



## wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 02:03 PM BST

PDB ID : 4UQJ  
EMDB ID: : EMD-2680  
Title : Cryo-EM density map of GluA2em in complex with ZK200775  
Authors : Meyerson, J.R.; Kumar, J.; Chittori, S.; Rao, P.; Pierson, J.; Bartesaghi, A.;  
Mayer, M.L.; Subramaniam, S.  
Deposited on : 2014-06-24  
Resolution : 10.40 Å(reported)  
Based on PDB ID : 3KG2

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.  
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/EMValidationReportHelp>

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MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk27241

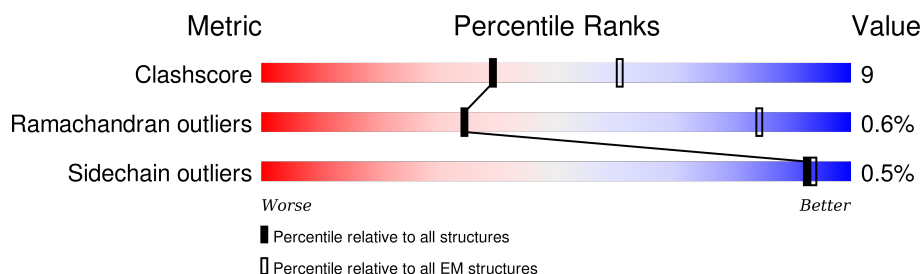
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 10.40 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	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	826	73% 17% 10%
1	B	826	73% 17% 10%
1	C	826	73% 17% 10%
1	D	826	72% 17% 10%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 21793 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.

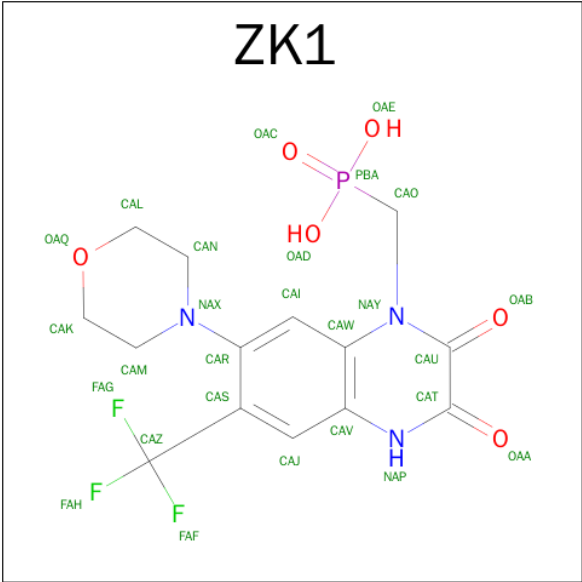
- Molecule 1 is a protein called GLUTAMATE RECEPTOR 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	746	Total	C	N	O	S	0	0
			5423	3497	858	1042	26		
1	B	746	Total	C	N	O	S	0	0
			5423	3497	858	1042	26		
1	C	746	Total	C	N	O	S	0	0
			5423	3497	858	1042	26		
1	D	745	Total	C	N	O	S	0	0
			5416	3492	857	1041	26		

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	GLU	ASN	ENGINEERED MUTATION	UNP P19491
A	382	LEU	VAL	ENGINEERED MUTATION	UNP P19491
A	589	ALA	CYS	ENGINEERED MUTATION	UNP P19491
A	758	LEU	VAL	VARIANT	UNP P19491
B	241	GLU	ASN	ENGINEERED MUTATION	UNP P19491
B	382	LEU	VAL	ENGINEERED MUTATION	UNP P19491
B	589	ALA	CYS	ENGINEERED MUTATION	UNP P19491
B	758	LEU	VAL	VARIANT	UNP P19491
C	241	GLU	ASN	ENGINEERED MUTATION	UNP P19491
C	382	LEU	VAL	ENGINEERED MUTATION	UNP P19491
C	589	ALA	CYS	ENGINEERED MUTATION	UNP P19491
C	758	LEU	VAL	VARIANT	UNP P19491
D	241	GLU	ASN	ENGINEERED MUTATION	UNP P19491
D	382	LEU	VAL	ENGINEERED MUTATION	UNP P19491
D	589	ALA	CYS	ENGINEERED MUTATION	UNP P19491
D	758	LEU	VAL	VARIANT	UNP P19491

- Molecule 2 is {[7-MORPHOLIN-4-YL-2,3-DIOXO-6-(TRIFLUOROMETHYL)-3,4-DIHYDROQUINOXALIN-1(2H)-YL]METHYL}PHOSPHONIC ACID (three-letter code: ZK1) (formula: C<sub>14</sub>H<sub>15</sub>F<sub>3</sub>N<sub>3</sub>O<sub>6</sub>P).



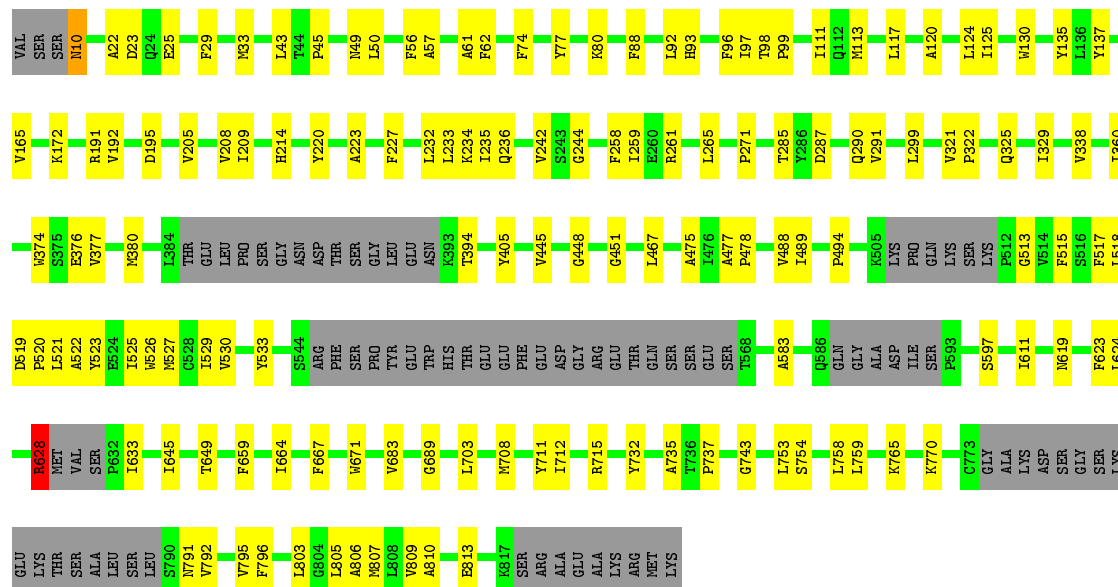
Mol	Chain	Residues	Atoms						AltConf
2	A	1	Total	C	F	N	O	P	0
			27	14	3	3	6	1	
2	B	1	Total	C	F	N	O	P	0
			27	14	3	3	6	1	
2	C	1	Total	C	F	N	O	P	0
			27	14	3	3	6	1	
2	D	1	Total	C	F	N	O	P	0
			27	14	3	3	6	1	





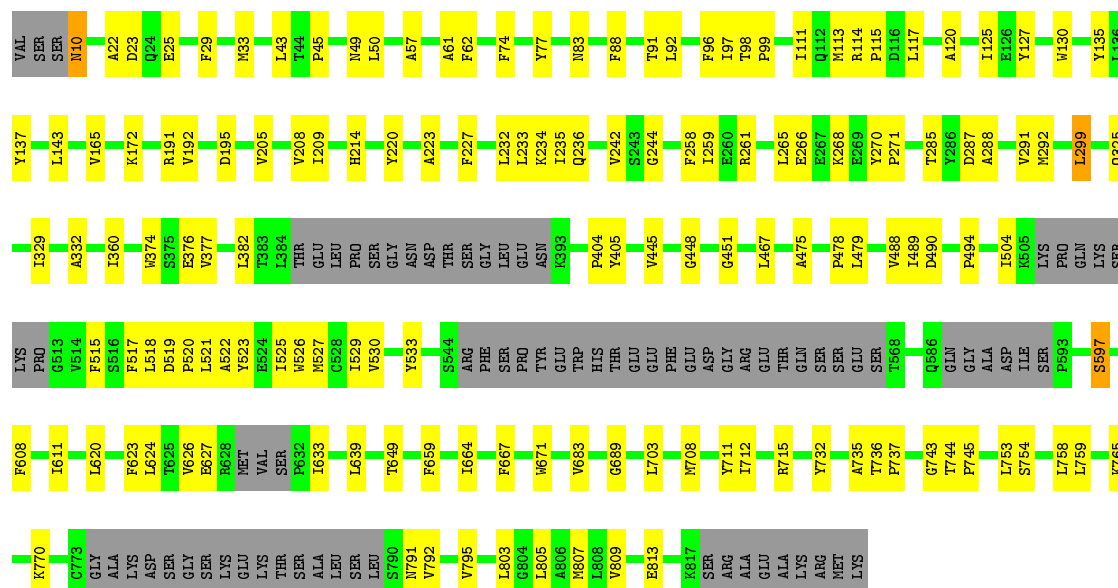
• Molecule 1: GLUTAMATE RECEPTOR 2

Chain C: 73% 17% 10%



• Molecule 1: GLUTAMATE RECEPTOR 2

Chain D: 72% 17% 10%



## 4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	INDIVIDUAL PARTICLES, Not provided	Depositor
Microscope	OTHER	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	Not provided	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	Not provided	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: ZK1

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  > 2$	RMSZ	# $ Z  > 2$
1	A	0.21	0/5539	0.41	3/7577 (0.0%)
1	B	0.21	0/5539	0.35	0/7577
1	C	0.21	0/5539	0.41	3/7577 (0.0%)
1	D	0.21	0/5531	0.35	0/7566
All	All	0.21	0/22148	0.38	6/30297 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	628	ARG	NE-CZ-NH2	-13.84	113.38	120.30
1	A	628	ARG	NE-CZ-NH2	-12.54	114.03	120.30
1	C	628	ARG	NE-CZ-NH1	11.79	126.19	120.30
1	A	628	ARG	NE-CZ-NH1	11.27	125.93	120.30
1	A	628	ARG	CD-NE-CZ	5.70	131.58	123.60
1	C	628	ARG	CD-NE-CZ	5.33	131.06	123.60

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	5423	0	4953	105	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	5423	0	4953	102	0
1	C	5423	0	4953	102	0
1	D	5416	0	4945	115	0
2	A	27	0	13	1	0
2	B	27	0	13	0	0
2	C	27	0	13	1	0
2	D	27	0	13	2	0
All	All	21793	0	19856	391	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 9.

All (391) 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:626:VAL:HB	1:C:628:ARG:HH21	1.24	1.01
1:B:649:THR:HG22	1:B:703:LEU:HB2	1.68	0.76
1:D:649:THR:HG22	1:D:703:LEU:HB2	1.68	0.76
1:C:649:THR:HG22	1:C:703:LEU:HB2	1.68	0.75
1:A:649:THR:HG22	1:A:703:LEU:HB2	1.68	0.74
1:C:611:ILE:HG21	1:D:795:VAL:HG21	1.70	0.73
1:A:628:ARG:NH1	1:D:627:GLU:H	1.91	0.68
1:C:521:LEU:HD22	1:C:526:TRP:CD2	2.32	0.64
1:B:521:LEU:HD23	1:B:525:ILE:HB	1.80	0.63
1:A:810:ALA:HB2	1:D:597:SER:OG	1.98	0.63
1:D:521:LEU:HD23	1:D:525:ILE:HB	1.80	0.63
1:B:521:LEU:HD22	1:B:526:TRP:CD2	2.35	0.62
1:A:475:ALA:HB3	1:A:735:ALA:HB3	1.82	0.62
1:C:475:ALA:HB3	1:C:735:ALA:HB3	1.82	0.62
1:A:521:LEU:HD23	1:A:525:ILE:HB	1.80	0.62
1:A:521:LEU:HD22	1:A:526:TRP:CD2	2.35	0.62
1:A:522:ALA:HB3	1:A:525:ILE:HG13	1.82	0.61
1:B:522:ALA:HB3	1:B:525:ILE:HG13	1.83	0.61
1:A:628:ARG:NH2	1:D:623:PHE:HD1	1.98	0.61
1:D:522:ALA:HB3	1:D:525:ILE:HG13	1.83	0.61
1:A:50:LEU:HD23	1:A:57:ALA:HB1	1.82	0.61
1:C:50:LEU:HD23	1:C:57:ALA:HB1	1.82	0.61
1:D:489:ILE:HD12	1:D:735:ALA:HB1	1.83	0.61
1:D:521:LEU:HD22	1:D:526:TRP:CD2	2.35	0.60
1:B:50:LEU:HD23	1:B:57:ALA:HB1	1.83	0.60
1:B:489:ILE:HD12	1:B:735:ALA:HB1	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:522:ALA:HB3	1:C:525:ILE:HG13	1.84	0.60
1:C:521:LEU:HD23	1:C:525:ILE:HB	1.83	0.59
1:D:50:LEU:HD23	1:D:57:ALA:HB1	1.83	0.59
1:D:235:ILE:HD13	1:D:242:VAL:HG21	1.85	0.59
1:D:517:PHE:HB2	1:D:791:ASN:OD1	2.02	0.59
1:A:795:VAL:HG21	1:D:611:ILE:HG21	1.84	0.59
1:A:628:ARG:HH12	1:D:627:GLU:H	1.50	0.59
1:B:10:ASN:N	1:B:10:ASN:HD22	2.01	0.58
1:B:597:SER:OG	1:C:810:ALA:HB2	2.04	0.58
1:C:10:ASN:HD22	1:C:10:ASN:N	2.01	0.57
1:B:235:ILE:HD13	1:B:242:VAL:HG21	1.85	0.57
1:A:235:ILE:HD13	1:A:242:VAL:HG21	1.85	0.57
1:C:99:PRO:HA	1:C:113:MET:HB2	1.87	0.57
1:A:93:HIS:ND1	1:A:322:PRO:HB3	2.19	0.57
1:C:235:ILE:HD13	1:C:242:VAL:HG21	1.85	0.57
1:C:521:LEU:HD22	1:C:526:TRP:CG	2.40	0.57
1:B:633:ILE:HD13	1:B:639:LEU:HD21	1.86	0.57
1:A:10:ASN:N	1:A:10:ASN:HD22	2.02	0.57
1:A:99:PRO:HA	1:A:113:MET:HB2	1.87	0.57
1:B:99:PRO:HA	1:B:113:MET:HB2	1.87	0.56
1:C:93:HIS:ND1	1:C:322:PRO:HB3	2.19	0.56
1:C:517:PHE:HB2	1:C:791:ASN:OD1	2.05	0.56
1:D:10:ASN:N	1:D:10:ASN:HD22	2.01	0.56
1:A:77:TYR:CE2	1:A:98:THR:HG21	2.40	0.56
1:C:227:PHE:CD1	1:C:244:GLY:HA3	2.41	0.56
1:B:517:PHE:HB2	1:B:791:ASN:OD1	2.06	0.56
1:D:633:ILE:HD13	1:D:639:LEU:HD21	1.86	0.56
1:D:299:LEU:HD11	1:D:332:ALA:HB2	1.88	0.56
1:B:521:LEU:HD22	1:B:526:TRP:CG	2.41	0.56
1:A:209:ILE:HA	1:A:214:HIS:CD2	2.41	0.56
1:B:77:TYR:CE2	1:B:98:THR:HG21	2.41	0.56
1:C:209:ILE:HA	1:C:214:HIS:CD2	2.41	0.56
1:B:376:GLU:HG3	1:B:377:VAL:HG13	1.88	0.56
1:B:227:PHE:CD1	1:B:244:GLY:HA3	2.41	0.56
1:B:209:ILE:HA	1:B:214:HIS:CD2	2.41	0.55
1:C:77:TYR:CE2	1:C:98:THR:HG21	2.40	0.55
1:D:99:PRO:HA	1:D:113:MET:HB2	1.87	0.55
1:A:227:PHE:CD1	1:A:244:GLY:HA3	2.41	0.55
1:C:376:GLU:HG3	1:C:377:VAL:HG13	1.88	0.55
1:A:521:LEU:HD22	1:A:526:TRP:CG	2.41	0.55
1:D:77:TYR:CE2	1:D:98:THR:HG21	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:GLU:HG3	1:A:377:VAL:HG13	1.88	0.55
1:D:209:ILE:HA	1:D:214:HIS:CD2	2.41	0.55
1:D:227:PHE:CD1	1:D:244:GLY:HA3	2.41	0.55
1:D:521:LEU:HD22	1:D:526:TRP:CG	2.41	0.55
1:A:765:LYS:O	1:A:770:LYS:HB2	2.07	0.55
1:D:376:GLU:HG3	1:D:377:VAL:HG13	1.88	0.55
1:B:765:LYS:O	1:B:770:LYS:HB2	2.07	0.55
1:D:753:LEU:HD22	1:D:758:LEU:HD13	1.89	0.55
1:C:520:PRO:HA	1:C:623:PHE:HE2	1.70	0.54
1:B:753:LEU:HD22	1:B:758:LEU:HD13	1.89	0.54
1:C:753:LEU:HD22	1:C:758:LEU:HD13	1.89	0.54
1:D:62:PHE:HE2	1:D:92:LEU:HD12	1.72	0.54
1:A:753:LEU:HD22	1:A:758:LEU:HD13	1.89	0.54
1:B:62:PHE:HE2	1:B:92:LEU:HD12	1.72	0.54
1:D:765:LYS:O	1:D:770:LYS:HB2	2.07	0.54
1:D:475:ALA:HB3	1:D:735:ALA:HB3	1.89	0.54
1:B:299:LEU:HD11	1:B:332:ALA:HB2	1.88	0.54
1:B:475:ALA:HB3	1:B:735:ALA:HB3	1.90	0.54
1:A:62:PHE:HE2	1:A:92:LEU:HD12	1.73	0.54
1:B:504:ILE:HG21	1:B:633:ILE:HD12	1.90	0.53
1:D:504:ILE:HG21	1:D:633:ILE:HD12	1.90	0.53
1:D:754:SER:HB3	1:D:759:LEU:HD12	1.90	0.53
1:A:754:SER:HB3	1:A:759:LEU:HD12	1.90	0.53
1:C:754:SER:HB3	1:C:759:LEU:HD12	1.90	0.53
1:C:62:PHE:HE2	1:C:92:LEU:HD12	1.73	0.53
1:C:765:LYS:O	1:C:770:LYS:HB2	2.07	0.53
1:B:626:VAL:HB	1:C:628:ARG:NH2	2.08	0.53
1:B:626:VAL:CB	1:C:628:ARG:HH21	2.08	0.53
1:D:405:TYR:CG	1:D:478:PRO:HG3	2.44	0.53
1:B:405:TYR:CG	1:B:478:PRO:HG3	2.44	0.53
1:A:628:ARG:HH11	1:D:626:VAL:HB	1.74	0.52
1:A:205:VAL:O	1:A:209:ILE:HG13	2.10	0.52
1:B:754:SER:HB3	1:B:759:LEU:HD12	1.90	0.52
1:D:405:TYR:CD2	1:D:478:PRO:HG3	2.45	0.52
1:A:405:TYR:CG	1:A:478:PRO:HG3	2.45	0.52
1:C:633:ILE:HD11	1:C:645:ILE:HD12	1.92	0.52
1:D:523:TYR:O	1:D:527:MET:HG2	2.10	0.52
1:A:611:ILE:HG21	1:B:795:VAL:HG21	1.92	0.52
1:B:205:VAL:O	1:B:209:ILE:HG13	2.10	0.52
1:A:523:TYR:O	1:A:527:MET:HG2	2.09	0.52
1:D:205:VAL:O	1:D:209:ILE:HG13	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:523:TYR:O	1:C:527:MET:HG2	2.10	0.51
1:A:633:ILE:HD11	1:A:645:ILE:HD12	1.92	0.51
1:A:405:TYR:CD2	1:A:478:PRO:HG3	2.45	0.51
1:C:405:TYR:CD2	1:C:478:PRO:HG3	2.45	0.51
1:B:611:ILE:HG21	1:C:795:VAL:HG21	1.93	0.51
1:C:515:PHE:HD1	1:C:518:LEU:HD12	1.76	0.51
1:B:523:TYR:O	1:B:527:MET:HG2	2.10	0.51
1:C:405:TYR:CG	1:C:478:PRO:HG3	2.45	0.51
1:C:205:VAL:O	1:C:209:ILE:HG13	2.10	0.51
1:B:405:TYR:CD2	1:B:478:PRO:HG3	2.45	0.50
1:A:608:PHE:HZ	1:B:796:PHE:CZ	2.30	0.50
1:A:515:PHE:HD1	1:A:518:LEU:HD12	1.76	0.50
1:D:74:PHE:CZ	1:D:285:THR:HG23	2.46	0.50
1:B:74:PHE:CZ	1:B:285:THR:HG23	2.46	0.50
1:D:43:LEU:O	1:D:45:PRO:HD3	2.11	0.50
1:D:515:PHE:HD1	1:D:518:LEU:HD12	1.76	0.50
1:B:43:LEU:O	1:B:45:PRO:HD3	2.11	0.50
1:A:56:PHE:CE2	1:B:91:THR:HG21	2.47	0.50
1:A:74:PHE:CZ	1:A:285:THR:HG23	2.47	0.50
1:A:494:PRO:HA	1:A:732:TYR:O	2.12	0.49
1:B:515:PHE:HD1	1:B:518:LEU:HD12	1.76	0.49
1:A:659:PHE:HB3	1:A:671:TRP:HB2	1.95	0.49
1:C:526:TRP:O	1:C:529:ILE:HG22	2.13	0.49
1:B:526:TRP:O	1:B:529:ILE:HG22	2.12	0.49
1:D:50:LEU:HD22	1:D:61:ALA:HB2	1.95	0.49
1:A:80:LYS:HA	1:B:83:ASN:ND2	2.27	0.49
1:C:659:PHE:HB3	1:C:671:TRP:HB2	1.95	0.49
1:A:517:PHE:HB2	1:A:791:ASN:OD1	2.12	0.49
1:C:50:LEU:HD22	1:C:61:ALA:HB2	1.95	0.49
1:A:445:VAL:HG13	1:A:448:GLY:HA2	1.95	0.49
1:C:445:VAL:HG13	1:C:448:GLY:HA2	1.95	0.49
1:D:659:PHE:HB3	1:D:671:TRP:HB2	1.95	0.49
1:C:74:PHE:CZ	1:C:285:THR:HG23	2.47	0.49
1:C:80:LYS:HA	1:D:83:ASN:ND2	2.27	0.49
1:B:50:LEU:HD22	1:B:61:ALA:HB2	1.94	0.49
1:D:445:VAL:HG13	1:D:448:GLY:HA2	1.95	0.49
1:C:56:PHE:CE2	1:D:91:THR:HG21	2.47	0.49
1:C:43:LEU:O	1:C:45:PRO:HD3	2.12	0.49
1:C:494:PRO:HA	1:C:732:TYR:O	2.12	0.49
1:C:619:ASN:HA	1:D:624:LEU:HD13	1.95	0.48
1:B:659:PHE:HB3	1:B:671:TRP:HB2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:494:PRO:HA	1:D:732:TYR:O	2.13	0.48
1:B:494:PRO:HA	1:B:732:TYR:O	2.14	0.48
1:B:536:VAL:HG22	1:C:803:LEU:HD21	1.94	0.48
1:A:526:TRP:O	1:A:529:ILE:HG22	2.13	0.48
1:A:628:ARG:CZ	1:D:623:PHE:HD1	2.27	0.48
1:A:489:ILE:HD12	1:A:735:ALA:HB1	1.94	0.48
1:A:50:LEU:HD22	1:A:61:ALA:HB2	1.95	0.48
1:A:43:LEU:O	1:A:45:PRO:HD3	2.12	0.48
1:B:664:ILE:HB	1:B:667:PHE:HD2	1.79	0.48
1:B:445:VAL:HG13	1:B:448:GLY:HA2	1.95	0.48
1:C:258:PHE:HD2	1:C:259:ILE:HD12	1.79	0.48
1:B:29:PHE:O	1:B:33:MET:HG2	2.14	0.48
1:A:258:PHE:HD2	1:A:259:ILE:HD12	1.79	0.48
1:B:539:VAL:HG13	1:C:807:MET:SD	2.53	0.48
1:C:195:ASP:HA	1:C:223:ALA:HB3	1.96	0.48
1:D:526:TRP:O	1:D:529:ILE:HG22	2.13	0.47
1:D:29:PHE:O	1:D:33:MET:HG2	2.14	0.47
1:A:124:LEU:HD13	1:A:380:MET:HE1	1.96	0.47
1:B:195:ASP:HA	1:B:223:ALA:HB3	1.96	0.47
1:C:521:LEU:HD22	1:C:526:TRP:CE2	2.49	0.47
1:C:489:ILE:HD12	1:C:735:ALA:HB1	1.95	0.47
1:C:711:TYR:O	1:C:715:ARG:HG2	2.14	0.47
1:A:664:ILE:HB	1:A:667:PHE:HD2	1.79	0.47
1:D:258:PHE:HD2	1:D:259:ILE:HD12	1.79	0.47
1:D:711:TYR:O	1:D:715:ARG:HG2	2.14	0.47
1:C:124:LEU:HD13	1:C:380:MET:HE1	1.96	0.47
1:A:711:TYR:O	1:A:715:ARG:HG2	2.14	0.47
1:B:97:ILE:HG13	1:B:111:ILE:HB	1.96	0.47
1:B:258:PHE:HD2	1:B:259:ILE:HD12	1.79	0.47
1:D:195:ASP:HA	1:D:223:ALA:HB3	1.96	0.47
1:C:664:ILE:HB	1:C:667:PHE:HD2	1.79	0.47
1:D:97:ILE:HG13	1:D:111:ILE:HB	1.97	0.47
1:D:530:VAL:O	1:D:533:TYR:HB3	2.15	0.47
1:A:195:ASP:HA	1:A:223:ALA:HB3	1.96	0.47
1:A:135:TYR:CE2	1:A:137:TYR:HB3	2.50	0.47
1:D:62:PHE:CE2	1:D:92:LEU:HD12	2.51	0.46
1:C:530:VAL:O	1:C:533:TYR:HB3	2.15	0.46
1:B:135:TYR:CE2	1:B:137:TYR:HB3	2.50	0.46
1:C:619:ASN:CA	1:D:624:LEU:HD13	2.46	0.46
1:A:62:PHE:CE2	1:A:92:LEU:HD12	2.51	0.46
1:B:711:TYR:O	1:B:715:ARG:HG2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:530:VAL:O	1:B:533:TYR:HB3	2.15	0.46
1:D:135:TYR:CE2	1:D:137:TYR:HB3	2.50	0.46
1:C:97:ILE:HG13	1:C:111:ILE:HB	1.96	0.46
1:D:22:ALA:HB1	1:D:25:GLU:HB2	1.97	0.46
1:C:135:TYR:CE2	1:C:137:TYR:HB3	2.50	0.46
1:A:521:LEU:CD2	1:A:525:ILE:HB	2.45	0.46
1:A:530:VAL:O	1:A:533:TYR:HB3	2.15	0.46
1:A:97:ILE:HG13	1:A:111:ILE:HB	1.97	0.46
1:B:117:LEU:HD12	1:B:120:ALA:HB3	1.98	0.46
1:A:628:ARG:HB2	1:D:626:VAL:HG21	1.98	0.46
1:B:490:ASP:HB2	1:B:736:THR:HG23	1.98	0.46
1:C:29:PHE:O	1:C:33:MET:HG2	2.15	0.46
1:D:23:ASP:HB3	1:D:271:PRO:HB2	1.98	0.46
1:B:208:VAL:HG12	1:B:214:HIS:HB3	1.97	0.46
1:A:209:ILE:CD1	1:A:234:LYS:HB2	2.46	0.46
1:A:22:ALA:HB1	1:A:25:GLU:HB2	1.98	0.46
1:C:209:ILE:CD1	1:C:234:LYS:HB2	2.46	0.46
1:C:208:VAL:HG12	1:C:214:HIS:HB3	1.97	0.46
1:A:29:PHE:O	1:A:33:MET:HG2	2.15	0.46
1:B:209:ILE:CD1	1:B:234:LYS:HB2	2.46	0.45
1:A:208:VAL:HG12	1:A:214:HIS:HB3	1.97	0.45
1:D:208:VAL:HG12	1:D:214:HIS:HB3	1.97	0.45
1:B:23:ASP:HB3	1:B:271:PRO:HB2	1.98	0.45
1:C:22:ALA:HB1	1:C:25:GLU:HB2	1.98	0.45
1:A:628:ARG:HH12	1:D:627:GLU:N	2.14	0.45
1:B:683:VAL:HG11	1:B:689:GLY:HA2	1.99	0.45
2:D:833:ZK1:HAI	2:D:833:ZK1:HAOA	1.82	0.45
1:D:664:ILE:HB	1:D:667:PHE:HD2	1.79	0.45
1:C:117:LEU:HD12	1:C:120:ALA:HB3	1.98	0.45
1:C:62:PHE:CE2	1:C:92:LEU:HD12	2.51	0.45
1:C:23:ASP:HB3	1:C:271:PRO:HB2	1.98	0.45
1:A:117:LEU:HD12	1:A:120:ALA:HB3	1.98	0.45
1:C:683:VAL:HG11	1:C:689:GLY:HA2	1.99	0.45
1:B:536:VAL:HG21	1:B:605:TRP:CE3	2.51	0.45
1:D:490:ASP:HB2	1:D:736:THR:HG23	1.98	0.45
1:B:601:VAL:HG23	1:C:806:ALA:CB	2.46	0.45
1:D:521:LEU:CD2	1:D:525:ILE:HB	2.46	0.45
1:A:519:ASP:HB2	1:A:520:PRO:HD3	1.98	0.45
1:C:521:LEU:CD2	1:C:525:ILE:HB	2.47	0.45
1:A:515:PHE:CD1	1:A:518:LEU:HD12	2.52	0.45
1:B:22:ALA:HB1	1:B:25:GLU:HB2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:VAL:HG22	1:B:165:VAL:O	2.17	0.45
1:B:521:LEU:HD22	1:B:526:TRP:CE2	2.52	0.45
1:B:525:ILE:HD13	1:C:792:VAL:HG21	1.99	0.45
1:B:519:ASP:HB2	1:B:520:PRO:HD3	1.98	0.44
1:C:261:ARG:O	1:C:265:LEU:HG	2.17	0.44
1:A:62:PHE:CE2	1:A:88:PHE:HB3	2.52	0.44
1:C:62:PHE:CE2	1:C:88:PHE:HB3	2.52	0.44
1:D:683:VAL:HG11	1:D:689:GLY:HA2	1.99	0.44
1:D:125:ILE:HG23	1:D:130:TRP:HB2	1.99	0.44
1:D:504:ILE:CG2	1:D:633:ILE:HD12	2.48	0.44
1:A:261:ARG:O	1:A:265:LEU:HG	2.17	0.44
1:A:23:ASP:HB3	1:A:271:PRO:HB2	1.98	0.44
1:D:261:ARG:O	1:D:265:LEU:HG	2.17	0.44
1:D:521:LEU:HD22	1:D:526:TRP:CE2	2.52	0.44
1:A:96:PHE:CE2	1:A:98:THR:HB	2.53	0.44
1:D:62:PHE:CE2	1:D:88:PHE:HB3	2.52	0.44
1:B:62:PHE:CE2	1:B:92:LEU:HD12	2.51	0.44
1:B:261:ARG:O	1:B:265:LEU:HG	2.17	0.44
1:D:519:ASP:HB2	1:D:520:PRO:HD3	1.98	0.44
1:D:620:LEU:HA	1:D:623:PHE:HD2	1.82	0.44
1:B:515:PHE:CD1	1:B:518:LEU:HD12	2.52	0.44
1:D:117:LEU:HD12	1:D:120:ALA:HB3	1.98	0.44
1:B:62:PHE:CE2	1:B:88:PHE:HB3	2.52	0.44
1:A:683:VAL:HG11	1:A:689:GLY:HA2	1.99	0.44
1:D:165:VAL:HG22	1:D:165:VAL:O	2.17	0.44
1:C:287:ASP:O	1:C:291:VAL:HG23	2.18	0.44
1:C:49:ASN:C	1:C:50:LEU:HD12	2.38	0.44
1:D:49:ASN:C	1:D:50:LEU:HD12	2.38	0.44
1:D:209:ILE:CD1	1:D:234:LYS:HB2	2.47	0.44
1:B:96:PHE:CE2	1:B:98:THR:HB	2.53	0.44
1:C:519:ASP:HB2	1:C:520:PRO:HD3	1.98	0.44
1:C:125:ILE:HG23	1:C:130:TRP:HB2	1.99	0.44
2:C:833:ZK1:HAOA	2:C:833:ZK1:HAI	1.82	0.44
1:A:125:ILE:HG23	1:A:130:TRP:HB2	1.99	0.44
1:B:130:TRP:CH2	1:B:191:ARG:HB3	2.53	0.44
1:C:96:PHE:CE2	1:C:98:THR:HB	2.53	0.44
1:D:96:PHE:CE2	1:D:98:THR:HB	2.53	0.44
1:C:130:TRP:CH2	1:C:191:ARG:HB3	2.53	0.44
1:A:130:TRP:CH2	1:A:191:ARG:HB3	2.53	0.44
1:C:521:LEU:HB3	1:C:526:TRP:CE2	2.53	0.43
1:B:521:LEU:CD2	1:B:525:ILE:HB	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:515:PHE:CD1	1:D:518:LEU:HD12	2.52	0.43
1:D:488:VAL:HG23	1:D:489:ILE:HG23	2.00	0.43
1:B:504:ILE:CG2	1:B:633:ILE:HD12	2.48	0.43
1:A:130:TRP:CE2	1:A:191:ARG:HD3	2.53	0.43
1:A:49:ASN:C	1:A:50:LEU:HD12	2.38	0.43
1:D:287:ASP:O	1:D:291:VAL:HG23	2.18	0.43
1:A:521:LEU:HD22	1:A:526:TRP:CE2	2.52	0.43
1:B:130:TRP:CE2	1:B:191:ARG:HD3	2.54	0.43
1:C:290:GLN:HG2	1:C:338:VAL:HG21	2.00	0.43
1:A:165:VAL:HG22	1:A:165:VAL:O	2.18	0.43
1:C:515:PHE:CD1	1:C:518:LEU:HD12	2.52	0.43
1:D:479:LEU:HD12	2:D:833:ZK1:OAA	2.19	0.43
1:D:130:TRP:CE2	1:D:191:ARG:HD3	2.54	0.43
1:A:287:ASP:O	1:A:291:VAL:HG23	2.18	0.43
1:A:290:GLN:HG2	1:A:338:VAL:HG21	2.00	0.43
1:A:628:ARG:NH1	1:D:623:PHE:HA	2.33	0.43
1:A:209:ILE:HD11	1:A:235:ILE:HG23	2.01	0.43
1:A:97:ILE:N	1:A:97:ILE:HD12	2.34	0.43
1:B:360:ILE:HD11	1:B:374:TRP:HB2	2.01	0.43
1:B:287:ASP:O	1:B:291:VAL:HG23	2.18	0.43
1:A:488:VAL:HG23	1:A:489:ILE:HG23	2.00	0.43
1:B:488:VAL:HG23	1:B:489:ILE:HG23	2.00	0.43
1:D:130:TRP:CH2	1:D:191:ARG:HB3	2.53	0.43
1:A:628:ARG:NH2	1:D:623:PHE:CD1	2.84	0.43
1:C:130:TRP:CE2	1:C:191:ARG:HD3	2.53	0.43
1:B:233:LEU:HD23	1:B:236:GLN:OE1	2.19	0.43
1:B:49:ASN:C	1:B:50:LEU:HD12	2.38	0.43
1:C:97:ILE:N	1:C:97:ILE:HD12	2.34	0.43
1:B:125:ILE:HG23	1:B:130:TRP:HB2	1.99	0.43
1:C:809:VAL:O	1:C:813:GLU:HG3	2.19	0.43
2:A:833:ZK1:HAI	2:A:833:ZK1:HAOA	1.82	0.43
1:C:488:VAL:HG23	1:C:489:ILE:HG23	2.00	0.42
1:A:233:LEU:HD23	1:A:236:GLN:OE1	2.19	0.42
1:B:209:ILE:HD11	1:B:235:ILE:HG23	2.01	0.42
1:D:809:VAL:O	1:D:813:GLU:HG3	2.19	0.42
1:C:165:VAL:O	1:C:165:VAL:HG22	2.18	0.42
1:D:97:ILE:HD12	1:D:97:ILE:N	2.34	0.42
1:C:360:ILE:HD11	1:C:374:TRP:HB2	2.01	0.42
1:D:521:LEU:HB3	1:D:526:TRP:CE2	2.54	0.42
1:B:97:ILE:N	1:B:97:ILE:HD12	2.34	0.42
1:B:809:VAL:O	1:B:813:GLU:HG3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:521:LEU:HB3	1:A:526:TRP:CE2	2.54	0.42
1:C:209:ILE:HD11	1:C:235:ILE:HG23	2.01	0.42
1:D:360:ILE:HD11	1:D:374:TRP:HB2	2.01	0.42
1:C:805:LEU:O	1:C:809:VAL:HG23	2.20	0.42
1:D:708:MET:O	1:D:712:ILE:HG12	2.20	0.42
1:B:521:LEU:HB3	1:B:526:TRP:CE2	2.55	0.42
1:B:608:PHE:HZ	1:C:796:PHE:CZ	2.38	0.42
1:C:233:LEU:HD23	1:C:236:GLN:OE1	2.19	0.42
1:C:525:ILE:HD13	1:D:792:VAL:HG21	2.00	0.42
1:C:533:TYR:CZ	1:C:583:ALA:HB1	2.54	0.42
1:A:708:MET:O	1:A:712:ILE:HG12	2.20	0.42
1:A:270:TYR:HA	1:A:271:PRO:HD2	1.87	0.42
1:A:360:ILE:HD11	1:A:374:TRP:HB2	2.01	0.42
1:D:209:ILE:HD11	1:D:235:ILE:HG23	2.01	0.41
1:A:744:THR:HB	1:A:745:PRO:HD3	2.02	0.41
1:A:795:VAL:HG12	1:D:608:PHE:CD1	2.54	0.41
1:D:805:LEU:O	1:D:809:VAL:HG23	2.20	0.41
1:A:809:VAL:O	1:A:813:GLU:HG3	2.19	0.41
1:A:795:VAL:HG12	1:D:608:PHE:HD1	1.85	0.41
1:C:98:THR:HA	1:C:99:PRO:HD3	1.80	0.41
1:A:533:TYR:CZ	1:A:583:ALA:HB1	2.55	0.41
1:B:232:LEU:O	1:B:236:GLN:HB2	2.20	0.41
1:B:805:LEU:O	1:B:809:VAL:HG23	2.21	0.41
1:C:232:LEU:O	1:C:236:GLN:HB2	2.20	0.41
1:D:233:LEU:HD23	1:D:236:GLN:OE1	2.19	0.41
1:C:708:MET:O	1:C:712:ILE:HG12	2.20	0.41
1:A:678:GLU:HA	1:A:679:PRO:C	2.41	0.41
1:A:805:LEU:O	1:A:809:VAL:HG23	2.21	0.41
1:D:467:LEU:HD22	1:D:737:PRO:HD3	2.03	0.41
1:D:744:THR:HB	1:D:745:PRO:HD3	2.03	0.41
1:C:624:LEU:O	1:C:628:ARG:HG3	2.21	0.41
1:B:608:PHE:HD1	1:C:795:VAL:HG12	1.85	0.41
1:C:803:LEU:O	1:C:807:MET:HG2	2.21	0.41
1:D:232:LEU:O	1:D:236:GLN:HB2	2.20	0.41
1:A:467:LEU:HD22	1:A:737:PRO:HD3	2.02	0.41
1:D:114:ARG:HA	1:D:115:PRO:HD3	1.92	0.41
1:A:624:LEU:O	1:A:628:ARG:HG3	2.20	0.41
1:A:321:VAL:HA	1:A:322:PRO:HD3	1.86	0.41
1:C:467:LEU:HD22	1:C:737:PRO:HD3	2.03	0.41
1:A:792:VAL:HG21	1:D:525:ILE:HD13	2.03	0.41
1:D:235:ILE:HG13	1:D:235:ILE:H	1.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:270:TYR:HA	1:D:271:PRO:HD2	1.87	0.41
1:A:519:ASP:N	1:A:520:PRO:CD	2.84	0.41
1:A:232:LEU:O	1:A:236:GLN:HB2	2.21	0.41
1:A:192:VAL:HB	1:A:220:TYR:CD1	2.56	0.41
1:D:288:ALA:O	1:D:292:MET:HG3	2.21	0.41
1:A:803:LEU:O	1:A:807:MET:HG2	2.21	0.41
1:D:127:TYR:CE1	1:D:382:LEU:HD21	2.56	0.41
1:A:375:SER:HB3	1:A:378:ASP:HB2	2.03	0.41
1:B:803:LEU:O	1:B:807:MET:HG2	2.21	0.41
1:D:192:VAL:HB	1:D:220:TYR:CD1	2.56	0.41
1:B:127:TYR:CE1	1:B:382:LEU:HD21	2.56	0.41
1:D:519:ASP:N	1:D:520:PRO:CD	2.84	0.40
1:C:519:ASP:N	1:C:520:PRO:CD	2.84	0.40
1:B:467:LEU:HD22	1:B:737:PRO:HD3	2.03	0.40
1:B:192:VAL:HB	1:B:220:TYR:CD1	2.56	0.40
1:B:288:ALA:O	1:B:292:MET:HG3	2.21	0.40
1:B:114:ARG:HA	1:B:115:PRO:HD3	1.92	0.40
1:B:533:TYR:HA	1:B:605:TRP:CH2	2.56	0.40
1:C:325:GLN:O	1:C:329:ILE:HG13	2.22	0.40
1:B:708:MET:O	1:B:712:ILE:HG12	2.20	0.40
1:B:758:LEU:HD23	1:B:758:LEU:O	2.22	0.40
1:A:758:LEU:HD23	1:A:758:LEU:O	2.21	0.40
1:D:404:PRO:HB3	1:D:711:TYR:CE1	2.56	0.40
1:B:231:ASP:HB3	1:B:234:LYS:HE2	2.04	0.40
1:C:321:VAL:HA	1:C:322:PRO:HD3	1.86	0.40
1:B:519:ASP:N	1:B:520:PRO:CD	2.84	0.40
1:D:803:LEU:O	1:D:807:MET:HG2	2.21	0.40
1:D:143:LEU:HD22	1:D:143:LEU:N	2.37	0.40
1:D:266:GLU:HG2	1:D:268:LYS:H	1.86	0.40
1:D:520:PRO:HA	1:D:623:PHE:HE2	1.87	0.40
1:A:628:ARG:CZ	1:D:623:PHE:HA	2.52	0.40
1:D:299:LEU:CD1	1:D:332:ALA:HB2	2.51	0.40
1:D:98:THR:HA	1:D:99:PRO:HD3	1.80	0.40
1:C:477:ALA:HB1	1:C:478:PRO:CD	2.52	0.40
1:B:404:PRO:HB3	1:B:711:TYR:CE1	2.56	0.40
1:C:192:VAL:HB	1:C:220:TYR:CD1	2.56	0.40
1:D:325:GLN:O	1:D:329:ILE:HG13	2.22	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	732/826 (89%)	683 (93%)	44 (6%)	5 (1%)	26	71
1	B	732/826 (89%)	683 (93%)	44 (6%)	5 (1%)	26	71
1	C	732/826 (89%)	682 (93%)	45 (6%)	5 (1%)	26	71
1	D	731/826 (88%)	683 (93%)	44 (6%)	4 (0%)	34	77
All	All	2927/3304 (89%)	2731 (93%)	177 (6%)	19 (1%)	34	74

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	172	LYS
1	B	172	LYS
1	C	172	LYS
1	D	172	LYS
1	A	513	GLY
1	B	513	GLY
1	C	513	GLY
1	A	597	SER
1	B	597	SER
1	C	597	SER
1	D	597	SER
1	A	743	GLY
1	C	743	GLY
1	B	743	GLY
1	D	743	GLY
1	A	451	GLY
1	B	451	GLY
1	C	451	GLY
1	D	451	GLY

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	527/705 (75%)	524 (99%)	3 (1%)	90	95
1	B	527/705 (75%)	525 (100%)	2 (0%)	93	96
1	C	527/705 (75%)	523 (99%)	4 (1%)	86	94
1	D	526/705 (75%)	524 (100%)	2 (0%)	93	96
All	All	2107/2820 (75%)	2096 (100%)	11 (0%)	92	96

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	299	LEU
1	A	394	THR
1	B	10	ASN
1	B	299	LEU
1	C	10	ASN
1	C	299	LEU
1	C	394	THR
1	C	628	ARG
1	D	10	ASN
1	D	299	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	619	ASN
1	B	83	ASN
1	B	619	ASN
1	D	83	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

4 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
2	ZK1	A	833	-	28,29,29	3.02	13 (46%)	35,45,45	1.53	3 (8%)
2	ZK1	B	833	-	28,29,29	3.02	13 (46%)	35,45,45	1.52	3 (8%)
2	ZK1	C	833	-	28,29,29	3.00	13 (46%)	35,45,45	1.53	3 (8%)
2	ZK1	D	833	-	28,29,29	3.02	13 (46%)	35,45,45	1.54	3 (8%)

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
2	ZK1	A	833	-	-	0/13/23/23	0/3/3/3
2	ZK1	B	833	-	-	0/13/23/23	0/3/3/3
2	ZK1	C	833	-	-	0/13/23/23	0/3/3/3
2	ZK1	D	833	-	-	0/13/23/23	0/3/3/3

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	833	ZK1	CAN-NAX	2.04	1.49	1.46
2	B	833	ZK1	CAN-NAX	2.06	1.49	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	833	ZK1	CAN-NAX	2.06	1.49	1.46
2	B	833	ZK1	PBA-CAO	2.09	1.86	1.81
2	A	833	ZK1	PBA-CAO	2.13	1.86	1.81
2	D	833	ZK1	CAN-NAX	2.13	1.50	1.46
2	D	833	ZK1	CAR-NAX	2.15	1.46	1.41
2	C	833	ZK1	PBA-CAO	2.15	1.86	1.81
2	C	833	ZK1	CAR-NAX	2.15	1.46	1.41
2	C	833	ZK1	CAM-NAX	2.15	1.50	1.46
2	A	833	ZK1	CAR-NAX	2.15	1.46	1.41
2	B	833	ZK1	CAR-NAX	2.16	1.46	1.41
2	D	833	ZK1	PBA-CAO	2.17	1.86	1.81
2	C	833	ZK1	CAV-NAP	2.19	1.38	1.35
2	D	833	ZK1	CAM-NAX	2.23	1.50	1.46
2	B	833	ZK1	CAM-NAX	2.25	1.50	1.46
2	B	833	ZK1	CAV-NAP	2.27	1.38	1.35
2	A	833	ZK1	CAM-NAX	2.27	1.50	1.46
2	D	833	ZK1	CAV-NAP	2.29	1.38	1.35
2	A	833	ZK1	CAV-NAP	2.32	1.39	1.35
2	A	833	ZK1	CAI-CAR	3.36	1.42	1.37
2	C	833	ZK1	CAI-CAR	3.39	1.42	1.37
2	D	833	ZK1	CAI-CAR	3.43	1.42	1.37
2	B	833	ZK1	CAI-CAR	3.49	1.42	1.37
2	B	833	ZK1	PBA-OAD	4.21	1.65	1.54
2	D	833	ZK1	PBA-OAD	4.22	1.65	1.54
2	C	833	ZK1	PBA-OAD	4.24	1.65	1.54
2	A	833	ZK1	PBA-OAD	4.25	1.65	1.54
2	C	833	ZK1	PBA-OAE	4.33	1.65	1.54
2	A	833	ZK1	PBA-OAE	4.34	1.65	1.54
2	B	833	ZK1	PBA-OAE	4.36	1.65	1.54
2	D	833	ZK1	PBA-OAE	4.39	1.65	1.54
2	C	833	ZK1	CAJ-CAS	4.53	1.44	1.37
2	B	833	ZK1	CAJ-CAS	4.57	1.44	1.37
2	D	833	ZK1	CAJ-CAS	4.58	1.44	1.37
2	A	833	ZK1	CAJ-CAS	4.59	1.44	1.37
2	C	833	ZK1	CAW-NAY	4.68	1.46	1.40
2	D	833	ZK1	CAW-NAY	4.79	1.46	1.40
2	A	833	ZK1	CAW-NAY	4.84	1.47	1.40
2	B	833	ZK1	CAW-NAY	4.85	1.47	1.40
2	B	833	ZK1	CAT-NAP	6.12	1.44	1.33
2	D	833	ZK1	CAT-NAP	6.13	1.44	1.33
2	C	833	ZK1	CAT-NAP	6.13	1.44	1.33
2	A	833	ZK1	CAT-NAP	6.18	1.44	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	833	ZK1	OAB-CAU	6.47	1.40	1.24
2	D	833	ZK1	OAB-CAU	6.48	1.40	1.24
2	C	833	ZK1	OAB-CAU	6.49	1.40	1.24
2	B	833	ZK1	OAB-CAU	6.49	1.40	1.24
2	C	833	ZK1	OAA-CAT	6.56	1.41	1.24
2	B	833	ZK1	OAA-CAT	6.56	1.41	1.24
2	D	833	ZK1	OAA-CAT	6.59	1.41	1.24
2	A	833	ZK1	OAA-CAT	6.62	1.41	1.24

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	833	ZK1	CAI-CAR-NAX	-2.62	118.92	122.62
2	C	833	ZK1	CAI-CAR-NAX	-2.55	119.02	122.62
2	D	833	ZK1	CAI-CAR-NAX	-2.55	119.02	122.62
2	B	833	ZK1	CAI-CAR-NAX	-2.55	119.02	122.62
2	B	833	ZK1	CAV-CAW-NAY	3.36	120.10	117.58
2	C	833	ZK1	CAV-CAW-NAY	3.42	120.14	117.58
2	A	833	ZK1	CAV-CAW-NAY	3.47	120.18	117.58
2	D	833	ZK1	CAV-CAW-NAY	3.61	120.28	117.58
2	A	833	ZK1	CAN-NAX-CAM	4.78	121.44	111.54
2	D	833	ZK1	CAN-NAX-CAM	4.79	121.45	111.54
2	B	833	ZK1	CAN-NAX-CAM	4.80	121.48	111.54
2	C	833	ZK1	CAN-NAX-CAM	4.84	121.57	111.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	833	ZK1	1	0
2	C	833	ZK1	1	0
2	D	833	ZK1	2	0

## 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.