:A:131:LYS:HD3	1:A:343:ASP:HA	1.83	0.59
1:A:530:SER:HA	1:A:769:LYS:HE3	1.83	0.59
2:B:235:GLU:O	2:B:239:ILE:HD12	2.02	0.58
3:D:207:MET:HE2	3:D:236:GLU:HG2	1.84	0.58
2:B:544:VAL:HG12	2:B:546:PRO:HD3	1.85	0.58
3:D:650:LEU:HB3	3:D:654:MET:HE1	1.85	0.58
1:A:824:ILE:HG21	3:D:634:VAL:HG11	1.85	0.58
1:A:125:MET:O	1:A:139:ARG:NH2	2.34	0.58
1:C:557:PRO:HB3	1:C:650:ASN:HB3	1.86	0.58
2:B:731:ASP:OD2	2:B:761:TYR:OH	2.21	0.57
1:C:164:ILE:HB	1:C:219:ILE:HG22	1.86	0.57
1:A:208:LEU:HB3	1:A:238:LEU:HD22	1.87	0.57
1:C:310:GLY:H	3:D:78:PRO:HG3	1.69	0.57
1:A:698:GLU:HA	2:B:432:ASN:HD21	1.69	0.57
1:A:732:ASP:OD2	5:A:908:GLY:N	2.38	0.57
1:C:536:GLN:NE2	1:C:756:SER:O	2.38	0.57
3:D:673:ARG:HG2	3:D:674:PRO:HD3	1.86	0.57
1:A:662:GLU:O	1:A:671:ARG:NH1	2.37	0.56
1:C:70:ASN:HD22	3:D:323:ASN:H	1.52	0.56
1:A:82:LEU:HG	1:A:90:ILE:HD12	1.87	0.56
1:A:363:GLN:NE2	1:A:365:GLY:O	2.38	0.56
1:C:234:ALA:HA	1:C:237:MET:HE2	1.86	0.56
2:B:277:ILE:HG22	2:B:365:VAL:HG22	1.88	0.56
1:C:354:MET:HE2	1:C:361:LEU:HB3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:THR:OG1	1:A:104:PRO:HD3	2.05	0.56
1:A:195:LEU:HB3	1:A:207:LEU:HD12	1.88	0.56
1:A:211:ALA:HA	1:A:214:LEU:HD12	1.88	0.56
1:A:287:VAL:HG11	1:A:340:PHE:HE2	1.71	0.56
3:D:277:LEU:HD23	3:D:373:TRP:HH2	1.71	0.56
2:B:338:MET:HB3	2:B:350:PHE:HD2	1.70	0.56
2:B:360:ARG:HB3	2:B:362:VAL:HG23	1.87	0.56
2:B:825:ALA:HA	2:B:828:MET:HE2	1.88	0.56
1:C:114:TYR:OH	1:C:308:CYS:O	2.22	0.55
1:C:674:ASN:ND2	3:D:743:ASP:O	2.38	0.55
3:D:389:TYR:HE2	3:D:450:PRO:HB2	1.71	0.55
3:D:643:LEU:HD12	7:D:905:JC9:C4	2.37	0.55
3:D:438:LYS:HB2	3:D:480:LEU:HD12	1.88	0.55
1:A:221:LEU:HB3	1:A:249:VAL:HG12	1.88	0.55
1:C:167:VAL:HG21	1:C:177:GLN:HB3	1.89	0.55
1:A:541:LEU:HB2	1:A:736:LEU:HD13	1.88	0.55
1:A:118:VAL:O	1:A:138:LEU:N	2.39	0.55
1:A:335:THR:HG22	1:A:336:GLY:H	1.72	0.55
3:D:410:THR:HG23	3:D:478:LEU:HD11	1.88	0.54
1:C:221:LEU:HB3	1:C:249:VAL:HG12	1.88	0.54
3:D:132:MET:O	3:D:145:GLN:NE2	2.40	0.54
3:D:84:ARG:O	3:D:88:LEU:HG	2.08	0.54
3:D:612:LEU:HD11	3:D:632:VAL:HG13	1.89	0.54
1:C:120:GLY:N	1:C:138:LEU:O	2.39	0.54
2:B:157:VAL:HG21	2:B:361:LEU:HD12	1.89	0.54
1:A:624:PRO:HD2	2:B:606:TRP:HH2	1.72	0.54
1:C:195:LEU:HB3	1:C:207:LEU:HD11	1.89	0.53
1:C:91:LEU:HD12	1:C:121:LEU:HD21	1.88	0.53
3:D:602:ILE:O	3:D:606:ILE:HG13	2.08	0.53
1:A:111:ALA:HB1	1:A:116:ILE:HB	1.89	0.53
1:A:119:LEU:HD21	1:A:325:LEU:HD21	1.90	0.53
2:B:51:LEU:HD23	2:B:54:LEU:HD21	1.90	0.53
1:C:463:LEU:HB2	1:C:514:VAL:HG11	1.91	0.53
1:C:70:ASN:ND2	3:D:323:ASN:H	2.07	0.53
3:D:59:PHE:HB3	3:D:62:LEU:HD12	1.91	0.53
2:B:237:VAL:HG21	2:B:269:PRO:HG2	1.91	0.52
1:C:74:MET:HE3	1:C:78:VAL:HG23	1.91	0.52
1:C:624:PRO:HB2	1:C:629:ALA:HB1	1.90	0.52
1:A:576:MET:HG3	1:A:632:LEU:HD21	1.92	0.52
2:B:410:THR:HG22	2:B:509:VAL:O	2.09	0.52
2:B:467:LEU:HD21	2:B:775:ILE:HG12	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:540:ILE:HG12	1:C:730:ILE:HG13	1.90	0.52
1:A:128:TYR:HB2	1:A:139:ARG:NH2	2.25	0.52
3:D:99:LEU:HD22	3:D:112:LEU:HD11	1.91	0.52
3:D:216:ILE:HG22	3:D:243:VAL:HG11	1.92	0.51
3:D:99:LEU:HD13	3:D:112:LEU:HD21	1.91	0.51
3:D:230:LEU:HB3	3:D:258:VAL:HG12	1.92	0.51
3:D:442:SER:OG	3:D:443:GLU:N	2.42	0.51
2:B:118:SER:HB3	2:B:142:SER:HB3	1.93	0.51
3:D:311:HIS:HB3	3:D:313:PHE:CE1	2.46	0.51
1:C:301:ILE:HA	1:C:317:THR:HG21	1.93	0.50
2:B:127:ILE:HB	2:B:291:ARG:HG2	1.93	0.50
2:B:433:THR:HG22	2:B:457:LYS:HB3	1.93	0.50
2:B:563:VAL:HA	2:B:566:LEU:HG	1.94	0.50
1:C:627:PHE:O	1:C:631:ILE:HD12	2.11	0.50
3:D:120:LEU:HD12	3:D:327:LYS:HB2	1.91	0.50
1:C:244:GLY:N	1:C:385:GLU:OE2	2.44	0.50
1:C:397:ARG:HA	1:C:473:THR:HG23	1.93	0.50
3:D:503:MET:SD	3:D:505:ARG:NH2	2.84	0.50
1:A:132:SER:HA	2:B:178:PRO:HB3	1.94	0.50
1:A:129:SER:OG	1:A:139:ARG:NH1	2.45	0.50
3:D:709:PHE:O	3:D:711:GLN:NE2	2.37	0.50
1:C:70:ASN:HB3	1:C:73:GLN:HB2	1.94	0.50
1:C:249:VAL:HB	1:C:253:GLU:HB2	1.94	0.50
1:C:106:PRO:HG3	3:D:114:PHE:CZ	2.46	0.49
1:A:212:ARG:NH1	1:A:238:LEU:O	2.46	0.49
1:A:315:TRP:CE3	1:A:317:THR:HB	2.47	0.49
1:C:254:ILE:HD13	1:C:268:GLY:HA3	1.94	0.49
1:C:566:VAL:O	1:C:570:VAL:HG23	2.13	0.49
2:B:518:ARG:HE	2:B:692:ARG:HH22	1.59	0.49
1:A:760:ILE:HD12	1:A:776:ILE:HD11	1.93	0.49
1:A:824:ILE:HG13	3:D:634:VAL:HG21	1.94	0.49
2:B:642:LEU:HD23	7:D:905:JC9:C12	2.42	0.49
1:A:813:MET:HE1	3:D:642:PHE:HB2	1.95	0.49
1:C:325:LEU:O	1:C:328:SER:OG	2.29	0.49
1:C:403:ILE:HG13	1:C:404:HIS:H	1.77	0.49
1:C:253:GLU:OE1	1:C:253:GLU:N	2.46	0.48
1:C:381:TRP:HB2	1:C:385:GLU:HB2	1.95	0.48
3:D:275:THR:HG23	3:D:371:ARG:HH22	1.77	0.48
1:A:664:ILE:HG22	1:A:666:GLY:H	1.78	0.48
1:C:119:LEU:HD21	1:C:325:LEU:HD21	1.95	0.48
1:C:212:ARG:HH21	1:C:238:LEU:HD11	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:463:ASP:OD2	3:D:796:TRP:NE1	2.40	0.48
2:B:42:HIS:HE1	2:B:49:ARG:NH2	2.11	0.48
1:C:103:THR:O	1:C:106:PRO:HD2	2.14	0.48
1:C:165:LEU:HG	1:C:167:VAL:HG23	1.94	0.48
1:C:467:ALA:HA	1:C:472:PHE:HB2	1.95	0.48
3:D:186:ILE:O	3:D:190:ILE:HG12	2.13	0.48
3:D:668:ASP:HA	3:D:672:GLN:HG3	1.96	0.48
2:B:375:VAL:HG22	2:B:388:ALA:HB3	1.95	0.48
3:D:613:VAL:HG23	3:D:635:TRP:HE1	1.78	0.48
3:D:615:ASN:ND2	7:D:905:JC9:C2	2.76	0.48
1:A:499:ASN:HD21	1:A:686:GLN:HG2	1.77	0.48
1:A:670:PRO:HG2	2:B:797:THR:HG21	1.95	0.48
1:A:502:MET:HE1	1:A:527:ILE:HG12	1.96	0.48
2:B:40:LEU:HD12	2:B:288:LEU:HD23	1.96	0.48
2:B:42:HIS:HB3	2:B:75:ASN:ND2	2.29	0.48
1:C:541:LEU:HB2	1:C:736:LEU:HD13	1.96	0.48
1:A:479:VAL:HG12	1:A:482:GLY:H	1.79	0.48
1:C:778:LYS:O	1:C:782:ASN:ND2	2.47	0.48
3:D:196:GLY:HA3	3:D:431:ARG:HD3	1.96	0.48
1:A:489:ARG:HA	1:A:496:LYS:HA	1.96	0.47
2:B:298:ILE:HG13	2:B:341:VAL:HG11	1.96	0.47
3:D:360:PRO:HD2	3:D:379:TRP:HZ3	1.80	0.47
1:C:631:ILE:HA	1:C:634:MET:HE2	1.95	0.47
1:C:86:GLN:HG2	1:C:306:ARG:HD2	1.96	0.47
3:D:157:MET:O	3:D:161:MET:HG3	2.14	0.47
2:B:462:ASP:OD2	2:B:795:TRP:NE1	2.48	0.47
2:B:539:ARG:HG2	2:B:746:LYS:HA	1.96	0.47
1:C:793:VAL:O	1:C:794:ARG:HG2	2.14	0.47
3:D:576:VAL:HA	3:D:579:TYR:CE2	2.50	0.47
2:B:204:ILE:HG21	2:B:218:GLN:HB3	1.97	0.47
1:C:403:ILE:HG23	1:C:405:GLN:H	1.79	0.47
3:D:42:VAL:HB	3:D:101:ASP:HA	1.96	0.47
1:A:635:VAL:HG21	2:B:827:ALA:HB1	1.96	0.47
3:D:468:ILE:O	3:D:472:VAL:HG22	2.14	0.47
1:A:672:LEU:HB3	1:A:702:MET:HE3	1.95	0.47
2:B:125:LEU:HA	2:B:145:PHE:O	2.14	0.47
2:B:174:THR:HG22	2:B:230:TYR:HB3	1.96	0.47
2:B:523:ASP:OD1	2:B:523:ASP:N	2.43	0.47
1:C:407:PRO:HG3	1:C:735:VAL:HA	1.97	0.47
2:B:539:ARG:HH21	2:B:748:VAL:HG11	1.80	0.47
2:B:206:LEU:HD12	2:B:239:ILE:HG12	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:744:CYS:HB3	1:C:798:CYS:HB3	1.51	0.46
2:B:249:THR:HB	2:B:273:PRO:HG3	1.97	0.46
1:C:30:GLY:H	1:C:87:VAL:HG13	1.79	0.46
1:A:45:ALA:HB1	1:A:285:VAL:HG21	1.97	0.46
1:A:351:TYR:O	1:A:367:TYR:N	2.47	0.46
2:B:737:TYR:OH	2:B:741:ARG:NH1	2.48	0.46
1:A:237:MET:SD	1:A:238:LEU:HD12	2.55	0.46
2:B:34:LEU:H	2:B:96:HIS:HD2	1.64	0.46
2:B:437:ARG:NH1	2:B:476:ASP:OD2	2.49	0.46
1:A:704:ARG:HD3	2:B:430:VAL:HG22	1.97	0.46
2:B:496:MET:HE2	2:B:512:LEU:HD22	1.98	0.46
1:C:541:LEU:HD21	1:C:746:LEU:HD13	1.97	0.46
1:C:668:ASN:OD1	1:C:673:ARG:NH2	2.44	0.46
3:D:497:MET:HE2	3:D:509:ALA:HB1	1.96	0.46
3:D:555:SER:HB3	3:D:560:VAL:HG13	1.98	0.46
1:C:115:ARG:HG3	1:C:314:ILE:HD12	1.98	0.46
3:D:217:GLN:OE1	3:D:247:VAL:HG22	2.16	0.46
3:D:73:MET:HE3	3:D:81:ILE:HG22	1.98	0.46
3:D:287:TYR:OH	3:D:295:ASP:OD2	2.29	0.46
1:C:32:VAL:N	1:C:91:LEU:O	2.35	0.46
1:C:71:ALA:HB1	3:D:114:PHE:HZ	1.80	0.46
1:C:501:MET:HE2	1:C:517:LEU:HD13	1.98	0.46
1:C:505:LEU:HD13	1:C:513:ILE:HG13	1.97	0.46
1:C:779:SER:HB2	1:C:785:MET:HE2	1.97	0.46
3:D:408:ILE:HG13	3:D:508:MET:HB2	1.98	0.46
2:B:530:GLU:OE1	1:C:755:ARG:NH2	2.40	0.46
3:D:287:TYR:CE2	3:D:292:ARG:HG2	2.52	0.46
3:D:678:SER:HB2	3:D:679:PRO:HD3	1.99	0.45
2:B:665:LEU:O	2:B:671:GLN:NE2	2.47	0.45
2:B:689:SER:N	6:B:903:GLU:OE1	2.49	0.45
1:A:249:VAL:HB	1:A:253:GLU:HB2	1.98	0.45
1:C:741:SER:OG	1:C:796:GLN:O	2.30	0.45
3:D:172:VAL:HG22	3:D:203:LEU:HD22	1.98	0.45
1:A:370:THR:O	1:A:371:HIS:ND1	2.50	0.45
3:D:156:VAL:HG21	3:D:362:LEU:HD12	1.98	0.45
2:B:489:VAL:O	2:B:492:VAL:HG12	2.17	0.45
2:B:513:THR:H	6:B:903:GLU:N	2.14	0.45
2:B:528:PHE:HD1	2:B:529:VAL:HG23	1.82	0.45
2:B:628:LYS:HB3	1:C:608:TRP:CZ3	2.52	0.45
2:B:438:LYS:O	2:B:452:VAL:N	2.48	0.45
3:D:299:ILE:HG23	3:D:342:VAL:HG21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:706:MET:O	1:A:710:ASN:N	2.49	0.45
1:A:830:LEU:HD13	1:A:833:ILE:HD12	1.99	0.45
1:C:539:THR:HG22	1:C:733:SER:HA	1.99	0.45
3:D:363:VAL:HG11	3:D:375:ARG:HD2	1.99	0.45
1:C:499:ASN:HD21	1:C:686:GLN:HG2	1.82	0.45
1:A:155:MET:SD	1:A:163:ILE:HG21	2.56	0.45
1:A:626:SER:O	1:A:630:ARG:N	2.46	0.45
3:D:738:TYR:HD2	3:D:739:MET:HE2	1.82	0.45
1:A:151:TRP:O	1:A:155:MET:HG3	2.17	0.44
2:B:42:HIS:CE1	2:B:49:ARG:NH2	2.85	0.44
1:C:174:ARG:O	1:C:177:GLN:HG2	2.17	0.44
1:C:578:TYR:HB2	1:C:603:LEU:HA	2.00	0.44
1:C:531:LYS:HA	1:C:531:LYS:HD3	1.83	0.44
1:C:570:VAL:HG22	1:C:636:TRP:CH2	2.52	0.44
1:C:751:GLU:HG2	1:C:752:LEU:N	2.32	0.44
1:A:42:PHE:CE2	1:A:62:ALA:HB1	2.52	0.44
1:A:138:LEU:HD22	1:A:325:LEU:HG	1.99	0.44
1:A:630:ARG:HG2	2:B:606:TRP:CZ3	2.53	0.44
2:B:42:HIS:HE1	2:B:49:ARG:HH22	1.65	0.44
3:D:410:THR:HG22	3:D:510:VAL:O	2.17	0.44
2:B:531:THR:OG1	2:B:532:GLY:N	2.51	0.44
3:D:134:MET:HG3	3:D:137:LYS:HG2	1.99	0.44
1:A:139:ARG:HG2	1:A:141:VAL:H	1.83	0.44
1:A:331:ALA:HA	1:A:337:ARG:HA	1.98	0.44
2:B:38:VAL:HG23	2:B:99:VAL:HB	1.99	0.44
3:D:102:ASP:OD1	3:D:102:ASP:N	2.48	0.44
1:A:133:ILE:HD11	2:B:136:ALA:HB3	1.99	0.43
1:A:573:VAL:HA	1:A:576:MET:HG2	2.00	0.43
1:A:662:GLU:HA	1:A:671:ARG:HH22	1.82	0.43
1:A:35:THR:OG1	1:A:38:HIS:ND1	2.37	0.43
2:B:34:LEU:O	2:B:68:ASN:N	2.51	0.43
2:B:628:LYS:HA	2:B:631:VAL:HG12	2.01	0.43
1:C:639:PHE:HA	1:C:642:ILE:HG12	1.99	0.43
1:A:540:ILE:HG12	1:A:730:ILE:HG12	1.99	0.43
1:A:634:MET:HE3	2:B:606:TRP:CZ3	2.53	0.43
2:B:387:HIS:ND1	2:B:387:HIS:O	2.51	0.43
3:D:458:LYS:HD2	3:D:796:TRP:HH2	1.83	0.43
2:B:278:SER:HB3	2:B:364:ILE:HG13	1.99	0.43
3:D:670:LYS:HG3	3:D:671:PHE:CD1	2.53	0.43
1:A:287:VAL:HG12	1:A:325:LEU:HD12	2.01	0.43
1:C:732:ASP:OD2	5:C:908:GLY:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:608:TRP:CE3	1:A:619:ILE:HD12	2.53	0.43
1:C:485:GLY:HA3	1:C:502:MET:HG2	2.00	0.43
3:D:422:VAL:HG11	3:D:454:LYS:HB3	2.00	0.43
1:C:302:THR:HG21	1:C:316:LYS:HB2	1.99	0.43
1:C:396:THR:HG22	1:C:473:THR:HG22	2.01	0.43
1:A:72:ILE:HG23	2:B:83:ILE:HD12	2.00	0.43
2:B:759:THR:OG1	2:B:760:GLY:N	2.52	0.43
1:C:152:PHE:HA	1:C:155:MET:HE3	2.00	0.43
3:D:142:MET:SD	3:D:332:SER:OG	2.76	0.43
3:D:243:VAL:O	3:D:247:VAL:HG23	2.19	0.43
1:A:143:PRO:HD2	1:A:146:HIS:ND1	2.34	0.42
3:D:171:ILE:HB	3:D:202:VAL:HG13	2.01	0.42
1:A:438:GLY:HA3	1:A:478:LEU:HB2	2.01	0.42
1:A:630:ARG:HH22	2:B:605:ILE:HG22	1.83	0.42
3:D:524:ASP:OD1	3:D:524:ASP:N	2.52	0.42
2:B:630:MET:SD	1:C:825:VAL:HG22	2.59	0.42
1:C:364:VAL:HG21	1:C:379:ILE:HG12	2.00	0.42
1:A:117:PRO:HA	1:A:136:SER:HB2	2.01	0.42
2:B:314:PRO:HG2	2:B:333:THR:HA	2.01	0.42
1:C:205:THR:O	1:C:209:MET:HG2	2.18	0.42
3:D:407:SER:O	3:D:507:TYR:N	2.52	0.42
1:C:283:ASP:OD2	1:C:335:THR:OG1	2.30	0.42
1:C:123:THR:HG21	1:C:137:PHE:HZ	1.85	0.42
3:D:41:LEU:HD11	3:D:70:LEU:HD12	2.02	0.42
1:A:164:ILE:HG13	1:A:216:ALA:HB3	2.01	0.42
1:A:486:THR:HG22	1:A:691:ILE:HD11	2.02	0.42
2:B:97:GLY:HA2	2:B:123:PRO:HG2	2.02	0.42
2:B:430:VAL:HG12	2:B:432:ASN:N	2.31	0.42
1:C:62:ALA:O	4:C:901:NAG:O6	2.33	0.42
1:A:554:PHE:CD2	1:A:555:MET:HG3	2.55	0.42
2:B:259:SER:HB2	2:B:280:SER:HA	2.02	0.42
1:C:655:LEU:HD12	1:C:656:VAL:HG23	2.02	0.42
2:B:650:ALA:O	2:B:654:ILE:HG12	2.19	0.42
1:C:502:MET:HE2	1:C:517:LEU:HD21	2.01	0.42
1:A:335:THR:HG22	1:A:336:GLY:N	2.35	0.41
1:C:125:MET:O	1:C:139:ARG:NH1	2.38	0.41
3:D:817:ASN:C	3:D:817:ASN:HD22	2.28	0.41
2:B:557:VAL:HG12	2:B:560:MET:H	1.84	0.41
2:B:612:VAL:HA	2:B:638:ALA:HB1	2.02	0.41
2:B:152:GLN:O	2:B:156:THR:HG23	2.20	0.41
2:B:265:THR:HG23	2:B:372:TRP:CD1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:436:CYS:HB2	2:B:456:CYS:H	1.85	0.41
1:C:604:SER:O	1:C:608:TRP:CD1	2.73	0.41
3:D:825:LEU:O	3:D:829:MET:HG2	2.20	0.41
1:C:124:ARG:HH12	1:C:142:PRO:HG2	1.85	0.41
1:C:404:HIS:CD2	1:C:411:VAL:HG22	2.55	0.41
3:D:97:VAL:HB	3:D:123:ILE:HG12	2.03	0.41
1:A:490:VAL:HB	1:A:493:SER:HB3	2.02	0.41
2:B:677:SER:OG	2:B:678:PRO:HD3	2.20	0.41
1:C:472:PHE:HE1	1:C:768:TRP:CE3	2.38	0.41
1:C:505:LEU:HD23	1:C:527:ILE:HG21	2.03	0.41
3:D:410:THR:CG2	3:D:478:LEU:HD11	2.50	0.41
1:A:570:VAL:HA	1:A:573:VAL:HG12	2.01	0.41
2:B:74:MET:HE1	2:B:81:SER:HB3	2.03	0.41
2:B:134:ILE:HD11	2:B:149:ALA:N	2.36	0.41
1:C:580:LEU:HD23	1:C:580:LEU:HA	1.85	0.41
1:A:656:VAL:HG13	1:A:657:LEU:HG	2.03	0.41
2:B:127:ILE:HD12	2:B:291:ARG:HB3	2.01	0.41
1:C:129:SER:HB2	1:C:139:ARG:HH21	1.86	0.41
1:A:139:ARG:NH1	1:A:143:PRO:HB3	2.36	0.41
1:A:407:PRO:HB3	1:A:738:PHE:CG	2.56	0.41
1:A:440:ASN:N	1:A:440:ASN:OD1	2.53	0.41
1:A:722:ARG:NH2	1:A:745:ASP:OD1	2.54	0.41
2:B:74:MET:HE2	2:B:76:ARG:HB2	2.03	0.41
2:B:103:ASP:OD1	2:B:103:ASP:N	2.51	0.41
2:B:628:LYS:HB3	1:C:608:TRP:HZ3	1.86	0.41
1:C:463:LEU:HD13	1:C:514:VAL:HG21	2.02	0.41
1:A:762:MET:HE3	1:A:762:MET:HB2	1.88	0.41
1:C:440:ASN:OD1	1:C:441:ASP:N	2.54	0.41
1:C:831:ILE:HD13	1:C:831:ILE:HA	1.95	0.41
3:D:190:ILE:HD12	3:D:197:TRP:CG	2.56	0.41
3:D:576:VAL:HA	3:D:579:TYR:CD2	2.56	0.41
1:A:541:LEU:HD22	1:A:736:LEU:HD22	2.03	0.41
1:C:79:CYS:HB3	1:C:308:CYS:HB3	1.68	0.41
1:C:124:ARG:HH22	1:C:351:TYR:HE1	1.69	0.41
3:D:365:ILE:HG22	3:D:373:TRP:HE3	1.86	0.41
3:D:519:ARG:HA	3:D:522:VAL:HG12	2.03	0.41
1:A:351:TYR:HB2	1:A:367:TYR:HB3	2.04	0.40
2:B:53:ASN:HA	2:B:69:VAL:HB	2.02	0.40
2:B:256:ILE:HG13	2:B:277:ILE:HD11	2.03	0.40
1:A:79:CYS:HB3	1:A:308:CYS:HB3	1.95	0.40
1:A:143:PRO:HD2	1:A:146:HIS:CE1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:829:ALA:O	2:B:833:ILE:HG12	2.21	0.40
1:C:626:SER:OG	1:C:627:PHE:N	2.55	0.40
2:B:206:LEU:HD13	2:B:215:THR:HB	2.02	0.40
2:B:301:THR:HG21	2:B:343:TRP:HB2	2.03	0.40
2:B:459:PHE:CE2	2:B:791:LEU:HB3	2.56	0.40
2:B:518:ARG:HH12	6:B:903:GLU:N	2.19	0.40
1:C:760:ILE:HD12	1:C:776:ILE:HD11	2.03	0.40
3:D:522:VAL:HG13	3:D:523:VAL:HG22	2.03	0.40
2:B:536:MET:HA	2:B:749:THR:HA	2.04	0.40
3:D:156:VAL:HG22	3:D:379:TRP:CD1	2.56	0.40
3:D:464:ILE:O	3:D:468:ILE:HG12	2.22	0.40
3:D:573:ALA:HA	3:D:576:VAL:HG12	2.03	0.40
1:A:824:ILE:HG21	3:D:634:VAL:HG21	2.02	0.40
1:C:29:ILE:HD13	1:C:29:ILE:HA	1.96	0.40
1:C:466:LEU:HB3	1:C:472:PHE:CE2	2.56	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	790/819 (96%)	771 (98%)	19 (2%)	0	100	100
1	C	801/819 (98%)	776 (97%)	25 (3%)	0	100	100
2	B	788/809 (97%)	763 (97%)	25 (3%)	0	100	100
3	D	787/808 (97%)	752 (96%)	35 (4%)	0	100	100
All	All	3166/3255 (97%)	3062 (97%)	104 (3%)	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	601/706 (85%)	601 (100%)	0	100	100
1	C	624/706 (88%)	624 (100%)	0	100	100
2	B	554/707 (78%)	554 (100%)	0	100	100
3	D	596/705 (84%)	596 (100%)	0	100	100
All	All	2375/2824 (84%)	2375 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	196	GLN
1	A	487	GLN
1	A	770	GLN
2	B	53	ASN
2	B	432	ASN
2	B	482	ASN
2	B	621	ASN
2	B	693	ASN
2	B	736	ASN
1	C	162	HIS
1	C	371	HIS
1	C	453	GLN
1	C	487	GLN
1	C	616	ASN
1	C	719	GLN
3	D	118	GLN
3	D	218	ASN
3	D	385	GLN
3	D	437	GLN
3	D	697	ASN
3	D	767	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 ⓘ

24 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	A	906	1	14,14,15	0.20	0	17,19,21	0.36	0
4	NAG	B	901	2	14,14,15	0.23	0	17,19,21	0.44	0
4	NAG	A	903	1	14,14,15	0.24	0	17,19,21	0.42	0
4	NAG	D	902	3	14,14,15	0.22	0	17,19,21	0.43	0
5	GLY	A	908	-	3,3,4	0.60	0	0,2,4	-	-
7	JC9	D	905	-	17,17,17	0.51	0	14,24,24	0.60	0
4	NAG	C	903	1	14,14,15	0.27	0	17,19,21	0.50	0
4	NAG	A	905	1	14,14,15	0.26	0	17,19,21	0.41	0
4	NAG	C	902	1	14,14,15	0.28	0	17,19,21	0.46	0
4	NAG	C	907	1	14,14,15	0.21	0	17,19,21	0.40	0
4	NAG	A	904	1	14,14,15	0.21	0	17,19,21	0.43	0
4	NAG	A	907	1	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	C	906	1	14,14,15	0.25	0	17,19,21	0.41	0
5	GLY	C	908	-	3,3,4	0.61	0	0,2,4	-	-
4	NAG	C	904	1	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	C	905	1	14,14,15	0.25	0	17,19,21	0.35	0
4	NAG	A	901	1	14,14,15	0.32	0	17,19,21	0.40	0
4	NAG	C	901	1	14,14,15	0.22	0	17,19,21	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GLU	D	904	-	7,8,9	0.87	0	4,9,11	0.98	0
4	NAG	A	902	1	14,14,15	0.22	0	17,19,21	0.44	0
6	GLU	B	903	-	7,8,9	0.87	0	4,9,11	1.01	0
4	NAG	B	902	2	14,14,15	0.20	0	17,19,21	0.42	0
4	NAG	D	903	3	14,14,15	0.21	0	17,19,21	0.43	0
4	NAG	D	901	3	14,14,15	0.21	0	17,19,21	0.42	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	906	1	-	2/6/23/26	0/1/1/1
4	NAG	B	901	2	-	4/6/23/26	0/1/1/1
4	NAG	A	903	1	-	2/6/23/26	0/1/1/1
4	NAG	D	902	3	-	2/6/23/26	0/1/1/1
5	GLY	A	908	-	-	0/0/1/2	-
7	JC9	D	905	-	-	1/6/23/23	0/2/2/2
4	NAG	C	903	1	-	2/6/23/26	0/1/1/1
4	NAG	A	905	1	-	1/6/23/26	0/1/1/1
4	NAG	C	902	1	-	0/6/23/26	0/1/1/1
4	NAG	C	907	1	-	2/6/23/26	0/1/1/1
4	NAG	A	904	1	-	2/6/23/26	0/1/1/1
4	NAG	A	907	1	-	2/6/23/26	0/1/1/1
4	NAG	C	906	1	-	2/6/23/26	0/1/1/1
5	GLY	C	908	-	-	0/0/1/2	-
4	NAG	C	904	1	-	2/6/23/26	0/1/1/1
4	NAG	C	905	1	-	2/6/23/26	0/1/1/1
4	NAG	A	901	1	-	0/6/23/26	0/1/1/1
4	NAG	C	901	1	-	2/6/23/26	0/1/1/1
6	GLU	D	904	-	-	2/6/7/9	-
4	NAG	A	902	1	-	2/6/23/26	0/1/1/1
6	GLU	B	903	-	-	0/6/7/9	-
4	NAG	B	902	2	-	2/6/23/26	0/1/1/1
4	NAG	D	903	3	-	1/6/23/26	0/1/1/1
4	NAG	D	901	3	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (37) torsion outliers are listed below:

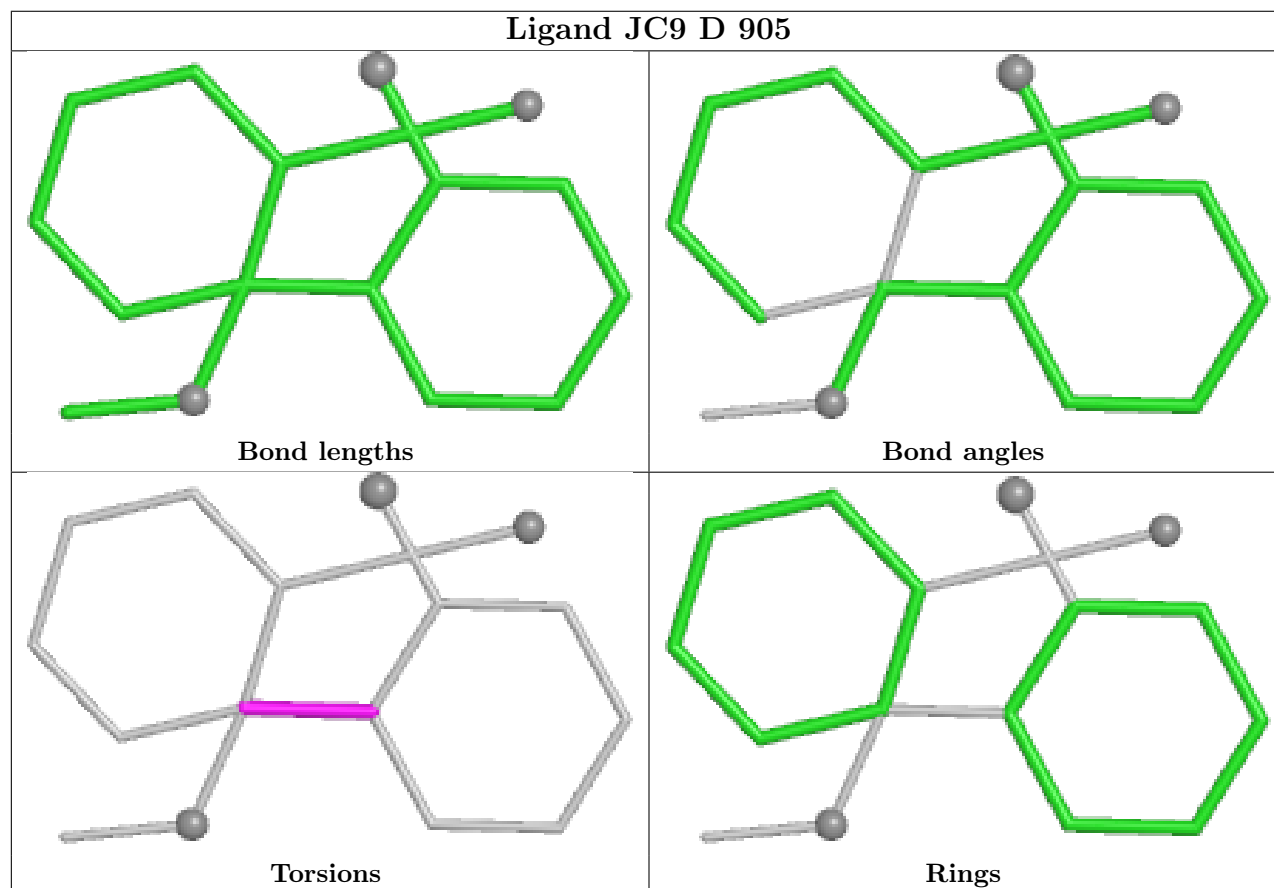
Mol	Chain	Res	Type	Atoms
7	D	905	JC9	N1-C1-C7-C9
4	C	901	NAG	O5-C5-C6-O6
4	B	902	NAG	C4-C5-C6-O6
4	A	903	NAG	O5-C5-C6-O6
4	C	903	NAG	O5-C5-C6-O6
4	A	903	NAG	C4-C5-C6-O6
4	A	902	NAG	O5-C5-C6-O6
4	A	907	NAG	O5-C5-C6-O6
4	C	904	NAG	O5-C5-C6-O6
4	C	901	NAG	C4-C5-C6-O6
4	A	906	NAG	O5-C5-C6-O6
4	B	902	NAG	O5-C5-C6-O6
4	C	905	NAG	O5-C5-C6-O6
4	A	902	NAG	C4-C5-C6-O6
4	A	907	NAG	C4-C5-C6-O6
4	D	901	NAG	C4-C5-C6-O6
4	B	901	NAG	C8-C7-N2-C2
4	B	901	NAG	O7-C7-N2-C2
4	D	902	NAG	C8-C7-N2-C2
4	D	902	NAG	O7-C7-N2-C2
4	C	903	NAG	C4-C5-C6-O6
4	C	907	NAG	O5-C5-C6-O6
4	C	905	NAG	C4-C5-C6-O6
4	A	906	NAG	C4-C5-C6-O6
4	D	901	NAG	O5-C5-C6-O6
4	C	904	NAG	C4-C5-C6-O6
4	C	906	NAG	O5-C5-C6-O6
4	A	904	NAG	O5-C5-C6-O6
4	A	904	NAG	C4-C5-C6-O6
4	D	903	NAG	O5-C5-C6-O6
4	C	907	NAG	C4-C5-C6-O6
4	A	905	NAG	C4-C5-C6-O6
6	D	904	GLU	OE1-CD-CG-CB
4	B	901	NAG	C4-C5-C6-O6
6	D	904	GLU	OE2-CD-CG-CB
4	B	901	NAG	O5-C5-C6-O6
4	C	906	NAG	C4-C5-C6-O6

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	908	GLY	1	0
7	D	905	JC9	4	0
5	C	908	GLY	1	0
4	C	901	NAG	1	0
6	B	903	GLU	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.