



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:14 PM GMT

PDB ID : 1EKJ
Title : THE X-RAY CRYSTALLOGRAPHIC STRUCTURE OF BETA CARBONIC ANHYDRASE FROM THE C3 DICOT PISUM SATIVUM
Authors : Kimber, M.S.; Pai, E.F.
Deposited on : 2000-03-08
Resolution : 1.93 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
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) : trunk26865

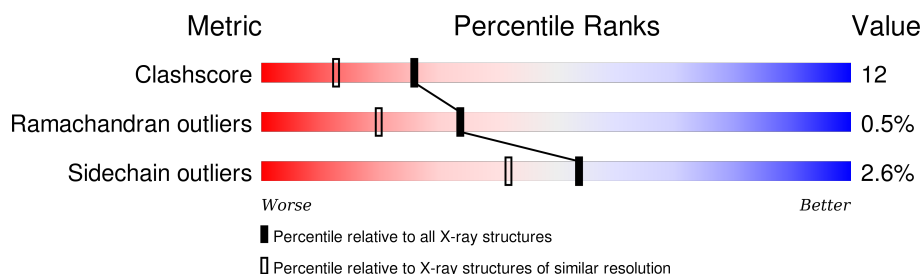
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	3095 (1.96-1.92)
Ramachandran outliers	100387	3062 (1.96-1.92)
Sidechain outliers	100360	3062 (1.96-1.92)


The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	221	
1	B	221	
1	C	221	
1	D	221	
1	E	221	
1	F	221	
1	G	221	

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Mol	Chain	Length	Quality of chain
1	H	221	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	3001	-	-	X	-
2	ACT	C	3007	-	-	X	-
2	ACT	E	3005	-	-	X	-
2	ACT	F	3008	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14070 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called BETA-CARBONIC ANHYDRASE.

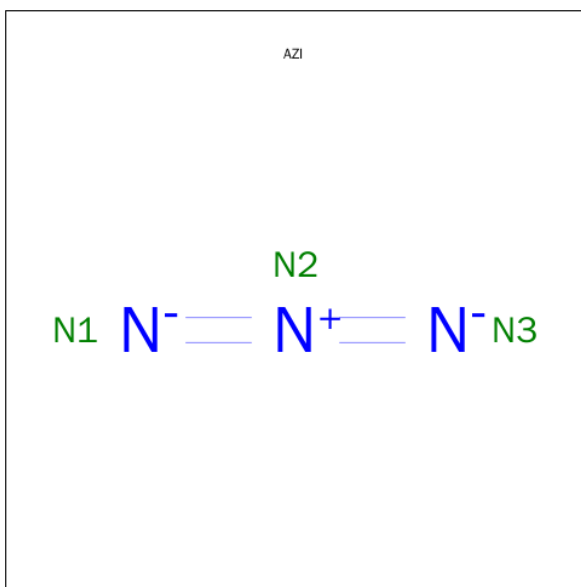
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	0	0
			1618	1051	264	297	6			
1	B	213	Total	C	N	O	S	0	0	0
			1640	1065	268	301	6			
1	C	217	Total	C	N	O	S	0	1	0
			1672	1083	273	309	7			
1	D	211	Total	C	N	O	S	0	1	0
			1630	1057	266	300	7			
1	E	212	Total	C	N	O	S	0	0	0
			1633	1060	267	300	6			
1	F	214	Total	C	N	O	S	0	0	0
			1648	1071	269	302	6			
1	G	220	Total	C	N	O	S	0	1	0
			1691	1093	276	315	7			
1	H	211	Total	C	N	O	S	0	1	0
			1630	1057	266	300	7			

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	G	1	Total	C	O	0	0
			4	2	2		
2	G	1	Total	C	O	0	0
			4	2	2		
2	H	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is AZIDE ION (three-letter code: AZI) (formula: N₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total N 3 3	0	0
3	D	1	Total N 3 3	0	0
3	E	1	Total N 3 3	0	0
3	H	1	Total N 3 3	0	0

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	G	1	Total Zn 1 1	0	0
4	D	1	Total Zn 1 1	0	0
4	E	1	Total Zn 1 1	0	0
4	H	1	Total Zn 1 1	0	0
4	B	1	Total Zn 1 1	0	0
4	C	1	Total Zn 1 1	0	0
4	A	1	Total Zn 1 1	0	0
4	F	1	Total Zn 1 1	0	0

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	1	Total	Cl	0	0
			1	1		
5	D	1	Total	Cl	0	0
			1	1		
5	E	1	Total	Cl	0	0
			1	1		
5	H	1	Total	Cl	0	0
			1	1		
5	B	1	Total	Cl	0	0
			1	1		
5	C	1	Total	Cl	0	0
			1	1		
5	A	1	Total	Cl	0	0
			1	1		
5	F	1	Total	Cl	0	0
			1	1		

- Molecule 6 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

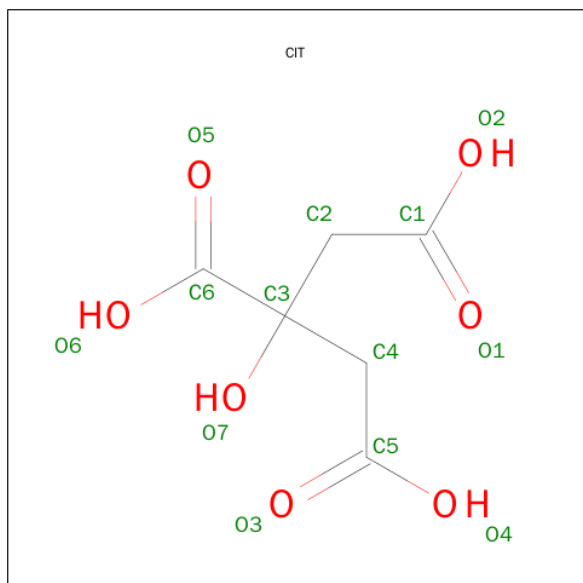
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	G	1	Total	Cu	0	0
			1	1		
6	B	1	Total	Cu	0	0
			1	1		
6	C	1	Total	Cu	0	0
			1	1		
6	E	1	Total	Cu	0	0
			1	1		

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	D	1	Total	C	O	0	0
			4	2	2		
7	E	1	Total	C	O	0	0
			4	2	2		
7	H	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 9 is water.

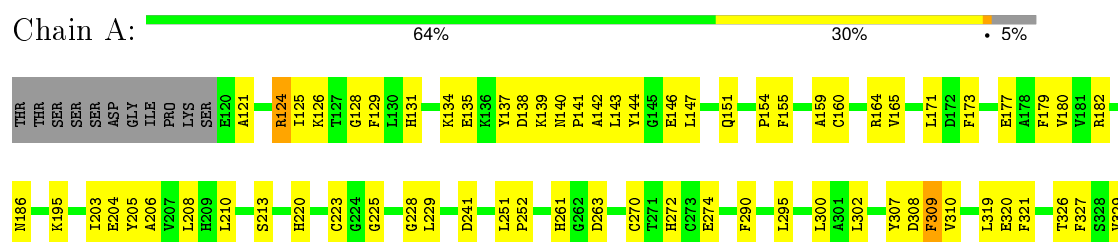
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	50	Total	O	0	0
			50	50		
9	B	43	Total	O	0	0
			43	43		
9	C	138	Total	O	0	0
			138	138		
9	D	131	Total	O	0	0
			131	131		
9	E	71	Total	O	0	0
			71	71		
9	F	84	Total	O	0	0
			84	84		
9	G	150	Total	O	0	0
			150	150		
9	H	144	Total	O	0	0
			144	144		

3 Residue-property plots [i](#)

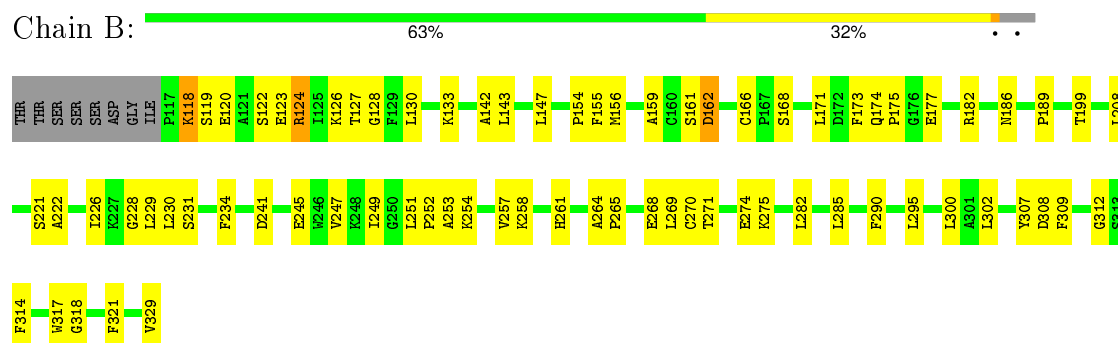
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

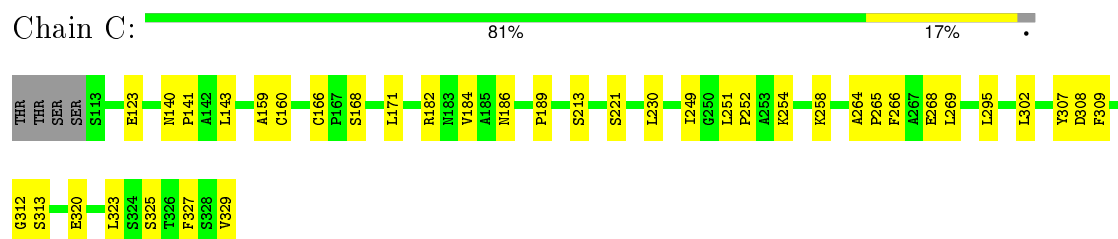
• Molecule 1: BETA-CARBONIC ANHYDRASE



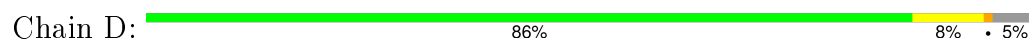
• Molecule 1: BETA-CARBONIC ANHYDRASE

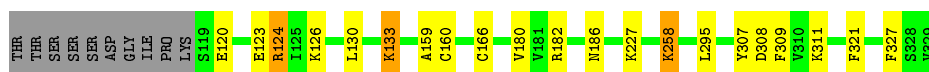


• Molecule 1: BETA-CARBONIC ANHYDRASE



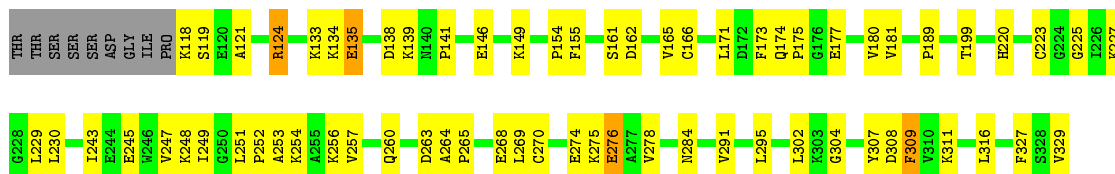
• Molecule 1: BETA-CARBONIC ANHYDRASE





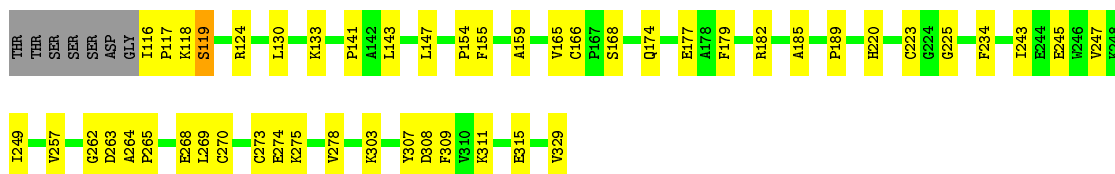
• Molecule 1: BETA-CARBONIC ANHYDRASE

Chain E: 66% 29% 5%



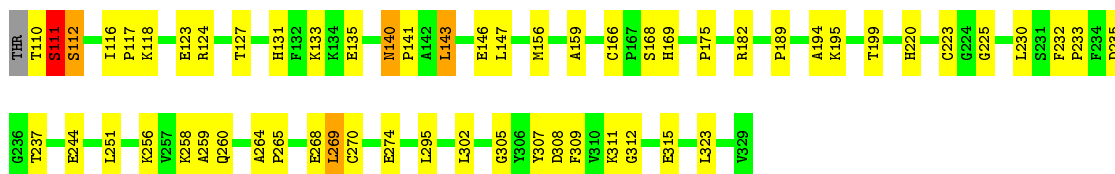
• Molecule 1: BETA-CARBONIC ANHYDRASE

Chain F: 75% 22% 3%



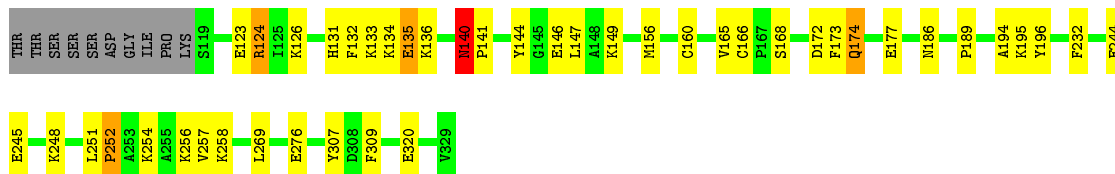
• Molecule 1: BETA-CARBONIC ANHYDRASE

Chain G: 73% 24% 3%



• Molecule 1: BETA-CARBONIC ANHYDRASE

Chain H: 76% 18% 6%



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	136.91Å 143.32Å 202.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.93	Depositor
% Data completeness (in resolution range)	87.5 (40.00-1.93)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.5	Depositor
R, R_{free}	0.229 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	14070	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AZI, ZN, CL, CIT, EDO, ACT, CU

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	0.46	0/1662	0.64	0/2250
1	B	0.43	0/1685	0.63	1/2280 (0.0%)
1	C	0.62	0/1717	0.72	0/2325
1	D	0.75	0/1674	0.80	0/2267
1	E	0.48	0/1677	0.66	0/2269
1	F	0.51	0/1693	0.66	1/2292 (0.0%)
1	G	0.67	0/1736	0.78	2/2351 (0.1%)
1	H	0.75	0/1674	0.77	1/2267 (0.0%)
All	All	0.60	0/13518	0.71	5/18301 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	156	MET	N-CA-C	-5.12	97.16	111.00
1	F	179	PHE	N-CA-C	-5.07	97.32	111.00
1	G	112	SER	N-CA-C	5.06	124.67	111.00
1	B	156	MET	N-CA-C	-5.03	97.42	111.00
1	H	156	MET	N-CA-C	-5.01	97.47	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1618	0	1593	57	0
1	B	1640	0	1619	59	0
1	C	1672	0	1646	31	0
1	D	1630	0	1603	16	0
1	E	1633	0	1611	53	0
1	F	1648	0	1629	37	0
1	G	1691	0	1664	47	0
1	H	1630	0	1604	44	0
2	A	8	0	6	3	0
2	C	8	0	6	2	0
2	E	4	0	3	2	0
2	F	4	0	3	2	0
2	G	8	0	6	2	0
2	H	4	0	3	0	0
3	A	3	0	0	0	0
3	D	3	0	0	0	0
3	E	3	0	0	0	0
3	H	3	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	E	1	0	0	0	0
6	G	1	0	0	0	0
7	A	4	0	6	0	0
7	D	4	0	6	0	0
7	E	4	0	6	1	0
7	H	4	0	6	0	0
8	D	13	0	4	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	50	0	0	0	0
9	B	43	0	0	4	0
9	C	138	0	0	3	0
9	D	131	0	0	2	0
9	E	71	0	0	2	0
9	F	84	0	0	4	0
9	G	150	0	0	12	0
9	H	144	0	0	4	0
All	All	14070	0	13024	325	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 12.

All (325) 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:G:111:SER:HA	1:G:117:PRO:HA	1.29	1.10
1:H:134:LYS:HG3	1:H:135:GLU:H	1.23	1.03
1:G:110:THR:O	1:G:111:SER:HB2	1.60	0.99
1:F:166:CYS:HB2	9:F:1130:HOH:O	1.63	0.99
1:H:146:GLU:O	1:H:149:LYS:HG2	1.66	0.94
1:E:166:CYS:HB2	9:E:1168:HOH:O	1.72	0.89
1:E:225:GLY:H	2:E:3005:ACT:H1	1.40	0.85
1:A:225:GLY:H	2:A:3001:ACT:H1	1.40	0.85
1:G:166[A]:CYS:SG	9:G:1185:HOH:O	2.35	0.84
1:B:142:ALA:HB3	9:B:1175:HOH:O	1.80	0.82
1:E:265:PRO:HD2	1:E:268:GLU:OE2	1.79	0.82
1:A:225:GLY:N	2:A:3001:ACT:H1	1.96	0.80
1:F:264:ALA:HB3	1:F:269:LEU:HD11	1.62	0.80
1:G:308:ASP:OD1	1:G:311:LYS:HD3	1.84	0.77
1:H:124:ARG:O	1:H:124:ARG:HD3	1.83	0.77
1:F:270:CYS:O	1:F:274:GLU:HG3	1.86	0.76
1:A:138:ASP:O	1:A:141:PRO:HD3	1.85	0.76
1:B:154:PRO:HG2	1:B:155:PHE:CD1	2.22	0.75
1:G:111:SER:HA	1:G:117:PRO:CA	2.12	0.74
1:H:124:ARG:C	1:H:124:ARG:HD3	2.08	0.74
1:E:295:LEU:HD21	1:E:302:LEU:HD13	1.69	0.74
1:C:166[B]:CYS:SG	9:C:1064:HOH:O	2.46	0.73
1:E:118:LYS:N	1:F:303:LYS:HZ1	1.85	0.73
1:D:308:ASP:OD2	1:D:311:LYS:HD3	1.89	0.72
1:E:189:PRO:HB3	1:E:249:ILE:HD13	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:TYR:CE2	2:A:3003:ACT:H3	2.25	0.71
1:B:166:CYS:HB2	9:B:1221:HOH:O	1.90	0.70
1:A:319:LEU:HD12	1:A:320:GLU:N	2.05	0.70
1:B:329:VAL:HG11	1:D:130:LEU:HD21	1.74	0.70
1:E:308:ASP:OD2	1:E:311:LYS:HD3	1.93	0.69
1:B:230:LEU:HD13	1:B:269:LEU:HD12	1.74	0.68
1:A:160:CYS:HB2	1:A:186:ASN:HD22	1.58	0.68
1:B:221:SER:HB3	1:B:308:ASP:HA	1.76	0.68
1:C:312:GLY:O	1:D:133:LYS:HE2	1.94	0.68
1:C:295:LEU:HD21	1:C:302:LEU:HD13	1.75	0.67
1:D:124:ARG:O	1:D:124:ARG:HD3	1.93	0.67
1:G:112:SER:HB3	9:G:1924:HOH:O	1.94	0.67
1:A:124:ARG:NH2	1:A:128:GLY:HA3	2.10	0.67
1:B:230:LEU:HA	1:B:254:LYS:HE3	1.76	0.66
1:B:270:CYS:O	1:B:274:GLU:HG3	1.96	0.66
1:B:182:ARG:HD2	9:B:1085:HOH:O	1.95	0.65
1:H:134:LYS:HG3	1:H:135:GLU:N	2.04	0.65
1:B:174:GLN:O	1:B:177:GLU:HB2	1.96	0.65
1:D:124:ARG:C	1:D:124:ARG:HD3	2.17	0.64
1:D:258:LYS:HE3	9:D:1386:HOH:O	1.97	0.64
1:E:118:LYS:HB3	1:E:121:ALA:HB3	1.79	0.64
1:B:282:LEU:HD23	1:B:285:LEU:HD12	1.78	0.64
1:G:140:ASN:HB2	9:G:1413:HOH:O	1.97	0.64
1:F:308:ASP:OD2	1:F:311:LYS:HD3	1.98	0.64
1:B:120:GLU:H	1:B:120:GLU:CD	2.00	0.64
1:E:307:TYR:CZ	1:E:309:PHE:HB3	2.32	0.64
1:E:245:GLU:O	1:E:248:LYS:HG2	1.98	0.63
1:A:121:ALA:O	1:A:125:ILE:HG13	1.99	0.63
1:H:166[B]:CYS:SG	9:H:1094:HOH:O	2.35	0.63
1:A:203:ILE:HG22	1:A:290:PHE:HE1	1.63	0.63
1:A:307:TYR:CZ	1:A:309:PHE:HB3	2.33	0.63
1:C:307:TYR:CZ	1:C:309:PHE:HB3	2.35	0.62
1:H:146:GLU:O	1:H:149:LYS:CG	2.46	0.62
1:A:164:ARG:HG2	1:B:147:LEU:HD22	1.82	0.61
1:C:189:PRO:HB3	1:C:249:ILE:CD1	2.31	0.61
1:F:116:ILE:N	1:F:117:PRO:HD2	2.16	0.61
1:A:295:LEU:HD21	1:A:302:LEU:HD13	1.81	0.61
1:F:225:GLY:H	2:F:3008:ACT:H1	1.66	0.61
1:E:154:PRO:HG2	1:E:155:PHE:CD1	2.36	0.61
1:H:146:GLU:OE1	1:H:149:LYS:HG3	2.01	0.60
1:G:124:ARG:NE	1:H:177:GLU:OE2	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:LYS:C	1:B:118:LYS:HD2	2.22	0.60
1:E:134:LYS:HG3	1:E:135:GLU:N	2.17	0.60
1:E:146:GLU:HA	1:E:149:LYS:HD3	1.83	0.59
1:A:140:ASN:HB3	1:A:143:LEU:CB	2.32	0.59
1:G:189:PRO:HD2	1:G:199:THR:HG21	1.83	0.59
1:G:110:THR:O	1:G:111:SER:CB	2.45	0.59
1:B:307:TYR:CE2	1:B:309:PHE:HB3	2.38	0.58
1:G:166[B]:CYS:SG	9:G:1185:HOH:O	2.52	0.58
1:C:254:LYS:HE2	1:C:258:LYS:HE3	1.84	0.58
1:C:168:SER:OG	1:D:166[A]:CYS:SG	2.58	0.58
1:G:312:GLY:O	1:H:133:LYS:HE3	2.03	0.58
1:G:168:SER:OG	1:H:166[A]:CYS:SG	2.49	0.58
1:G:307:TYR:CE2	1:G:309:PHE:HB3	2.39	0.58
1:F:264:ALA:HB3	1:F:269:LEU:CD1	2.31	0.58
1:E:189:PRO:HB3	1:E:249:ILE:CD1	2.34	0.57
1:B:230:LEU:HD23	1:B:254:LYS:HG3	1.85	0.57
1:B:118:LYS:HD3	1:B:123:GLU:HB2	1.86	0.57
1:F:133:LYS:HD3	1:F:133:LYS:O	2.03	0.57
1:C:189:PRO:HB3	1:C:249:ILE:HD13	1.85	0.57
1:E:124:ARG:HD2	1:E:124:ARG:O	2.04	0.57
1:A:171:LEU:HD12	1:A:173:PHE:HE1	1.69	0.57
1:A:124:ARG:HH21	1:A:128:GLY:HA3	1.69	0.57
1:B:171:LEU:HD12	1:B:173:PHE:HE1	1.70	0.57
1:C:327:PHE:CD1	1:C:329:VAL:HG22	2.40	0.57
1:C:329:VAL:O	1:C:329:VAL:HG23	2.05	0.57
1:B:295:LEU:HD21	1:B:302:LEU:HD13	1.85	0.57
1:A:142:ALA:O	1:A:146:GLU:HG2	2.04	0.56
1:B:154:PRO:HG2	1:B:155:PHE:HD1	1.66	0.56
1:H:245:GLU:O	1:H:248:LYS:HG2	2.06	0.56
1:E:189:PRO:HD2	1:E:199:THR:HG21	1.87	0.56
1:B:265:PRO:HD2	1:B:268:GLU:OE2	2.04	0.56
1:A:144:TYR:CD2	1:A:147:LEU:HD12	2.41	0.56
1:G:110:THR:N	9:G:1117:HOH:O	2.39	0.56
1:B:123:GLU:O	1:B:127:THR:HG23	2.06	0.56
1:H:174:GLN:HG2	1:H:177:GLU:OE1	2.06	0.55
1:A:154:PRO:HG2	1:A:155:PHE:CD1	2.41	0.55
1:H:168:SER:O	1:H:172:ASP:N	2.38	0.55
1:E:174:GLN:HB3	1:E:175:PRO:HD2	1.87	0.55
1:E:134:LYS:HG3	1:E:135:GLU:H	1.72	0.55
1:H:160:CYS:HB2	1:H:186:ASN:HD22	1.71	0.55
1:A:270:CYS:O	1:A:274:GLU:HG3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:LYS:HD3	1:A:135:GLU:OE1	2.07	0.54
1:B:307:TYR:CZ	1:B:309:PHE:HB3	2.41	0.54
1:H:168:SER:HA	1:H:173:PHE:HB2	1.89	0.54
9:F:1748:HOH:O	2:G:3009:ACT:H1	2.06	0.54
1:H:146:GLU:HA	1:H:146:GLU:OE1	2.08	0.54
1:A:329:VAL:O	1:A:329:VAL:HG12	2.07	0.54
1:E:270:CYS:O	1:E:274:GLU:HG3	2.07	0.54
1:C:265:PRO:HD2	1:C:268:GLU:OE1	2.08	0.53
1:G:146:GLU:HG2	9:G:1678:HOH:O	2.07	0.53
1:A:137:TYR:CE2	1:B:312:GLY:HA3	2.44	0.53
1:F:265:PRO:HG2	1:F:268:GLU:HB2	1.91	0.53
1:A:295:LEU:HD13	1:A:321:PHE:HB3	1.91	0.53
1:E:275:LYS:O	1:E:278:VAL:HB	2.08	0.53
1:E:264:ALA:HB3	1:E:269:LEU:CD2	2.38	0.53
1:B:329:VAL:OXT	1:C:313:SER:HB2	2.08	0.53
1:B:257:VAL:HG13	1:B:258:LYS:N	2.24	0.53
1:F:141:PRO:HA	9:F:1736:HOH:O	2.09	0.53
1:A:129:PHE:CE1	1:B:314:PHE:HB2	2.43	0.53
1:C:123:GLU:HG3	9:C:1385:HOH:O	2.08	0.53
1:E:133:LYS:O	1:E:133:LYS:HD3	2.09	0.52
1:H:258:LYS:NZ	9:H:1166:HOH:O	2.42	0.52
1:B:189:PRO:HB3	1:B:249:ILE:HD13	1.90	0.52
1:A:251:LEU:N	1:A:252:PRO:HD2	2.24	0.52
1:C:166[A]:CYS:HB2	9:C:1064:HOH:O	2.08	0.52
1:C:251:LEU:N	1:C:252:PRO:HD2	2.25	0.52
1:A:171:LEU:HB2	1:A:173:PHE:CE1	2.45	0.52
1:G:123:GLU:O	1:G:127:THR:HG23	2.10	0.52
1:E:251:LEU:N	1:E:252:PRO:HD2	2.24	0.52
1:E:165:VAL:O	1:E:165:VAL:HG23	2.09	0.52
1:G:264:ALA:HB3	1:G:269:LEU:HD13	1.92	0.52
1:E:229:LEU:HD13	1:E:247:VAL:O	2.10	0.52
1:A:220:HIS:CE1	1:A:223:CYS:HA	2.45	0.51
1:B:226:ILE:O	1:B:229:LEU:HB3	2.10	0.51
1:B:251:LEU:N	1:B:252:PRO:HD2	2.25	0.51
1:H:131:HIS:C	1:H:131:HIS:ND1	2.63	0.51
1:H:132:PHE:HE1	1:H:136:LYS:HD2	1.75	0.51
1:A:319:LEU:HD12	1:A:320:GLU:H	1.75	0.51
1:G:233:PRO:HG2	1:G:235:ASP:OD1	2.11	0.51
1:C:230:LEU:HA	1:C:254:LYS:HE3	1.93	0.51
1:H:256:LYS:HE3	1:H:276:GLU:OE2	2.12	0.50
1:E:329:VAL:HG12	1:E:329:VAL:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:225:GLY:N	2:F:3008:ACT:H1	2.25	0.50
1:G:131:HIS:NE2	1:G:135:GLU:HG3	2.26	0.50
1:B:329:VAL:O	1:B:329:VAL:HG12	2.11	0.50
1:B:208:LEU:HD21	1:B:290:PHE:HB2	1.93	0.50
1:C:230:LEU:HD23	1:C:254:LYS:HG3	1.94	0.50
1:F:329:VAL:HG22	1:H:126:LYS:NZ	2.25	0.50
1:H:257:VAL:HG12	1:H:269:LEU:HD22	1.92	0.50
1:D:227:LYS:HD3	8:D:3101:CIT:H42	1.92	0.50
1:H:254:LYS:HE3	1:H:258:LYS:HE3	1.94	0.50
1:F:143:LEU:HD12	1:F:147:LEU:HD13	1.93	0.50
1:H:232:PHE:CD1	1:H:244:GLU:HG2	2.46	0.49
1:F:265:PRO:O	1:F:268:GLU:N	2.40	0.49
1:A:134:LYS:HG3	1:A:135:GLU:H	1.76	0.49
1:F:234:PHE:HZ	1:F:247:VAL:HG23	1.77	0.49
1:D:123:GLU:OE2	1:D:126:LYS:HD2	2.12	0.49
1:F:119:SER:HB2	9:F:1259:HOH:O	2.12	0.49
1:A:138:ASP:C	1:A:140:ASN:H	2.16	0.49
1:B:261:HIS:HB3	1:B:264:ALA:HB2	1.95	0.49
1:G:270:CYS:O	1:G:274:GLU:HG3	2.12	0.49
1:A:327:PHE:C	1:A:327:PHE:CD1	2.86	0.49
1:A:140:ASN:HB3	1:A:143:LEU:HB3	1.94	0.49
1:E:154:PRO:HG2	1:E:155:PHE:HD1	1.77	0.49
1:C:221:SER:HB3	1:C:308:ASP:HA	1.94	0.49
1:A:160:CYS:N	1:A:186:ASN:HB3	2.28	0.49
1:E:220:HIS:CE1	1:E:223:CYS:HA	2.48	0.48
1:A:140:ASN:HB3	1:A:143:LEU:HB2	1.95	0.48
1:G:237:THR:HA	9:G:1759:HOH:O	2.12	0.48
1:B:318:GLY:O	1:C:323:LEU:HA	2.13	0.48
1:C:307:TYR:CE2	1:C:309:PHE:HB3	2.48	0.48
1:G:307:TYR:CZ	1:G:309:PHE:HB3	2.47	0.48
1:B:189:PRO:HD2	1:B:199:THR:HG21	1.96	0.48
1:C:159:ALA:O	1:C:182:ARG:HA	2.13	0.48
1:E:274:GLU:O	1:E:278:VAL:HG23	2.13	0.48
1:B:300:LEU:HB2	9:B:1706:HOH:O	2.14	0.48
1:B:228:GLY:HA2	1:B:231:SER:OG	2.14	0.48
1:H:307:TYR:CZ	1:H:309:PHE:HB3	2.48	0.48
1:G:111:SER:OG	1:G:117:PRO:HB3	2.14	0.47
1:E:134:LYS:CD	1:E:135:GLU:OE1	2.62	0.47
1:G:230:LEU:HD12	1:G:270:CYS:SG	2.54	0.47
1:C:160:CYS:HB2	1:C:186:ASN:HD22	1.79	0.47
1:D:327:PHE:CD1	1:D:327:PHE:C	2.87	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:256:LYS:HD3	1:E:276:GLU:HG2	1.97	0.47
1:A:151:GLN:HG3	1:A:210:LEU:HD22	1.96	0.47
1:B:317:TRP:CB	1:C:325:SER:HA	2.45	0.47
1:G:308:ASP:CG	1:G:311:LYS:HD3	2.34	0.47
1:H:132:PHE:CE1	1:H:136:LYS:HD2	2.50	0.47
1:E:256:LYS:HD3	1:E:276:GLU:CG	2.45	0.47
1:B:126:LYS:O	1:B:130:LEU:HG	2.15	0.47
1:E:124:ARG:HD2	1:E:124:ARG:C	2.33	0.47
1:E:307:TYR:CE2	1:E:309:PHE:HB3	2.49	0.47
1:E:284:ASN:OD1	7:E:3304:EDO:H11	2.15	0.47
1:A:131:HIS:O	1:A:135:GLU:HG2	2.15	0.46
1:B:234:PHE:HZ	1:B:247:VAL:HG23	1.80	0.46
1:G:256:LYS:HE2	1:G:260:GLN:OE1	2.15	0.46
1:B:154:PRO:HD2	1:B:177:GLU:O	2.15	0.46
1:C:160:CYS:SG	2:C:3007:ACT:H2	2.55	0.46
1:C:140:ASN:N	1:C:141:PRO:CD	2.78	0.46
1:B:253:ALA:O	1:B:257:VAL:HG12	2.16	0.46
1:E:133:LYS:C	1:E:133:LYS:HD3	2.35	0.46
1:H:131:HIS:O	1:H:131:HIS:ND1	2.49	0.46
1:E:230:LEU:HD23	1:E:254:LYS:HG3	1.97	0.46
1:B:317:TRP:HA	1:C:325:SER:HA	1.97	0.46
1:E:291:VAL:O	1:E:295:LEU:HG	2.16	0.45
1:B:133:LYS:HD3	1:B:133:LYS:O	2.15	0.45
1:B:264:ALA:HB1	1:B:268:GLU:OE2	2.16	0.45
1:A:131:HIS:CE1	1:A:135:GLU:HG3	2.51	0.45
1:F:257:VAL:HG11	1:F:273:CYS:HA	1.97	0.45
1:B:119:SER:HB3	1:B:122:SER:OG	2.17	0.45
1:F:189:PRO:HB3	1:F:249:ILE:CD1	2.47	0.45
1:F:269:LEU:CD1	1:F:269:LEU:N	2.78	0.45
1:C:264:ALA:HB1	1:C:265:PRO:HD2	1.98	0.45
1:A:261:HIS:CD2	1:A:272:HIS:NE2	2.85	0.45
1:H:146:GLU:C	1:H:149:LYS:HG2	2.33	0.45
1:A:307:TYR:CE2	1:A:309:PHE:HB3	2.51	0.45
1:B:189:PRO:HB3	1:B:249:ILE:CD1	2.46	0.45
1:G:116:ILE:CD1	9:G:1278:HOH:O	2.65	0.45
1:G:311:LYS:CD	1:G:311:LYS:N	2.80	0.45
1:G:295:LEU:HD21	1:G:302:LEU:HG	1.98	0.45
1:G:159:ALA:O	1:G:182:ARG:HA	2.17	0.45
1:F:174:GLN:O	1:F:177:GLU:HB2	2.17	0.45
1:H:174:GLN:O	1:H:177:GLU:HB2	2.17	0.45
1:D:160:CYS:HB2	1:D:186:ASN:HD22	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:259:ALA:HB2	9:G:1596:HOH:O	2.15	0.45
1:E:227:LYS:HE2	9:E:1688:HOH:O	2.17	0.44
1:F:116:ILE:N	1:F:117:PRO:CD	2.80	0.44
1:E:256:LYS:HE3	1:E:260:GLN:OE1	2.16	0.44
1:G:225:GLY:H	2:G:3002:ACT:H1	1.81	0.44
1:A:195:LYS:HB3	1:A:195:LYS:HE2	1.72	0.44
1:E:276:GLU:OE1	1:E:276:GLU:HA	2.18	0.44
1:G:143:LEU:HD11	1:G:175:PRO:HG2	1.99	0.44
1:A:180:VAL:HG12	1:A:182:ARG:HG3	2.00	0.44
1:B:174:GLN:HB3	1:B:175:PRO:HD2	1.99	0.44
1:A:154:PRO:HG2	1:A:155:PHE:HD1	1.81	0.44
1:B:282:LEU:HA	1:B:285:LEU:HD12	2.00	0.43
1:E:243:ILE:O	1:E:247:VAL:HG22	2.18	0.43
1:G:195:LYS:HG2	1:H:194:ALA:O	2.18	0.43
1:D:180:VAL:HG12	1:D:182:ARG:HG3	2.00	0.43
1:B:143:LEU:HD21	1:B:175:PRO:HG2	1.99	0.43
1:B:124:ARG:NH2	1:B:128:GLY:HA3	2.34	0.43
1:H:144:TYR:HA	1:H:147:LEU:HD12	2.00	0.43
1:F:308:ASP:CG	1:F:311:LYS:HD3	2.38	0.43
1:B:271:THR:O	1:B:275:LYS:HG3	2.19	0.43
1:E:161:SER:O	1:E:162:ASP:C	2.57	0.43
1:A:213:SER:O	1:A:300:LEU:HA	2.19	0.43
1:E:134:LYS:HD2	1:E:135:GLU:OE1	2.18	0.43
1:D:307:TYR:CZ	1:D:309:PHE:HB3	2.53	0.43
1:G:147:LEU:HD23	1:G:147:LEU:HA	1.80	0.43
1:B:124:ARG:O	1:B:124:ARG:NE	2.47	0.43
1:B:168:SER:HA	1:B:173:PHE:HB2	2.01	0.43
1:F:130:LEU:HD23	1:F:130:LEU:HA	1.83	0.43
1:G:232:PHE:CG	1:G:244:GLU:HG2	2.53	0.43
1:D:159:ALA:O	1:D:182:ARG:HA	2.18	0.43
1:A:179:PHE:CD2	1:A:206:ALA:HB2	2.54	0.42
1:H:134:LYS:HG3	1:H:135:GLU:HG2	2.01	0.42
1:A:177:GLU:OE2	1:B:124:ARG:HD2	2.19	0.42
1:E:304:GLY:O	1:E:316:LEU:HD12	2.19	0.42
1:B:159:ALA:C	1:B:186:ASN:HB3	2.39	0.42
1:G:258:LYS:HE3	9:G:1389:HOH:O	2.19	0.42
1:A:126:LYS:HG3	1:C:329:VAL:HG21	2.01	0.42
1:E:171:LEU:HB2	1:E:173:PHE:CD1	2.53	0.42
1:C:266:PHE:O	1:C:269:LEU:HB2	2.19	0.42
1:E:174:GLN:O	1:E:177:GLU:HB2	2.19	0.42
1:D:295:LEU:HD13	1:D:321:PHE:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:166[A]:CYS:SG	9:H:1932:HOH:O	2.62	0.42
1:F:185:ALA:HB2	1:F:225:GLY:HA3	2.00	0.42
1:G:141:PRO:HA	9:G:1490:HOH:O	2.19	0.42
1:F:159:ALA:O	1:F:182:ARG:HA	2.20	0.42
1:F:165:VAL:O	1:F:165:VAL:HG23	2.19	0.42
1:B:321:PHE:CD2	1:B:321:PHE:C	2.93	0.42
1:G:305:GLY:HA2	1:G:315:GLU:O	2.19	0.42
1:F:269:LEU:HD12	1:F:269:LEU:N	2.35	0.41
1:A:124:ARG:HH21	1:A:128:GLY:CA	2.32	0.41
1:H:251:LEU:N	1:H:252:PRO:HD2	2.35	0.41
1:G:194:ALA:O	1:H:195:LYS:HG2	2.19	0.41
1:F:307:TYR:CZ	1:F:309:PHE:HB3	2.55	0.41
1:F:275:LYS:O	1:F:278:VAL:HB	2.20	0.41
1:A:326:THR:O	1:A:327:PHE:HB3	2.20	0.41
1:A:171:LEU:HB2	1:A:173:PHE:CD1	2.55	0.41
1:E:253:ALA:O	1:E:257:VAL:HG22	2.20	0.41
1:A:165:VAL:HG23	1:A:165:VAL:O	2.19	0.41
1:E:138:ASP:O	1:E:141:PRO:HD3	2.21	0.41
1:A:295:LEU:HD23	1:A:300:LEU:O	2.21	0.41
1:H:307:TYR:CE2	1:H:309:PHE:HB3	2.56	0.41
1:A:204:GLU:HG3	1:A:208:LEU:HD12	2.02	0.41
1:A:228:GLY:O	1:A:229:LEU:C	2.58	0.41
1:E:225:GLY:N	2:E:3005:ACT:H1	2.22	0.41
1:F:133:LYS:HD3	1:F:133:LYS:C	2.40	0.41
1:F:220:HIS:CE1	1:F:223:CYS:HA	2.55	0.41
1:F:124:ARG:HD2	1:F:124:ARG:O	2.21	0.41
1:G:166[B]:CYS:SG	1:G:169:HIS:HD2	2.44	0.41
1:D:311:LYS:NZ	9:D:1559:HOH:O	2.54	0.41
1:A:137:TYR:HA	1:A:144:TYR:CE1	2.56	0.41
1:C:184:VAL:HG13	2:C:3007:ACT:H1	2.03	0.41
1:E:180:VAL:HG12	1:E:181:VAL:N	2.35	0.41
1:E:327:PHE:C	1:E:327:PHE:CD1	2.94	0.41
1:E:173:PHE:HD2	1:E:177:GLU:HB3	1.85	0.41
1:G:116:ILE:HD12	9:G:1278:HOH:O	2.20	0.41
1:B:161:SER:O	1:B:162:ASP:C	2.59	0.41
1:H:165:VAL:O	1:H:166[A]:CYS:C	2.59	0.40
1:F:154:PRO:HG2	1:F:155:PHE:CD1	2.57	0.40
1:F:166:CYS:SG	1:F:168:SER:HB2	2.62	0.40
1:A:159:ALA:C	1:A:186:ASN:HB3	2.42	0.40
1:A:129:PHE:CZ	1:B:314:PHE:HB2	2.56	0.40
1:H:232:PHE:CG	1:H:244:GLU:HG2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:189:PRO:HB3	1:F:249:ILE:HD13	2.03	0.40
1:G:220:HIS:CE1	1:G:223:CYS:HA	2.55	0.40
1:H:140:ASN:N	1:H:141:PRO:CD	2.84	0.40
1:G:265:PRO:O	1:G:268:GLU:HB2	2.22	0.40
1:H:146:GLU:HA	1:H:149:LYS:HG2	2.03	0.40
1:A:308:ASP:OD1	1:A:310:VAL:HB	2.21	0.40
1:G:251:LEU:HD23	1:G:251:LEU:HA	1.89	0.40
1:C:171:LEU:HA	1:C:171:LEU:HD23	1.83	0.40
1:H:320:GLU:HG3	9:H:1907:HOH:O	2.22	0.40
1:H:146:GLU:OE1	1:H:149:LYS:NZ	2.50	0.40
1:F:243:ILE:O	1:F:247:VAL:HG22	2.21	0.40
1:H:189:PRO:HB2	1:H:196:TYR:CD1	2.57	0.40
1:G:133:LYS:O	1:G:133:LYS:HD3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	208/221 (94%)	190 (91%)	17 (8%)	1 (0%)	34	20
1	B	211/221 (96%)	192 (91%)	17 (8%)	2 (1%)	21	9
1	C	216/221 (98%)	209 (97%)	7 (3%)	0	100	100
1	D	210/221 (95%)	205 (98%)	5 (2%)	0	100	100
1	E	210/221 (95%)	194 (92%)	15 (7%)	1 (0%)	34	20
1	F	212/221 (96%)	200 (94%)	10 (5%)	2 (1%)	21	9
1	G	219/221 (99%)	206 (94%)	12 (6%)	1 (0%)	34	20
1	H	210/221 (95%)	199 (95%)	9 (4%)	2 (1%)	19	7
All	All	1696/1768 (96%)	1595 (94%)	92 (5%)	9 (0%)	34	20

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	111	SER
1	A	139	LYS
1	F	119	SER
1	B	222	ALA
1	H	135	GLU
1	B	162	ASP
1	E	139	LYS
1	F	262	GLY
1	H	140	ASN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	170/180 (94%)	166 (98%)	4 (2%)	57	46
1	B	173/180 (96%)	169 (98%)	4 (2%)	58	48
1	C	177/180 (98%)	174 (98%)	3 (2%)	68	61
1	D	172/180 (96%)	168 (98%)	4 (2%)	58	48
1	E	172/180 (96%)	166 (96%)	6 (4%)	43	28
1	F	174/180 (97%)	170 (98%)	4 (2%)	58	48
1	G	180/180 (100%)	174 (97%)	6 (3%)	45	31
1	H	172/180 (96%)	167 (97%)	5 (3%)	50	36
All	All	1390/1440 (96%)	1354 (97%)	36 (3%)	54	42

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	124	ARG
1	A	241	ASP
1	A	263	ASP
1	A	309	PHE
1	B	118	LYS

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Mol	Chain	Res	Type
1	B	124	ARG
1	B	241	ASP
1	B	245	GLU
1	C	143	LEU
1	C	213	SER
1	C	320	GLU
1	D	120	GLU
1	D	124	ARG
1	D	133	LYS
1	D	258	LYS
1	E	119	SER
1	E	124	ARG
1	E	135	GLU
1	E	263	ASP
1	E	276	GLU
1	E	309	PHE
1	F	118	LYS
1	F	245	GLU
1	F	263	ASP
1	F	315	GLU
1	G	111	SER
1	G	118	LYS
1	G	140	ASN
1	G	143	LEU
1	G	269	LEU
1	G	323	LEU
1	H	123	GLU
1	H	124	ARG
1	H	140	ASN
1	H	174	GLN
1	H	252	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	HIS
1	A	186	ASN
1	A	260	GLN
1	B	186	ASN
1	B	260	GLN
1	C	169	HIS
1	C	186	ASN

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Mol	Chain	Res	Type
1	C	260	GLN
1	D	186	ASN
1	F	209	HIS
1	G	169	HIS
1	H	186	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 ⓘ

Of 38 ligands modelled in this entry, 20 are monoatomic - leaving 18 for Mogul analysis.

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	ACT	A	3001	4	1,3,3	2.65	1 (100%)	0,3,3	0.00	-
2	ACT	A	3003	4	1,3,3	2.65	1 (100%)	0,3,3	0.00	-
3	AZI	A	3203	-	0,2,2	0.00	-	0,1,1	0.00	-
7	EDO	A	3303	-	3,3,3	0.79	0	2,2,2	0.49	0
2	ACT	C	3004	-	1,3,3	2.92	1 (100%)	0,3,3	0.00	-
2	ACT	C	3007	-	1,3,3	3.14	1 (100%)	0,3,3	0.00	-
8	CIT	D	3101	-	3,12,12	3.84	1 (33%)	3,17,17	0.44	0
3	AZI	D	3202	-	0,2,2	0.00	-	0,1,1	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	EDO	D	3302	-	3,3,3	0.84	0	2,2,2	1.08	0
2	ACT	E	3005	4	1,3,3	1.88	0	0,3,3	0.00	-
3	AZI	E	3204	-	0,2,2	0.00	-	0,1,1	0.00	-
7	EDO	E	3304	-	3,3,3	0.75	0	2,2,2	0.51	0
2	ACT	F	3008	-	1,3,3	2.84	1 (100%)	0,3,3	0.00	-
2	ACT	G	3002	4	1,3,3	3.04	1 (100%)	0,3,3	0.00	-
2	ACT	G	3009	-	1,3,3	2.27	1 (100%)	0,3,3	0.00	-
2	ACT	H	3006	4	1,3,3	2.04	1 (100%)	0,3,3	0.00	-
3	AZI	H	3201	-	0,2,2	0.00	-	0,1,1	0.00	-
7	EDO	H	3301	-	3,3,3	1.22	0	2,2,2	0.60	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
2	ACT	A	3001	4	-	0/0/0/0	0/0/0/0
2	ACT	A	3003	4	-	0/0/0/0	0/0/0/0
3	AZI	A	3203	-	-	0/0/0/0	0/0/0/0
7	EDO	A	3303	-	-	0/1/1/1	0/0/0/0
2	ACT	C	3004	-	-	0/0/0/0	0/0/0/0
2	ACT	C	3007	-	-	0/0/0/0	0/0/0/0
8	CIT	D	3101	-	-	0/6/16/16	0/0/0/0
3	AZI	D	3202	-	-	0/0/0/0	0/0/0/0
7	EDO	D	3302	-	-	0/1/1/1	0/0/0/0
2	ACT	E	3005	4	-	0/0/0/0	0/0/0/0
3	AZI	E	3204	-	-	0/0/0/0	0/0/0/0
7	EDO	E	3304	-	-	0/1/1/1	0/0/0/0
2	ACT	F	3008	-	-	0/0/0/0	0/0/0/0
2	ACT	G	3002	4	-	0/0/0/0	0/0/0/0
2	ACT	G	3009	-	-	0/0/0/0	0/0/0/0
2	ACT	H	3006	4	-	0/0/0/0	0/0/0/0
3	AZI	H	3201	-	-	0/0/0/0	0/0/0/0
7	EDO	H	3301	-	-	0/1/1/1	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	D	3101	CIT	O7-C3	-6.65	1.32	1.43
2	H	3006	ACT	CH3-C	2.04	1.51	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	3009	ACT	CH3-C	2.27	1.52	1.48
2	A	3001	ACT	CH3-C	2.65	1.52	1.48
2	A	3003	ACT	CH3-C	2.65	1.52	1.48
2	F	3008	ACT	CH3-C	2.84	1.52	1.48
2	C	3004	ACT	CH3-C	2.92	1.52	1.48
2	G	3002	ACT	CH3-C	3.04	1.53	1.48
2	C	3007	ACT	CH3-C	3.14	1.53	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	3001	ACT	2	0
2	A	3003	ACT	1	0
2	C	3007	ACT	2	0
8	D	3101	CIT	1	0
2	E	3005	ACT	2	0
7	E	3304	EDO	1	0
2	F	3008	ACT	2	0
2	G	3002	ACT	1	0
2	G	3009	ACT	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.